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# Interim Groundwater Monitoring Report (2019-2020)

*Union Pacific Railroad Houston Wood Preserving Works Facility*

*Post-Closure Care Permit No HW-50343/Industrial SWR No. 31547*

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## 1.0 INTRODUCTION

This Interim Groundwater Monitoring Report (IGMR) was prepared by Golder Associates Inc. (Golder) on behalf of Union Pacific Railroad (UPRR) to summarize the site-wide groundwater sampling activities conducted in July 2019 and the additional groundwater assessment activities and site-wide sampling event in January-March 2020 at the UPRR Houston Wood Preserving Works Facility (the Site) located at 4910 Liberty Road, Houston, Texas. The January-March 2020 site-wide groundwater sampling included the sampling of the additional monitoring wells that were installed in February through March 2020 to further assess groundwater at the Site, as discussed in Section 2.0 below. The analytical data from the recently installed wells are included in this IGMR.

Details on the additional groundwater investigations, groundwater monitoring activities, groundwater elevation data and analytical results are provided in the following sections. In addition, information regarding off-site property notification in accordance with Texas Risk Reduction Program (TRRP) Rules 350.55(a) is also provided.

## 2.0 ADDITIONAL GROUNDWATER INVESTIGATION ACTIVITIES

### 2.1 Assessment Methods

As detailed in the Conceptual Response Action Plan (RAP) dated July 9, 2019 (Golder, 2019b) in response to the TCEQ 4<sup>th</sup> Technical Notice of Deficiency (TNOD) dated April 11 2019, UPRR proposed to install additional monitoring wells to refine the lateral and vertical extent of the dissolved phase plume in the A-Transmissive Zone (A-TZ), B-Cohesive Zone/ B-Transmissive Zone (B-TZ/B-CZ), and C-Transmissive Zone (C-TZ). The TCEQ reviewed the proposed monitoring well locations and requested additional monitoring wells be installed as detailed in a TCEQ September 6, 2019 letter. Following a meeting with the TCEQ on September 23, 2019 discussing the proposed locations, UPRR prepared a response letter dated October 23, 2019 to the TCEQ confirming the number and location of monitoring wells to be installed. The TCEQ issued a letter dated December 13, 2019 approving UPRR to proceed with the planned monitoring well installation activities. The following monitoring wells were installed during the period from February through March 2020 and were sampled in March 2020 (locations shown on Figure 1):

- Nine A-TZ wells, MW-47A, MW-84A, MW-88A, MW-91A, MW-94A, MW-95A, MW-97A, MW-98A, and MW-60AR (replacement well);
- Ten B-TZ/B-CZ wells - MW-50B, MW-54B, MW-60B, MW-61B, MW-76B, MW-88B, MW-92B, MW-93B, MW-96B, and MW-98B; and
- Two C-TZ wells - MW-70C and MW 99C.

After each monitoring well was installed, the wells were developed to remove sediment from the wells.

Permanent monitoring wells were surveyed by a licensed, professional surveyor to Texas State Plane coordinates (NAD 27, Texas South Central, U.S. Feet). A summary of the monitoring well construction for wells located at the Site is provided on Table A-1 of Attachment A. Monitoring well boring logs for the monitoring wells installed January through March 2020 are provided in Attachment A.

### 2.2 Data Quality

Samples collected from the Site were analyzed in accordance with the guidelines of U.S. Environmental Protection Agency (EPA) SW-846, Test Methods for Evaluating Solid Waste-Physical/Chemical Methods. The procedures for laboratory analysis, with any modifications, are further documented in the laboratory standard operating procedures, which are maintained at the laboratory, and are listed in the laboratory's quality assurance

plan. Analyses of groundwater samples for these sampling events were performed by ALS Laboratory Group in Houston, Texas. Data obtained from field and laboratory measurements were reviewed for conformance to project requirements, ensuring the lowest method quantitation limit (MQL) was used in the evaluation.

Analytical data from the groundwater samples collected during the groundwater investigations from 2019 through 2020 were reviewed for adherence to established quality assurance/quality control (QA/QC) criteria, and Data Usability Summaries (DUSs) were prepared following TCEQ publication RG-366/TRRP-13 (TCEQ, 2010) to demonstrate the quality of the laboratory analytical data and present any deviations from the established QA/QC criteria. Copies of the analytical data reports and details of the DUSs are provided in Attachment B. For data included in this IGMR, the groundwater data are considered usable for the purpose of evaluating chemicals of concern (COCs) in the environmental media to assess the affected property based on the COCs and establish TRRP protective concentration level exceedance (PCLE) zones.

### 3.0 GROUNDWATER MONITORING ACTIVITIES

Golder, on behalf of UPRR, conducted groundwater monitoring activities at the Site in July 2019 and January-March 2020 to further evaluate the lateral and vertical extent COCs in groundwater and continue to evaluate groundwater COC trends at the Site. Groundwater samples were collected using low-flow sampling techniques described in the EPA guidance document Low-Flow (Minimal Drawdown) Ground Water Sampling Procedures (EPA, 1996). Each sample was analyzed for the site-specific COC list (volatile organic compounds (VOCs) by EPA Method 8260, semi-volatile organic compounds (SVOCs) by EPA Method 8270C, and arsenic by EPA Method 6020).

Prior to sampling, the depth to water and the thickness or presence of non-aqueous phase liquids (NAPL) was measured using an oil-water interface probe. The wells were then purged at a low flow rate using a peristaltic pump with dedicated tubing with the pump intake near the lower portion of the screened interval, unless dense NAPL (DNAPL) was present. Purging was accomplished in such a way as to minimize disturbance of sediments at the bottom of the well, and therefore minimize turbidity of the water samples. During well purging, field parameters, including specific conductance, pH, temperature, dissolved oxygen, oxidation/reduction potential (redox), and turbidity were monitored with field meters. Field meters were calibrated before sampling each day, using the manufacturer's recommended procedures. Odor and color of the purge water were also noted on the groundwater sampling record. Each monitoring well was purged until the field parameters pH, specific conductance, and temperature had stabilized, or until the well purged dry. Groundwater sampling records from July 2019 through March 2020 are provided in Attachment C.

After purging, groundwater samples were collected from the discharge of the peristaltic pump following low-flow sampling techniques. In July 2019, none of the wells with detectable DNAPL were sampled. However, as requested by the TCEQ in a letter dated September 6, 2019 (Additional Comments to UPRR's July 10, 2019 Response to 4<sup>th</sup> TNOD), wells with detectable DNAPL were sampled during the January–March 2020 sampling event. For wells with DNAPL measured, the pump intake was raised to at least 2 to 3 feet above the top of the measured DNAPL.

Since dedicated tubing was used in each well, no equipment rinse sample was collected. Sampling information (i.e., sample time, bottle sets, sampler name, use of filter, etc.) was recorded on the groundwater sampling forms. Groundwater samples were placed in coolers and delivered to ALS Laboratory Group in Houston, Texas for analysis. The samples were analyzed in accordance with EPA protocol for the analytical methods requested.



Analytical laboratory reports and DUSs for both sampling events are provided in Attachment B.

### 3.1 July 2019 Site-Wide Sampling event

The July 2019 sampling event was conducted between July 9 and July 31, 2019, in which 97 groundwater monitoring wells were sampled from the following groundwater bearing units (GWBUs):

- Forty (40) A-TZ monitoring wells;
- Thirty-two (32) B-TZ/B-CZ monitoring wells;
- Twenty-one (21) C-TZ monitoring wells; and
- Four (4) D-TZ monitoring wells.

Monitoring wells containing detectable DNAPL were not sampled during this monitoring event.

### 3.2 January-March 2020 Site-Wide Sampling Event

The January-March 2020 sampling event was conducted between January 9 and March 20, 2020 (January-March 2020 sampling event), and included the sampling of monitoring wells installed in February and March 2020. During the January-March 2020 sampling event, 127 groundwater monitoring wells were sampled from the following GWBUs:

- Forty-eight (48) A-TZ monitoring wells;
- Forty-eight (48) B-TZ/B-CZ monitoring wells;
- Twenty-seven (27) C-TZ monitoring wells; and
- Four (4) D-TZ monitoring wells.

Monitoring wells containing detectable DNAPL were sampled during this monitoring event.

## 4.0 GROUNDWATER ELEVATION DATA

Prior to purging and sampling, the static depth to groundwater in wells was measured from the designated surveyed measuring point with an interface probe to the nearest one hundredth of one foot (0.01 ft). Groundwater elevations for the monitoring wells measured in July 2019 and January 2020 are summarized on Table 1. Potentiometric surface maps for each of the four transmissive zones, A-TZ, B-TZ/B-CZ, C-TZ, and D-TZ, are presented on Figures 5A-1a through and 5A-4a for the July 2019 sampling event, and on Figures 5A-1b through and 5A-4b for the January-March 2020 sampling event. Groundwater flow directions are described in the following sections.

### 4.1 A-TZ Wells

Groundwater in the A-TZ generally flows from southwest to east-northeast across the Site at a gradient ranging from approximately 0.002 ft/ft to 0.016 ft/ft (approximate average of 0.004 ft/ft), with a groundwater divide on the east side of the Site just west of the Lockwood Road Bridge (Figures 5A-1a and 5A-1b). As discussed in the Updated APAR Addendum (PBW, 2010), the 60-in wastewater line runs north to south just west of the Lockwood Bridge and appears to intersect the A-TZ. Groundwater flow in the A-TZ flows to the east on the west side of the wastewater line and flows to the west on the east side of the wastewater line.

The highest groundwater elevation in the A-TZ during the July 2019 and January 2020 event was near the west side of the Site at Solid Waste Management Unit (SWMU) 1 (45.35 feet relative to the City of Houston Vertical Datum (HVD) (MW-08, July 2019) and 45.47 ft HVD (MW-02, January 2020)), with the lowest elevations near the east side of the Site along Lockwood Drive (34.06 feet HVD (MW-18A, July 2019) and 33.86 ft HVD (MW-18A, January 2020)) near the area where the wastewater line is located.

## 4.2 B-CZ/B-TZ Wells

Groundwater in the B-TZ generally flows from west to east-northeast across the Site at gradients ranging from 0.003 ft/ft to 0.02 ft/ft, with an average of approximately 0.004 ft/ft. As shown on Figures 5A-2a and 5A-2b, there is a piezometric high near the west perimeter of the Site near SWMU 1, similar to the flow conditions in the A-TZ. Off-site to the north, the groundwater flow conditions in the B-TZ are more variable. In July 2019 and January 2020, groundwater flow was to the east, north, or northwest depending on the location. For the two monitoring events, the highest groundwater elevation in the B-TZ was 45.70 feet HVD (P-12 (July 2019), and lowest elevation in the B-TZ wells was 36.09 feet HVD (MW-80B, July 2019).

As discussed in the Response Action Plan (RAP) (Attachment 1A - APAR Addendum, PBW, 2014), there is a lateral change in the hydrogeology of the B-TZ as it pinches out from west to east into the B-CZ clay unit, where groundwater is encountered in very thin carbonate seams (typically less than 0.1 feet thick) within the clay unit at relatively the same elevation as the B-TZ. The B-CZ, where these thin carbonate seams are found, was classified as a Class 3 groundwater resource that was approved by the TCEQ in the 3<sup>rd</sup> TNOD letter dated April 10, 2017.

Groundwater flow in the B-CZ during the gauging events show variable flow direction with flow to the east-northeast on the east portion of the Site. Groundwater potentiometric elevations from the off-site wells north of the Site (i.e., MW-36B and MW-59B) indicate flow to the north toward MW-67B (Figures 5A-2a and 5A-2b). There is also a groundwater elevation high near off-site wells MW-36B and MW-71B showing flow radiating from that area (Figures 5A-2a and 5A-2b). The highest groundwater elevation in the B-CZ in July 2019 was 43.97 feet HVD (MW-71B), and the lowest elevation in the B-CZ wells was 34.20 feet HVD (MW-49B) (Figure 5A-2a)

## 4.3 C-TZ Wells

Groundwater flow direction in the C-TZ is generally from northeast to southwest across the Site (Figures 5A-3a and 5A-3b) with a gradient about 0.001 ft/ft. The flow direction is similar to previous gauging events in the C-TZ. Groundwater elevations measured in July 2019 and January 2020 events ranged from a high of approximately 29.72 feet HVD (MW-53C, July 2019) to 25.52 feet HVD (MW-86C, January 2020). There are two areas showing relatively low potentiometric elevations, near MW-23C and in the vicinity of MW-44C, MW-45C, and MW-46C. Both areas, where current DNAPL recovery efforts are being conducted (i.e., monthly recovery events), have potentiometric elevations about 1 to 3 feet lower than the surrounding potentiometric elevations (Figures 5A-3a (July 2019) and 5A-3b (January 2020)).

## 4.4 D-TZ Wells

Groundwater flow direction in the D-TZ appears to be from the southeast to northwest during the July 2019 event (Figure 5A-4a) and from the east to west during the January 2020 event (Figure 5A-4b) at gradients ranging from 0.0015 ft/ft to 0.0038 ft/ft. Groundwater elevations range from a high of -34.75 feet HVD (MW-66D (January 2020) to a low of -38.84 feet HVD (MW-59D, January 2020).

## 5.0 GROUNDWATER ANALYTICAL RESULTS

Groundwater analytical data were compared to the TCEQ TRRP Residential Groundwater PCLs or Residential Assessment Levels (RALs) (the lesser value between <sup>GW</sup>GW<sub>Ing</sub> and <sup>Air</sup>GW<sub>Inh-V</sub> PCLs), dated November 2019, assuming the source area greater than 0.5 acre in size (30-acre source area). Critical PCLs (cPCLs) to assess PCLE Zones were then established assuming residential or commercial/industrial (C/I) land use depending on the location of the monitoring well (i.e., on-site wells compared to C/I PCLs and off-site wells compared to residential PCLs). The PCLE Zones were established based on the COC concentrations for both the July 2019 and January-March 2020 sampling events. Semi-annual groundwater monitoring data from January 2009 through March 2020 were used to construct the following summary tables and evaluate trends in COC concentrations in groundwater:

- Table 1 – Groundwater Measurements
- Table 2 – Analytical Results Summary – Class 2 Groundwater
- Table 3 – Analytical Results Summary – Class 3 Groundwater
- Table 4 - Analytical Results Summary – A-TZ Wells – Vapor Intrusion Screening Level (VISL) Comparison (2019-2020)

COCs evaluated were the site-specific COCs identified in the RCRA Facility Investigation (RFI) Work Plan (IC, 1994) prepared for the Site. In addition to the 34 site-specific COCs, groundwater samples from selected wells were analyzed for vinyl chloride by EPA Method 8260 and samples from the site-wide monitoring wells were analyzed for arsenic by EPA Method 6020 based on previous groundwater sampling events. Comparing the maximum groundwater analytical data from the July 2019 and January 2020 groundwater sampling events to cPCLs, concentrations of 18 target COCs (plus arsenic) exceeded their respective RALs:

### Volatile Organic Compounds (VOCs)

- Benzene (A-TZ, B-CZ/B-TZ, C-TZ)

### Semi-Volatile Organic Compounds (SVOCs)

- 2,4-Dimethylphenol (A-TZ, B-TZ)
- 2,6-Dinitrotoluene (C-TZ)
- 2-Methylnaphthalene (A-TZ, B-CZ/B-TZ, & C-TZ)
- *Acenaphthene (B-CZ/B-TZ only\*)*
- *Acenaphthylene (B-CZ/B-TZ only\*)*
- *Anthracene (B-CZ only\*)*
- Benzo(a)anthracene (B-CZ/B-TZ)
- Benzo(a)pyrene (A-TZ, B-CZ/B-TZ, & C-TZ)
- *Chrysene (B-TZ only\*)*
- Dibenzofuran (A-TZ, B-CZ/B-TZ, & C-TZ)
- *Fluoranthene (B-CZ/B-TZ only\*)*
- *Fluorene (B-CZ/B-TZ only\*)*
- Naphthalene (A-TZ, B-CZ/B-TZ, & C-TZ)
- Pentachlorophenol (C-TZ)
- Phenanthrene (B-CZ/B-TZ only)

Volatile Organic Compounds (VOCs)

Semi-Volatile Organic Compounds (SVOCs)

- Phenol (A-TZ, B-CZ)
- Pyrene (B-CZ/B-TZ only\*)

\* - COC only detected in wells with DNAPL present

As noted above, seven SVOC concentrations (acenaphthene, acenaphthylene, anthracene, chrysene, fluoranthene, fluorene, and pyrene) were detected above RALs in wells MW-12B, MW-32B, MW-49B, and MW-70B. Each of these wells contained detectable DNAPL during the sampling event. As a result, these concentrations would tend to overestimate the dissolved concentration in the groundwater if some DNAPL was entrained in the sample during sample collection. As an example, naphthalene concentrations were detected in MW-12B at 760 mg/L in January 2020, which is more than an order of magnitude greater than the solubility of naphthalene in water at 41 mg/L, thereby demonstrating the entrainment of some naphthalene-containing NAPL within the sample collected from this well.

### 5.1 July 2019 and January-March 2020 COC Concentrations

The spatial distribution of the COCs exceeding RALs in each GWBU from the July 2019 and January-March 2020 monitoring events are presented on the following figures (provided in Attachment D and E, respectively for each event):

- Figure 5B-1 for unit A-TZ,
- Figure 5B-2 for B-TZ/B-CZ,
- Figure 5B-3 for C-TZ, and
- Figure 5B-4 for D-TZ.

In addition to the figures listed above, individual COC concentration maps for the most commonly detected groundwater COCs at the Site (primary COCs –VOC benzene, and SVOCs 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene) were prepared for the A-TZ, B-CZ/B-TZ, and C-TZ units at the Site using the July 2019 (Attachment D) and January-March 2020 (Attachment E) data. The figures (Figure 5B-5 through 5B-19) are listed below:

GWBU	A-TZ	B-CZ/B-TZ	C-TZ
COC	Figure Number		
Benzene	5B-5	5B-10	5B-15
2,4 – Dimethylphenol	5B-6	5B-11	5B-16
2-Methylnaphthalene	5B-7	5B-12	5B-17
Dibenzofuran	5B-8	5B-13	5B-18
Naphthalene	5B-9	5B-14	5B-19

For each COC concentration map, a cumulative groundwater PCLE Zone based on the combined PCLE Zones for each individual COC in each GWBU is shown on the figures.



Monitoring wells in the four GWBUs were also sampled and analyzed for arsenic. Arsenic concentrations from the July 2019 and January-March 2020 sampling events were detected above the TCEQ TRRP RAL and cPCL in the A-TZ and B-CZ/B-CZ zones (Tables 2 and 3). Arsenic isopleth maps were generated for the A-TZ, B-CZ/B-TZ, and C-TZ based on the 2019 and 2020 data (Figure 5B-20 through Figure 5B-22 in Attachment D (July 2019) and E (January-March 2020)). Details of the distribution of the COCs detected in each transmissive zone are discussed below.

### 5.1.1 A-TZ Wells

#### VOCs

During the July 2019 and January-March 2020 groundwater monitoring events, benzene concentrations were the only VOCs that were detected above its RAL and cPCL of 0.005 mg/L in A-TZ wells. Benzene concentrations detected above the RAL were predominantly located on the eastern portion of the Site and in the northern portion of the Englewood Intermodal Yard (Figure 5B-5 (Attachments D and E)).

Except for the benzene concentration detected at off-site well MW-26A at 0.00036J mg/L (July 2019 event), an order of magnitude below the RAL, none of the off-site A-TZ wells to the north had benzene concentrations detected above the sample detection limit (SDL) during the two sampling events. Therefore, benzene concentrations are delineated to the RAL to the north. Benzene concentrations are also delineated to the RAL to the east of the Site in wells MW-47A, MW-59A, MW-60AR, and MW-61A; delineated to the RAL to the south at MW-50A, MW-51A, MW-97A, and MW-98A ; and delineated to the RAL to the west at wells MW-05, MW-09, MW-12A, MW-15A, and MW-88A based on the analytical data from both events. Based on the analytical results, the PCLE Zone for benzene in the A-TZ is within the Site boundary except to the north (along Liberty Road) and east in which it extends onto the City of Houston right-of-way (ROW) (Figure 5B-5, Attachments D and E)).

For the wells installed in February-March 2020, benzene concentrations were below SDLs in the eight new A-TZ wells MW-47A, MW-84A, MW-88A, MW-91A, MW-94A, MW-95A, MW-97A, MW-98A, and in replacement well MW-60AR sampled during the March 2020 sampling event.

Other VOCs (1,2-dichloroethane, chlorobenzene, ethylbenzene, toluene, vinyl chloride, and xylenes) were detected above SDLs in the A-TZ wells sampled during the July 2019 and January-March 2020 sampling events, however, none of the results exceeded their respective RALs.

#### SVOCs

Similar to the benzene detections, SVOCs were detected above the applicable RALs in A-TZ wells located generally on the eastern portion of the Site and in the northern portion of the Englewood Intermodal Yard. The primary SVOCs detected in the A-TZ above RALs included 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene (Figures 5B-6 through 5B-9, Attachments D and E). Analytical data from the July 2019 and January-March 2020 sampling events indicated that concentrations of SVOCs 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene were delineated in the A-TZ to the north, east, south and west below their respective RALs similar to the benzene concentrations. Dibenzofuran was detected above the RAL in one A-TZ well (MW-12A) on the western side of the Site during both sampling events but below the cPCL. Dibenzofuran concentrations were below RALs in the off-site well to the northwest (MW-38A) of MW-12A during both events.

For both sampling events, the groundwater PCLE Zone for the primary SVOCs that exceeded their cPCLs generally falls within the benzene PCLE Zone. Therefore, similar to the benzene PCLE Zone, none of the SVOC

concentrations detected in A-TZ wells offsite exceeded their respective RALs. The PCLE Zones for the individual SVOCs are predominately within the Site boundary except to the north (along Liberty Road) and east in which it extends onto the City of Houston ROW.

For the newly installed wells in February-March 2020, SVOCs 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene concentrations were below RALs in the eight new A-TZ wells MW-47A, MW-84A, MW-88A, MW-91A, MW-94A, MW-95A, MW-97A, MW-98A, and in replacement well MW-60AR during the March 2020 sampling event.

Except for benzo(a)pyrene and phenol, concentrations for the other SVOCs analyzed were either detected below RALs or below the SDL in the A-TZ wells during the July 2019 and January-March 2020 sampling events. Benzo(a)pyrene at MW-57A and phenol at MW-17 were detected at concentrations greater than the respective RALs in July 2019. Benzo(a)pyrene at MW-57A and MW-79A was detected at concentrations greater than the RAL in January-March 2020. Trace DNAPL was observed in MW-57A in early 2019 but has not been detected since then.

### **Arsenic**

Arsenic concentrations have generally fluctuated in A-TZ wells between sampling events (Table 2). Of the A-TZ wells sampled in July 2019, arsenic concentrations in 13 A-TZ wells exceeded the RAL where only 7 of those same wells sampled in January 2020 had arsenic concentrations detected above the RAL. The July 2019 arsenic concentrations in A-TZ wells generally exceed the RAL along the northern boundary of the Site (MW-13, MW-15A, MW-17, and MW-18A), offsite to the north (MW-26A, MW-32AR, MW-33A, MW-35A, MW-44A, and MW-68A), in the southwestern portion of the Site (MW-04 and MW-05), and one well in the central portion of the Englewood Intermodal Yard (MW-77A) (Figure 5B-20, Attachment D). The January-March 2020 arsenic concentrations in A-TZ wells exceed the RAL along the northern boundary of the Site (MW-13, MW-15A, MW-18A, and MW-57A), offsite to the north (MW-26A, MW-35A, and MW-68A), one well offsite to the west (MW-38A), and at one well in the central portion of the Englewood intermodal yard (MW-77A) (Figure 5B-20, Attachment E). For the newly installed wells in February-March 2020, arsenic concentrations were below the RAL in the eight new A-TZ wells MW-47A, MW-84A, MW-88A, MW-91A, MW-94A, MW-95A, MW-97A, MW-98A, and in replacement well MW-60AR during the March 2020 sampling event. Based on the January-March 2020 groundwater data with the newly installed wells, the arsenic concentrations are delineated to the RAL in the A-TZ except near MW-13 on the northwest corner of the Site.

Arsenic concentrations in the A-TZ wells do not appear to correlate with the elevated concentrations of the primary COCs. Many wells with elevated arsenic concentrations (exceeding RAL) do not have exceedances of other COC concentrations, especially north of the Site. For example, arsenic concentrations were highest at MW-26A and MW-13 in July 2019, but all other COC concentrations were detected below RALs at MW-26A and at MW-13 (Table 2). Similar to July 2019, January-March 2020 arsenic concentrations were highest at MW-13, and all other COC concentrations in MW-13 were below RALs (Table 2). Dissolved concentrations of arsenic in groundwater at the Site are likely associated with naturally-occurring arsenic in the aquifer matrix that is reduced from less mobile arsenate species to more mobile arsenite species under reducing conditions caused by the natural degradation of petroleum hydrocarbons (Cozzarelli, et al., 2015).

### **EPA Vapor Intrusion Screening Level Evaluation**

Groundwater analytical data from July 2019 and January-March 2020 for the off-site monitoring wells completed in the A-TZ north, west and east of the Site were also compared to values developed using the U.S. Environmental Protection Agency's (EPA's) vapor intrusion screening level (VISL) calculator (EPA, 2015, EPA,

2019). The VISL calculator was used to calculate conservative, non-site specific, risk-based potential vapor intrusion (VI) screening values for the identified COCs in the A-TZ. The EPA tool calculates the COC concentration in groundwater, based on certain default conditions, at which the COC is not expected to pose an unacceptable VI risk and, as such, can be eliminated from further consideration.

#### EPA VISL Calculator

The VISL calculator (EPA, 2019) provides a screening level based on several basic inputs, including a residential or commercial exposure scenario, target hazard quotient, target carcinogenic risk, and groundwater temperature. For this evaluation, the selected inputs were – residential scenario, hazard quotient of 0.1, carcinogenic risk of  $10^{-5}$  (consistent with the TRRP criteria), and groundwater temperature of 25°C. Based on these inputs, the calculated groundwater screening concentrations were estimated. These values were compared with the July 2019 and January-March 2020 groundwater sample COC concentrations for the Site-specific COCs (Table 4). A summary of the VISL values relative to the maximum detected COC concentrations from the two sampling events are provided in the following table.

COC	VISL Screening Level for Elimination from Further Consideration (mg/L)	Maximum Detected COC Concentrations in Off-Site A-TZ Wells July 2019 & January-March 2020 Samples (mg/L)
Benzene	0.014	0.00036J (MW-26A-July 2019)
Chlorobenzene	0.041	<0.0003
Ethylbenzene	0.035	<0.0003
Methylene Chloride	0.471	<0.001
Toluene	1.92	<0.0002
Xylenes	0.0385	<0.0003
Benzo(a)anthracene	0.344	0.00012 (MW-25A-Jan 2020)
Naphthalene	0.0174	0.0026 (MW-27A – July 2019)
Nitrobenzene	0.715	<0.000025

EPA's VISL calculator uses a conservative default attenuation factor of 0.001 for the attenuation of vapors between the groundwater source and the overlying receptor. This attenuation factor is not adjusted for the more rapid attenuation observed in petroleum hydrocarbon vapors and is therefore likely overly conservative for these compounds. Nevertheless, as indicated above, the COC concentrations from the July 2019 and January-March 2020 off-site A-TZ wells were below the conservative EPA VISL-calculated screening levels by orders of magnitude indicating the VI pathway from the shallow groundwater is incomplete.

## 5.1.2 B-CZ/B-TZ Wells

### VOCs

During the July 2019 and January-March 2020 groundwater monitoring events, benzene concentrations were the only VOCs that were detected above its RAL of 0.005 mg/L (Class 2 groundwater PCL) in B-TZ wells and 0.5 mg/L (Class 3 groundwater PCL) in the B-CZ wells. Benzene concentrations detected above the two RALs in July 2019 and January-March 2020 were predominantly located in area extending from the northeastern portion of the Englewood Intermodal Yard, through the eastern portion of the Site, and off-site to the north (Figure 5B-10 (Attachment D and E)). There is also a smaller area where benzene concentrations exceed the RAL on the western portion of the Site.

Benzene concentrations in the B-CZ are delineated to the Class 3 groundwater RAL to the north and east by MW-33BR, MW-59B, MW-60B, MW-61B, MW-63B, MW-67B, and MW-71B (also shown on Figure 5B-10 (Attachment D and E)). Benzene concentrations in the B-TZ in the eastern portion of the Site are delineated to the RAL to the west by MW-15B, MW-82B, MW-88B, MW-89B, MW-90B, MW-92B, and MW-93B and to the south by MW-50B, MW-76B, MW-80B, MW-81B, and MW-98B. Benzene concentrations in the smaller area on the west side of the Site are delineated to the RAL to the north by MW-39B, to the east by MW-42B, to the south by P-11, and to the east by MW-62B, TW-41B, and MW-38B. Benzene concentrations were below RALs in the ten new B-CZ/B-TZ wells during the March 2020 sampling event: MW-50B, MW-54B, MW-60B, MW-61B, MW-76B, MW-88B, MW-92B, MW-93B, MW-96B, and MW-98B.

The PCLE Zone for benzene based on land use (i.e., C/I for on-Site wells and Residential for off-Site wells) in the B-CZ/B-TZ (Class 3 and 2 groundwater, respectively) for both the July 2019 and January-March 2020 events encompasses the northeastern portion of the Englewood Intermodal Yard, eastern portion of the Site, and extends off-site to the north of the Site (referred to as the "Main" PCLE Zone). The PCLE Zone for the smaller area on the west side of the Site appears to be within the Site property boundary (referred to as the "West" PCLE Zone).

Other VOCs (ethylbenzene, toluene, vinyl chloride, and xylenes) were detected above SDLs in the B-CZ/B-TZ wells sampled during the July 2019 and January-March 2020 sampling events; however, none of the results exceeded their respective RALs.

### SVOCs

Similar to the benzene concentrations, primary SVOCs 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene were detected above their respective RALs in B-CZ/B-TZ wells located generally in the area extending from the northeastern portion of the Englewood Intermodal Yard, through the eastern portion of the Site, and off-site to the north (Figures 5B-11 through 5B-14 (Attachment D and E)). A smaller area on the western portion of the Site also had RAL exceedances for 2-methylnaphthalene, dibenzofuran, and naphthalene.

Analytical data from the July 2019 and January-March 2020 sampling events indicated that concentrations of SVOCs 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene were delineated in the B-CZ/B-TZ to the north, east, south and west below their respective RALs for both areas. For the newly installed wells in February-March 2020, 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene concentrations were below RALs in the ten new B-CZ/B-TZ wells during the March 2020 sampling event: MW-50B, MW-54B, MW-60B, MW-61B, MW-76B, MW-88B, MW-92B, MW-93B, MW-96B, and MW-98B.



For both sampling events, the groundwater PCLE Zones based on the land use and groundwater classification cPCLs for the primary SVOCs either fall within or are very similar to the benzene PCLE Zones for both the Main and West PCLE Zones.

Except for benzo(a)pyrene in MW-68B and phenol in MW-74B, all other SVOCs were either detected below RALs or below the SDL in the B-CZ/B-TZ wells sampled in July 2019. During the January-March 2020 sampling event, wells with DNAPL were sampled as requested by the TCEQ in a letter dated September 6, 2019. The following other SVOCs were detected above RALs in January-March 2020 in MW-12B, MW-32B, MW-41B, MW-49B, MW-68B, MW-70B, MW-74B, and MW-75B: acenaphthene, acenaphthylene, anthracene, benzo(a)anthracene, benzo(a)pyrene, chrysene, fluoranthene, fluorene, phenanthrene, phenol, and pyrene. Except for MW-74B, DNAPL was detected in each of these B-CZ/B-TZ wells that had SVOC exceedances during the January-March 2020 sampling event. Well MW-74B has not had detectable DNAPL since installation; however, DNAPL was noted in the boring log for this well in the B-CZ/B-TZ interval during monitoring well installation.

### **Arsenic**

Arsenic concentrations in the B-TZ in July 2019 exceeded the Class 2 groundwater RAL (0.01 mg/L) in seven wells at the Site in two areas: the western portion of the Site (MW-40B, TW-41B, and P-11) and offsite north of the Site (MW-15B, MW-68B, MW-83B, and MW-84B). Arsenic concentrations in the B-TZ in January-March 2020 exceeded the RAL in the same two areas but in 11 wells: western portion of the Site (MW-40B, TW-41B, and P-11, MW-12B, MW-41B, and MW-62B) but also extended offsite to the west (MW-22BR and MW-38B), and offsite to the north (MW-35B and MW-83B).

Arsenic levels in B-CZ/B-TZ wells north of the Site do not appear to correlate with the elevated concentrations of the primary COCs in the B-CZ/B-TZ in the same area. Concentrations of primary COCs (benzene and SVOCs) were above RALs in July 2019 and January-March 2020 in the eastern portion of the Site, where arsenic concentrations were below the arsenic RAL in that same area. The arsenic concentrations to the north of the Site from the January-March 2020 sampling event are delineated to the RAL to the north, east, south, and west (Figure 5B-21 (Attachment E)).

B-TZ wells on the western side of the Site had both primary COC and arsenic concentrations above the RALs. Based on the groundwater data from the January-March 2020 sampling event, arsenic concentrations are delineated to the RAL in the B-TZ except to the west near MW-38B and MW-62B (Figure 5B-21 (Attachment E)). Arsenic concentrations in MW-38B and MW-62B have fluctuated above and below the RAL over the past two years. In June 2018, arsenic concentrations in MW-38B (0.0386 mg/L) and MW-68B (0.0112 mg/L) exceeded the RAL, whereas in January and July 2019, arsenic concentrations in MW-38B (<0.0004 mg/L and 0.000553 mg/L J, respectively) and MW-62B (<0.0004 mg/L and 0.00194 mg/L J, respectively) were below the RAL. All other site-related COC concentrations were below respective RALs in MW-38B and MW-62B during the past two years.

Arsenic concentrations were below RALs in the ten new B-CZ/B-TZ wells sampled during the March 2020 sampling event: MW-50B, MW-54B, MW-60B, MW-61B, MW-76B, MW-88B, MW-92B, MW-93B, MW-96B, and MW-98B.

### 5.1.3 C-TZ Wells

#### VOCs

During July 2019 and January-March 2020 groundwater monitoring events, benzene was the only VOC that had concentrations detected above its RAL in C-TZ wells. Benzene concentrations were detected above the RAL in wells located in the northeast portion of the Site (MW-17C and MW-18C (July 2019), and MW-18C (Jan-Mar 2020)) and offsite to the north and northeast (only MW-46C and MW-70C in Jan-Mar 2020) (Figure 5B-15 (Attachment D and E)). DNAPL was historically observed in wells MW-44C, MW-45C, and MW-46C, but has not been detected since June 2019 in MW-44C and MW-46C, and since October 2015 in MW-45C. These three wells were sampled in January 2020, and benzene was detected above the RAL in only MW-46C. Benzene was detected above the RAL in newly installed well MW-70C north of the Site and below the SDL in the other newly installed C-TZ well MW-99C northeast of the Site.

Benzene concentrations are delineated to the RAL to the east of the Site by MW-54C; to the north by MW-28C, MW-53C, MW-87C, and MW-99C; to the west by MW-15C, MW-17C and MW-68C; and to the south by MW-19C, MW-47C, MW-48C, and MW-88C. The PCLE Zone for benzene in the C-TZ is in the northeast portion of the Site and extends offsite to the north (along and across Liberty Road (MW-70C)) and east in which it extends onto the City of Houston ROW (MW-46C) (Figure 5B-15 (Attachments D and E)). The PCLE Zone is delineated based on the same wells listed above.

Other VOCs (chlorobenzene, ethylbenzene, toluene, and xylenes) were detected above SDLs in the C-TZ wells sampled during the July 2019 and January-March 2020 sampling events, however, none of the results exceeded their respective RALs.

#### SVOCs

Following a similar pattern for the benzene detections above its RAL in the C-TZ, SVOCs were detected above the applicable RALs in C-TZ wells located in the northeastern portion of the Site and offsite to the north and northeast. The primary SVOCs detected in the C-TZ above RALs include 2-methylnaphthalene, dibenzofuran, and naphthalene (Figures 5B-16 through 5B-19 (Attachment D and E)). Analytical data from the July 2019 and January-March 2020 sampling events indicated that concentrations of SVOCs 2-methylnaphthalene, dibenzofuran, and naphthalene were delineated in the C-TZ to the north (MW-53C, MW-87C, MW-28C, and MW-99C), east (MW-54C), south (MW-19C, MW-47C, MW-48C, and MW-88C), and west (MW-15C, MW-17C and MW-68C) below their respective RALs similar to the benzene concentrations.

For both sampling events, the groundwater PCLE Zone for the primary SVOCs generally falls within the extent of the benzene PCLE Zone. Therefore, similar to the benzene PCLE Zone, the PCLE Zones for the individual SVOCs are predominately within the Site boundary except to the north (along and across Liberty Road (MW-70C)) and east where it extends onto the City of Houston ROW (MW-25C and MW-46C).

For the newly installed wells in February-March 2020, SVOCs 2,4-dimethylphenol, 2-methylnaphthalene, and dibenzofuran concentrations were at or below RALs in the two new C-TZ wells, MW-70C and MW-99C, during the March 2020 sampling event. The naphthalene concentration in MW-70C was above the RAL in March 2020.

Other SVOCs (2,6-dinitrotoluene, benzo(a)pyrene, and pentachlorophenol) were detected at concentrations that exceeded their RALs in the C-TZ during the July 2019 and January-March 2020 sampling events. The 2,6-dinitrotoluene concentration in MW-53C was detected above the RAL (0.001 mg/L) in July 2019 but below the SDL during the January-March 2020 sampling event. The reported 2,6-dinitrotoluene concentration in MW-21C was below the SDL (0.00004 mg/L) during both sampling events but detected above the RAL at 0.003 mg/L in the

field duplicate collected at MW-21C during the January-March 2020 sampling event (qualified with a J-flag). Benzo(a)pyrene was detected above the RAL (0.0002 mg/L) at well MW-25C in July 2019, but below the SDL during the January-March 2020 sampling event. In January-March 2020, benzo(a)pyrene was detected above the RAL at wells MW-18C, MW-23C, MW-44C, MW-45C, and MW-46C. As previously mentioned, DNAPL was historically observed in wells MW-23C, MW-44C, MW-45C, and MW-46C, and these wells were not sampled in July 2019. Pentachlorophenol was detected above the RAL (0.001 mg/L) in well MW-18C in July 2019 and January 2020 and in off-Site well MW-25C in January 2020. Except for the unverified 2,6-dinitrotoluene concentration in MW-21C, the PCLE Zones for these additional SVOCs are within the benzene and primary SVOC PCLE Zones discussed above.

### **Arsenic**

Arsenic was not detected above the RAL in any of the C-TZ wells in July 2019 or January-March 2020.

## **5.1.4 D-TZ Wells**

### **VOCs**

In July 2019 and January-March 2020, no VOCs were detected above RALs in D-TZ wells. All VOCs were below the SDL.

### **SVOCs**

In July 2019 and January-March 2020, no SVOCs were detected above RALs in D-TZ wells. All SVOCs were either below the SDL or detected (estimated – J-flagged) multiple orders of magnitude below respective RALs.

### **Arsenic**

Arsenic was not detected above the RAL in any of the D-TZ wells in July 2019 or January-March 2020.

## **5.2 COC Concentration Trends**

Graphs of primary COC concentrations (benzene, 2,4-dimethylphenol, 2-methylnaphthalene, dibenzofuran, and naphthalene) over time at each monitoring well are provided in Attachment F. Trends of the primary COC concentrations at monitoring wells in each GWBU from 2009 (or date of installation) to 2020 with at least four sampling events, including samples collected from wells containing detectable DNAPL, were preliminarily evaluated by performing Mann-Kendall trend test. The Mann-Kendall test is a non-parametric statistical test that makes no distributional assumptions of data and is used to analyze linear increasing or decreasing trends in data over time (GSI, 2012). The temporal trend module of the Monitoring and Remediation Optimization System (MAROS) software (developed by GSI Environmental Inc. (GSI) on behalf of Air Force Civil Engineer Center (AFCEC)) was used to run the Mann-Kendall tests at individual wells over time (MAROS version 3 (GSI, 2012)). The Mann-Kendall test was applied to concentrations from each primary COC at each well, comprised of concentrations for detects and detection limits for non-detects.

Using the Mann-Kendall statistic (MK (S)), the trends in the data can then be evaluated. Increases in constituent concentrations over time are indicated by positive values, and decreases in constituent concentrations are indicated by negative values. The MK (S) indicates the direction of the trend and the confidence factor (p-value) indicates the strength of the trend.

- Mann-Kendall test results with a p-value less than 0.05 indicate a statistically significant increasing or decreasing trend (95% level of confidence).
- Results with a p-value greater than 0.05 but less than 0.1 indicate a “probably” increasing or “probably” decreasing trend. Mann-Kendall test results with a p-value greater than 0.10 indicate either a “stable” trend or “no trend”.
- A “stable” trend occurs when data from a constituent in a well has a coefficient of variation less than 1, appears to follow a decreasing trend, and has a p-value from the Mann-Kendall test greater than 0.10.
- In all other cases when the p-value from the Mann-Kendall test is greater than 0.10, the constituent and well is considered to have “no trend” (GSI, 2012).

The preliminary results of the Mann-Kendall analysis and output from the software are provided in Attachment G. The Mann-Kendall test results generally indicated that primary COC concentrations are either stable, decreasing, probably decreasing, or exhibit no trend in most of the monitoring wells at the Site. However, the following Mann-Kendall test results preliminarily indicate an increasing/probably increasing trend for certain primary COC concentrations in selected monitoring wells:

#### A-TZ:

- Results indicate a probably increasing trend in naphthalene in on-site well MW-51A and a probably increasing trend in 2-methylnaphthalene concentrations in on-site well MW-03. However, in both these wells, primary COC concentrations are multiple orders of magnitude lower than applicable PCLs with a majority of the datasets comprising non-detects or estimate values (J-flagged).

#### B-CZ/B-TZ:

- Results indicate an increasing trend in benzene, 2,4-dimethylphenol, 2-methylnaphthalene, and dibenzofuran and probably increasing trends in naphthalene concentrations in on-site well MW-49B. DNAPL has been observed in well MW-49B.
- Results indicate an increasing trend in 2-methylnaphthalene and naphthalene in off-site well MW-70B. No trend was determined for the other primary COCs at MW-70B. DNAPL has also been observed in MW-70B.
- Results indicate an increasing trend in 2-methylnaphthalene, dibenzofuran, and naphthalene and probably increasing trend in benzene in on-site well TW-41B. Test well TW-41B is located between DNAPL recovery wells MW-12B and MW-41B, which both have had DNAPL in-depth thicknesses as thick as 15 to 26 feet. TW-41B is a 4-inch stainless steel well installed in 2009 to serve as a possible DNAPL recovery well given the relatively thick DNAPL noted in the two nearby wells. No DNAPL has been detected in TW-41 since installation.
- Results indicate a probably increasing trend in benzene concentrations in off-site well MW-83B. Since installation in 2018, the benzene concentration in MW-83B has been above the RAL and has ranged from 0.018 mg/L to 0.032 mg/L with a standard deviation of 0.006 mg/L. The highest concentration was detected in January 2019 at 0.032 mg/L, and has decreased to 0.021 mg/L in July 2019 and January 2020. None of the other primary COCs showed increasing trends in this well (decreasing or probably decreasing trends (2-methylnaphthalene, dibenzofuran, and naphthalene) and not trend (2,4-dimethylphenol).
- Results indicate an increasing trend and probably increasing trend in naphthalene concentrations in off-site well MW-90B and on-site well MW-81B, respectively. However, in both these wells, all primary COC concentrations are multiple orders of magnitude lower than applicable PCLs with a majority of the datasets comprising non-detects or estimates (J-flagged).

**C-TZ:**

- Results indicate an increasing trend in 2,4-dimethylphenol concentrations in on-site well MW-23C. DNAPL has been observed in well MW-23C. Mann-Kendall test results indicate no trend or a stable trend for the other primary COC concentrations in MW-23C.
- Results indicate an increasing trend in naphthalene concentrations in on-site MW-86C. However, primary COC concentrations are multiple orders of magnitude lower than applicable PCLs with a majority of the dataset comprising non-detects or estimated values (J-flagged).

Of the 104 wells evaluated, 10 wells listed above showed an increasing or probably increasing trend in primary COC concentrations. DNAPL has been observed in 3 of those 10 wells. For 5 of the 10 wells, the majority of the datasets comprised of non-detects or estimates (J-flagged).

### 5.3 Naphthalene Concentration in MW-11B

As detailed in the Corrective Action Monitoring Report - 2019 Second Semi-Annual Event dated January 17, 2020 (Golder, 2020), naphthalene concentrations were detected in point of compliance (POC) well MW-11B completed in the B-TZ above the Groundwater Protective Standards (GWPS) established in the approved Compliance Plan (CP) Table III for the Closed Surface Impoundment (Solid Waste Management Unit (SWMU) 1) during the second semi-annual monitoring period of 2019.

Based on the analytical results from the July 2019 monitoring event, naphthalene concentrations were detected at POC well MW-11B at 0.70 mg/L, above the GWPS of 0.49 mg/L. As allowed under the CP Section VI.D3, monitoring well MW-11B was resampled on July 30, 2019 for naphthalene to confirm the initial GWPS exceedance. The resampled results indicated that naphthalene concentrations were detected at 1.1 mg/L. Therefore, the initial GWPS exceedance at MW-11B for naphthalene was verified with the resampling event and the POC well was considered to be non-compliant. The TCEQ was notified of the verified exceedance in a letter dated September 5, 2019. An additional groundwater sample from POC well MW-11B was collected in October 2019. The naphthalene concentration in MW-11B in October 2019 was detected at 0.6 mg/L, also above the GWPS. Naphthalene concentrations in groundwater samples from all other SWMU 1 monitoring wells in A-TZ and B-TZ have not exceeded the established CP PCLs since July 2005 and are considered compliant. Following submittal of the Corrective Action Monitoring Report - 2019 Second Semi-Annual Event, the TCEQ issued a comment letter dated March 18, 2020 with the following comment:

*“Based on our review of the above-referenced Report, the TCEQ concurs that the report generally fulfills the reporting requirements of Provision VII.C.2 of the above-referenced Permit with the following comment: UPRR should finish their evaluation of naphthalene concentration trends in well MW-11B, present their findings in the April 30, 2020 Initial Groundwater Sampling Interim Report, and assess whether WMU No. 1 remains in the corrective action program or switches to compliance monitoring.”*

Since submittal of the Corrective Action Monitoring Report - 2019 Second Semi-Annual Event (Golder, 2020), UPRR conducted the first semi-annual corrective action monitoring event in January 2020. The groundwater analytical data from the January 2020 sampling event indicated that the naphthalene concentrations in POC well MW-11B decreased to less than the SDL (<0.002 mg/L). These results indicate that the well is now compliant since the analytical results are less than the GWSP established in CP Table III.

In response to the TCEQ comment, the historical naphthalene concentrations at MW-11B were evaluated. A summary table and a graph of historical naphthalene concentrations in MW-11B from 1994 through 2020 (including the January 2020 sample result) are provided in Attachment H. The historical naphthalene

concentrations show a relatively consistent seasonal pattern of higher naphthalene concentrations detected during the second semi-annual monitoring events typically conducted in July compared to the first semi-annual monitoring events typically conducted in January (Table 1 and Graph 1, Attachment H). From January 2009 through January 2020, naphthalene concentrations during the first semi-annual monitoring event for each year have been below SDLs at MW-11B except for January 2014 when naphthalene concentrations were detected slightly above the SDL at 0.000382J mg/L. In comparison, naphthalene concentrations in MW-11B have been detected above the SDL during each second semi-annual event dating back to 1994, with the recent GWPS exceedances detected in July and October 2019.

UPRR is scheduled to conduct the second semi-annual monitoring event in July 2020, and will assess if POC well MW-11B remains compliant under the current corrective action monitoring program. As detailed in the Section XI (CP) of the RCRA Permit Renewal application (initial submittal-December 2014 through Revision 4 -June 2017), UPRR proposed to revise the CP to move the SWMU 1 to compliance monitoring. UPRR will evaluate the proposed switch to compliance monitoring following review of the second semi-annual 2020 sampling event.

## 6.0 CONCLUSIONS

The following conclusions were made for each GWBU based on the groundwater sampling analytical events from the site-wide events conducted in July 2019 and January-March 2020:

- **A-TZ:**
  - Site-specific VOCs and SVOCs are delineated to applicable RALs in the A-TZ at the Site. The combined PCLE Zone in the A-TZ is within the Site boundary except to the north (along Liberty Road) and east in which it extends onto the City of Houston ROW.
  - Based on the January-March 2020 groundwater data including data from the wells installed in February-March 2020, arsenic concentrations are delineated to the RAL in the A-TZ on-site and off-site except near MW-13 on the northwest corner of the Site. Arsenic concentrations in the A-TZ wells do not appear to correlate with the elevated concentrations of the primary COCs, and are likely associated with naturally-occurring arsenic mobilizing under reducing conditions caused by the natural degradation of petroleum hydrocarbons.
  - Site-specific COC concentrations detected in the off-site A-TZ wells during the July 2019 and January-March 2020 sampling events were below the conservative EPA VISL-calculated screening levels, indicating the off-site VI pathway from groundwater is incomplete.
- **B-CZ/B-TZ:**
  - Site-specific VOCs and SVOCs are delineated to applicable RALs in the B-CZ and B-TZ at the Site. The combined PCLE Zone based on land use (i.e., C/I for on-Site wells and Residential for off-Site wells) in the B-CZ/B-TZ (Class 3 and 2 groundwater, respectively) for both the July 2019 and January-March 2020 events encompasses the northeastern portion of the Englewood Intermodal Yard, eastern portion of the Site, and extends off-site to the north of the Site (referred to as the “Main” PCLE Zone). The PCLE Zone for the smaller area on the west side of the Site appears to be within the Site property boundary (referred to as the “West” PCLE Zone).
  - Based on the January-March 2020 groundwater data, arsenic concentrations that exceeded the RAL in the wells to the north of the Site were delineated to the RAL based on the results of the wells nearby. The arsenic concentrations that exceed the RAL on the west side of the Site are delineated to the RAL in the B-TZ except to the west near MW-38B and MW-62B.



- **C-TZ:**
  - Site-specific VOCs and SVOCs are delineated to applicable RALs in the C-TZ at the Site. The combined PCLE Zone in the C-TZ is in the northeast portion of the Site and extends offsite to the north and to the east onto the City of Houston ROW.
  - Arsenic concentrations were not detected above the RAL in any of the C-TZ wells during the July 2019 or January-March 2020 sampling events.
- **D-TZ:**
  - No site-specific VOC, SVOC, or arsenic concentrations were detected above RALs in D-TZ wells.

## 7.0 ADDITIONAL GROUNDWATER SAMPLING EVENT

An additional groundwater sampling event will be conducted for the monitoring wells installed in February/March 2020 as agreed upon during the meeting with the TCEQ on December 19, 2019 and detailed in the UPRR response letter dated February 7, 2020. The second sampling event is scheduled for late May 2020. Data collected during the additional sampling event will be used to confirm initial results from the March 2020 sampling event.

## 8.0 NOTIFICATIONS

With the additional groundwater data collected in July 2019 and January-March 2020 being submitted to the TCEQ in this IGMR, property owners identified where the groundwater PCLE Zone extends off-site onto property not owned by UPRR and onto the City of Houston will be notified of the availability of information related to the groundwater monitoring at the Site as required under 30 Texas Administrative Code §350.55(b). The list and map showing the off-site property owners that will be notified is provided in Attachment I.

## 9.0 REFERENCES

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**TABLES**

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-01A	47.92	9/2/1993	6.96			40.99
	47.92	12/21/1993	3.28			44.67
	47.92	3/24/1994	3.95			44
	47.92	6/22/1994	5.30			42.65
	47.92	9/28/1994	7.10			40.85
	47.92	10/13/1994	7.26			40.69
	47.92	1/24/1995	2.63			45.32
	47.92	4/11/1995	2.61			45.34
	47.92	7/11/1995	4.78			43.17
	47.92	1/23/1996	5.67			42.28
	47.92	7/19/1996	7.84			40.11
	47.92	9/17/1996	8.33			39.62
	47.92	10/31/1996	6.90			41.05
	47.92	11/22/1996	8.63			39.32
	47.92	12/27/1996	5.50			42.45
	47.92	1/22/1997	3.41			44.54
	47.92	2/21/1997	2.68			45.27
	47.92	3/25/1997	2.96			44.99
	47.92	4/23/1997	4.27			43.68
	47.92	4/24/1997	4.47			43.48
	47.92	5/13/1997	2.91			45.04
	47.92	6/20/1997	4.88			43.07
	47.92	6/25/1997	2.59			45.36
	47.92	7/1/1997	4.04			43.91
	47.92	7/24/1997	6.80			41.15
	47.92	8/16/1997	7.84			40.11
	47.92	8/22/1997	9.52			38.43
	47.92	9/25/1997	6.02			41.93
	47.92	10/22/1997	4.89			43.06
	47.92	11/25/1997	4.88			43.07
	47.92	12/19/1997	4.26			43.69
	47.92	1/20/1998	3.10			44.85
	47.92	3/3/1998	2.87			45.08
	47.92	3/18/1998	2.68			45.27
	47.92	4/24/1998	6.73			41.22
	47.92	5/21/1998	6.89			41.06
	47.92	7/30/1998	7.96			39.99
	47.92	8/25/1998	6.87			41.08
	47.92	9/21/1998	4.70			43.25
	47.92	10/26/1998	5.98			41.97
	47.92	11/23/1998	4.11			43.84
	47.92	1/29/1999	3.01			44.94
	47.92	2/26/1999	3.20			44.75
	47.92	3/16/1999	3.71			44.24
	47.92	4/29/1999	3.93			44.02
	47.92	6/1/1999	3.98			43.97
	47.92	7/30/1999	4.31			43.64
	47.92	8/27/1999	4.11			43.84
	47.92	9/27/1999	9.67			38.28
	47.92	10/29/1999	10.67			37.28
	47.92	12/29/1999	10.00			37.95
	47.92	2/4/2000	12.71			35.24
	47.92	2/25/2000	9.10			38.85
	47.92	3/27/2000	7.38			40.57
	47.92	4/7/2000	7.00			40.95
	47.92	5/31/2000	7.15			40.8
	47.92	6/1/2000	7.00			40.95
	47.92	7/28/2000	7.11			40.84
	47.92	8/30/2000	10.33			37.62
	47.92	9/19/2000	11.56			36.39
	47.92	10/27/2000	9.01			38.94
	47.92	11/21/2000	8.49			39.46
	47.92	5/1/2001	6.60			41.35
	47.92	10/1/2001	6.85			41.1
	47.92	3/11/2002	3.31			44.64
	47.92	9/23/2002	3.23			44.72
	47.92	3/10/2003	2.48			45.44
	47.92	9/23/2003	4.29			43.63

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-01A	47.92	3/15/2004	3.49			44.43
	47.92	9/13/2004	8.26			39.66
	47.92	7/18/2005	3.73			44.19
	47.92	1/4/2006	8.54			39.38
	47.92	7/27/2006	3.10			44.82
	47.92	1/23/2007	2.26			45.66
	47.92	3/7/2007	2.36			45.56
	47.92	7/27/2007	4.05			43.87
	47.92	1/28/2008	2.51			45.41
	47.92	7/16/2008	7.21			40.71
	47.92	1/22/2009	6.21			41.71
	47.92	7/22/2009	6.96			40.96
	47.92	1/8/2010	3.07			44.85
	47.92	7/12/2010	3.87			44.05
	47.88	1/12/2011	3.63			44.25
	47.88	7/13/2011	9.94			37.94
	47.88	1/27/2012	3.19			44.69
	47.88	7/10/2013	9.96			37.92
	47.88	1/8/2014	5.21			42.67
	47.88	7/2/2014	6.81			41.07
	47.88	1/7/2015	2.36			45.52
	47.88	8/10/2015	4.11			43.77
	47.90	1/12/2016	2.49			45.41
	47.90	7/7/2016	5.42			42.48
	47.90	1/12/2017	4.29			43.61
	47.90	7/12/2017	6.19			41.71
	47.90	1/3/2018	6.47			41.43
	47.90	7/18/2018	5.88			42.02
	47.90	1/3/2019	5.96			41.94
	47.90	7/1/2019	2.85			45.05
47.90	1/14/2020	2.71			45.19	
MW-02	47.97	9/2/1993	7.45			40.58
	47.97	12/21/1993	2.58			45.45
	47.97	3/24/1994	4.08			43.95
	47.97	6/22/1994	5.85			42.18
	47.97	9/28/1994	7.05			40.98
	47.97	10/13/1994	7.69			40.34
	47.97	1/24/1995	2.12			45.91
	47.97	4/11/1995	2.53			45.5
	47.97	7/11/1995	5.34			42.69
	47.97	1/23/1996	5.69			42.34
	47.97	7/19/1996	8.28			39.75
	47.97	9/17/1996	8.84			39.19
	47.97	10/31/1996	7.11			40.92
	47.97	11/22/1996	8.99			39.04
	47.97	12/27/1996	5.42			42.61
	47.97	1/22/1997	3.08			44.95
	47.97	2/21/1997	2.60			45.43
	47.97	3/25/1997	2.98			45.05
	47.97	4/23/1997	4.60			43.43
	47.97	4/24/1997	4.78			43.25
	47.97	5/13/1997	2.89			45.14
	47.97	6/20/1997	5.45			42.58
	47.97	6/25/1997	2.59			45.44
	47.97	7/1/1997	4.48			43.55
	47.97	7/24/1997	7.42			40.61
	47.97	8/16/1997	8.42			39.61
	47.97	8/22/1997	9.20			38.83
	47.97	9/25/1997	4.53			43.5
	47.97	10/22/1997	4.95			43.08
	47.97	11/25/1997	4.97			43.06
47.97	12/19/1997	4.33			43.7	
47.97	1/20/1998	3.05			44.98	
47.97	3/3/1998	2.88			45.15	
47.97	3/18/1998	2.66			45.37	
47.97	4/24/1998	7.09			40.94	
47.97	5/21/1998	7.00			41.03	
47.97	7/30/1998	8.11			39.92	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-02	47.97	8/25/1998	7.33			40.7
	47.97	9/21/1998	4.18			43.85
	47.97	10/26/1998	6.85			41.18
	47.97	11/23/1998	4.63			43.4
	47.97	1/29/1999	3.51			44.52
	47.97	2/26/1999	3.61			44.42
	47.97	3/16/1999	3.55			44.48
	47.97	4/29/1999	3.76			44.27
	47.97	6/1/1999	3.76			44.27
	47.97	7/30/1999	4.61			43.42
	47.97	8/27/1999	3.96			44.07
	47.97	9/27/1999	10.12			37.91
	47.97	10/29/1999	11.33			36.7
	47.97	12/29/1999	10.66			37.37
	47.97	2/4/2000	13.19			34.84
	47.97	2/25/2000	9.57			38.46
	47.97	3/27/2000	7.73			40.3
	47.97	4/7/2000	7.30			40.73
	47.97	5/31/2000	7.33			40.7
	47.97	6/1/2000	7.31			40.72
	47.97	7/28/2000	7.35			40.68
	47.97	8/30/2000	10.55			37.48
	47.97	9/19/2000	11.93			36.1
	47.97	10/27/2000	9.04			38.99
	47.97	11/21/2000	8.66			39.37
	47.97	5/1/2001	6.91			41.12
	47.97	10/1/2001	8.22			39.81
	47.97	3/11/2002	3.33			44.7
	47.97	9/23/2002	3.16			44.87
	47.97	3/10/2003	2.54			45.43
	47.97	9/23/2003	3.29			44.68
	47.97	3/15/2004	2.87			45.1
	47.97	9/13/2004	8.71			39.26
	47.97	7/18/2005	2.98			44.99
	47.97	1/4/2006	8.77			39.2
	47.97	7/27/2006	2.87			45.1
	47.97	1/23/2007	2.34			45.63
	47.97	3/7/2007	2.23			45.74
	47.97	7/27/2007	4.40			43.57
	47.97	1/28/2008	2.42			45.55
47.97	7/16/2008	7.72			40.25	
47.97	1/22/2009	6.31			41.66	
47.97	7/22/2009	7.56			40.41	
47.97	1/8/2010	3.91			44.06	
47.97	7/12/2010	4.37			43.6	
48.00	1/12/2011	3.63			44.37	
48.00	7/13/2011	10.28			37.72	
48.00	1/27/2012	2.67			45.33	
48.00	7/10/2013	10.58			37.42	
48.00	1/8/2014	5.47			42.53	
48.00	7/2/2014	7.51			40.49	
48.00	1/7/2015	2.41			45.59	
48.00	8/10/2015	4.96			43.04	
47.89	1/12/2016	2.91			44.98	
47.89	7/7/2016	6.12			41.77	
47.89	1/12/2017	4.62			43.27	
47.89	7/12/2017	6.82			41.07	
47.89	1/3/2018	6.87			41.02	
47.89	7/18/2018	6.62				
47.89	1/3/2019	6.31			41.58	
47.89	7/1/2019	3.09				
47.89	1/14/2020	2.42			45.47	
MW-03	48.34	9/2/1993	8.17			40.17
	48.34	12/21/1993	3.81			44.53
	48.34	3/24/1994	4.74			43.6
	48.34	6/22/1994	6.35			41.99
	48.34	9/28/1994	7.56			40.78
48.34	10/13/1994	8.21			40.13	



**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-03	48.34	1/24/1995	3.18			45.16
	48.34	4/11/1995	3.22			45.12
	48.34	7/11/1995	7.90			40.44
	48.34	1/23/1996	6.27			42.07
	48.34	7/19/1996	8.77			39.57
	48.34	9/17/1996	9.31			39.03
	48.34	10/31/1996	7.61			40.73
	48.34	11/22/1996	9.48			38.86
	48.34	12/27/1996	6.14			42.2
	48.34	1/22/1997	5.68			42.66
	48.34	2/21/1997	3.13			45.21
	48.34	3/25/1997	3.48			44.86
	48.34	4/23/1997	5.17			43.17
	48.34	4/24/1997	5.25			43.09
	48.34	5/13/1997	3.41			44.93
	48.34	6/20/1997	5.91			42.43
	48.34	6/25/1997	3.11			45.23
	48.34	7/11/1997	4.91			43.43
	48.34	7/24/1997	7.90			40.44
	48.34	8/16/1997	8.91			39.43
	48.34	8/22/1997	9.65			38.69
	48.34	9/25/1997	6.96			41.38
	48.34	10/22/1997	5.50			42.84
	48.34	11/25/1997	5.55			42.79
	48.34	12/19/1997	5.10			43.24
	48.34	1/20/1998	3.58			44.76
	48.34	3/3/1998	3.37			44.97
	48.34	3/18/1998	3.16			45.18
	48.34	4/24/1998	7.54			40.8
	48.34	5/21/1998	7.50			40.84
	48.34	7/30/1998	8.44			39.9
	48.34	8/25/1998	7.56			40.78
	48.34	9/21/1998	5.28			43.06
	48.34	10/26/1998	6.96			41.38
	48.34	11/23/1998	5.11			43.23
	48.34	1/29/1999	4.21			44.13
	48.34	2/26/1999	4.32			44.02
	48.34	3/16/1999	4.16			44.18
	48.34	4/29/1999	4.33			44.01
	48.34	6/1/1999	4.39			43.95
	48.34	7/30/1999	5.88			42.46
	48.34	8/27/1999	4.57			43.77
	48.34	9/27/1999	10.48			37.86
	48.34	10/29/1999	11.61			36.73
	48.34	12/29/1999	10.11			38.23
	48.34	2/4/2000	13.22			35.12
	48.34	2/25/2000	9.14			39.2
	48.34	3/27/2000	8.06			40.28
	48.34	4/7/2000	7.64			40.7
	48.34	5/31/2000	7.70			40.64
	48.34	6/1/2000	7.66			40.68
	48.34	7/28/2000	7.71			40.63
	48.34	8/30/2000	10.59			37.75
	48.34	9/19/2000	12.29			36.05
	48.34	10/27/2000	9.09			39.25
	48.34	11/21/2000	9.11			39.23
	48.34	5/1/2001	7.26			41.08
	48.34	10/1/2001	7.57			40.77
	48.34	3/11/2002	7.40			40.94
	48.34	9/23/2002	4.60			43.74
	48.34	3/10/2003	2.89			45.45
	48.34	9/23/2003	3.74			44.6
	48.34	3/15/2004	3.27			45.07
	48.34	9/13/2004	9.03			39.31
	48.34	7/18/2005	3.94			44.4
	48.34	1/4/2006	9.13			39.21
	48.34	7/27/2006	3.30			45.04
	48.34	3/7/2007	2.62			45.72

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-03	48.34	7/27/2007	3.74			44.6
	48.34	1/30/2008	2.85			45.49
	48.34	7/16/2008	7.96			40.38
	48.34	2/4/2009	7.18			41.16
	48.34	7/24/2009	7.63			40.71
	48.34	1/8/2010	5.06			43.28
	48.34	7/12/2010	3.86			44.48
	48.34	1/12/2011	3.71			44.63
	48.34	7/12/2011	6.42			41.92
	48.34	1/26/2012	--			
	48.34	7/9/2012	4.06			44.28
	48.34	1/7/2013	5.09			43.25
	48.34	7/22/2013	8.24			40.1
	48.34	1/7/2014	8.09			40.25
	48.34	7/15/2014	8.78			39.56
	48.34	1/5/2015	7.06			41.28
	48.34	2/11/2018	5.29			43.05
	48.34	3/11/2018	5.72			42.62
	48.34	5/14/2018	5.61			42.73
	48.34	7/2/2018	5.93			42.41
48.34	1/3/2019	5.03			43.31	
48.34	7/9/2019	5.57			42.77	
48.34	1/8/2020	3.06			45.28	
MW-04	49.85	9/2/1993	8.57			41.28
	49.85	12/21/1993	5.42			44.43
	49.85	3/24/1994	5.85			44
	49.85	6/22/1994	6.77			43.08
	49.85	9/28/1994	8.18			41.67
	49.85	10/13/1994	8.93			40.92
	49.85	1/24/1995	4.72			45.13
	49.85	4/11/1995	4.57			45.28
	49.85	7/11/1995	6.47			43.38
	49.85	1/23/1996	7.85			42
	49.85	7/19/1996	9.62			40.23
	49.85	9/17/1996	10.09			39.76
	49.85	10/31/1996	7.93			41.92
	49.85	11/22/1996	10.62			39.23
	49.85	12/27/1996	8.06			41.79
	49.85	1/22/1997	6.07			43.78
	49.85	2/21/1997	4.86			44.99
	49.85	3/25/1997	5.16			44.69
	49.85	4/23/1997	6.25			43.6
	49.85	4/24/1997	6.45			43.4
	49.85	5/13/1997	5.07			44.78
	49.85	6/20/1997	6.69			43.16
	49.85	6/25/1997	4.68			45.17
	49.85	7/1/1997	5.91			43.94
	49.85	7/24/1997	8.61			41.24
	49.85	8/16/1997	9.62			40.23
	49.85	8/22/1997	10.35			39.5
	49.85	9/25/1997	8.13			41.72
	49.85	10/22/1997	7.23			42.62
	49.85	11/25/1997	7.25			42.6
	49.85	12/19/1997	6.76			43.09
	49.85	1/20/1998	5.40			44.45
	49.85	3/3/1998	5.00			44.85
49.85	3/18/1998	4.82			45.03	
49.85	4/24/1998	8.63			41.22	
49.85	5/21/1998	9.30			40.55	
49.85	7/30/1998	10.19			39.66	
49.85	8/25/1998	9.05			40.8	
49.85	9/21/1998	7.05			42.8	
49.85	10/26/1998	8.12			41.73	
49.85	11/23/1998	6.01			43.84	
49.85	1/29/1999	5.19			44.66	
49.85	2/26/1999	5.22			44.63	
49.85	3/16/1999	6.21			43.64	
49.85	4/29/1999	6.33			43.52	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-04	49.85	6/1/1999	6.39			43.46
	49.85	7/30/1999	7.79			42.06
	49.85	8/27/1999	6.51			43.34
	49.85	9/27/1999	11.32			38.53
	49.85	10/29/1999	12.21			37.64
	49.85	12/29/1999	11.52			38.33
	49.85	2/4/2000	14.33			35.52
	49.85	2/25/2000	10.63			39.22
	49.85	3/27/2000	9.38			40.47
	49.85	4/7/2000	9.09			40.76
	49.85	5/31/2000	9.13			40.72
	49.85	6/1/2000	9.10			40.75
	49.85	7/28/2000	9.18			40.67
	49.85	8/30/2000	12.17			37.68
	49.85	9/19/2000	13.39			36.46
	49.85	10/27/2000	10.69			39.16
	49.85	11/21/2000	9.61			40.24
	49.85	5/1/2001	8.41			41.44
	49.85	10/1/2001	8.68			41.17
	49.85	3/11/2002	5.41			44.44
	49.85	9/23/2002	5.29			44.56
	49.85	3/10/2003	4.36			45.49
	49.85	9/23/2003	5.28			44.57
	49.85	3/15/2004	4.80			45.05
	49.85	9/13/2004	9.80			40.05
	49.85	7/18/2005	5.84			44.01
	49.85	1/4/2006	10.48			39.37
	49.85	7/27/2006	5.30			44.55
	49.85	3/7/2007	4.10			45.75
	49.85	7/27/2007	5.36			44.49
	49.85	1/29/2008	4.18			45.67
	49.85	7/16/2008	8.66			41.19
	49.85	2/4/2009	8.93			40.92
	49.85	7/24/2009	9.27			40.58
	49.85	1/8/2010	6.34			43.51
	49.85	7/12/2010	5.02			44.83
	49.85	1/12/2011	5.26			44.59
	49.85	7/12/2011	8.06			41.79
	49.85	1/26/2012	--			
	49.85	7/9/2012	3.74			46.11
49.85	1/7/2013	4.62			45.23	
49.85	7/22/2013	7.59			42.26	
49.85	1/7/2014	7.16			42.69	
49.85	7/15/2014	7.62			42.23	
49.85	1/5/2015	6.12			43.73	
49.85	8/10/2015	4.26			45.59	
49.85	1/13/2016	3.92			45.93	
49.85	7/6/2016	4.31			45.54	
49.85	1/12/2017	4.67			45.18	
49.85	7/6/2017	5.12			44.73	
49.85	9/5/2017	5.01			44.84	
49.85	2/11/2018	5.12			44.73	
49.85	3/11/2018	5.67			44.18	
49.85	5/14/2018	6.06			43.79	
49.85	7/2/2018	6.42			43.43	
49.85	1/3/2019	5.52			44.33	
49.85	7/9/2019	6.02			43.83	
49.85	1/8/2020	4.64			45.21	
MW-05	49.24	9/2/1993	4.90			44.34
	49.24	12/21/1993	2.21			47.03
	49.24	3/24/1994	2.30			46.94
	49.24	6/22/1994	2.80			46.44
	49.24	9/28/1994	3.90			45.34
	49.24	10/13/1994	5.05			44.19
	49.24	1/24/1995	1.36			47.88
	49.24	4/11/1995	3.90			45.34
	49.24	7/11/1995	5.33			43.91
	49.24	1/23/1996	7.42			41.82

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-05	49.24	7/19/1996	8.61			40.63
	49.24	9/17/1996	9.01			40.23
	49.24	10/31/1996	7.84			41.4
	49.24	11/22/1996	9.68			39.56
	49.24	12/27/1996	7.66			41.58
	49.24	1/22/1997	5.89			43.35
	49.24	2/21/1997	4.45			44.79
	49.24	3/25/1997	4.65			44.59
	49.24	4/23/1997	5.53			43.71
	49.24	4/24/1997	5.68			43.56
	49.24	5/13/1997	4.39			44.85
	49.24	6/20/1997	5.67			43.57
	49.24	6/25/1997	3.97			45.27
	49.24	7/1/1997	5.06			44.18
	49.24	7/24/1997	7.46			41.78
	49.24	8/16/1997	8.57			40.67
	49.24	8/22/1997	9.20			40.04
	49.24	9/25/1997	7.28			41.96
	49.24	10/22/1997	6.70			42.54
	49.24	11/25/1997	6.70			42.54
	49.24	12/19/1997	6.26			42.98
	49.24	1/20/1998	5.05			44.19
	49.24	3/4/1998	4.54			44.7
	49.24	3/18/1998	4.36			44.88
	49.24	4/24/1998	7.67			41.57
	49.24	5/21/1998	8.80			40.44
	49.24	7/30/1998	9.90			39.34
	49.24	8/25/1998	8.86			40.38
	49.24	9/21/1998	6.59			42.65
	49.24	10/26/1998	7.77			41.47
	49.24	11/23/1998	5.79			43.45
	49.24	1/29/1999	4.88			44.36
	49.24	2/26/1999	4.96			44.28
	49.24	3/16/1999	5.81			43.43
	49.24	4/29/1999	5.91			43.33
	49.24	6/1/1999	5.99			43.25
	49.24	7/30/1999	7.00			42.24
	49.24	8/27/1999	6.13			43.11
	49.24	9/27/1999	10.17			39.07
	49.24	10/29/1999	11.65			37.59
	49.24	12/29/1999	10.90			38.34
	49.24	2/4/2000	13.77			35.47
	49.24	2/25/2000	9.46			39.78
	49.24	3/27/2000	8.62			40.62
	49.24	4/7/2000	8.20			41.04
	49.24	5/31/2000	8.26			40.98
	49.24	6/1/2000	8.21			41.03
	49.24	7/28/2000	8.26			40.98
	49.24	8/30/2000	11.33			37.91
	49.24	9/19/2000	12.33			36.91
	49.24	10/27/2000	9.94			39.3
	49.24	11/21/2000	9.21			40.03
	49.24	5/1/2001	7.47			41.77
	49.24	10/1/2001	7.79			41.45
	49.24	3/11/2002	4.92			44.32
	49.24	9/23/2002	4.76			44.48
	49.24	3/10/2003	3.77			45.47
	49.24	9/23/2003	4.61			44.63
	49.24	3/15/2004	4.22			45.02
	49.24	9/13/2004	8.58			40.66
	49.24	7/18/2005	5.61			43.63
	49.24	1/4/2006	9.76			39.48
	49.24	7/27/2006	4.85			44.39
	49.24	3/7/2007	5.94			43.3
	49.24	7/27/2007	4.53			44.71
	49.24	1/29/2008	3.71			45.53
	49.24	7/15/2008	7.77			41.47
	49.24	2/4/2009	8.33			40.91

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-05	49.24	7/24/2009	8.67			40.57
	49.24	1/8/2010	6.06			43.18
	49.24	7/12/2010	4.86			44.38
	49.24	1/12/2011	5.06			44.18
	49.24	7/12/2011	10.96			38.28
	49.24	2/2/2012	4.9			44.34
	49.24	7/9/2012	4.61			44.63
	49.24	1/7/2013	7.58			41.66
	49.24	7/22/2013	10.44			38.8
	49.24	1/7/2014	6.92			42.32
	49.24	7/16/2014	8.46			40.78
	49.24	1/5/2015	5.96			43.28
	49.24	8/10/2015	4.13			45.11
	49.24	1/13/2016	3.76			45.48
	49.24	7/7/2016	3.94			45.30
	49.24	1/12/2017	4.31			44.93
	49.24	7/6/2017	4.84			44.40
	49.24	9/5/2017	4.71			44.53
	49.24	2/11/2018	5.56			43.68
	49.24	3/11/2018	5.98			43.26
49.24	5/14/2018	6.57			42.67	
49.24	7/2/2018	6.83			42.41	
49.24	1/3/2019	6.07			43.17	
49.24	7/9/2019	6.56			42.68	
49.24	1/7/2020	4.69			44.55	
MW-07	48.86	9/2/1993	8.09			40.77
	48.86	12/21/1993	4.60			44.26
	48.86	3/24/1994	5.06			43.8
	48.86	6/22/1994	6.03			42.83
	48.86	9/28/1994	7.52			41.34
	48.86	10/13/1994	8.13			40.73
	48.86	1/24/1995	3.81			45.05
	48.86	4/11/1995	3.41			45.45
	48.86	7/11/1995	5.74			43.12
	48.86	1/23/1996	6.99			41.87
	48.86	7/19/1996	8.89			39.97
	48.86	9/17/1996	9.41			39.45
	48.86	10/31/1996	8.04			40.82
	48.86	11/22/1996	9.94			38.92
	48.86	12/27/1996	7.30			41.56
	48.86	1/22/1997	5.25			43.61
	48.86	2/21/1997	4.00			44.86
	48.86	3/25/1997	4.32			44.54
	48.86	4/23/1997	5.51			43.35
	48.86	4/24/1997	5.67			43.19
	48.86	5/13/1997	4.26			44.6
	48.86	6/20/1997	6.00			42.86
	48.86	6/25/1997	3.86			45
	48.86	7/1/1997	5.21			43.65
	48.86	7/24/1997	7.99			40.87
	48.86	8/16/1997	8.92			39.94
	48.86	8/22/1997	9.72			39.14
	48.86	9/25/1997	7.50			41.36
	48.86	10/22/1997	6.48			42.38
	48.86	11/25/1997	6.50			42.36
	48.86	12/19/1997	6.12			42.74
	48.86	1/20/1998	4.52			44.34
	48.86	3/4/1998	4.14			44.72
	48.86	3/18/1998	3.94			44.92
	48.86	4/24/1998	7.85			41.01
	48.86	5/21/1998	8.61			40.25
48.86	7/30/1998	9.54			39.32	
48.86	8/25/1998	8.63			40.23	
48.86	9/21/1998	6.34			42.52	
48.86	10/26/1998	7.56			41.3	
48.86	11/23/1998	5.91			42.95	
48.86	1/29/1999	4.71			44.15	
48.86	2/26/1999	4.76			44.1	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-07	48.86	3/16/1999	5.32			43.54
	48.86	4/29/1999	5.41			43.45
	48.86	6/1/1999	5.49			43.37
	48.86	7/30/1999	6.98			41.88
	48.86	8/27/1999	5.61			43.25
	48.86	9/27/1999	10.64			38.22
	48.86	10/29/1999	11.56			37.3
	48.86	12/29/1999	9.90			38.96
	48.86	2/4/2000	14.21			34.65
	48.86	2/25/2000	8.86			40
	48.86	3/27/2000	8.62			40.24
	48.86	4/7/2000	8.15			40.71
	48.86	5/31/2000	8.21			40.65
	48.86	6/1/2000	8.22			40.64
	48.86	7/28/2000	8.29			40.57
	48.86	8/30/2000	11.55			37.31
	48.86	9/19/2000	12.65			36.21
	48.86	10/27/2000	10.00			38.86
	48.86	11/21/2000	9.46			39.4
	48.86	5/1/2001	7.64			41.22
	48.86	10/1/2001	8.00			40.86
	48.86	3/11/2002	4.56			44.3
	48.86	9/23/2002	4.69			44.17
	48.86	3/10/2003	3.52			45.34
	48.86	9/23/2003	4.70			44.16
	48.86	3/15/2004	3.89			44.97
	48.86	9/13/2004	9.04			39.82
	48.86	7/18/2005	5.27			43.59
	48.86	1/4/2006	9.91			38.95
	48.86	7/27/2006	4.60			44.26
	48.86	1/23/2007	3.46			45.4
	48.86	3/7/2007	3.82			45.04
	48.86	7/27/2007	4.94			43.92
	48.86	1/29/2008	3.39			45.47
	48.86	7/16/2008	7.94			40.92
	48.86	1/22/2009	7.49			41.37
	48.86	7/24/2009	NM			
	48.86	1/8/2010	4.02			44.84
	48.86	7/12/2010	4.72			44.14
	48.92	1/12/2011	4.56			44.36
	48.92	7/12/2011	10.91			38.01
	48.92	1/27/2012	3.86			45.06
48.92	7/10/2013	10.62			38.30	
48.92	1/8/2014	6.42			42.50	
48.92	7/3/2014	7.61			41.31	
48.92	1/7/2015	3.46			45.46	
48.92	8/10/2015	5.01			43.91	
48.91	1/12/2016	3.09			45.82	
48.91	7/7/2016	6.72			42.19	
48.91	1/12/2017	5.81			43.10	
48.91	7/12/2017	7.71			41.20	
48.91	1/3/2018	7.87			41.04	
48.91	7/19/2018	6.93			41.98	
48.91	1/3/2019	7.32			41.59	
48.91	7/1/2019	3.93			44.98	
48.91	1/13/2020	4.29			44.62	
MW-08	49.33	9/2/1993	8.18			41.19
	49.33	12/21/1993	5.02			44.35
	49.33	3/24/1994	5.53			43.84
	49.33	6/22/1994	6.38			42.99
	49.33	9/28/1994	7.72			41.65
	49.33	10/13/1994	8.43			40.94
	49.33	1/24/1995	4.15			45.22
	49.33	4/11/1995	4.02			45.35
	49.33	7/11/1995	5.95			43.42
	49.33	1/23/1996	7.20			42.17
	49.33	7/19/1996	9.06			40.31
	49.33	9/17/1996	9.51			39.86



**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-08	49.33	10/31/1996	7.99			41.38
	49.33	11/22/1996	9.98			39.39
	49.33	12/27/1996	7.24			42.13
	49.33	1/22/1997	5.25			44.12
	49.33	2/21/1997	4.21			45.16
	49.33	3/25/1997	4.48			44.89
	49.33	4/23/1997	5.61			43.76
	49.33	4/24/1997	5.76			43.61
	49.33	5/13/1997	4.45			44.92
	49.33	6/20/1997	6.09			43.28
	49.33	6/25/1997	4.56			44.81
	49.33	7/1/1997	5.06			44.31
	49.33	7/24/1997	7.97			41.4
	49.33	8/16/1997	8.05			41.32
	49.33	8/22/1997	9.73			39.64
	49.33	9/25/1997	7.57			41.8
	49.33	10/22/1997	6.43			42.94
	49.33	11/25/1997	6.48			42.89
	49.33	12/19/1997	5.22			44.15
	49.33	1/20/1998	4.70			44.67
	49.33	3/4/1998	4.38			44.99
	49.33	3/18/1998	4.18			45.19
	49.33	4/24/1998	8.00			41.37
	49.33	5/21/1998	8.45			40.92
	49.33	7/30/1998	9.33			40.04
	49.33	8/25/1998	8.46			40.91
	49.33	9/21/1998	6.31			43.06
	49.33	10/26/1998	7.66			41.71
	49.33	11/23/1998	5.96			43.41
	49.33	1/29/1999	4.80			44.57
	49.33	2/26/1999	4.89			44.48
	49.33	3/16/1999	5.45			43.92
	49.33	4/29/1999	5.66			43.71
	49.33	6/1/1999	5.66			43.71
	49.33	7/30/1999	7.20			42.17
	49.33	8/27/1999	5.85			43.52
	49.33	9/27/1999	10.78			38.59
	49.33	10/29/1999	11.76			37.61
	49.33	12/29/1999	11.03			38.34
	49.33	2/4/2000	14.66			34.71
	49.33	2/25/2000	10.33			39.04
	49.33	3/27/2000	8.75			40.62
	49.33	4/7/2000	8.37			41
	49.33	5/31/2000	8.40			40.97
	49.33	6/1/2000	8.36			41.01
	49.33	7/28/2000	8.40			40.97
	49.33	8/30/2000	11.29			38.08
	49.33	9/19/2000	12.82			36.55
	49.33	10/27/2000	12.63			36.74
	49.33	11/21/2000	9.64			39.73
	49.33	5/1/2001	7.83			41.54
	49.33	10/1/2001	8.05			41.32
	49.33	3/11/2002	4.75			44.62
	49.33	9/23/2002	4.69			44.68
	49.33	3/10/2003	3.84			45.49
	49.33	9/23/2003	4.73			44.6
	49.33	3/15/2004	4.31			45.02
	49.33	9/13/2004	9.31			40.02
	49.33	7/18/2005	5.32			44.01
	49.33	1/4/2006	10.63			38.7
	49.33	7/27/2006	4.79			44.54
	49.33	1/22/2007	3.81			45.52
	49.33	3/7/2007	3.96			45.37
	49.33	7/27/2007	5.06			44.27
	49.33	1/29/2008	3.71			45.62
	49.33	7/16/2008	8.32			41.01
	49.33	1/22/2009	7.71			41.62
	49.33	7/24/2009	NM			

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-08	49.33	1/8/2010	4.17			45.16
	49.33	7/12/2010	4.96			44.37
	49.33	1/12/2011	5.32			44.01
	49.33	7/12/2011	11.24			38.09
	49.33	1/27/2012	4.68			44.65
	49.33	7/10/2013	11.07			38.26
	49.33	1/8/2014	6.87			42.46
	49.33	7/3/2014	8.16			41.17
	49.33	1/7/2015	3.82			45.51
	49.33	8/10/2015	5.06			44.27
	49.33	1/12/2016	3.87			45.46
	49.33	7/7/2016	6.44			42.89
	49.33	1/12/2017	5.82			43.51
	49.33	7/12/2017	7.92			41.41
	49.33	1/3/2018	8.02			41.31
	49.33	7/19/2018	7.22			42.11
	49.33	1/3/2019	7.52			41.81
	49.33	7/1/2019	3.98			45.35
49.33	1/13/2020	4.45			44.88	
MW-09	49.26	9/2/1993	7.43			41.86
	49.26	12/21/1993	4.89			44.4
	49.26	3/24/1994	4.92			44.37
	49.26	6/22/1994	5.51			43.78
	49.26	9/28/1994	6.90			42.39
	49.26	10/13/1994	7.66			41.63
	49.26	1/24/1995	4.10			45.19
	49.26	4/11/1995	3.74			45.55
	49.26	7/11/1995	5.08			44.21
	49.26	1/23/1996	7.09			42.2
	49.26	7/19/1996	8.27			41.02
	49.26	9/17/1996	8.58			40.71
	49.26	10/31/1996	7.27			42.02
	49.26	11/22/1996	9.17			40.12
	49.26	12/27/1996	7.05			42.24
	49.26	1/22/1997	5.42			43.87
	49.26	2/21/1997	4.09			45.2
	49.26	3/25/1997	4.17			45.12
	49.26	4/23/1997	5.05			44.24
	49.26	4/24/1997	5.21			44.08
	49.26	5/13/1997	4.16			45.13
	49.26	6/20/1997	5.32			43.97
	49.26	6/25/1997	3.80			45.49
	49.26	7/1/1997	4.57			44.72
	49.26	7/24/1997	7.03			42.26
	49.26	8/16/1997	8.26			41.03
	49.26	8/22/1997	8.67			40.62
	49.26	9/25/1997	6.99			42.3
	49.26	10/22/1997	6.10			43.19
	49.26	11/25/1997	6.12			43.17
	49.26	12/19/1997	5.62			43.67
	49.26	1/20/1998	4.60			44.69
	49.26	3/4/1998	4.15			45.14
	49.26	3/18/1998	4.02			45.27
	49.26	4/24/1998	7.32			41.97
	49.26	5/21/1998	8.10			41.19
	49.26	7/30/1998	9.12			40.17
	49.26	8/25/1998	8.41			40.88
	49.26	9/21/1998	6.11			43.18
	49.26	10/26/1998	7.61			41.68
49.26	11/23/1998	5.43			43.86	
49.26	1/29/1999	4.60			44.69	
49.26	2/26/1999	4.68			44.61	
49.26	3/16/1999	5.46			43.83	
49.26	4/29/1999	5.66			43.63	
49.26	6/1/1999	5.66			43.63	
49.26	7/30/1999	7.11			42.18	
49.26	8/27/1999	5.86			43.43	
49.26	9/27/1999	9.81			39.48	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-09	49.26	10/29/1999	10.63			38.66
	49.26	12/29/1999	9.99			39.3
	49.26	2/4/2000	12.44			36.85
	49.26	2/25/2000	8.88			40.41
	49.26	3/27/2000	8.22			41.07
	49.26	4/7/2000	8.10			41.19
	49.26	5/31/2000	8.15			41.14
	49.26	6/1/2000	8.00			41.29
	49.26	7/28/2000	8.11			41.18
	49.26	8/30/2000	11.10			38.19
	49.26	9/19/2000	11.91			37.38
	49.26	10/27/2000	9.84			39.45
	49.26	11/21/2000	8.89			40.4
	49.26	5/1/2001	7.16			42.13
	49.26	10/1/2001	7.39			41.9
	49.26	3/11/2002	4.61			44.68
	49.26	9/23/2002	4.45			44.84
	49.26	3/10/2003	3.59			45.67
	49.26	9/23/2003	4.31			44.95
	49.26	3/15/2004	4.18			45.08
	49.26	9/13/2004	8.39			40.87
	49.26	7/18/2005	5.53			43.73
	49.26	1/4/2006	9.46			39.8
	49.26	7/27/2006	4.85			44.41
	49.26	3/7/2007	5.58			43.68
	49.26	7/27/2007	3.78			45.48
	49.26	1/29/2008	3.52			45.74
	49.26	7/15/2008	7.04			42.22
	49.26	2/4/2009	8.01			41.25
	49.26	7/24/2009	8.34			40.92
	49.26	1/8/2010	5.89			43.37
	49.26	7/12/2010	4.32			44.94
	49.26	1/12/2011	4.61			44.65
	49.26	7/12/2011	10.71			38.55
	49.26	1/26/2012	4.73			44.53
	49.26	7/9/2012	4.23			45.03
	49.26	1/7/2013	6.73			42.53
	49.26	7/22/2013	9.16			40.1
	49.26	1/7/2014	8.72			40.54
	49.26	7/16/2014	8.17			41.09
49.26	1/5/2015	8.01			41.25	
49.26	8/10/2015	6.17			43.09	
49.26	1/13/2016	5.81			43.45	
49.26	7/6/2016	6.14			43.12	
49.26	1/12/2017	6.71			42.55	
49.26	7/6/2017	7.09			42.17	
49.26	9/5/2017	7.06			42.20	
49.26	2/11/2018	5.16			44.10	
49.26	3/11/2018	6.01			43.25	
49.26	5/14/2018	6.21			43.05	
49.26	7/2/2018	6.67			42.59	
49.26	1/3/2019	5.61			43.65	
49.26	7/9/2019	6.21			43.05	
49.26	1/7/2020	4.11			45.15	
MW-10A	49.86	9/28/1994	8.69			41.21
	49.86	10/13/1994	9.36			40.54
	49.86	1/24/1995	4.62			45.28
	49.86	4/11/1995	4.60			45.3
	49.86	7/11/1995	7.00			42.9
	49.86	1/23/1996	7.74			42.16
	49.86	7/19/1996	9.98			39.92
	49.86	9/17/1996	10.54			39.36
	49.86	10/31/1996	7.94			41.96
	49.86	11/22/1996	10.82			39.08
	49.86	12/27/1996	7.81			42.09
	49.86	1/22/1997	5.45			44.45
	49.86	2/21/1997	4.63			45.27
	49.86	3/25/1997	5.01			44.89

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-10A	49.86	4/23/1997	6.39			43.51
	49.86	4/24/1997	6.58			43.32
	49.86	5/13/1997	4.93			44.97
	49.86	6/20/1997	7.08			42.82
	49.86	6/25/1997	4.58			45.32
	49.86	7/1/1997	6.13			43.77
	49.86	7/24/1997	9.11			40.79
	49.86	8/16/1997	10.10			39.8
	49.86	8/22/1997	10.81			39.09
	49.86	9/25/1997	8.47			41.43
	49.86	10/22/1997	7.02			42.88
	49.86	11/25/1997	7.05			42.85
	49.86	12/19/1997	6.89			43.01
	49.86	1/20/1998	5.10			44.8
	49.86	3/3/1998	4.87			45.03
	49.86	3/18/1998	4.65			45.25
	49.86	4/24/1998	8.84			41.06
	49.86	5/21/1998	9.10			40.8
	49.86	7/30/1998	10.23			39.67
	49.86	8/25/1998	9.11			40.79
	49.86	9/21/1998	6.82			43.08
	49.86	10/26/1998	8.19			41.71
	49.86	11/23/1998	6.12			43.78
	49.86	1/29/1999	5.61			44.29
	49.86	2/26/1999	5.69			44.21
	49.86	3/16/1999	5.91			43.99
	49.86	4/29/1999	6.11			43.79
	49.86	6/1/1999	6.10			43.8
	49.86	7/30/1999	7.70			42.2
	49.86	8/27/1999	6.31			43.59
	49.86	9/27/1999	11.73			38.17
	49.86	10/29/1999	12.69			37.21
	49.86	12/29/1999	12.00			37.9
	49.86	2/4/2000	14.30			35.6
	49.86	2/25/2000	11.44			38.46
	49.86	3/27/2000	9.57			40.33
	49.86	4/7/2000	9.27			40.63
	49.86	5/31/2000	9.31			40.59
	49.86	6/1/2000	9.10			40.8
	49.86	7/28/2000	9.30			40.6
	49.86	8/30/2000	12.09			37.81
	49.86	9/19/2000	13.70			36.2
	49.86	10/27/2000	10.69			39.21
	49.86	11/21/2000	10.49			39.41
	49.86	5/1/2001	8.64			41.26
	49.86	10/1/2001	8.93			40.97
	49.86	3/11/2002	5.30			44.6
	49.86	9/23/2002	5.19			44.71
	49.86	3/10/2003	4.43			45.43
	49.86	9/23/2003	5.31			44.55
	49.86	3/15/2004	4.69			45.17
	49.86	9/13/2004	10.30			39.56
	49.86	7/18/2005	5.57			44.29
	49.86	1/4/2006	9.68			40.18
	49.86	7/27/2006	5.01			44.85
	49.86	1/23/2007	4.29			45.57
	49.86	3/7/2007	4.13			45.73
	49.86	7/27/2007	6.03			43.83
	49.86	1/28/2008	4.22			45.64
	49.86	7/16/2008	9.31			40.55
	49.86	1/22/2009	8.27			41.59
	49.86	7/24/2009	NM			
	49.86	1/8/2010	4.64			45.22
	49.86	7/12/2010	5.23			44.63
	49.82	1/12/2011	5.72			44.10
	49.82	7/12/2011	12.07			37.75
	49.82	7/13/2011	11.96			37.86
	49.82	1/27/2012	4.88			44.94

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-10A	49.82	7/10/2013	12.07			37.75
	49.82	1/8/2014	7.33			42.49
	49.82	7/2/2014	8.92			40.90
	49.82	1/7/2015	4.26			45.56
	49.82	8/10/2015	6.02			43.80
	49.83	1/12/2016	4.41			45.42
	49.83	7/7/2016	7.36			42.47
	49.83	1/12/2017	6.69			43.14
	49.83	7/12/2017	8.23			41.60
	49.83	1/3/2018	8.63			41.20
	49.83	7/18/2018	7.97			41.86
	49.83	1/3/2019	8.09			41.74
	49.83	7/1/2019	4.69			45.14
	49.83	1/14/2020	4.68			45.15
MW-10B	49.94	9/28/1994	8.77			41.2
	49.94	10/13/1994	9.45			40.52
	49.94	1/24/1995	4.72			45.25
	49.94	4/11/1995	4.72			45.25
	49.94	7/11/1995	7.13			42.84
	49.94	1/23/1996	7.84			42.13
	49.94	7/19/1996	10.27			39.7
	49.94	9/17/1996	10.64			39.33
	49.94	10/31/1996	8.01			41.96
	49.94	11/22/1996	10.93			39.04
	49.94	12/27/1996	7.99			41.98
	49.94	1/22/1997	5.72			44.25
	49.94	2/21/1997	4.78			45.19
	49.94	3/25/1997	5.13			44.84
	49.94	4/23/1997	6.52			43.45
	49.94	4/24/1997	6.71			43.26
	49.94	5/13/1997	5.09			44.88
	49.94	6/20/1997	7.21			42.76
	49.94	6/25/1997	4.71			45.26
	49.94	7/1/1997	6.27			43.7
	49.94	7/24/1997	9.15			40.82
	49.94	8/16/1997	10.19			39.78
	49.94	8/22/1997	10.92			39.05
	49.94	9/25/1997	8.69			41.28
	49.94	10/22/1997	7.18			42.79
	49.94	11/25/1997	7.21			42.76
	49.94	12/19/1997	6.56			43.41
	49.94	1/20/1998	5.25			44.72
	49.94	3/3/1998	5.00			44.97
	49.94	3/18/1998	4.79			45.18
	49.94	4/24/1998	8.95			41.02
	49.94	5/21/1998	9.30			40.67
	49.94	7/30/1998	10.30			39.67
	49.94	8/25/1998	9.20			40.77
	49.94	9/21/1998	7.06			42.91
	49.94	10/26/1998	8.31			41.66
	49.94	11/23/1998	6.25			43.72
	49.94	1/29/1999	5.71			44.26
	49.94	2/26/1999	5.76			44.21
	49.94	3/16/1999	6.05			43.92
	49.94	4/29/1999	6.10			43.87
49.94	6/1/1999	6.10			43.87	
49.94	7/30/1999	7.61			42.36	
49.94	8/27/1999	6.33			43.64	
49.94	9/27/1999	11.90			38.07	
49.94	10/29/1999	12.60			37.37	
49.94	12/29/1999	12.10			37.87	
49.94	2/4/2000	14.29			35.68	
49.94	2/25/2000	11.15			38.82	
49.94	3/27/2000	9.67			40.3	
49.94	4/7/2000	9.32			40.65	
49.94	5/31/2000	9.38			40.59	
49.94	6/1/2000	9.21			40.76	
49.94	7/28/2000	9.33			40.64	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-10B	49.94	8/30/2000	12.11			37.86
	49.94	9/19/2000	13.77			36.2
	49.94	10/27/2000	10.63			39.34
	49.94	11/21/2000	10.64			39.33
	49.94	5/1/2001	8.75			41.22
	49.94	10/1/2001	9.12			40.85
	49.94	3/11/2002	5.47			44.5
	49.94	9/23/2002	5.40			44.57
	49.94	3/10/2003	4.59			45.35
	49.94	9/23/2003	5.58			44.36
	49.94	3/15/2004	5.78			44.16
	49.94	9/13/2004	10.41			39.53
	49.94	7/18/2005	5.97			43.97
	49.94	1/4/2006	10.75			39.19
	49.94	7/27/2006	5.73			44.21
	49.94	1/23/2007	4.45			45.49
	49.94	3/7/2007	4.61			45.33
	49.94	7/27/2007	6.15			43.79
	49.94	1/28/2008	4.44			45.5
	49.94	7/16/2008	9.42			40.52
	49.94	1/22/2009	8.39			41.55
	49.94	7/24/2009	NM			
	49.94	1/8/2010	4.91			45.03
	49.94	7/12/2010	5.33			44.61
	49.95	1/12/2011	5.96			43.99
	49.95	7/13/2011	12.07			37.88
	49.95	1/27/2012	5.02			44.93
	49.95	7/10/2013	12.18			37.77
	49.95	1/8/2014	7.46			42.49
	49.95	7/2/2014	8.96			40.99
	49.95	1/7/2015	4.46			45.49
	49.95	8/10/2015	6.14			43.81
49.96	1/12/2016	4.64			45.32	
49.96	7/7/2016	7.62			42.34	
49.96	1/12/2017	6.57			43.39	
49.96	7/12/2017	8.33			41.63	
49.96	1/3/2018	8.71			41.25	
49.96	7/18/2018	8.04			41.92	
49.96	1/3/2019	8.16			41.80	
49.96	7/1/2019	4.82			45.14	
49.96	1/14/2020	4.92			45.04	
MW-11A	50.05	9/28/1994	8.66			41.38
	50.05	10/13/1994	9.35			40.69
	50.05	1/24/1995	4.88			45.16
	50.05	4/11/1995	4.81			45.23
	50.05	7/11/1995	6.67			43.37
	50.05	1/23/1996	8.01			42.03
	50.05	7/19/1996	10.09			39.95
	50.05	9/17/1996	10.56			39.48
	50.05	10/31/1996	8.16			41.88
	50.05	11/22/1996	10.98			39.06
	50.05	12/27/1996	8.21			41.83
	50.05	1/22/1997	6.06			43.98
	50.05	2/21/1997	4.98			45.06
	50.05	3/25/1997	5.32			44.72
	50.05	4/23/1997	6.59			43.45
	50.05	4/24/1997	6.77			43.27
	50.05	5/13/1997	5.31			44.73
	50.05	6/20/1997	7.15			42.89
	50.05	6/25/1997	4.88			45.16
	50.05	7/1/1997	6.29			43.75
	50.05	7/24/1997	9.12			40.92
	50.05	8/16/1997	10.11			39.93
50.05	8/22/1997	10.82			39.22	
50.05	9/25/1997	8.70			41.34	
50.05	10/22/1997	7.40			42.64	
50.05	11/25/1997	7.41			42.63	
50.05	12/19/1997	6.10			43.94	



**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-11A	50.05	1/20/1998	5.49			44.55
	50.05	3/3/1998	5.16			44.88
	50.05	3/18/1998	4.96			45.08
	50.05	4/24/1998	8.98			41.06
	50.05	5/21/1998	9.40			40.64
	50.05	7/30/1998	10.56			39.48
	50.05	8/25/1998	9.32			40.72
	50.05	9/21/1998	7.28			42.76
	50.05	10/26/1998	8.43			41.61
	50.05	11/23/1998	6.41			43.63
	50.05	1/29/1999	5.31			44.73
	50.05	2/26/1999	5.39			44.65
	50.05	3/16/1999	6.32			43.72
	50.05	4/29/1999	6.51			43.53
	50.05	6/1/1999	6.57			43.47
	50.05	7/30/1999	8.00			42.04
	50.05	8/27/1999	6.79			43.25
	50.05	9/27/1999	11.73			38.31
	50.05	10/29/1999	12.81			37.23
	50.05	12/29/1999	12.11			37.93
	50.05	2/4/2000	14.33			35.71
	50.05	2/25/2000	11.10			38.94
	50.05	3/27/2000	9.66			40.38
	50.05	4/7/2000	9.40			40.64
	50.05	5/31/2000	9.50			40.54
	50.05	6/1/2000	9.30			40.74
	50.05	7/28/2000	9.47			40.57
	50.05	8/30/2000	12.44			37.6
	50.05	9/19/2000	13.74			36.3
	50.05	10/27/2000	11.01			39.03
	50.05	11/21/2000	10.69			39.35
	50.05	5/1/2001	8.78			41.26
	50.05	10/1/2001	9.12			40.93
	50.05	3/11/2002	5.59			44.45
	50.05	9/23/2002	5.60			44.44
	50.05	3/10/2003	4.66			45.39
	50.05	9/23/2003	5.73			44.32
	50.05	3/15/2004	4.99			45.06
	50.05	9/13/2004	10.28			39.77
	50.05	7/18/2005	6.66			43.39
	50.05	1/5/2006	10.85			39.2
	50.05	7/27/2006	5.02			45.03
	50.05	1/23/2007	4.54			45.51
	50.05	3/7/2007	4.26			45.79
	50.05	7/27/2007	6.09			43.96
	50.05	1/28/2008	4.46			45.59
	50.05	7/16/2008	9.25			40.8
	50.05	1/22/2009	8.57			41.48
	50.05	7/24/2009	NM			
	50.05	1/8/2010	4.97			45.08
	50.05	7/12/2010	5.51			44.54
	50.07	1/12/2011	6.21			43.86
	50.07	7/12/2011	12.02			38.05
	50.07	1/27/2012	5.31			44.76
	50.07	7/10/2013	12.01			38.06
	50.07	1/8/2014	7.46			42.61
	50.07	7/2/2014	9.02			41.05
	50.07	1/7/2015	4.58			45.49
	50.07	8/10/2015	6.11			43.96
	50.16	1/12/2016	4.71			45.45
	50.16	7/7/2016	7.61			42.55
	50.16	1/12/2017	8.47			41.69
	50.16	7/12/2017	8.46			41.70
	50.16	1/3/2018	8.94			41.22
	50.16	7/18/2018	8.09			42.07
	50.16	1/3/2019	8.48			41.68
	50.16	7/1/2019	5.06			45.10
	50.16	1/14/2020	5.11			45.05

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-11B	50.18	9/28/1994	8.92			41.27
	50.18	10/13/1994	9.59			40.6
	50.18	1/24/1995	5.04			45.15
	50.18	4/11/1995	5.01			45.18
	50.18	7/11/1995	7.23			42.96
	50.18	1/23/1996	8.20			41.99
	50.18	7/19/1996	8.92			41.27
	50.18	9/17/1996	10.83			39.36
	50.18	10/31/1996	9.34			40.85
	50.18	11/22/1996	11.23			38.96
	50.18	12/27/1996	8.45			41.74
	50.18	1/22/1997	6.28			43.91
	50.18	2/21/1997	5.16			45.03
	50.18	3/25/1997	5.51			44.68
	50.18	4/23/1997	6.81			43.38
	50.18	4/24/1997	6.99			43.2
	50.18	5/13/1997	5.46			44.73
	50.18	6/20/1997	7.40			42.79
	50.18	6/25/1997	5.06			45.13
	50.18	7/1/1997	6.52			43.67
	50.18	7/24/1997	9.36			40.83
	50.18	8/16/1997	10.36			39.83
	50.18	8/22/1997	11.11			39.08
	50.18	9/25/1997	8.96			41.23
	50.18	10/22/1997	7.61			42.58
	50.18	11/25/1997	7.63			42.56
	50.18	12/19/1997	7.11			43.08
	50.18	1/20/1998	5.70			44.49
	50.18	3/3/1998	5.35			44.84
	50.18	3/18/1998	5.14			45.05
	50.18	4/24/1998	9.19			41
	50.18	5/21/1998	9.61			40.58
	50.18	7/30/1998	10.72			39.47
	50.18	8/25/1998	9.48			40.71
	50.18	9/21/1998	7.49			42.7
	50.18	10/26/1998	8.57			41.62
	50.18	11/23/1998	6.32			43.87
	50.18	2/26/1999	5.32			44.87
	50.18	3/16/1999	6.49			43.7
	50.18	4/29/1999	6.66			43.53
	50.18	6/1/1999	6.66			43.53
	50.18	7/30/1999	8.12			42.07
	50.18	8/27/1999	6.88			43.31
	50.18	9/27/1999	12.04			38.15
	50.18	10/29/1999	13.00			37.19
	50.18	12/29/1999	12.33			37.86
	50.18	2/4/2000	15.61			34.58
	50.18	2/25/2000	11.49			38.7
	50.18	3/27/2000	9.93			40.26
	50.18	4/7/2000	9.54			40.65
	50.18	5/31/2000	9.61			40.58
	50.18	6/1/2000	9.51			40.68
	50.18	7/28/2000	9.60			40.59
	50.18	8/30/2000	12.76			37.43
	50.18	9/19/2000	13.97			36.22
	50.18	10/27/2000	11.23			38.96
	50.18	11/21/2000	10.88			39.31
	50.18	5/1/2001	5.97			44.22
	50.18	10/1/2001	9.33			40.86
	50.18	3/1/2002	5.80			44.39
	50.18	9/23/2002	5.79			44.4
	50.18	3/10/2003	4.85			45.33
	50.18	9/23/2003	5.95			44.23
	50.18	3/15/2004	5.16			45.02
	50.18	9/13/2004	10.53			39.65
	50.18	7/18/2005	5.45			44.73
	50.18	1/4/2006	11.01			39.17
	50.18	7/27/2006	5.26			44.92

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-11B	50.18	1/23/2007	4.13			46.05
	50.18	3/7/2007	4.42			45.76
	50.18	7/27/2007	6.29			43.89
	50.18	1/28/2008	4.69			45.49
	50.18	7/16/2008	9.49			40.69
	50.18	1/22/2009	8.72			41.46
	50.18	7/24/2009	NM			
	50.18	1/8/2010	5.15			45.03
	50.18	7/12/2010	5.67			44.51
	50.23	1/12/2011	6.37			43.86
	50.23	7/12/2011	12.23			38.00
	50.23	1/27/2012	5.38			44.85
	50.23	7/10/2013	12.22			38.01
	50.23	1/8/2014	7.82			42.41
	50.23	7/2/2014	9.14			41.09
	50.23	1/7/2015	4.79			45.44
	50.23	8/10/2015	6.27			43.96
	50.24	1/12/2016	4.99			45.25
	50.24	7/7/2016	7.59			42.65
	50.24	1/12/2017	8.54			41.70
	50.24	7/12/2017	8.49			41.75
	50.24	1/3/2018	9.04			41.20
	50.24	7/18/2018	8.24			42.00
	50.24	1/3/2019	8.57			41.67
50.24	7/1/2019	5.21			45.03	
50.24	1/14/2020	5.3			44.94	
MW-12A	49.96	3/25/1997	5.52			44.44
	49.96	4/23/1997	6.51			43.45
	49.96	4/24/1997	6.66			43.3
	49.96	5/13/1997	5.47			44.49
	49.96	6/20/1997	6.81			43.15
	49.96	9/25/1997	8.08			41.88
	49.96	10/22/1997	7.10			42.86
	49.96	11/25/1997	7.12			42.84
	49.96	12/19/1997	6.96			43
	49.96	1/20/1998	5.69			44.27
	49.96	3/4/1998	4.52			45.44
	49.96	3/18/1998	5.28			44.68
	49.96	4/24/1998	8.70			41.26
	49.96	5/21/1998	9.10			40.86
	49.96	8/25/1998	10.05			39.91
	49.96	9/21/1998	7.11			42.85
	49.96	10/26/1998	9.11			40.85
	49.96	11/23/1998	6.01			43.95
	49.96	1/29/1999	5.44			44.52
	49.96	2/26/1999	5.52			44.44
	49.96	3/16/1999	6.21			43.75
	49.96	4/29/1999	6.38			43.58
	49.96	6/1/1999	6.31			43.65
	49.96	7/30/1999	7.88			42.08
	49.96	8/27/1999	6.56			43.4
	49.96	9/27/1999	11.61			38.35
	49.96	10/29/1999	12.79			37.17
	49.96	11/18/1999	13.18			36.78
	49.96	12/29/1999	12.03			37.93
	49.96	2/4/2000	15.43			34.53
	49.96	2/25/2000	11.34			38.62
	49.96	3/27/2000	9.22			40.74
	49.96	4/7/2000	8.80			41.16
49.96	5/31/2000	8.84			41.12	
49.96	6/1/2000	8.81			41.15	
49.96	7/28/2000	8.87			41.09	
49.96	8/30/2000	11.76			38.2	
49.96	9/19/2000	13.22			36.74	
49.96	10/27/2000	10.54			39.42	
49.96	11/21/2000	10.16			39.8	
49.96	5/1/2001	8.60			41.36	
49.96	10/1/2001	8.73			41.23	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-12A	49.96	3/11/2002	6.01			43.95
	49.96	9/23/2002	5.87			44.09
	49.96	3/10/2003	5.37			44.59
	49.96	9/23/2003	5.96			44
	49.96	3/15/2004	5.54			44.42
	49.96	9/13/2004	10.30			39.66
	49.96	7/18/2005	7.01			42.95
	49.96	1/4/2006	10.57			39.39
	49.96	7/27/2006	6.60			43.36
	49.96	3/7/2007	6.94			43.02
	49.96	7/27/2007	5.79			44.17
	49.96	1/30/2008	5.29			44.67
	49.96	7/15/2008	9.19			40.77
	49.96	2/4/2009	8.81			41.15
	49.96	7/24/2009	9.13			40.83
	49.96	1/8/2010	5.47			44.49
	49.96	7/12/2010	9.72			40.24
	49.96	1/12/2011	5.59			44.37
	49.96	7/12/2011	12.46			37.5
	49.96	1/26/2012	5.78			44.18
	49.96	7/9/2012	5.96			44
	49.96	1/7/2013	9.04			40.92
	49.96	7/22/2013	11.64			38.32
	49.96	1/7/2014	7.38			42.58
	49.96	7/16/2014	9.82			40.14
	49.96	1/5/2015	6.46			43.50
	49.96	8/10/2015	5.26			44.70
	49.96	1/13/2016	4.67			45.29
	49.96	7/6/2016	4.96			45.00
	49.96	1/12/2017	5.67			44.29
49.96	7/6/2017	6.03			43.93	
49.96	9/5/2017	5.86			44.10	
49.96	2/11/2018	6.48			43.48	
49.96	3/11/2018	7.12			42.84	
49.96	5/14/2018	8.92			41.04	
49.96	1/3/2019	8.37			41.59	
49.96	7/9/2019	8.02			41.94	
49.96	1/7/2020	6.04			43.92	
MW-12B	50.02	3/25/1997	5.60			44.42
	50.02	4/23/1997	6.64			43.38
	50.02	4/24/1997	6.74			43.28
	50.02	5/13/1997	5.55			44.47
	50.02	6/20/1997	7.01			43.01
	50.02	9/25/1997	8.32			41.7
	50.02	10/22/1997	7.25			42.77
	50.02	11/25/1997	7.29			42.73
	50.02	12/19/1997	6.86			43.16
	50.02	1/20/1998	5.88			44.14
	50.02	3/4/1998	5.64	44.08	1.72	44.38
	50.02	3/18/1998	5.38	44.07	1.73	44.64
	50.02	4/9/1998	7.87		0.98	42.15
	50.02	4/16/1998	8.31		1.35	41.71
	50.02	4/24/1998	8.72	43.82	1.98	41.3
	50.02	5/8/1998	NM		0.50	
	50.02	5/12/1998	NM		0.50	
	50.02	5/21/1998	10.48			39.54
	50.02	5/25/1998	NM		1.00	
	50.02	6/9/1998	NM		1.00	
	50.02	6/16/1998	NM		1.20	
	50.02	6/26/1998	NM		1.50	
	50.02	7/2/1998	NM		1.50	
	50.02	7/10/1998	NM		2.00	
	50.02	7/14/1998	NM		2.00	
	50.02	7/23/1998	NM		2.00	
	50.02	8/5/1998	NM		2.00	
	50.02	8/13/1998	NM		2.00	
	50.02	8/18/1998	NM		2.00	
	50.02	8/25/1998	10.22			39.8

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-12B	50.02	9/15/1998	NM		2.00	
	50.02	9/21/1998	7.73			42.29
	50.02	9/30/1998	NM		4.00	
	50.02	10/8/1998	NM		4.00	
	50.02	10/16/1998	NM		4.00	
	50.02	10/26/1998	8.88			41.14
	50.02	11/6/1998	NM		4.00	
	50.02	11/13/1998	NM		1.49	
	50.02	11/19/1998	NM		4.00	
	50.02	11/23/1998	6.11			43.91
	50.02	12/16/1998	NM		4.00	
	50.02	1/7/1999	NM		4.00	
	50.02	1/15/1999	NM		4.00	
	50.02	1/22/1999	NM		4.00	
	50.02	1/26/1999	NM		4.00	
	50.02	1/29/1999	5.70			44.32
	50.02	2/4/1999	NM		4.00	
	50.02	2/9/1999	NM		3.00	
	50.02	2/26/1999	5.83	39.95	5.85	44.19
	50.02	3/16/1999	6.30	43.60	2.20	43.72
	50.02	4/29/1999	6.44	38.90	6.90	43.58
	50.02	5/21/1999	7.40	36.90	8.90	42.62
	50.02	5/27/1999	7.38	36.90	8.90	42.64
	50.02	6/1/1999	6.40	37.90	7.90	43.62
	50.02	6/10/1999	7.36	36.90	8.90	42.66
	50.02	7/30/1999	7.98			42.04
	50.02	8/27/1999	6.61	38.90	6.90	43.41
	50.02	9/27/1999	11.71	42.34	3.46	38.31
	50.02	10/29/1999	12.76	41.84	3.96	37.26
	50.02	11/18/1999	13.22			36.8
	50.02	12/29/1999	12.01	41.84	3.96	38.01
	50.02	2/4/2000	13.22	41.84	3.96	36.8
	50.02	2/25/2000	11.44	41.84	3.96	38.58
	50.02	3/27/2000	NM			
	50.02	4/7/2000	8.73	41.81	3.99	41.29
	50.02	5/31/2000	8.77	41.81	3.99	41.25
	50.02	6/1/2000	8.73	41.81	3.99	41.29
	50.02	7/28/2000	8.77	41.89	3.91	41.25
	50.02	8/30/2000	11.66	41.82	3.98	38.36
	50.02	9/19/2000	13.33	40.89	4.91	36.69
	50.02	10/27/2000	11.75	41.80	4.00	38.27
	50.02	11/21/2000	10.64	43.48	2.32	39.38
	50.02	5/1/2001	8.71	43.46	2.34	41.31
	50.02	10/1/2001	8.37		15.00	41.65
	50.02	3/14/2002	6.37	36.99	8.81	43.65
	50.02	9/23/2002	6.10	40.03	5.77	43.92
	50.02	3/10/2003	5.45			44.57
	50.02	9/24/2003	6.29	39.85	5.95	43.73
	50.02	3/15/2004	5.63			44.39
	50.02	9/13/2004	10.44	38.72	7.08	39.58
	50.02	7/18/2005	7.14	38.40	7.40	42.88
	50.02	1/4/2006	10.75	35.98	9.82	39.27
	50.02	7/27/2006	6.07	35.74	10.06	43.95
	50.02	3/7/2007	6.96	34.60	11.20	43.06
	50.02	7/27/2007	5.36	33.45	12.35	44.66
	50.02	1/31/2008	5.75	33.34	12.46	44.27
	50.02	7/15/2008	9.38	38.88	6.92	40.64
	50.02	2/4/2009	8.89	38.14	7.66	41.13
	50.02	7/24/2009	9.18	38.51	7.29	40.84
	50.02	1/8/2010	6.81	37.46	8.34	43.21
50.02	5/27/2010	7.29	39.5	6.30	42.73	
50.02	6/28/2010	7.39	44.1	1.70	42.63	
50.02	7/12/2010	7.47	44.25	1.55	42.55	
50.02	8/31/2010	7.26	45.42	0.38	42.76	
50.02	1/12/2011	7.01	45.39	0.41	43.01	
50.02	7/12/2011	10.09	45.39	0.41	39.93	
50.02	3/8/2012	6.87	40.2	5.60	43.15	
50.02	7/9/2012	7.16	40.1	5.70	42.86	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-12B	50.02	1/7/2013	9.17	39.86	5.94	40.85
	50.02	7/22/2013	11.16	39.04	6.76	38.86
	50.02	1/7/2014	11.34	45.12	0.68	38.68
	50.02	7/15/2014	10.59	44.89	0.91	39.43
	50.02	1/5/2015	10.06	44.91	1.29	39.96
	50.02	8/10/2015	7.39	46.1	0.10	42.63
	50.02	1/13/2016	6.06	45.79	0.41	43.96
	50.02	7/6/2016	6.29	45.72	0.48	43.73
	50.02	1/12/2017	7.02	45.81	0.39	43.00
	50.02	7/6/2017	7.01	45.71	1.89	43.01
	50.02	9/5/2017	7.03	45.6	2.00	42.99
	50.02	2/7/2018	7.13	45.87	0.33	42.89
	50.02	3/11/2018	7.42	45.96	0.24	42.6
	50.02	5/14/2018	8.59	45.91	0.29	41.43
	50.02	1/3/2019	7.96	45.87	0.33	42.06
	50.02	7/9/2019	6.67	45.06	1.14	43.35
	50.02	1/8/2020	5.66	27.12	19.08	44.36
MW-12C	50.14	5/13/1997	39.34			10.8
	50.14	6/20/1997	38.94			11.2
	50.14	9/25/1997	36.70			13.44
	50.14	10/22/1997	36.09			14.05
	50.14	11/25/1997	36.13			14.01
	50.14	12/19/1997	35.34			14.8
	50.14	1/20/1998	32.60			17.54
	50.14	3/4/1998	31.56			18.58
	50.14	3/18/1998	31.64			18.5
	50.14	4/24/1998	31.06			19.08
	50.14	5/21/1998	38.20			11.94
	50.14	8/25/1998	31.00			19.14
	50.14	9/21/1998	29.86			20.28
	50.14	10/26/1998	30.12			20.02
	50.14	11/23/1998	28.38			21.76
	50.14	1/29/1999	27.61			22.53
	50.14	2/26/1999	27.69			22.45
	50.14	3/16/1999	28.00			22.14
	50.14	4/29/1999	28.21			21.93
	50.14	6/1/1999	28.20			21.94
	50.14	7/30/1999	29.80			20.34
	50.14	8/27/1999	28.41			21.73
	50.14	9/27/1999	29.20			20.94
	50.14	10/29/1999	29.78			20.36
	50.14	11/18/1999	30.17			19.97
	50.14	12/29/1999	29.09			21.05
	50.14	2/4/2000	29.66			20.48
	50.14	2/25/2000	30.32			19.82
	50.14	3/27/2000	28.91			21.23
	50.14	4/7/2000	27.40			22.74
	50.14	5/31/2000	27.44			22.7
	50.14	6/1/2000	27.43			22.71
	50.14	7/28/2000	27.45			22.69
	50.14	8/30/2000	33.61			16.53
	50.14	9/19/2000	30.03			20.11
	50.14	10/27/2000	33.94			16.2
50.14	11/21/2000	29.12			21.02	
50.14	5/1/2001	26.85			23.29	
50.14	10/1/2001	26.85			23.29	
50.14	3/11/2002	25.59			24.55	
50.14	9/23/2002	26.57			23.57	
50.14	3/10/2003	24.85			25.29	
50.14	9/23/2003	26.06			24.08	
50.14	3/15/2004	24.31			25.83	
50.14	9/13/2004	26.15			23.99	
50.14	7/18/2005	26.23			23.91	
50.14	1/4/2006	22.26			27.88	
50.14	7/27/2006	25.28			24.86	
50.14	3/7/2007	23.78			26.36	
50.14	7/27/2007	22.05			28.09	
50.14	1/30/2008	22.69			27.45	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-12C	50.14	7/15/2008	24.41			25.73
	50.14	2/4/2009	24.59			25.55
	50.14	7/24/2009	24.91			25.23
	50.14	1/8/2010	23.03			27.11
	50.14	7/12/2010	23.91			26.23
	50.14	1/12/2011	23.76			26.38
	50.14	7/12/2011	25.98			24.16
	50.14	1/26/2012	25.76			24.38
	50.14	7/9/2012	24.59			25.55
	50.14	1/7/2013	26.04			24.1
	50.14	7/22/2013	27.09			23.05
	50.14	1/7/2014	26.52			23.62
	50.14	7/16/2014	25.15			24.99
	50.14	1/5/2015	26.01			24.13
	50.14	8/10/2015	24.26			25.88
	50.14	1/13/2016	23.83			26.31
	50.14	7/6/2016	24.13			26.01
	50.14	1/12/2017	24.49			25.65
	50.14	7/6/2017	24.88			25.26
	50.14	9/5/2017	24.84			25.30
50.14	2/11/2018	25.13			25.01	
50.14	3/11/2018	24.04			26.1	
50.14	4/14/2018	25.96			24.18	
50.14	1/3/2019	25.34			24.8	
50.14	7/9/2019	23.73			26.41	
50.14	1/7/2020	24.61			25.53	
MW-13	50.65	3/25/1997	9.43			41.22
	50.65	4/23/1997	9.87			40.78
	50.65	4/24/1997	9.92			40.73
	50.65	5/13/1997	9.30			41.35
	50.65	6/20/1997	10.11			40.54
	50.65	9/25/1997	10.75			39.9
	50.65	10/22/1997	10.09			40.56
	50.65	11/25/1997	10.11			40.54
	50.65	12/19/1997	10.01			40.64
	50.65	1/20/1998	9.32			41.33
	50.65	3/4/1998	9.23			41.42
	50.65	3/18/1998	8.90			41.75
	50.65	4/24/1998	10.74			39.82
	50.65	5/21/1998	12.11			38.54
	50.65	8/25/1998	12.00			38.56
	50.65	9/21/1998	10.13			40.43
	50.65	10/26/1998	11.15			39.41
	50.65	11/23/1998	9.22			41.34
	50.65	1/29/1999	8.00			42.65
	50.65	2/26/1999	8.11			42.54
	50.65	3/16/1999	9.51			41.14
	50.65	4/29/1999	9.79			40.86
	50.65	6/1/1999	9.70			40.95
	50.65	7/30/1999	11.01			39.64
	50.65	8/27/1999	9.96			40.69
	50.65	9/27/1999	12.84			37.81
	50.65	10/29/1999	13.88			36.77
	50.65	11/17/1999	14.00			36.65
	50.65	12/29/1999	13.08			37.57
	50.65	2/4/2000	15.61			35.04
	50.65	2/25/2000	12.17			38.48
	50.65	3/27/2000	10.95			39.7
	50.65	4/7/2000	10.51			40.14
	50.65	5/31/2000	10.57			40.08
50.65	6/1/2000	10.51			40.14	
50.65	7/28/2000	10.54			40.11	
50.65	8/30/2000	13.63			37.02	
50.65	9/19/2000	14.57			36.08	
50.65	10/27/2000	11.11			39.54	
50.65	11/21/2000	11.44			39.21	
50.65	5/1/2001	10.70			39.95	
50.65	10/1/2001	10.31			40.34	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-13	50.65	3/11/2002	9.62			41.03
	50.65	9/23/2002	9.17			41.48
	50.65	3/10/2003	9.17			41.48
	50.65	9/23/2003	9.14			41.51
	50.65	3/15/2004	9.30			41.35
	50.65	9/13/2004	11.98			38.67
	50.65	7/18/2005	10.25			40.4
	50.65	1/4/2006	12.03			38.62
	50.65	7/27/2006	8.82			41.83
	50.65	3/7/2007	9.95			40.7
	50.65	7/27/2007	8.90			41.75
	50.65	1/30/2008	8.85			41.8
	50.65	7/15/2008	10.89			39.76
	50.65	2/4/2009	10.59			40.06
	50.65	7/23/2009	11.07			39.58
	50.65	1/8/2010	9.22			41.43
	50.65	7/12/2010	11.12			39.53
	50.65	1/12/2011	8.89			41.76
	50.65	7/12/2011	12.96			37.69
	50.65	1/26/2012	9.31			41.34
	50.65	7/9/2012	9.14			41.51
	50.65	1/7/2013	10.68			39.97
	50.65	7/22/2013	12.13			38.52
	50.65	1/7/2014	10.13			40.52
	50.65	7/16/2014	11.04			39.61
	50.65	1/5/2015	9.34			41.31
	50.65	8/10/2015	7.67			42.98
	50.65	1/13/2016	7.01			43.64
	50.65	7/6/2016	7.39			43.26
	50.65	1/12/2017	7.81			42.84
	50.65	7/6/2017	7.96			42.69
	50.65	9/5/2017	9.01			41.64
	50.65	2/11/2018	9.58			41.07
50.65	3/11/2018	10.09			40.56	
50.65	5/14/2018	10.96			39.69	
50.65	1/3/2019	10.52			40.13	
50.65	7/9/2019	10.63			40.02	
50.65	1/7/2020	9.42			41.23	
MW-14	50.66	3/25/1997	7.71			42.95
	50.66	4/23/1997	8.31			42.35
	50.66	4/24/1997	8.34			42.32
	50.66	5/13/1997	7.83			42.83
	50.66	6/20/1997	8.64			42.02
	50.66	9/25/1997	9.95			40.71
	50.66	10/22/1997	8.89			41.77
	50.66	11/25/1997	8.86			41.8
	50.66	12/19/1997	8.62			42.04
	50.66	1/20/1998	8.08			42.58
	50.66	3/4/1998	7.72			42.94
	50.66	3/18/1998	7.66			43
	50.66	4/24/1998	9.75			40.91
	50.66	5/21/1998	11.00			39.66
	50.66	8/25/1998	12.00			38.66
	50.66	9/21/1998	9.41			41.25
	50.66	10/26/1998	11.10			39.56
	50.66	11/23/1998	8.08			42.58
	50.66	1/29/1999	7.10			43.56
	50.66	2/26/1999	7.21			43.45
	50.66	3/16/1999	8.74			41.92
	50.66	4/29/1999	8.93			41.73
	50.66	6/1/1999	8.92			41.74
	50.66	7/30/1999	10.44			40.22
	50.66	8/27/1999	9.21			41.45
	50.66	9/27/1999	12.56			38.1
	50.66	10/29/1999	13.56			37.1
	50.66	11/17/1999	13.63			37.03
50.66	12/29/1999	12.88			37.78	
50.66	2/4/2000	14.22			36.44	



**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-14	50.66	2/25/2000	11.73			38.93
	50.66	3/27/2000	10.54			40.12
	50.66	4/7/2000	10.14			40.52
	50.66	5/31/2000	10.17			40.49
	50.66	6/1/2000	10.13			40.53
	50.66	7/28/2000	10.17			40.49
	50.66	8/30/2000	13.22			37.44
	50.66	9/19/2000	14.27			36.39
	50.66	10/27/2000	11.56			39.1
	50.66	11/21/2000	11.17			39.49
	50.66	5/1/2001	9.71			40.95
	50.66	10/1/2001	10.64			40.02
	50.66	3/11/2002	8.45			42.21
	50.66	9/23/2002	7.90			42.76
	50.66	3/10/2003	8.59			42.07
	50.66	9/23/2003	7.70			42.96
	50.66	3/15/2004	7.96			42.7
	50.66	9/13/2004	11.05			39.61
	50.66	7/18/2005	9.55			41.11
	50.66	1/4/2006	11.83			38.83
	50.66	7/27/2006	7.80			42.86
	50.66	3/7/2007	8.96			41.7
	50.66	7/27/2007	8.01			42.65
	50.66	1/30/2008	7.66			43
	50.66	7/15/2008	10.41			40.25
	50.66	2/4/2009	10.27			40.39
	50.66	7/23/2009	10.67			39.99
	50.66	1/8/2010	8.24			42.42
	50.66	7/12/2010	10.54			40.12
	50.66	1/12/2011	18.09			32.57
	50.66	7/12/2011	12.93			37.73
	50.66	1/26/2012	8.57			42.09
	50.66	7/9/2012	8.61			42.05
	50.66	1/7/2013	10.46			40.2
	50.66	7/22/2013	11.91			38.75
	50.66	1/7/2014	9.39			41.27
	50.66	7/16/2014	10.58			40.08
	50.66	1/5/2015	8.79			41.87
	50.66	8/10/2015	6.34			44.32
	50.66	1/13/2016	5.79			44.87
50.66	7/6/2016	6.06			44.60	
50.66	1/12/2017	6.59			44.07	
50.66	7/6/2017	6.92			43.74	
50.66	9/5/2017	6.83			43.83	
50.66	2/11/2018	8.66			42.00	
50.66	3/11/2018	8.99			41.67	
50.66	5/14/2018	10.09			40.57	
50.66	1/3/2019	9.37			41.29	
50.66	7/9/2019	9.57			41.09	
50.66	1/7/2020	8.39			42.27	
MW-15A	50.41	3/25/1997	8.22			42.19
	50.41	4/23/1997	8.28			42.13
	50.41	4/24/1997	8.51			41.9
	50.41	5/13/1997	8.06			42.35
	50.41	6/20/1997	8.64			41.77
	50.41	9/25/1997	9.75			40.66
	50.41	10/22/1997	9.09			41.32
	50.41	11/25/1997	9.13			41.28
	50.41	12/19/1997	8.89			41.52
	50.41	1/20/1998	8.35			42.06
	50.41	3/4/1998	8.09			42.32
	50.41	3/18/1998	7.98			42.43
	50.41	4/24/1998	9.57			40.84
	50.41	5/21/1998	11.10			39.31
	50.41	8/25/1998	11.78			38.63
	50.41	9/21/1998	9.59			40.82
	50.41	10/26/1998	10.69			39.72
50.41	11/23/1998	8.46			41.95	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-15A	50.41	1/29/1999	7.11			43.3
	50.41	2/26/1999	7.23			43.18
	50.41	3/16/1999	9.17			41.24
	50.41	4/29/1999	9.29			41.12
	50.41	6/1/1999	9.29			41.12
	50.41	7/30/1999	10.83			39.58
	50.41	8/27/1999	9.39			41.02
	50.41	9/27/1999	12.02			38.39
	50.41	10/29/1999	13.11			37.3
	50.41	11/17/1999	13.44			36.97
	50.41	12/29/1999	12.49			37.92
	50.41	2/4/2000	15.71			34.7
	50.41	2/25/2000	11.34			39.07
	50.41	3/27/2000	10.66			39.75
	50.41	4/7/2000	10.20			40.21
	50.41	5/31/2000	10.23			40.18
	50.41	6/1/2000	10.22			40.19
	50.41	7/28/2000	10.23			40.18
	50.41	8/30/2000	13.34			37.07
	50.41	9/19/2000	14.01			36.4
	50.41	10/27/2000	11.77			38.64
	50.41	11/21/2000	11.09			39.32
	50.41	5/1/2001	9.85			40.56
	50.41	10/1/2001	9.73			40.68
	50.41	3/11/2002	8.81			41.6
	50.41	9/23/2002	8.21			42.2
	50.41	3/10/2003	7.76			42.65
	50.41	9/23/2003	7.87			42.54
	50.41	3/15/2004	7.94			42.47
	50.41	9/13/2004	10.72			39.69
	50.41	7/18/2005	9.33			41.08
	50.41	1/4/2006	11.66			38.75
	50.41	7/27/2006	7.92			42.49
	50.41	3/7/2007	9.19			41.22
	50.41	7/27/2007	7.88			42.53
	50.41	1/30/2008	8.02			42.39
	50.41	7/15/2008	10.26			40.15
	50.41	2/4/2009	10.59			39.82
	50.41	7/23/2009	11.01			39.4
	50.41	1/8/2010	8.64			41.77
50.41	7/12/2010	10.81			39.6	
50.41	1/12/2011	8.77			41.64	
50.41	7/12/2011	12.78			37.63	
50.41	1/26/2012	9.29			41.12	
50.41	7/9/2012	5.92			44.49	
50.41	1/7/2013	10.77			39.64	
50.41	7/22/2013	12.21			38.2	
50.41	1/7/2014	9.85			40.56	
50.41	7/16/2014	10.65			39.76	
50.41	1/5/2015	9.07			41.34	
50.41	8/10/2015	6.49			43.92	
50.41	1/13/2016	5.79			44.62	
50.41	7/6/2016	6.21			44.20	
50.41	1/12/2017	6.82			43.59	
50.41	7/6/2017	7.47			42.94	
50.41	9/5/2017	7.43			42.98	
50.41	2/11/2018	8.89			41.52	
50.41	3/11/2018	9.23			41.18	
50.41	5/14/2018	10.18			40.23	
50.41	1/3/2019	9.41			41	
50.41	7/9/2019	9.03			41.38	
50.41	1/7/2020	8.57			41.84	
MW-15B	50.20	1/26/2012	10.13			40.07
	50.20	7/9/2012	8.32			41.88
	50.20	1/7/2013	10.71			39.49
	50.20	7/22/2013	11.97			38.23
	50.20	1/7/2014	9.81			40.39
50.20	7/15/2014	10.36			39.84	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-15B	50.20	1/5/2015	9.26			40.94
	50.20	8/10/2015	7.29			42.91
	50.20	1/13/2016	6.81			43.39
	50.20	7/6/2016	7.56			42.64
	50.20	1/12/2017	8.09			42.11
	50.20	7/6/2017	8.61			41.59
	50.20	9/5/2017	8.56			41.64
	50.20	2/11/2018	8.74			41.46
	50.20	3/11/2018	9.09			41.11
	50.20	5/14/2018	9.91			40.29
	50.20	1/3/2019	9.4			40.8
	50.20	7/9/2019	9.23			40.97
	50.20	1/7/2020	8.39			41.81
	MW-15C	50.01	5/13/1997	33.46		
50.01		6/20/1997	34.18			15.83
50.01		9/25/1997	33.77			16.24
50.01		10/22/1997	32.89			17.12
50.01		11/25/1997	32.95			17.06
50.01		12/19/1997	32.01			18
50.01		1/20/1998	29.90			20.11
50.01		3/4/1998	28.56			21.45
50.01		3/18/1998	28.53			21.48
50.01		4/24/1998	28.46			21.55
50.01		5/21/1998	35.00			15.01
50.01		8/25/1998	29.30			20.71
50.01		9/21/1998	28.15			21.86
50.01		10/26/1998	28.11			21.9
50.01		11/23/1998	26.50			23.51
50.01		1/29/1999	25.44			24.57
50.01		2/26/1999	25.51			24.5
50.01		3/16/1999	26.11			23.9
50.01		4/29/1999	26.33			23.68
50.01		6/1/1999	26.39			23.62
50.01		7/30/1999	27.99			22.02
50.01		8/27/1999	26.51			23.5
50.01		9/27/1999	27.46			22.55
50.01		10/29/1999	28.26			21.75
50.01		11/17/1999	28.55			21.46
50.01		12/29/1999	27.61			22.4
50.01		2/4/2000	28.11			21.9
50.01		2/25/2000	28.23			21.78
50.01		3/27/2000	27.45			22.56
50.01		4/7/2000	26.11			23.9
50.01		5/31/2000	26.13			23.88
50.01		6/1/2000	26.03			23.98
50.01		7/28/2000	26.14			23.87
50.01		8/30/2000	29.11			20.9
50.01		9/19/2000	28.67			21.34
50.01		10/27/2000	27.64			22.37
50.01		11/21/2000	27.56			22.45
50.01		5/1/2001	25.24			24.77
50.01		10/1/2001	25.40			24.61
50.01		3/11/2002	24.17			25.84
50.01		9/23/2002	25.35			24.66
50.01		3/10/2003	23.52			26.49
50.01		9/23/2003	24.88			25.13
50.01		3/15/2004	22.97			27.04
50.01		9/13/2004	24.80			25.21
50.01		7/18/2005	25.17			24.84
50.01		1/4/2006	26.23			23.78
50.01	7/27/2006	24.31			25.7	
50.01	3/7/2007	22.76			27.25	
50.01	7/27/2007	21.03			28.98	
50.01	1/30/2008	21.80			28.21	
50.01	7/15/2008	23.63			26.38	
50.01	2/4/2009	23.73			26.28	
50.01	7/23/2009	23.96			26.05	
50.01	1/8/2010	21.88			28.13	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-15C	50.01	7/12/2010	23.08			26.93
	50.01	1/12/2011	23.04			26.97
	50.01	7/12/2011	25.09			24.92
	50.01	1/26/2012	24.37			25.64
	50.01	7/9/2012	24.41			25.6
	50.01	1/7/2013	25.21			24.8
	50.01	7/22/2013	26.10			23.91
	50.01	1/7/2014	25.26			24.75
	50.01	7/16/2014	24.15			25.86
	50.01	1/5/2015	25.34			24.67
	50.01	8/10/2015	22.74			27.27
	50.01	1/13/2016	21.92			28.09
	50.01	7/6/2016	22.26			27.75
	50.01	1/12/2017	22.69			27.32
	50.01	7/6/2017	23.31			26.70
	50.01	9/5/2017	23.29			26.72
	50.01	2/11/2018	23.63			26.38
	50.01	3/11/2018	22.47			27.54
	50.01	5/14/2018	23.33			26.68
	50.01	1/3/2019	23.87			26.14
50.01	7/9/2019	22.38			27.63	
50.01	1/7/2020	23.12			26.89	
MW-16	51.51	3/25/1997	7.41			44.1
	51.51	4/23/1997	8.44			43.07
	51.51	4/24/1997	8.52			42.99
	51.51	5/13/1997	8.29			43.22
	51.51	6/20/1997	8.41			43.1
	51.51	9/25/1997	10.71			40.8
	51.51	10/22/1997	9.53			41.98
	51.51	11/25/1997	9.55			41.96
	51.51	12/19/1997	9.10			42.41
	51.51	1/20/1998	8.60			42.91
	51.51	3/4/1998	8.13			43.38
	51.51	3/18/1998	8.59			42.92
	51.51	4/24/1998	9.96			41.55
	51.51	5/21/1998	11.43			40.08
	51.51	7/30/1998	12.56			38.95
	51.51	8/25/1998	11.53			39.98
	51.51	9/21/1998	9.81			41.7
	51.51	10/26/1998	10.44			41.07
	51.51	11/23/1998	8.98			42.53
	51.51	1/29/1999	7.12			44.39
	51.51	2/26/1999	7.23			44.28
	51.51	3/16/1999	10.06			41.45
	51.51	4/29/1999	10.16			41.35
	51.51	6/1/1999	10.16			41.35
	51.51	7/30/1999	11.76			39.75
	51.51	8/27/1999	10.33			41.18
	51.51	9/27/1999	11.79			39.72
	51.51	10/29/1999	12.93			38.58
	51.51	11/17/1999	13.71			37.8
	51.51	12/29/1999	12.20			39.31
	51.51	2/4/2000	15.11			36.4
	51.51	2/25/2000	11.10			40.41
	51.51	3/27/2000	11.48			40.03
	51.51	4/7/2000	11.09			40.42
	51.51	5/31/2000	11.11			40.4
	51.51	6/1/2000	11.00			40.51
	51.51	7/28/2000	11.11			40.4
	51.51	8/30/2000	13.10			38.41
	51.51	9/19/2000	14.83			36.68
	51.51	10/27/2000	11.66			39.85
51.51	11/21/2000	11.29			40.22	
51.51	5/1/2001	9.92			41.59	
51.51	10/1/2001	9.93			41.58	
51.51	3/11/2002	9.12			42.39	
51.51	9/23/2002	8.65			42.86	
51.51	3/10/2003	7.74			43.77	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-16	51.51	9/23/2003	8.48			43.03
	51.51	3/15/2004	8.09			43.42
	51.51	9/13/2004	10.38			41.13
	51.51	7/18/2005	10.42			41.09
	51.51	1/4/2006	12.48			39.03
	51.51	7/27/2006	9.37			42.14
	51.51	3/7/2007	9.66			41.85
	51.51	7/27/2007	7.85			43.66
	51.51	1/31/2008	8.42	25.40	3.40	43.09
	51.51	7/15/2008	10.16			41.35
	51.51	2/5/2009	11.93			39.58
	51.51	7/23/2009	12.67			38.84
	51.51	1/8/2010	8.66			42.85
	51.51	7/12/2010	10.31			41.2
	51.51	1/12/2011	9.89			41.62
	51.51	7/12/2011	12.98			38.53
	51.51	1/26/2012	9.92			41.59
	51.51	7/9/2012	9.68			41.83
	51.51	1/7/2013	11.41			40.1
	51.51	7/22/2013	12.39			39.12
51.51	1/7/2014	12.02			39.49	
51.51	7/15/2014	9.69			41.82	
51.51	1/5/2015	11.07			40.44	
51.51	8/10/2015	9.42			42.09	
MW-17	50.92	3/25/1997	9.97			40.95
	50.92	4/23/1997	10.41			40.51
	50.92	4/24/1997	10.51			40.41
	50.92	5/13/1997	10.32			40.6
	50.92	6/20/1997	11.07			39.85
	50.92	9/25/1997	12.39			38.53
	50.92	10/22/1997	11.19			39.73
	50.92	11/25/1997	11.21			39.71
	50.92	12/19/1997	11.01			39.91
	50.92	1/20/1998	10.25			40.67
	50.92	3/4/1998	9.93			40.99
	50.92	3/18/1998	9.94			40.98
	50.92	4/9/1998	11.32			39.6
	50.92	4/16/1998	11.52			39.4
	50.92	4/24/1998	11.80			39.12
	50.92	5/8/1998	NM			
	50.92	5/12/1998	NM			
	50.92	5/21/1998	13.30			37.62
	50.92	5/25/1998	NM			
	50.92	6/9/1998	NM			
	50.92	6/16/1998	NM			
	50.92	6/26/1998	NM			
	50.92	7/2/1998	NM			
	50.92	7/10/1998	NM			
	50.92	7/14/1998	NM			
	50.92	7/23/1998	NM			
	50.92	8/5/1998	NM			
	50.92	8/13/1998	NM			
	50.92	8/25/1998	13.78			37.14
	50.92	9/15/1998	NM			
	50.92	9/21/1998	11.49			39.43
	50.92	9/30/1998	NM			
	50.92	10/8/1998	NM			
	50.92	10/16/1998	NM			
	50.92	10/26/1998	12.22			38.7
	50.92	11/6/1998	NM			
	50.92	11/13/1998	NM			
	50.92	11/19/1998	NM			
	50.92	11/23/1998	10.21			40.71
	50.92	12/16/1998	NM			
50.92	1/7/1999	NM				
50.92	1/15/1999	NM				
50.92	1/22/1999	NM				
50.92	1/26/1999	NM				

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-17	50.92	1/29/1999	10.88			40.04
	50.92	2/4/1999	NM			
	50.92	2/9/1999	NM			
	50.92	2/26/1999	10.93			39.99
	50.92	3/16/1999	11.18			39.74
	50.92	4/29/1999	11.00			39.92
	50.92	5/21/1999	11.25			39.67
	50.92	5/27/1999	11.31			39.61
	50.92	6/1/1999	11.07			39.85
	50.92	6/10/1999	11.28			39.64
	50.92	7/30/1999	12.67			38.25
	50.92	8/27/1999	11.27			39.65
	50.92	9/27/1999	14.67			36.25
	50.92	10/29/1999	15.11			35.81
	50.92	11/17/1999	16.08			34.84
	50.92	12/29/1999	14.43			36.49
	50.92	2/4/2000	17.21			33.71
	50.92	2/25/2000	13.63			37.29
	50.92	3/27/2000	13.08	32.60	0.70	37.84
	50.92	4/7/2000	12.63	32.30	1.00	38.29
	50.92	5/31/2000	12.67	32.30	1.00	38.25
	50.92	6/1/2000	12.61	32.30	1.00	38.31
	50.92	7/28/2000	12.69	32.30	1.00	38.23
	50.92	8/30/2000	15.56			35.36
	50.92	9/19/2000	16.24	32.20	1.10	34.68
	50.92	10/27/2000	14.10			36.82
	50.92	11/21/2000	13.12			37.8
	50.92	5/1/2001	11.82	32.44	0.86	39.1
	50.92	10/1/2001	12.55	32.30	1.00	38.37
	50.92	3/14/2002	10.91	31.79	1.51	40.01
	50.92	9/23/2002	10.48			40.44
	50.92	3/10/2003	9.76			41.16
	50.92	9/24/2003	10.59	32.85	0.45	40.33
	50.92	3/15/2004	10.15			40.77
	50.92	9/13/2004	13.09			37.83
	50.92	7/18/2005	12.06	32.90	0.40	38.86
	50.92	1/4/2006	13.90	32.90	0.40	37.02
	50.92	7/27/2006	10.71	33.28	0.02	40.21
	50.92	3/7/2007	10.91	33.00	0.30	40.01
	50.92	7/27/2007	9.33	33.02	0.28	41.59
	50.92	1/31/2008	10.00	31.17	2.13	40.92
	50.92	7/15/2008	12.95	33.08	0.23	37.97
	50.92	2/4/2009	12.64	Trace	Trace	38.28
	50.92	1/8/2010	10.62			40.3
	50.92	7/12/2010	12.96			37.96
	50.92	7/12/2010	12.96			37.96
	50.92	1/12/2011	11.06			39.86
	50.92	7/12/2011	14.93			35.99
	50.92	1/26/2012	11.2			39.72
	50.92	7/9/2012	11.02			39.9
	50.92	1/7/2013	13.14			37.78
	50.92	7/22/2013	14.62			36.3
	50.92	1/7/2014	12.36			38.56
	50.92	7/15/2014	12.54			38.38
	50.92	1/5/2015	11.71			39.21
	50.92	8/10/2015	9.61			41.31
	50.92	1/13/2016	9.02			41.90
	50.92	7/6/2016	9.47			41.45
	50.92	1/12/2017	10.06			40.86
	50.92	7/6/2017	10.62			40.30
	50.92	9/5/2017	10.51			40.41
	50.92	2/11/2018	10.76			40.16
	50.92	3/11/2018	11.21			39.71
	50.92	5/14/2018	12.21			38.71
	50.92	1/3/2019	11.72			39.2
	50.92	7/9/2019	11.66			39.26
	50.92	1/7/2020	8.23			42.69
MW-17C	50.17	3/15/2004	22.75			27.42

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-17C	50.17	9/13/2004	24.56			25.61
	50.17	7/18/2005	25.02			25.15
	50.17	1/4/2006	26.07			24.1
	50.17	7/27/2006	24.15			26.02
	50.17	3/7/2007	22.51			27.66
	50.17	7/27/2007	20.93			29.24
	50.17	1/30/2008	21.74			28.43
	50.17	7/15/2008	23.65			26.52
	50.17	2/4/2009	23.72			26.45
	50.17	7/23/2009	24.08			26.09
	50.17	1/8/2010	21.98			28.19
	50.17	7/12/2010	23.03			27.14
	50.17	1/12/2011	23.16			27.01
	50.17	7/12/2011	25.11			25.06
	50.17	1/26/2012	24.27			25.9
	50.17	7/9/2012	24.32			25.85
	50.17	1/7/2013	24.76			25.41
	50.17	7/22/2013	25.89			24.28
	50.17	1/7/2014	25.06			25.11
	50.17	7/15/2014	23.98			26.19
	50.17	1/5/2015	24.62			25.55
	50.17	8/10/2015	22.47			27.70
	50.17	1/13/2016	21.81			28.36
	50.17	7/6/2016	22.16			28.01
	50.17	1/12/2017	22.67			27.50
	50.17	7/6/2017	23.09			27.08
	50.17	9/5/2017	23.01			27.16
	50.17	2/11/2018	23.11			27.06
	50.17	3/11/2018	22.21			27.96
	50.17	5/14/2018	23.02			27.15
	50.17	1/3/2019	22.71			27.46
	50.17	7/9/2019	22.14			28.03
50.17	1/7/2020	23.21			26.96	
MW-18A	51.57	3/25/1997	15.41			36.16
	51.57	4/23/1997	15.80			35.77
	51.57	5/13/1997	14.92			36.65
	51.57	6/20/1997	16.02			35.55
	51.57	9/25/1997	15.15			36.42
	51.57	10/22/1997	16.38			35.19
	51.57	11/25/1997	16.37			35.2
	51.57	12/19/1997	16.11			35.46
	51.57	1/20/1998	15.49			36.08
	51.57	3/4/1998	15.19			36.38
	51.57	3/18/1998	14.28			37.29
	51.57	4/24/1998	17.53			34.04
	51.57	5/21/1998	18.41			33.16
	51.57	7/30/1998	18.59			32.98
	51.57	8/25/1998	16.95			34.62
	51.57	9/21/1998	16.39			35.18
	51.57	10/26/1998	15.77			35.8
	51.57	11/23/1998	16.26			35.31
	51.57	1/29/1999	17.02			34.55
	51.57	2/26/1999	17.11			34.46
	51.57	4/29/1999	16.01			35.56
	51.57	6/1/1999	16.11			35.46
	51.57	7/30/1999	17.55			34.02
	51.57	8/27/1999	16.39			35.18
	51.57	9/27/1999	19.13			32.44
	51.57	10/29/1999	20.50			31.07
	51.57	11/17/1999	21.63			29.94
	51.57	12/29/1999	19.83			31.74
	51.57	2/4/2000	23.71			27.86
	51.57	2/25/2000	18.80			32.77
	51.57	3/27/2000	17.98			33.59
	51.57	4/7/2000	17.61			33.96
51.57	5/31/2000	17.65			33.92	
51.57	6/1/2000	17.60			33.97	
51.57	7/28/2000	17.67			33.9	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-18A	51.57	8/30/2000	20.30			31.27
	51.57	9/19/2000	19.54			32.03
	51.57	10/27/2000	18.75			32.82
	51.57	11/21/2000	16.52			35.05
	51.57	5/1/2001	17.91	27.85	7.94	33.66
	51.57	10/1/2001	17.47			34.1
	51.57	3/11/2002	16.68			34.89
	51.57	9/23/2002	15.30			36.27
	51.57	3/10/2003	15.77			35.8
	51.57	9/23/2003	25.08			26.49
	51.57	3/15/2004	15.58			35.99
	51.57	9/13/2004	18.32			33.25
	51.57	7/18/2005	14.88			36.69
	51.57	1/4/2006	17.96			33.61
	51.57	7/27/2006	14.15			37.42
	51.57	3/7/2007	17.32			34.25
	51.57	7/27/2007	15.22			36.35
	51.57	1/30/2008	15.63			35.94
	51.57	7/15/2008	17.43			34.14
	51.57	2/5/2009	18.67			32.9
	51.57	7/23/2009	19.03			32.54
	51.57	1/8/2010	16.51			35.06
	51.57	7/12/2010	18.11			33.46
	51.57	1/12/2011	15.82			35.75
	51.57	7/12/2011	19.02			32.55
	51.57	1/26/2012	16.9			34.67
	51.57	7/9/2012	15.06			36.51
	51.57	1/7/2013	18.39			33.18
	51.57	7/22/2013	18.74			32.83
	51.57	1/7/2014	18.06			33.51
	51.57	7/16/2014	18.14			33.43
	51.57	1/5/2015	17.39			34.18
51.57	8/10/2015	15.02			36.55	
51.57	1/13/2016	14.36			37.21	
51.57	7/6/2016	14.71			36.86	
51.57	1/12/2017	15.09			36.48	
51.57	7/6/2017	15.59			35.98	
51.57	9/5/2017	15.49			36.08	
51.57	2/11/2018	16.62			34.95	
51.57	3/11/2018	17.12			34.45	
51.57	5/14/2018	17.71			33.86	
51.57	1/3/2019	17.52			34.05	
51.57	7/9/2019	17.51			34.06	
51.57	1/7/2020	17.71			33.86	
MW-18C	51.47	5/13/1997	29.45			22.02
	51.47	6/20/1997	30.37			21.1
	51.47	9/25/1997	31.53			19.94
	51.47	10/22/1997	30.71			20.76
	51.47	11/25/1997	30.75			20.72
	51.47	12/19/1997	30.10			21.37
	51.47	1/20/1998	28.30			23.17
	51.47	3/4/1998	27.03			24.44
	51.47	3/18/1998	26.81			24.66
	51.47	4/9/1998	27.04			24.43
	51.47	4/16/1998	27.03			24.44
	51.47	4/24/1998	27.25			24.22
	51.47	5/8/1998	NM			
	51.47	5/12/1998	NM			
	51.47	5/21/1998	27.68			23.79
	51.47	5/25/1998	NM			
	51.47	6/9/1998	NM			
	51.47	6/16/1998	NM			
	51.47	6/26/1998	NM			
	51.47	7/2/1998	NM			
	51.47	7/10/1998	NM			
	51.47	7/14/1998	NM			
	51.47	7/23/1998	NM			
	51.47	7/30/1998	28.40			23.07



**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)	
MW-18C	51.47	8/5/1998	NM				
	51.47	8/13/1998	NM				
	51.47	8/25/1998	28.88			22.59	
	51.47	9/15/1998	NM				
	51.47	9/21/1998	27.94			23.53	
	51.47	9/30/1998	NM				
	51.47	10/8/1998	NM				
	51.47	10/16/1998	NM				
	51.47	10/26/1998	27.62			23.85	
	51.47	11/6/1998	NM				
	51.47	11/11/1998	26.85			24.62	
	51.47	11/19/1998	NM				
	51.47	11/23/1998	26.21			25.26	
	51.47	12/16/1998	NM				
	51.47	1/7/1999	NM				
	51.47	1/15/1999	NM				
	51.47	1/22/1999	NM				
	51.47	1/26/1999	NM				
	51.47	1/29/1999	25.36			26.11	
	51.47	2/4/1999	NM				
	51.47	2/9/1999	NM				
	51.47	2/26/1999	25.41			26.06	
	51.47	4/29/1999	26.33			25.14	
	51.47	5/21/1999	25.75			25.72	
	51.47	5/27/1999	25.76			25.71	
	51.47	6/1/1999	26.38			25.09	
	51.47	6/10/1999	25.68			25.79	
	51.47	7/30/1999	25.61			25.86	
	51.47	8/27/1999	26.51			24.96	
	51.47	9/27/1999	27.28			24.19	
	51.47	10/29/1999	27.95			23.52	
	51.47	11/17/1999	28.42			23.05	
	51.47	12/29/1999	27.26			24.21	
	51.47	2/4/2000	27.84			23.63	
	51.47	2/25/2000	27.83			23.64	
	51.47	3/27/2000	27.48			23.99	
	51.47	4/7/2000	25.80			25.67	
	51.47	5/31/2000	25.83			25.64	
	51.47	6/1/2000	25.81			25.66	
	51.47	7/28/2000	25.86			25.61	
	51.47	8/30/2000	28.42			23.05	
	51.47	9/19/2000	28.77		80.44	0.97	22.7
	51.47	10/27/2000	28.69				22.78
	51.47	11/21/2000	27.67				23.8
	51.47	5/1/2001	25.20				26.27
	51.47	10/1/2001	25.59				25.8
	51.47	3/14/2002	24.35				27.12
	51.47	9/25/2002	25.45				26.02
	51.47	3/10/2003	23.60				27.87
	51.47	9/24/2003	25.15				26.32
	51.47	3/15/2004	24.23				27.24
	51.47	9/13/2004	25.12		78.22	1.70	26.35
	51.47	7/18/2005	25.50		66.20	0.30	25.97
	51.47	1/4/2006	26.71				24.76
	51.47	7/27/2006	24.80				26.67
	51.47	3/7/2007	23.11				28.36
	51.47	7/27/2007	24.80				26.67
	51.47	1/30/2008	22.64				28.83
	51.47	7/15/2008	24.43				27.04
	51.47	2/5/2009	24.34				27.13
51.47	7/23/2009	24.61				26.86	
51.47	1/8/2010	22.56				28.91	
51.47	7/12/2010	23.77				27.7	
51.47	1/12/2011	24.03				27.44	
51.47	7/12/2011	25.87				25.6	
51.47	1/26/2012	26.82				24.65	
51.47	7/9/2012	24.82				26.65	
51.47	1/7/2013	25.61				25.86	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-18C	51.47	7/22/2013	26.76			24.71
	51.47	1/7/2014	25.68			25.79
	51.47	7/16/2014	24.60			26.87
	51.47	1/5/2015	25.02			26.45
	51.47	8/10/2015	23.41			28.06
	51.47	1/13/2016	22.76			28.71
	51.47	7/6/2016	23.12			28.35
	51.47	1/12/2017	23.73			27.74
	51.47	7/6/2017	24.13			27.34
	51.47	9/5/2017	24.08			27.39
	51.47	2/11/2018	23.7			27.77
	51.47	3/11/2018	22.88			28.59
	51.47	5/14/2018	23.47			28.00
	51.47	1/3/2019	23.01			28.46
	51.47	7/9/2019	23.08			28.39
	51.47	1/7/2020	23.54			27.93
	MW-19C	53.05	11/23/1998	28.84		
53.05		1/29/1999	28.21			24.84
53.05		2/26/1999	28.28			24.77
53.05		3/16/1999	28.31			24.74
53.05		4/29/1999	28.56			24.49
53.05		6/1/1999	28.48			24.57
53.05		7/30/1999	30.00			23.05
53.05		8/27/1999	28.61			24.44
53.05		9/27/1999	29.72			23.33
53.05		10/29/1999	30.46			22.59
53.05		11/17/1999	30.76			22.29
53.05		12/29/1999	29.44			23.61
53.05		2/4/2000	30.22			22.83
53.05		2/25/2000	29.93			23.12
53.05		3/27/2000	29.80			23.25
53.05		4/7/2000	28.40			24.65
53.05		5/31/2000	28.44			24.61
53.05		6/1/2000	28.33			24.72
53.05		7/28/2000	28.37			24.68
53.05		8/30/2000	29.99			23.06
53.05		9/19/2000	30.97			22.08
53.05		10/27/2000	28.49			24.56
53.05		11/21/2000	29.88			23.17
53.05		5/1/2001	27.61	71.55	3.56	25.44
53.05		10/1/2001	27.84			25.21
53.05		3/11/2002	26.68			26.37
53.05		9/23/2002	27.66			25.39
53.05		3/10/2003	25.77			27.28
53.05		9/23/2003	27.21			25.84
53.05		3/15/2004	25.36			27.69
53.05		9/13/2004	27.20			25.85
53.05		7/18/2005	27.71			25.34
53.05		1/4/2006	28.78			24.27
53.05		7/27/2006	26.91			26.14
53.05		3/7/2007	25.22			27.83
53.05		7/27/2007	23.71			29.34
53.05		1/31/2008	24.57			28.48
53.05		7/15/2008	26.38			26.67
53.05		2/4/2009	26.44			26.61
53.05		7/23/2009	26.81			26.24
53.05		1/9/2010	24.47			28.58
53.05		7/12/2010	25.67			27.38
53.05		1/12/2011	25.86			27.19
53.05	7/12/2011	27.81			25.24	
53.05	1/26/2012	26.74			26.31	
53.05	7/9/2012	27.26			25.79	
53.05	1/7/2013	27.73			25.32	
53.05	7/22/2013	28.58			24.47	
53.05	1/7/2014	27.71			25.34	
53.05	7/15/2014	26.65			26.40	
53.05	1/5/2015	27.34			25.71	
53.05	8/10/2015	25.21			27.84	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-19C	53.05	1/13/2016	24.68			28.37
	53.05	7/6/2016	NM			
	53.05	2/11/2018	21.74			31.31
	53.05	3/11/2018	24.74			28.31
	53.05	5/14/2018	25.72			27.33
	53.05	1/3/2019	25.03			28.02
	53.05	7/9/2019	24.58			28.47
	53.05	1/8/2020	25.71			27.34
MW-20A	50.43	11/23/1998	8.31			42.116
	50.43	1/29/1999	8.70			41.726
	50.43	2/26/1999	8.81			41.616
	50.43	3/16/1999	9.26			41.166
	50.43	4/29/1999	9.33			41.096
	50.43	6/1/1999	9.30			41.126
	50.43	7/30/1999	10.91			39.516
	50.43	8/27/1999	9.56			40.866
	50.43	9/27/1999	10.79			39.636
	50.43	10/29/1999	11.96			38.466
	50.43	11/17/1999	13.06			37.366
	50.43	12/29/1999	11.11			39.316
	50.43	2/4/2000	14.89			35.536
	50.43	2/25/2000	10.33			40.096
	50.43	3/27/2000	10.79			39.636
	50.43	4/7/2000	10.41			40.016
	50.43	5/31/2000	10.46			39.966
	50.43	6/1/2000	10.41			40.016
	50.43	7/28/2000	10.47			39.956
	50.43	8/30/2000	12.56			37.866
	50.43	9/19/2000	13.68			36.746
	50.43	10/27/2000	11.01			39.416
	50.43	11/21/2000	10.64			39.786
	50.43	5/1/2001	9.40			41.03
	50.43	10/1/2001	10.42			40.01
	50.43	3/11/2002	8.59			41.836
	50.43	9/23/2002	8.51			41.916
	50.43	3/10/2003	7.42			43.006
	50.43	9/23/2003	7.95			42.476
	50.43	3/15/2004	7.72			42.706
	50.43	9/13/2004	10.22			40.206
	50.43	7/18/2005	9.88			40.546
	50.43	1/4/2006	11.72			38.706
	50.43	7/27/2006	8.59			41.836
	50.43	3/7/2007	8.91			41.516
	50.43	7/27/2007	7.63			42.796
	50.43	1/30/2008	7.91			42.516
	50.43	7/15/2008	10.05			40.376
	50.43	2/4/2009	10.18			40.246
	50.43	7/23/2009	10.47			39.956
	50.43	1/9/2010	8.23			42.196
	50.43	7/12/2010	10.62			39.806
	50.43	1/12/2011	8.76			41.666
	50.43	7/12/2011	12.53			37.896
	50.43	1/26/2012	11.61			38.816
	50.43	7/9/2012	9.18			41.246
	50.43	1/7/2013	10.66			39.766
50.43	7/22/2013	12.17			38.256	
50.43	1/7/2014	11.62			38.806	
50.43	7/15/2014	9.83			40.60	
50.43	1/5/2015	11.09			39.34	
50.43	8/10/2015	9.34			41.09	
50.43	7/6/2017	8.12			42.31	
50.43	9/6/2017	8.06			42.37	
50.43	2/11/2018	9.22			41.21	
50.43	3/11/2018	9.03			41.396	
50.43	5/14/2018	9.89			40.536	
50.43	1/3/2019	9.26			41.17	
50.43	7/9/2019	9.04			41.386	
50.43	1/7/2020	9.06			41.366	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-20A	50.43					50.426
	50.43					50.426
	50.43					50.426
MW-21C	49.05	11/23/1998	27.83			21.223
	49.05	1/29/1999	27.11			21.943
	49.05	2/26/1999	27.26			21.793
	49.05	3/16/1999	27.42			21.633
	49.05	4/29/1999	27.99			21.063
	49.05	6/1/1999	27.80			21.253
	49.05	7/30/1999	29.00			20.053
	49.05	8/27/1999	27.99			21.063
	49.05	9/27/1999	28.43			20.623
	49.05	10/29/1999	29.12			19.933
	49.05	11/18/1999	29.25			19.803
	49.05	12/29/1999	10.89			38.163
	49.05	2/4/2000	28.94			20.113
	49.05	2/25/2000	11.43			37.623
	49.05	3/27/2000	28.13			20.923
	49.05	4/7/2000	26.79			22.263
	49.05	5/31/2000	26.83			22.223
	49.05	6/1/2000	26.83			22.223
	49.05	7/28/2000	26.88			22.173
	49.05	8/30/2000	29.91			19.143
	49.05	9/19/2000	29.15			19.903
	49.05	10/27/2000	30.21			18.843
	49.05	11/21/2000	28.33			20.723
	49.05	5/1/2001	26.01			23.04
	49.05	10/1/2001	26.05			23
	49.05	3/11/2002	24.80			24.253
	49.05	9/23/2002	25.50			23.553
	49.05	3/10/2003	23.82			25.233
	49.05	9/23/2003	25.08			23.973
	49.05	3/15/2004	23.48			25.573
	49.05	9/13/2004	25.44			23.613
	49.05	7/18/2005	25.33			23.723
	49.05	1/4/2006	26.44			22.613
	49.05	7/27/2006	24.55			24.503
	49.05	3/7/2007	22.91			26.143
	49.05	7/27/2007	21.29			27.763
	49.05	1/29/2008	22.09			26.963
	49.05	7/15/2008	23.31			25.743
	49.05	2/4/2009	24.03			25.023
	49.05	7/24/2009	24.29			24.763
	49.05	1/9/2010	21.89			27.163
49.05	7/12/2010	23.01			26.043	
49.05	1/12/2011	23.21			25.843	
49.05	7/12/2011	25.09			23.963	
49.05	1/26/2012	24.48			24.573	
49.05	7/9/2012	23.39			25.663	
49.05	1/7/2013	25.17			23.883	
49.05	7/22/2013	26.49			22.563	
49.05	1/7/2014	25.94			23.113	
49.05	7/15/2014	24.61			24.44	
49.05	1/5/2015	25.31			23.74	
49.05	8/10/2015	23.37			25.68	
49.05	1/13/2016	22.71			26.34	
49.05	7/6/2016	23.04			26.01	
49.05	1/12/2017	23.59			25.46	
49.05	7/6/2017	24.02			25.03	
49.05	9/5/2017	23.96			25.09	
49.05	2/11/2018	24.08			24.97	
49.05	3/11/2018	23.07			25.98	
49.05	5/14/2018	23.97			25.08	
49.05	1/3/2019	23.17			25.88	
49.05	7/9/2019	23.11			25.943	
49.05	1/7/2020	23.46			25.59	
MW-22A	46.07	11/23/1998	NM			
	46.07	1/29/1999	2.10			43.969

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-22A	46.07	2/26/1999	2.21			43.859
	46.07	3/16/1999	2.65			43.419
	46.07	4/29/1999	2.71			43.359
	46.07	6/1/1999	2.68			43.389
	46.07	7/30/1999	4.12			41.949
	46.07	8/27/1999	2.81			43.259
	46.07	9/27/1999	8.53			37.539
	46.07	10/29/1999	10.23			35.839
	46.07	11/18/1999	9.92			36.149
	46.07	12/29/1999	9.56			36.509
	46.07	2/4/2000	12.31			33.759
	46.07	2/25/2000	8.72			37.349
	46.07	3/27/2000	6.30			39.769
	46.07	4/7/2000	6.03			40.039
	46.07	5/31/2000	6.12			39.949
	46.07	6/1/2000	6.00			40.069
	46.07	7/28/2000	6.13			39.939
	46.07	8/30/2000	9.09			36.979
	46.07	9/19/2000	10.12			35.949
	46.07	10/27/2000	8.64			37.429
	46.07	11/21/2000	7.69			38.379
	46.07	5/1/2001	5.15			40.92
	46.07	10/1/2001	5.49			40.58
	46.07	3/11/2002	2.34			43.729
	46.07	9/23/2002	2.11			43.959
	46.07	3/10/2003	1.68			44.389
	46.07	9/23/2003	2.30			43.769
	46.07	3/15/2004	2.05			44.019
	46.07	9/14/2004	6.89			39.179
	46.07	7/18/2005	3.65			42.419
	46.07	1/6/2006	7.29			38.779
	46.07	7/27/2006	1.65			44.419
	46.07	3/7/2007	NM			
	46.07	7/27/2007	2.84			43.229
	46.07	1/29/2008	1.05			45.019
	46.07	7/14/2008	5.33			40.739
	46.07	2/3/2009	5.24			40.829
	46.07	7/23/2009	5.91			40.159
	46.07	1/9/2010	1.32			44.749
	46.07	7/12/2010	6.52			39.549
	46.07	1/12/2011	3.21			42.859
	46.07	7/11/2011	8.39			37.679
	46.07	1/27/2012	0.98			45.089
	46.07	7/10/2012	1.74			44.326
	46.07	1/8/2013	3.09			42.979
46.07	7/22/2013	NM				
46.07	1/7/2014	3.81			42.26	
46.07	7/15/2014	3.22			42.85	
46.07	1/5/2015	NM				
46.07	8/10/2015	NM				
46.07	1/13/2016	NM				
46.07	7/6/2016	NM				
46.07	1/12/2017	NM				
46.07	7/6/2017	NM				
46.07	9/5/2017	NM		REPLACED		
MW-22AR	45.56	2/11/2018	3.43			42.13
	45.56	3/11/2018	2.24			43.32
	45.56	5/14/2018	4.41			41.15
	45.56	7/2/2018	4.48			41.08
	45.56	1/3/2019	3.67			41.89
	45.56	7/9/2019	3.96			41.60
	45.56	2/10/2020	1.55			44.01
MW-22B	45.86	11/23/1998	2.25			43.606
	45.86	1/29/1999	2.28			43.576
	45.86	2/26/1999	2.34			43.516
	45.86	3/16/1999	2.42			43.436
	45.86	4/29/1999	2.56			43.296

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-22B	45.86	6/1/1999	2.60			43.256
	45.86	7/30/1999	4.31			41.546
	45.86	8/27/1999	2.83			43.026
	45.86	9/27/1999	8.45			37.406
	45.86	10/29/1999	10.11			35.746
	45.86	11/18/1999	9.75			36.106
	45.86	12/29/1999	9.43			36.426
	45.86	2/4/2000	12.56			33.296
	45.86	2/25/2000	8.63			37.226
	45.86	3/27/2000	6.00			39.856
	45.86	4/7/2000	5.64			40.216
	45.86	5/31/2000	5.69			40.166
	45.86	6/1/2000	5.61			40.246
	45.86	7/28/2000	5.67			40.186
	45.86	8/30/2000	8.57			37.286
	45.86	9/19/2000	9.94			35.916
	45.86	10/27/2000	7.03			38.826
	45.86	11/21/2000	7.63			38.226
	45.86	5/1/2001	4.93			40.93
	45.86	10/1/2001	5.40			40.46
	45.86	3/11/2002	1.75			44.106
	45.86	9/23/2002	2.11			43.746
	45.86	3/10/2003	1.02			44.836
	45.86	9/23/2003	2.99			42.866
	45.86	3/15/2004	1.20			44.656
	45.86	9/14/2004	NM			
	45.86	7/18/2005	NM			
	45.86	1/6/2006	7.05			38.806
	45.86	7/27/2006	1.58			44.276
	45.86	3/7/2007	NM			
	45.86	7/27/2007	2.85			43.006
	45.86	1/29/2008	0.85			45.006
	45.86	7/14/2008	5.45			40.406
	45.86	2/3/2009	4.78			41.076
	45.86	7/23/2009	5.39			40.466
	45.86	1/9/2010	3.27			42.586
	45.86	7/12/2010	6.21			39.646
	45.86	1/12/2011	0.37			45.486
	45.86	7/11/2011	8.32			37.536
	45.86	1/27/2012	0.06			45.796
	45.86	7/10/2012	1.27			44.586
	45.86	1/8/2013	NM			
45.86	7/22/2013	NM				
45.86	1/7/2014	4.14			41.716	
45.86	7/15/2014	3.79			42.07	
45.86	1/5/2015	3.87			41.99	
45.86	8/10/2015	2.62			43.24	
45.86	1/13/2016	2.09			43.77	
45.86	7/6/2016	NM				
45.86	1/12/2017	NM				
45.86	7/6/2017	NM				
45.86	9/5/2017	NM		REPLACED		
MW-22BR	45.71	2/11/2018	4.14			41.57
	45.71	3/12/2018	3.29			42.42
	45.71	5/14/2018	5.27			40.44
	45.71	7/2/2018	5.39			40.32
	45.71	1/3/2019	4.29			41.42
	45.71	7/9/2019	4.41			41.3
	45.71	2/10/2020	1.44			44.27
MW-23C	51.91	11/23/1998	27.41			24.504
	51.91	1/29/1999	26.80			25.114
	51.91	2/26/1999	26.88			25.034
	51.91	3/16/1999	26.93			24.984
	51.91	4/29/1999	27.09			24.824
	51.91	6/1/1999	27.00			24.914
	51.91	7/30/1999	29.55			22.364
51.91	8/27/1999	27.29			24.624	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-23C	51.91	9/27/1999	28.40			23.514
	51.91	10/29/1999	29.11			22.804
	51.91	11/17/1999	29.49			22.424
	51.91	12/29/1999	28.46			23.454
	51.91	2/4/2000	28.96			22.954
	51.91	2/25/2000	28.96			22.954
	51.91	3/27/2000	28.61			23.304
	51.91	4/7/2000	27.10			24.814
	51.91	5/31/2000	27.15			24.764
	51.91	6/1/2000	27.11			24.804
	51.91	7/28/2000	27.15			24.764
	51.91	8/30/2000	29.96			21.954
	51.91	9/19/2000	29.77			22.144
	51.91	10/27/2000	28.44			23.474
	51.91	11/21/2000	28.61			23.304
	51.91	5/1/2001	26.26			25.65
	51.91	10/1/2001	26.50			25.41
	51.91	3/11/2002	25.33			26.584
	51.91	9/23/2002	26.43			25.484
	51.91	3/10/2003	24.53			27.384
	51.91	9/23/2003	25.95			25.964
	51.91	3/15/2004	24.15			27.764
	51.91	9/13/2004	25.97			25.944
	51.91	7/18/2005	26.46			25.454
	51.91	1/4/2006	27.53			24.384
	51.91	3/7/2007	23.96			27.954
	51.91	7/27/2007	22.41			29.504
	51.91	1/31/2008	23.22	75.98	1.71	28.694
	48.89 <sup>1</sup>	2/4/2009	22.11	72.05	1.47	26.78
	48.89 <sup>1</sup>	7/23/2009	22.93	73.01	0.51	25.961
	48.89 <sup>1</sup>	1/9/2010	20.29	71.8	1.72	28.601
	48.89 <sup>1</sup>	5/27/2010	22.81	71.5	2.02	26.081
	48.89 <sup>1</sup>	6/28/2010	22.93	72.15	1.37	25.961
	48.89 <sup>1</sup>	7/12/2010	21.41	72.4	1.12	27.481
	48.89 <sup>1</sup>	8/31/2010	21.61	72.65	0.87	27.281
	48.89 <sup>1</sup>	1/12/2011	21.7	71.25	1.45	27.191
	48.89	7/12/2011	23.11	70.65	2.05	25.782
	48.89	1/26/2012	22.81	71.57	1.13	26.082
	48.89	7/9/2012	22.31	71.45	1.25	26.582
	48.89	1/7/2013	23.32	71.06	1.64	25.572
	48.89	7/22/2013	24.38			24.512
	48.89	1/7/2014	23.51	70.8	2.30	25.38
	48.89	7/15/2014	24.06	70.96	2.14	24.83
48.89	1/5/2015	22.47	71.72	1.08	26.42	
48.89	8/10/2015	19.34	72.17	0.63	29.55	
48.89	1/13/2016	23.16	71.91	0.89	25.73	
48.89	7/6/2016	23.09	71.56	1.24	25.80	
48.89	1/12/2017	23.74	71.81	0.99	25.15	
54.16	7/6/2017	23.61	77.27	0.53	30.55	
54.16	9/5/2017	23.67	77.29	0.51	30.49	
54.16	2/7/2018	23.86	77.46	0.34	30.30	
54.16	3/11/2018	23.99	77.41	0.39	30.17	
54.16	5/14/2018	25.02	77.49	0.31	29.14	
54.16	1/3/2019	24.29	77.31	0.49	29.87	
54.16	7/9/2019	24.42	76.99	0.81	29.74	
54.16	1/9/2020	26.59	75.3	1.50	27.57	
MW-24A	45.79	3/27/2000	7.87			37.92
	45.79	4/7/2000	7.63			38.16
	45.79	5/31/2000	7.65			38.14
	45.79	6/1/2000	7.43			38.36
	45.79	7/28/2000	7.60			38.19
	45.79	8/30/2000	10.44			35.35
	45.79	9/19/2000	10.57			35.22
	45.79	10/27/2000	NM			NM
	45.79	11/21/2000	7.09			38.7
	45.79	5/1/2001	6.72			39.07
	45.79	10/1/2001	7.81			37.98
45.79	3/11/2002	3.91			41.88	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)	
MW-24A	45.79	9/23/2002	5.04			40.75	
	45.79	3/10/2003	2.76			43.03	
	45.79	9/23/2003	4.66			41.13	
	45.79	3/15/2004	3.10			42.69	
	45.79	9/14/2004	8.24			37.55	
	45.79	7/18/2005	6.03			39.76	
	45.79	1/6/2006	8.93			36.86	
	45.79	7/27/2006	4.21			41.58	
	45.79	3/7/2007	3.86			41.93	
	45.79	1/30/2008	NM			NM	
	MW-24AR	45.65	2/5/2009	5.18			40.47
		45.65	7/23/2009	7.36			38.29
		45.65	1/9/2010	3.72			41.93
		45.65	7/12/2010	4.29			41.36
		45.65	1/13/2011	3.58			42.07
		45.65	7/11/2011	6.38			39.27
		45.65	1/27/2012	4.59			41.06
		45.65	7/10/2012	4.38			41.27
		45.65	1/8/2013	5.59			40.06
		45.65	7/23/2013	10.14	71.06		35.51
45.65		1/8/2014	7.11			38.54	
45.65		1/5/2015	NM			NM	
45.65		7/6/2016	NM			NM	
MW-24B		46.06	3/27/2000	11.91			34.15
		46.06	4/7/2000	11.60			34.46
	46.06	5/31/2000	11.63			34.43	
	46.06	6/1/2000	11.51			34.55	
	46.06	7/28/2000	11.69			34.37	
	46.06	8/30/2000	13.91			32.15	
	46.06	9/19/2000	14.72			31.34	
	46.06	10/27/2000	12.44			33.62	
	46.06	11/21/2000	11.38			34.68	
	46.06	5/1/2001	10.71			35.35	
	46.06	10/1/2001	11.75			34.31	
	46.06	3/11/2002	9.01			37.05	
	46.06	9/23/2002	9.69			36.37	
	46.06	3/10/2003	7.83			38.23	
	46.06	9/23/2003	8.98			37.08	
	46.06	3/15/2004	7.33			38.73	
	46.06	9/14/2004	9.24			36.82	
	46.06	7/18/2005	9.54			36.52	
	46.06	1/6/2006	11.86			34.2	
	46.06	7/27/2006	10.50			35.56	
	46.06	3/7/2007	8.88			37.18	
	46.06	7/27/2007	9.85			36.21	
	46.06	1/28/2008	7.37			38.69	
	46.06	7/14/2008	11.41			34.65	
	46.06	2/3/2009	11.18			34.88	
	46.06	7/23/2009	12.26			33.8	
	46.06	1/9/2010	9.89			36.17	
	46.06	7/12/2010	12.82			33.24	
	46.06	1/13/2011	11.1			34.96	
	46.06	7/11/2011	14.09			31.97	
	46.06	1/27/2012	11.36			34.7	
	46.06	7/10/2012	10.49			35.57	
46.06	1/8/2013	12.96			33.1		
46.06	7/23/2013	8.49			37.57		
46.06	1/5/2015	NM			NM		
MW-24C	46.05	3/27/2000	25.77			20.28	
	46.05	4/7/2000	24.27			21.78	
	46.05	5/31/2000	24.30			21.75	
	46.05	6/1/2000	24.22			21.83	
	46.05	7/28/2000	24.26			21.79	
	46.05	8/30/2000	27.34			18.71	
	46.05	9/19/2000	26.59			19.46	
	46.05	10/27/2000	27.64			18.41	
	46.05	11/21/2000	25.43			20.62	
46.05	5/1/2001	23.90			22.15		



**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-24C	46.05	10/1/2001	23.71			22.34
	46.05	3/11/2002	22.40			23.65
	46.05	9/23/2002	23.04			23.01
	46.05	3/10/2003	21.71			24.34
	46.05	9/23/2003	23.04			23.01
	46.05	3/15/2004	21.45			24.6
	46.05	9/14/2004	22.45			23.6
	46.05	7/18/2005	22.19			23.86
	46.05	1/6/2006	23.57			22.48
	46.05	7/27/2006	22.61			23.44
	46.05	3/7/2007	21.07			24.98
	46.05	7/27/2007	19.62			26.43
	46.05	1/28/2008	19.43			26.62
	46.05	7/14/2008	20.63			25.42
	46.05	2/3/2009	21.68			24.37
	46.05	7/23/2009	23.07			22.98
	46.05	1/9/2010	20.46			25.59
	46.05	7/12/2010	20.44			25.61
	46.05	1/13/2011	20.26			25.79
	46.05	7/11/2011	21.59			24.46
46.05	1/27/2012	21.23			24.82	
46.05	7/10/2012	20.81			25.24	
46.05	1/8/2013	22.42			23.63	
46.05	7/23/2013	23.81			22.24	
46.05	1/5/2015	NM				
MW-25A	44.65	3/27/2000	9.15			35.5
	44.65	4/7/2000	8.79			35.86
	44.65	5/31/2000	8.81			35.84
	44.65	6/1/2000	8.86			35.79
	44.65	7/28/2000	8.84			35.81
	44.65	8/30/2000	11.43			33.22
	44.65	9/19/2000	11.12			33.53
	44.65	10/27/2000	10.09			34.56
	44.65	11/21/2000	8.10			36.55
	44.65	5/1/2001	8.94			35.71
	44.65	10/1/2001	8.81			35.84
	44.65	3/11/2002	7.23			37.42
	44.65	9/23/2002	5.65			39
	44.65	3/10/2003	5.84			38.81
	44.65	9/23/2003	5.35			39.3
	44.65	3/15/2004	5.75			38.9
	44.65	9/14/2004	7.00			37.65
	44.65	7/18/2005	6.42			38.23
	44.65	1/6/2006	9.29			35.36
	44.65	7/27/2006	5.10			39.55
	44.65	3/7/2007	4.76			39.89
	44.65	7/27/2007	4.22			40.43
	44.65	1/28/2008	4.25			40.4
	44.65	7/14/2008	8.59			36.06
	44.65	2/3/2009	8.90			35.75
	44.65	7/23/2009	8.71			35.94
	44.65	1/9/2010	6.84			37.81
	44.65	7/12/2010	7.78			36.87
	44.65	1/12/2011	6.26			38.39
	44.65	7/11/2011	10.22			34.43
	44.65	1/27/2012	5.24			39.41
	44.65	7/10/2012	4.56			40.09
44.65	1/8/2013	8.62			36.03	
44.65	7/23/2013	9.37			35.28	
44.65	1/8/2014	8.92			35.73	
44.65	7/16/2014	8.61			36.04	
44.65	1/5/2015	8.71			35.94	
44.65	8/10/2015	6.94			37.71	
44.65	1/13/2016	6.07			38.58	
44.65	7/6/2016	6.62			38.03	
44.65	1/12/2017	6.98			37.67	
44.65	7/6/2017	7.31			37.34	
44.65	9/5/2017	7.16			37.49	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-25A	44.65	2/11/2018	5.71			38.94
	44.65	3/12/2018	6.06			38.59
	44.65	5/14/2018	7.49			37.16
	44.65	1/3/2019	6.84			37.81
	44.65	7/9/2019	6.77			37.88
	44.65	1/7/2020	8.01			36.64
MW-25C	44.49	3/27/2000	19.92			24.57
	44.49	4/7/2000	19.50			24.99
	44.49	5/31/2000	19.56			24.93
	44.49	6/1/2000	19.51			24.98
	44.49	7/28/2000	19.54			24.95
	44.49	8/30/2000	22.14			22.35
	44.49	9/19/2000	21.30	66.73	0.90	23.19
	44.49	10/27/2000	20.63			23.86
	44.49	11/21/2000	27.63			16.86
	44.49	5/1/2001	18.14			26.35
	44.49	10/1/2001	18.29		0.40	26.2
	44.49	3/14/2002	17.39	64.32	4.13	27.1
	44.49	9/23/2002	17.81	61.41	6.00	26.68
	44.49	3/10/2003	16.73			27.76
	44.49	9/23/2003	22.35			22.14
	44.49	3/15/2004	16.15			28.34
	44.49	9/14/2004	17.00	60.14	2.56	27.49
	44.49	7/18/2005	15.57			28.92
	44.49	1/6/2006	18.49			26
	44.49	7/27/2006	15.32	60.64	2.03	29.17
	44.49	3/7/2007	15.87	59.82	2.18	28.62
	44.49	7/27/2007	14.25	60.61	1.04	30.24
	44.49	1/28/2008	14.91	60.88	0.67	29.58
	44.49	7/14/2008	17.24	60.95	0.60	27.25
	44.49	2/3/2009	15.97	TRACE	TRACE	28.52
	44.49	7/23/2009	16.39			28.1
	44.49	1/9/2010	13.68	61.45	0.65	30.81
	44.49	5/27/2010	16.09			28.4
	44.49	6/28/2010	16.26			28.23
	44.49	7/12/2010	16.05			28.44
	44.49	8/31/2010	16.21			28.28
	44.49	1/12/2011	16.29			28.2
	44.49	7/11/2011	18.81			25.68
	44.49	1/27/2012	17.29			27.2
	44.49	7/10/2012	16.53			27.96
	44.49	1/8/2013	18.34			26.15
	44.49	7/23/2013	18.74			25.75
	44.49	1/8/2014	18.23			26.26
	44.49	7/16/2014	18.66			25.83
	44.49	1/5/2015	17.81			26.68
	44.49	8/10/2015	16.09			28.40
	44.49	1/13/2016	15.61			28.88
	44.49	7/6/2016	16.02			28.47
	44.49	1/12/2017	16.64			27.85
44.49	7/5/2017	16.84			27.65	
44.49	9/5/2017	16.81			27.68	
44.49	2/11/2018	15.27			29.22	
44.49	3/12/2018	15.63			28.86	
44.49	5/14/2018	16.02			28.47	
44.49	1/3/2019	15.29			29.2	
44.49	7/9/2019	15.86			28.63	
44.49	1/7/2020	16.72			27.77	
MW-26A	44.62	3/27/2000	7.40			37.22
	44.62	4/7/2000	6.99			37.63
	44.62	5/31/2000	7.10			37.52
	44.62	6/1/2000	7.00			37.62
	44.62	7/28/2000	7.11			37.51
	44.62	8/30/2000	9.69			34.93
	44.62	9/19/2000	11.43			33.19
	44.62	10/27/2000	8.11			36.51
	44.62	11/21/2000	8.24			36.38
44.62	5/1/2001	6.01			38.61	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-26A	44.62	10/1/2001	6.34			38.28
	44.62	3/11/2002	4.05			40.57
	44.62	9/23/2002	4.29			40.33
	44.62	3/10/2003	2.84			41.78
	44.62	9/23/2003	4.84			39.78
	44.62	3/15/2004	3.30			41.32
	44.62	9/14/2004	6.80			37.82
	44.62	7/18/2005	6.72			37.9
	44.62	1/6/2006	9.34			35.28
	44.62	7/27/2006	4.42			40.2
	44.62	3/7/2007	4.70			39.92
	44.62	7/27/2007	3.98			40.64
	44.62	1/29/2008	2.37			42.25
	44.62	7/14/2008	7.87			36.75
	44.62	2/3/2009	6.89			37.73
	44.62	7/23/2009	7.88			36.74
	44.62	1/9/2010	4.31			40.31
	44.62	7/12/2010	8.12			36.5
	44.62	1/13/2011	2.38			42.24
	44.62	7/11/2011	10.27			34.35
	44.62	1/27/2012	3.09			41.53
	44.62	7/10/2012	2.77			41.85
	44.62	1/8/2013	7.27			37.35
	44.62	7/23/2013	9.72			34.9
	44.62	1/8/2014	6.33			38.29
	44.62	7/16/2014	7.64			36.98
	44.62	1/5/2015	5.74			38.88
	44.62	8/10/2015	4.03			40.59
	44.62	1/13/2016	3.41			41.21
	44.62	7/6/2016	3.72			40.90
	44.62	1/12/2017	4.92			39.70
	44.62	7/5/2017	5.34			39.28
44.62	9/5/2017	5.27			39.35	
44.62	2/11/2018	4.43			40.19	
44.62	3/12/2018	4.77			39.85	
44.62	5/14/2018	6.61			38.01	
44.62	1/3/2019	6.06			38.56	
44.62	7/9/2019	6.31			38.31	
44.62	1/7/2020	4.78			39.84	
MW-27A	44.90	5/1/2001	6.41			38.49
	44.90	10/1/2001	5.31			39.59
	44.90	3/11/2002	4.21			40.69
	44.90	9/23/2002	3.31			41.59
	44.90	3/10/2003	4.05			40.85
	44.90	9/23/2003	3.24			41.66
	44.90	3/15/2004	2.99			41.91
	44.90	9/14/2004	5.09			39.81
	44.90	7/18/2005	4.45			40.45
	44.90	1/6/2006	4.55			40.35
	44.90	7/27/2006	4.26			40.64
	44.90	3/7/2007	3.01			41.89
	44.90	7/27/2007	2.12			42.92
	44.90	1/28/2008	1.88			43.16
	44.90	7/14/2008	4.57			40.47
	44.90	2/3/2009	4.27			40.77
	44.90	7/23/2009	4.36			40.68
	44.90	1/9/2010	3.69			41.35
	44.90	7/12/2010	5.31			39.73
	44.90	1/12/2011	3.76			41.28
	44.90	7/12/2011	6.72			38.32
	44.90	1/26/2012				NM
	44.90	7/10/2012	well covered			NM
	44.90	1/7/2013	well covered			NM
	44.90	7/23/2013	NM			NM
	44.90	8/10/2015	NM			NM
	44.90	2/11/2018	4.21			40.69
	44.90	3/12/2018	4.59			40.31
44.90	5/14/2018	5.06			39.84	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-27A	44.90	1/3/2019	NM			NM
	44.90	7/9/2019	7.21			37.69
	44.90	1/8/2020	7.22			37.68
MW-27C	45.04	5/1/2001	17.82			27.22
	45.04	10/1/2001	17.82			27.22
	45.04	3/11/2002	16.36			28.68
	45.04	9/23/2002	16.49			28.55
	45.04	3/10/2003	18.68			26.36
	45.04	9/23/2003	16.89			28.15
	45.04	3/15/2004	14.35			30.69
	45.04	9/14/2004	14.49			30.55
	45.04	7/18/2005	16.12			28.92
	45.04	1/6/2006	18.07			26.97
	45.04	7/27/2006	17.13			27.91
	45.04	3/7/2007	15.47			29.57
	44.90	7/27/2007	14.85			30.05
	45.04	1/28/2008	14.31			30.73
	45.04	7/14/2008	17.51			27.53
	45.04	2/3/2009	15.76			29.28
	45.04	7/23/2009	16.38			28.66
	45.04	1/9/2010	14.82			30.22
	45.04	7/12/2010	16.12			28.92
	45.04	1/12/2011	15.84			29.2
	45.04	7/11/2011	18.17			26.87
	45.04	1/27/2012	17.14			27.9
	45.04	7/10/2012	16.56			28.48
	45.04	1/8/2013	17.04			28
	45.04	7/23/2013	18.61			26.43
	45.04	1/8/2014	18.12			26.92
	45.04	7/16/2014	16.94			28.10
	45.04	1/5/2015	17.74			27.30
	45.04	8/10/2015	15.71			29.33
	45.04	1/13/2016	15.04			30.00
	45.04	7/6/2016	15.32			29.72
45.04	1/12/2017	15.91			29.13	
45.04	7/5/2017	16.39			28.65	
45.04	9/5/2017	16.36			28.68	
45.04	2/11/2018	16.59			28.45	
45.04	3/12/2018	16.97			28.07	
45.04	5/14/2018	15.89			29.15	
45.04	1/3/2019	14.32			30.72	
45.04	7/9/2019	15.61			29.43	
45.04	1/8/2020	16.93			28.11	
MW-28A	43.86	5/1/2001	7.45			36.41
	43.86	10/1/2001	8.26			35.6
	43.86	3/11/2002	4.90			38.96
	43.86	9/23/2002	5.71			38.15
	43.86	3/10/2003	3.11			40.75
	43.86	9/23/2003	5.81			38.05
	43.86	9/14/2004	9.34			34.52
	43.86	7/18/2005	7.52			36.34
	43.86	1/6/2006	9.32			34.54
	43.86	7/27/2006	5.54			38.32
	43.86	3/7/2007	5.06			38.8
	43.86	7/27/2007	2.86			41
	43.86	1/29/2008	2.61			41.25
	43.86	7/14/2008	8.74			35.12
	43.86	2/3/2009	8.36			35.5
	43.86	7/23/2009	8.94			34.92
	43.86	1/9/2010	4.54			39.32
	43.86	7/12/2010	8.66			35.2
	43.86	1/12/2011	3.87			39.99
	43.86	7/11/2011	11.43			32.43
43.86	1/27/2012	2.66			41.2	
43.86	7/10/2012	4.52			39.34	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-28A	43.86	1/8/2013	8.11			35.75
	43.86	7/23/2013	10.78			33.08
	43.86	1/8/2014	7.71			36.15
	43.86	7/16/2014	8.19			35.67
	43.86	1/5/2015	7.21			36.65
	43.86	8/10/2015	5.72			38.14
	43.86	1/13/2016	5.09			38.77
	43.86	7/6/2016	5.42			38.44
	43.86	1/12/2017	5.89			37.97
	43.86	7/5/2017	6.13			37.73
	43.86	9/5/2017	6.06			37.80
	43.86	2/11/2018	5.31			38.55
	43.86	3/12/2018	5.61			38.25
	43.86	5/14/2018	6.02			37.84
	43.86	1/3/2019	5.41			38.45
	43.86	7/9/2019	6.52			37.34
43.86	1/7/2020	6.68			37.18	
MW-28C	43.96	5/1/2001	17.14			26.82
	43.96	10/1/2001	17.51			26.45
	43.96	3/11/2002	16.29			27.67
	43.96	9/23/2002	17.75			26.21
	43.96	3/10/2003	15.84			28.12
	43.96	9/23/2003	17.48			26.48
	43.96	3/15/2004	15.56			28.4
	43.96	9/14/2004	17.20			26.76
	43.96	7/18/2005	16.60			27.36
	43.96	1/6/2006	17.61			26.35
	43.96	7/27/2006	17.73			26.23
	43.96	3/7/2007	15.59			28.37
	43.96	7/27/2007	12.90			31.06
	43.96	1/29/2008	14.35			29.61
	43.96	7/14/2008	16.26			27.7
	43.96	2/3/2009	16.03			27.93
	43.96	7/23/2009	16.53			27.43
	43.96	1/9/2010	14.89			29.07
	43.96	7/12/2010	15.89			28.07
	43.96	1/12/2011	18.37			25.59
	43.96	7/11/2011	18.16			25.8
	43.96	1/27/2012	16.12			27.84
	43.96	7/10/2012	16.79			27.17
	43.96	1/8/2013	17.62			26.34
	43.96	7/23/2013	18.87			25.09
	43.96	1/8/2014	17.59			26.37
	43.96	7/16/2014	16.98			26.98
	43.96	1/5/2015	16.84			27.12
	43.96	8/10/2015	14.39			29.57
	43.96	1/13/2016	13.72			30.24
	43.96	7/6/2016	14.03			29.93
	43.96	1/12/2017	14.64			29.32
43.96	7/5/2017	14.88			29.08	
43.96	9/5/2017	14.89			29.07	
43.96	2/11/2018	17.33			26.63	
43.96	3/12/2018	14.73			29.23	
43.96	5/14/2018	16.59			27.37	
43.96	1/3/2019	15.88			28.08	
43.96	7/9/2019	15.03			28.93	
43.96	1/7/2020	15.56			28.4	
MW-29A	46.59	5/1/2001	5.01			41.58
	46.59	10/1/2001	5.38			41.21
	46.59	3/11/2002	1.51			45.08
	46.59	9/23/2002	1.65			44.94
	46.59	3/10/2003	1.42			45.17
	46.59	9/23/2003	1.50			45.09
	46.59	3/15/2004	1.85			44.74
	46.59	9/14/2004	6.35			40.24
	46.59	7/18/2005	3.12			43.47
	46.59	1/6/2006	6.57			40.02

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-29A	46.59	7/27/2006	1.44			45.15
	46.59	3/7/2007	1.95			44.64
	46.59	7/27/2007	2.49			44.1
	46.59	1/28/2008	1.28			45.31
	46.59	7/14/2008	4.14			42.45
	46.59	2/3/2009	3.50			43.09
	46.59	7/23/2009	4.09			42.5
	46.59	1/9/2010	1.76			44.83
	46.59	7/12/2010	3.62			42.97
	46.59	1/13/2011	3.07			43.52
	46.59	7/11/2011	7.14			39.45
	46.59	7/10/2012	4.17			42.42
	46.59	1/8/2013	4.91			41.68
	46.59	7/23/2013	--			--
	Plugged					NM
MW-29B	46.26	5/1/2001	19.01			27.25
	46.26	10/1/2001	19.41			26.85
	46.26	3/11/2002	18.04			28.22
	46.26	9/23/2002	18.82			27.44
	46.26	3/10/2003	17.21			29.05
	46.26	9/23/2003	18.09			28.17
	46.26	3/15/2004	17.10			29.16
	46.26	9/14/2004	17.76			28.5
	46.26	7/18/2005	18.11			28.15
	46.26	1/6/2006	18.83			27.43
	46.26	7/27/2006	18.41			27.85
	46.26	3/7/2007	17.21			29.05
	46.26	7/27/2007	15.49			30.77
	46.26	1/28/2008	15.32			30.94
	46.26	7/14/2008	18.23			28.03
	46.26	2/3/2009	17.72			28.54
	46.26	7/23/2009	16.19			30.07
	46.26	1/9/2010	16.02			30.24
	46.26	7/12/2010	19.29			26.97
	46.26	1/13/2011	17.73			28.53
	46.26	7/11/2011	20.06			26.2
46.26	7/10/2012	9.71			36.55	
46.26	1/8/2013	9.92			36.34	
	Plugged					
MW-29C	46.46	5/1/2001	25.51			20.95
	46.46	10/1/2001	25.04			21.42
	46.46	3/11/2002	23.51			22.95
	46.46	9/23/2002	24.10			22.36
	46.46	3/10/2003	22.71			23.75
	46.46	9/23/2003	23.48			22.98
	46.46	3/15/2004	22.24			24.22
	46.46	9/14/2004	24.12			22.34
	46.46	7/18/2005	23.75			22.71
	46.46	1/6/2006	25.12			21.34
	46.46	7/27/2006	23.35			23.11
	46.46	3/7/2007	22.38			24.08
	46.46	7/27/2007	20.42			26.04
	46.46	1/28/2008	21.08			25.38
	46.46	7/14/2008	22.38			24.08
	46.46	2/3/2009	22.86			23.6
	46.46	7/23/2009	22.81			23.65
	46.46	1/9/2010	20.71			25.75
	46.46	7/12/2010	21.32			25.14
	46.46	1/13/2011	20.39			26.07
	46.46	7/11/2011	23.17			23.29
46.46	7/10/2012	20.69			25.77	
46.46	1/8/2013	21.27			25.19	
46.46	7/23/2013	--			--	
	Plugged					
MW-30A	50.45	3/15/2004	9.71			40.74
	50.45	9/13/2004	12.76			37.69
	50.45	7/18/2005	11.80			38.65
	50.45	1/4/2006	13.52			36.93

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-30A	50.45	7/27/2006	10.45			40
	50.45	3/7/2007	10.98			39.47
	50.45	7/27/2007	9.49			40.96
	50.45	1/30/2008	9.62			40.83
	50.45	7/15/2008	12.52			37.93
	50.45	2/4/2009	13.01			37.44
	50.45	7/23/2009	13.71			36.74
	50.45	1/9/2010	10.87			39.58
	50.45	7/12/2010	12.61			37.84
	50.45	1/12/2011	10.06			40.39
	50.45	7/12/2011	14.76			35.69
	50.45	1/26/2012	10.78			39.67
	50.45	7/9/2012	11.13			39.32
	50.45	1/8/2013	12.91			37.54
	50.45	7/23/2013	14.16			36.29
	50.45	1/8/2014	13.81			36.64
	50.45	7/15/2014	12.10			38.35
50.45	1/5/2015	13.22			37.23	
	50.45	8/10/2015	12.16	Plugged and Abandoned		38.29
MW-31A	52.08	3/15/2004	10.97			41.11
	52.08	9/13/2004	13.00			39.08
	52.08	7/18/2005	13.05			39.03
	52.08	1/4/2006	14.77			37.31
	52.08	7/27/2006	11.83			40.25
	52.08	3/7/2007	12.43			39.65
	52.08	7/27/2007	10.83			41.25
	52.08	1/31/2008	10.99			41.09
	52.08	7/15/2008	13.68			38.4
	52.08	2/4/2009	14.23			37.85
	52.08	7/23/2009	14.73			37.35
	52.08	1/9/2010	12.31			39.77
	52.08	7/12/2010	14.06			38.02
	52.08	1/12/2011	11.62			40.46
	52.08	7/12/2011	15.92			36.16
	52.08	1/26/2012	12.24			39.84
	52.08	7/9/2012	12.79			39.29
52.08	1/8/2013	14.14			37.94	
52.08	7/23/2013	16.24			35.84	
52.08	1/8/2014	15.96			36.12	
52.08	7/15/2014	13.19			38.89	
52.08	1/5/2015	15.16			36.92	
	52.08	8/10/2015	12.76	Plugged and Abandoned		39.32
MW-32A	43.77	3/15/2004	1.00			42.77
	43.77	9/14/2004	6.03	29.00	3.48	37.74
	43.77	7/18/2005	5.82	26.56	5.92	37.95
	43.77	1/6/2006	6.93	24.92	7.57	36.84
	43.77	7/27/2006	12.96	25.71	6.74	30.81
	43.77	3/7/2007	4.03	25.26	7.19	39.74
	43.77	7/27/2007	1.95	30.76	1.70	41.82
	43.77	1/28/2008	2.18			41.59
	43.77	7/14/2008	6.14	26.25	6.20	37.63
	43.77	2/3/2009	5.71	26.29	6.16	38.06
	43.77	7/23/2009	6.29	26.51	5.94	37.48
	43.77	1/9/2010	3.55	25.41	7.04	40.22
	43.77	5/27/2010	5.86	26.2	6.25	37.91
	43.77	6/28/2010	6.02	29.1	3.35	37.75
	43.77	7/12/2010	6.12	29.45	3.00	37.65
	43.77	8/31/2010	5.43	30.67	1.78	38.34
	43.77	1/13/2011	2.63	29.15	3.30	41.14
43.77	7/11/2011	5.92	28.82	3.63	37.85	
MW-32AR	Plugged					37.85
	44.56	1/27/2012	3.22			41.34
	44.56	7/10/2012	3.73			40.83
	44.56	1/8/2013	6.64			37.92
	44.56	7/23/2013	9.42			35.14
	44.56	1/8/2014	5.64			38.92
44.56	7/16/2014	6.74			37.82	
	44.56	1/5/2015				

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-32AR	44.56	8/10/2015	3.18			41.38
	44.56	1/13/2016	2.66			41.90
	44.56	7/6/2016	3.14			41.42
	44.56	1/12/2017	3.67			40.89
	44.56	7/5/2017	4.16			40.40
	44.56	9/6/2017	4.03			40.53
	44.56	2/11/2018	4.06			40.50
	44.56	3/12/2018	5.02			39.54
	44.56	5/14/2018	5.91			38.65
	44.56	1/3/2019	5.42			39.14
	44.56	7/9/2019	6.41			38.15
	44.56	1/8/2020	3.88			40.68
MW-32B	44.41	1/27/2012	3.11	30.52	5.77	41.3
	44.41	7/10/2012	3.81	30.16	6.13	40.6
	44.41	1/8/2013	6.34	30.02	6.38	38.07
	44.41	7/23/2013	7.14			37.27
	44.41	1/8/2014	6.72	34.82	1.58	37.69
	44.41	7/16/2014	6.72	34.29	2.11	37.69
	44.41	1/5/2015	6.02	35.77	0.63	38.39
	44.41	8/10/2015	4.41	36.09	0.31	40.00
	44.41	1/13/2016	3.61	36.07	0.33	40.80
	44.41	7/6/2016	3.91	35.96	0.44	40.50
	44.41	1/12/2017	4.83	36.02	0.38	39.58
	44.41	7/5/2017	4.86	36.13	0.27	39.55
	44.41	9/6/2017	4.78	36.24	3.67	39.63
	44.41	2/7/2018	5.16	36.21	0.19	39.25
	44.41	3/12/2018	5.41	36.13	0.27	39.00
	44.41	5/15/2018	6.47	36.21	0.19	37.94
	44.41	1/3/2019	6.09	36.29	0.11	38.32
	44.41	7/9/2019	5.61	35.89	0.51	38.8
44.41	1/8/2020	2.71			41.7	
MW-33A	44.25	3/15/2004	3.90			40.35
	44.25	9/14/2004	7.85			36.4
	44.25	7/18/2005	6.35			37.9
	44.25	1/6/2006	8.00			36.25
	44.25	7/27/2006	4.73			39.52
	44.25	3/7/2007	5.22			39.03
	44.25	7/27/2007	3.48			40.77
	44.25	1/29/2008	3.34			40.91
	44.25	7/14/2008	7.42	25.19	0.03	36.83
	44.25	2/3/2009	7.28			36.97
	44.25	7/23/2009	7.63			36.62
	44.25	1/9/2010	4.79			39.46
	44.25	7/12/2010	7.61			36.64
	44.25	1/13/2011	3.19			41.06
	44.25	7/11/2011	9.87			34.38
	44.25	1/27/2012	2.69			41.56
	44.25	7/10/2012	3.86			40.39
	44.25	1/8/2013	6.76			37.49
	44.25	7/23/2013	9.83			34.42
	44.25	1/8/2014	6.71			37.54
	44.25	7/16/2014	7.09			37.16
	44.25	1/5/2015	5.02			39.23
	44.25	8/10/2015	4.09			40.16
	44.25	1/13/2016	3.51			40.74
	44.25	7/6/2016	3.89			40.36
	44.25	1/12/2017	5.01			39.24
	44.25	7/5/2017	5.59			38.66
	44.25	9/6/2017	5.51			38.74
44.25	2/11/2018	4.38			39.87	
44.25	3/12/2018	4.86			39.39	
44.25	5/14/2018	6.42			37.83	
44.25	1/3/2019	5.77			38.48	
44.25	7/9/2019	5.09			39.16	
44.25	1/20/2020	4.41			39.84	
MW-33B	44.35	3/7/2007	4.21			40.04
	44.35	7/27/2007	3.72			40.53
	44.35	1/29/2008	2.37	39.12	3.37	41.88



**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)	
MW-33B	44.35	7/14/2008	5.74	37.44	5.05	38.51	
	44.35	2/3/2009	9.28	36.91	5.58	34.97	
	44.35	7/23/2009	NM			NM	
	44.35	1/9/2010	4.61	35.21	7.28	39.74	
	44.35	5/27/2010	6.82			37.53	
	44.35	6/28/2010	6.91			37.44	
	44.35	7/12/2010	7.02			37.33	
	44.35	8/31/2010	7.22			37.13	
	44.35	1/13/2011	3.11	29.7	0.30	41.24	
	44.35	7/11/2011	10.19	29.75	0.25	34.16	
	44.35	1/5/2015	NM			NM	
	MW-33BR	44.35	1/27/2012	4.07			40.28
		44.35	7/10/2012	2.59			41.76
		44.35	1/8/2013	3.86			40.49
		44.35	7/23/2013	9.68			34.67
		44.35	1/8/2014	7.41			36.94
		44.35	7/16/2014	6.72			37.63
		44.35	1/5/2015	5.22			39.13
		44.35	8/10/2015	3.96			40.39
		44.35	1/13/2016	3.22			41.13
		44.35	7/6/2016	3.71			40.64
		44.35	1/12/2017	4.74			39.61
		44.35	7/5/2017	5.19			39.16
		44.35	9/6/2017	4.99			39.36
44.35		2/11/2018	4.74			39.61	
44.35		3/12/2018	5.19			39.16	
44.35		5/14/2018	6.03			38.32	
44.35		1/3/2019	5.18			39.17	
44.35		7/9/2019	5.92			38.43	
44.35		1/8/2020	5.06			39.29	
MW-34C		45.31	3/15/2004	17.40			27.91
		45.31	9/14/2004	18.82			26.49
		45.31	7/18/2005	19.41	65.29	7.19	25.9
		45.31	1/6/2006	20.54	65.27	8.38	24.77
		45.31	7/27/2006	18.55	63.84	8.61	26.76
	45.31	4/9/2007	16.34	62.06	10.39	28.97	
	45.31	7/27/2007	NM				
	45.31	1/29/2008	16.32			28.99	
	45.31	7/15/2008	18.13	43.49	29.01	27.18	
	45.31	2/5/2009	18.08	61.79	10.71	27.23	
	45.31	7/23/2009	NM				
	45.31	1/9/2010	16.41	69.20	3.30	28.9	
	45.31	7/12/2010	NM				
	45.31	1/12/2011	16.41	64.90		28.9	
	45.31	7/11/2011	19.08	65.26		26.23	
	45.31	2/8/2012	18.41			26.9	
	45.31	7/10/2012	NM				
	45.31	1/8/2013	NM				
	45.31	7/23/2013	NM				
	MW-34CR	46.47	7/16/2014	19.17			27.30
		46.47	1/5/2015	19.01			27.46
		46.47	8/10/2015	17.39			29.08
		46.47	1/13/2016	15.99			30.48
		46.47	7/6/2016	16.06			30.41
46.47		1/12/2017	16.94			29.53	
46.47		7/5/2017	17.01			29.46	
46.47		9/6/2017	17.11			29.36	
46.47		2/11/2018	18.19			28.28	
46.47		3/12/2018	18.52			27.95	
46.47		5/14/2018	18.26			28.21	
46.47		1/3/2019	18.26			28.21	
46.47		7/9/2019	NM			damaged	
46.47		3/31/2020	17.49			28.98	
MW-35A	44.75	3/7/2007	3.49			41.82	
	44.75	7/27/2007	3.05			42.26	
	44.75	1/29/2008	1.82			43.49	
	44.75	7/14/2008	6.21			39.1	
	44.75	2/3/2009	5.54			39.77	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-35A	44.75	7/23/2009	5.76			39.55
	44.75	1/9/2010	4.14			41.17
	44.75	7/12/2010	6.04			39.27
	44.75	1/13/2011	2.46			42.85
	44.75	7/11/2011	8.44			36.87
	44.75	1/27/2012	1.35			43.96
	44.75	7/10/2012	2.33			42.98
	44.75	1/8/2013	5.37			39.94
	44.75	7/23/2013	9.18			36.13
	44.75	1/8/2014	5.06			40.25
	44.75	7/15/2014	6.51			38.24
	44.75	1/5/2015	4.22			40.53
	44.75	8/10/2015	3.68			41.07
	44.75	1/13/2016	3.08			41.67
	44.75	7/6/2016	3.34			41.41
	44.75	1/12/2017	3.87			40.88
	44.75	7/5/2017	4.41			40.34
	44.75	9/6/2017	NM			
	44.75	2/11/2018	3.69			41.06
	44.75	3/11/2018	4.06			40.69
44.75	5/14/2018	8.71			36.04	
44.75	1/3/2019	8.06			36.69	
44.75	7/9/2019	7.92			36.83	
44.75	1/8/2020	4.41			40.34	
MW-35B	44.83	3/7/2007	3.31			41.52
	44.83	7/27/2007	3.29			41.54
	44.83	1/29/2008	1.95			42.88
	44.83	7/14/2008	6.40			38.43
	44.83	2/3/2009	5.79			39.04
	44.83	7/23/2009	6.42			38.41
	44.83	1/9/2010	3.51			41.32
	44.83	7/12/2010	6.39			38.44
	44.83	1/13/2011	2.96			41.87
	44.83	7/11/2011	8.67			36.16
	44.83	1/27/2012	1.59			43.24
	44.83	7/10/2012	2.74			42.09
	44.83	1/8/2013	6.09			38.74
	44.83	7/23/2013	9.22			35.61
	44.83	1/8/2014	5.31			39.52
	44.83	7/15/2014	6.75			38.08
	44.83	1/5/2015	4.81			40.02
	44.83	8/10/2015	3.97			40.86
	44.83	1/13/2016	3.26			41.57
	44.83	7/6/2016	3.57			41.26
44.83	1/12/2017	4.06			40.77	
44.83	7/5/2017	4.66			40.17	
44.83	9/6/2017	NM				
44.83	2/11/2018	4.06			40.77	
44.83	3/11/2018	4.31			40.52	
44.83	5/14/2018	6.11			38.72	
44.83	1/3/2019	5.33			39.5	
44.83	7/9/2019	5.62			39.21	
44.83	1/8/2020	4.67			40.16	
MW-36A	44.53	3/7/2007	8.71			35.82
	44.53	7/27/2007	6.54			37.99
	44.53	1/29/2008	5.59			38.94
	44.53	7/14/2008	9.33			35.2
	44.53	2/3/2009	10.69			33.84
	44.53	7/23/2009	12.03			32.5
	44.53	1/9/2010	9.23			35.3
	44.53	7/12/2010	9.14			35.39
	44.53	1/13/2011	8.62			35.91
	44.53	7/11/2011	12.16			32.37
	44.53	1/27/2012	6.82			37.71
	44.53	7/10/2012	6.68			37.85
	44.53	1/8/2013	7.61			36.92
	44.53	7/23/2013	11.36			33.17
	44.53	1/8/2014	9.23			35.3

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-36A	44.53	7/16/2014	8.62			35.91
	44.53	1/5/2015	8.67			35.86
	44.53	8/10/2015	6.47			38.06
	44.53	1/13/2016	5.79			38.74
	44.53	7/6/2016	6.13			38.40
	44.53	1/12/2017	6.58			37.95
	44.53	7/5/2017	7.01			37.52
	44.53	9/6/2017	6.92			37.61
	44.53	2/11/2018	7.77			36.76
	44.53	3/11/2018	8.06			36.47
	44.53	5/14/2018	8.92			35.61
	44.53	1/3/2019	8.22			36.31
	44.53	7/9/2019	8.32			36.21
	44.53	1/7/2020	8.83			35.7
MW-36B	44.07	7/12/2010	1.32			42.75
	44.07	1/13/2011	9.71			34.36
	44.07	7/11/2011	11.57			32.5
	44.07	1/27/2012	0.46			43.61
	44.07	7/10/2012	6.64			37.43
	44.07	1/8/2013	6.71			37.36
	44.07	7/23/2013	9.39			34.68
	44.07	1/8/2014	4.09			39.98
	44.07	7/16/2014	3.61			40.46
	44.07	1/5/2015	3.21			40.86
	44.07	8/10/2015	1.46			42.61
	44.07	1/13/2016	1.06			43.01
	44.07	7/6/2016	4.06			40.01
	44.07	1/12/2017	4.59			39.48
	44.07	7/5/2017	4.72			39.35
	44.07	9/6/2017	4.41			39.66
	44.07	2/11/2018	0.32			43.75
	44.07	3/11/2018	1.81			42.26
	44.07	5/14/2018	1.62			42.45
	44.07	1/3/2019	1.09			42.98
44.07	7/9/2019	1.86			42.21	
44.07	1/7/2020	0.39			43.68	
MW-36D	44.33	7/12/2010	85.39			-41.06
	44.33	1/13/2011	85.03			-40.7
	44.33	7/11/2011	85.33			-41
	44.33	1/27/2012	85.62			-41.29
	44.33	7/10/2012	85.17			-40.84
	44.33	1/8/2013	85.37			-41.04
	44.33	7/23/2013	85.93			-41.6
	44.33	1/8/2014	85.32			-40.99
	44.33	7/16/2014	84.77			-40.44
	44.33	1/5/2015	85.01			-40.68
	44.33	8/10/2015	84.67			-40.34
	44.33	1/13/2016	84.29			-39.96
	44.33	7/6/2016	84.42			-40.09
	44.33	1/12/2017	84.73			-40.40
	44.33	7/5/2017	84.89			-40.56
	44.33	9/6/2017	84.86			-40.53
	44.33	2/11/2018	82.59			-38.26
	44.33	3/11/2018	82.77			-38.44
	44.33	5/14/2018	83.09			-38.76
44.33	1/3/2019	82.51			-38.18	
44.33	7/9/2019	82.91			-38.58	
44.33	1/7/2020	82.06			-37.73	
MW-38A	46.39	3/7/2007	3.26			43.13
	46.39	7/27/2007	3.08			43.31
	46.39	1/29/2008	1.85			44.54
	46.39	7/14/2008	5.84			40.55
	46.39	2/3/2009	5.15			41.24
	46.39	7/23/2009	5.06			41.33
	46.39	1/9/2010	2.27			44.12
	46.39	7/12/2010	6.42			39.97
	46.39	1/13/2011	1.76			44.63
46.39	7/11/2011	8.16			38.23	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-38A	46.39	1/27/2012	1.8			44.59
	46.39	7/10/2012	2.52			43.87
	46.39	1/8/2013	4.62			41.77
	46.39	7/23/2013	8.34			38.05
	46.39	1/8/2014	4.77			41.62
	46.39	7/15/2014	6.20			40.19
	46.39	1/5/2015	4.16			42.23
	46.39	8/10/2015	3.61			42.78
	46.39	1/13/2016	3.02			43.37
	46.39	7/6/2016	3.42			42.97
	46.39	1/12/2017	4.01			42.38
	46.39	7/5/2017	4.21			42.18
	46.39	9/6/2017	4.12			42.27
	46.39	2/11/2018	2.61			43.78
	46.39	3/11/2018	4.12			42.27
	46.39	5/14/2018	5.41			40.98
	46.39	1/3/2019	4.66			41.73
46.39	7/9/2019	4.13			42.26	
46.39	1/8/2020	1.51			44.88	
MW-38B	45.51	3/15/2004	1.07			44.44
	45.51	9/14/2004	6.10			39.41
	45.51	7/18/2005	2.41			43.1
	45.51	1/6/2006	6.33			39.18
	45.51	7/27/2006	1.27			44.24
	45.51	3/7/2007	2.38			43.13
	45.51	7/27/2007	2.25			43.26
	45.51	1/29/2008	0.61			44.9
	45.51	7/14/2008	4.86			40.65
	45.51	2/3/2009	4.33			41.18
	45.51	7/23/2009	4.47			41.04
	45.51	1/9/2010	1.44			44.07
	45.51	7/12/2010	5.72			39.79
	45.51	1/13/2011	0.68			44.83
	45.51	7/11/2011	7.82			37.69
	45.51	1/27/2012	0.85			44.66
	45.51	7/10/2012	0.74			44.77
	45.51	1/8/2013	3.97			41.54
	45.51	7/23/2013	7.51			38
	45.51	1/8/2014	3.47			42.04
	45.51	7/15/2014	5.50			40.01
	45.51	1/5/2015	3.07			42.44
	45.51	8/10/2015	2.17			43.34
	45.51	1/13/2016	2.41			43.10
	45.51	7/6/2016	2.96			42.55
	45.51	1/12/2017	3.81			41.70
	45.51	7/5/2017	4.07			41.44
45.51	9/6/2017	3.91			41.60	
45.51	2/11/2018	2.02			43.49	
45.51	3/11/2018	3.22			42.29	
45.51	5/14/2018	4.62			40.89	
45.51	1/3/2019	3.79			41.72	
45.51	7/9/2019	3.26			42.25	
45.51	1/8/2020	3.79			41.72	
MW-39B	49.58	3/15/2004	5.48			44.1
	49.58	9/13/2004	10.02			39.56
	49.58	7/18/2005	7.21			42.37
	49.58	1/4/2006	10.37			39.21
	49.58	7/27/2006	6.08			43.5
	49.58	3/7/2007	6.91			42.67
	49.58	7/27/2007	5.74			43.84
	49.58	1/30/2008	6.34			43.24
	49.58	7/15/2008	8.96			40.62
	49.58	2/4/2009	8.60			40.98
	49.58	7/24/2009	9.13			40.45
	49.58	1/8/2010	5.61			43.97
	49.58	7/12/2010	9.31			40.27
	49.58	1/12/2011	5.64			43.94
	49.58	7/12/2011	11.97			37.61

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-39B	49.58	1/26/2012	5.84			43.74
	49.58	7/9/2012	5.77			43.81
	49.58	1/7/2013	8.68			40.9
	49.58	7/22/2013	11.17			38.41
	49.58	1/7/2014	7.23			42.35
	49.58	7/16/2014	9.46			40.12
	49.58	1/5/2015	6.71			42.87
	49.58	8/10/2015	4.82			44.76
	49.58	1/13/2016	4.17			45.41
	49.58	7/6/2016	4.26			45.32
	49.58	1/12/2017	5.61			43.97
	49.58	7/5/2017	5.87			43.71
	49.58	9/6/2017	5.66			43.92
	49.58	2/11/2018	6.09			43.49
	49.58	3/11/2018	7.04			42.54
	49.58	5/14/2018	8.73			40.85
	49.58	1/3/2019	7.97			41.61
	49.58	7/9/2019	7.47			42.11
49.58	1/7/2020	6.02			43.56	
MW-40B	49.59	3/15/2004	5.46			44.13
	49.59	9/13/2004	9.72			39.87
	49.59	7/18/2005	7.19			42.4
	49.59	1/4/2006	10.25			39.34
	49.59	7/27/2006	6.18			43.41
	49.59	3/7/2007	6.81			42.78
	49.59	7/27/2007	5.00			44.59
	49.59	1/30/2008	5.23			44.36
	49.59	7/15/2008	8.76			40.83
	49.59	2/4/2009	8.57			41.02
	49.59	7/24/2009	9.06			40.53
	49.59	1/8/2010	5.37			44.22
	49.59	7/12/2010	9.17			40.42
	49.59	1/12/2011	5.81			43.78
	49.59	7/12/2011	11.46			38.13
	49.59	1/26/2012	5.68			43.91
	49.59	7/9/2012	5.74			43.85
	49.59	1/7/2013	8.63			40.96
	49.59	7/22/2013	11.06			38.53
	49.59	1/7/2014	7.24			42.35
	49.59	7/16/2014	9.27			40.32
	49.59	1/5/2015	7.02			42.57
	49.59	8/10/2015	5.02			44.57
	49.59	1/13/2016	4.39			45.20
	49.59	7/6/2016	4.67			44.92
	49.59	1/12/2017	5.22			44.37
	49.59	7/5/2017	5.77			43.82
	49.59	9/6/2017	5.71			43.88
49.59	2/11/2018	6.21			43.38	
49.59	3/11/2018	6.82			42.77	
49.59	5/14/2018	8.44			41.15	
49.59	1/3/2019	7.91			41.68	
49.59	7/9/2019	7.23			42.36	
49.59	1/7/2020	6.17			43.42	
MW-41B	49.37	3/15/2004	4.66			44.71
	49.37	9/13/2004	9.76	35.01	9.80	39.61
	49.37	7/18/2005	5.96	32.23	12.58	43.41
	49.37	1/4/2006	10.03	32.21	12.60	39.34
	49.37	7/27/2006	5.65	29.55	15.26	43.72
	49.37	3/7/2007	4.41	29.13	15.68	44.96
	49.37	7/27/2007	5.27	12.00	32.81	44.1
	49.37	2/22/2008	5.04	25.14	19.67	44.7
	49.37	7/15/2008	8.87	25.09	19.72	40.5
	49.37	2/4/2009	8.93	23.79	21.02	40.44
	49.37	7/24/2009	9.46	23.91	20.90	39.91
	49.37	1/8/2010	5.92	23.65	21.16	43.45
	49.37	5/27/2010	6.13	25.45	19.36	43.24
	49.37	6/28/2010	6.21	38.2	6.61	43.16
	49.37	7/12/2010	6.32	38.45	6.36	43.05

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-41B	49.37	8/31/2010	6.26	39.22	5.59	43.11
	49.37	1/12/2011	6.02	39.6	5.21	43.35
	49.37	7/12/2011	8.86	39.75	5.06	40.51
	49.37	3/8/2012	6.31	20.67	24.14	43.06
	49.37	7/9/2012	8.23			41.14
	49.37	1/7/2013	9.09	41.13	3.68	40.28
	49.37	7/22/2013	10.31	39.29	5.52	39.06
	49.37	1/7/2014	9.06	39.17	5.64	40.31
	49.37	7/15/2014	8.62	37.86	6.95	40.75
	49.37	1/5/2015	8.26	39.02	5.79	41.11
	49.37	8/10/2015	6.01	40.39	4.42	43.36
	49.37	1/13/2016	5.51	39.91	4.90	43.86
	49.37	7/6/2016	5.72	40.01	4.80	43.65
	49.37	1/12/2017	6.39	40.56	4.25	42.98
	49.37	7/6/2017	6.34	40.57	4.24	43.03
	49.37	9/6/2017	6.36	40.62	4.19	43.01
	49.37	2/7/2018	6.97	40.76	4.05	42.40
	49.37	3/11/2018	7.21	40.63	4.18	42.16
	49.37	5/14/2018	8.71	40.82	3.99	40.66
	49.37	7/2/2018	8.97	40.96	3.85	40.4
49.37	1/3/2019	8.22	40.83	3.98	41.15	
49.37	7/9/2019	7.57	40.86	1.74	41.8	
49.37	8/1/2019	7.46	39.42	5.39	41.91	
49.37	1/8/2020	4.2	40.34	2.26	45.17	
MW-42B	50.52	3/7/2007	7.31			43.21
	50.52	7/27/2007	5.74			44.78
	50.52	1/30/2008	6.62			43.9
	50.52	7/15/2008	8.73			41.79
	50.52	2/4/2009	9.32			41.2
	50.52	7/24/2009	9.61			40.91
	50.52	1/8/2010	6.02			44.5
	50.52	7/12/2010	7.13			43.39
	50.52	1/12/2011	6.33			44.19
	50.52	7/12/2011	11.76			38.76
	50.52	1/26/2012	6.62			43.9
	50.52	7/9/2012	6.81			43.71
	50.52	1/7/2013	9.23			41.29
	50.52	7/22/2013	11.08			39.44
	50.52	1/7/2014	8.02			42.5
	50.52	7/15/2014	7.37			43.15
	50.52	1/5/2015	7.31			43.21
	50.52	8/10/2015	5.67			44.85
	50.52	1/13/2016	4.92			45.60
	50.52	7/6/2016	5.36			45.16
50.52	1/12/2017	5.94			44.58	
50.52	7/6/2017	6.27			44.25	
50.52	9/6/2017	6.39			44.13	
50.52	2/11/2018	6.84			43.68	
50.52	3/11/2018	7.12			43.40	
50.52	5/14/2018	8.76			41.76	
50.52	7/2/2018	8.99			41.53	
50.52	1/3/2019	8.02			42.50	
50.52	7/9/2019	7.42			43.1	
50.52	1/7/2020	6.97			43.55	
MW-44A	45.11	3/7/2007	10.86			34.25
	45.11	7/27/2007	7.46			37.65
	45.11	1/30/2008	8.44			36.67
	45.11	7/14/2008	10.75			34.36
	45.11	2/3/2009	12.55			32.56
	45.11	7/23/2009	12.76			32.35
	45.11	1/9/2010	10.23			34.88
	45.11	7/12/2010	11.24			33.87
	45.11	1/12/2011	9.63			35.48
	45.11	7/11/2011	12.59			32.52
	45.11	1/27/2012	9.27			35.84
	45.11	7/10/2012	10.11			35
	45.11	1/8/2013	11.01			34.1
45.11	7/23/2013	12.24			32.87	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-44A	45.11	1/8/2014	11.91			33.2
	45.11	7/16/2014	11.32			33.79
	45.11	1/5/2015	11.27			33.84
	45.11	8/10/2015	9.71			35.40
	45.11	1/13/2016	9.11			36.00
	45.11	7/6/2016	9.26			35.85
	45.11	1/12/2017	9.71			35.40
	45.11	7/5/2017	10.06			35.05
	45.11	9/6/2017	9.94			35.17
	45.11	2/11/2018	8.79			36.32
	45.11	3/11/2018	9.83			35.28
	45.11	5/14/2018	9.91			35.20
	45.11	1/3/2019	9.23			35.88
	45.11	7/9/2019	8.67			36.44
45.11	1/8/2020	10.18			34.93	
MW-44C	45.03	3/15/2004	17.54			27.49
	45.03	9/14/2004	18.35			26.68
	45.03	7/18/2005	18.90	64.77	5.35	26.13
	45.03	1/6/2006	20.03	66.50	5.37	25
	45.03	7/27/2006	18.47	63.35	6.75	26.56
	45.03	3/7/2007	16.02	62.30	7.75	29.01
	45.03	7/27/2007	14.83	65.45	5.50	30.2
	45.03	1/29/2008	15.95			29.08
	45.03	7/14/2008	17.91	64.95	6.18	27.12
	45.03	2/3/2009	16.72	64.15	6.98	28.31
	45.03	7/23/2009	17.12	64.05	6.75	27.91
	45.03	1/9/2010	15.57	63.81	6.99	29.46
	45.03	5/27/2010	16.67	64.7	6.10	28.36
	45.03	6/28/2010	16.77	67.85	2.95	28.26
	45.03	7/12/2010	16.91	70.35	0.45	28.12
	45.03	8/31/2010	16.89	70.63	0.17	28.14
	45.03	1/12/2011	16.77	70.05	0.75	28.26
	45.03	7/11/2011	19.31	70.05	0.75	25.72
	45.03	1/27/2012	17.91	63.88	6.92	27.12
	45.03	7/10/2012	17.61	63.7	7.10	27.42
	45.03	1/8/2013	19.02	62.94	7.86	26.01
	45.03	7/23/2013	20.36	70.26	0.54	24.67
	45.03	1/8/2014	19.67	70.42	0.38	25.36
	45.03	7/16/2014	18.72	69.31	1.49	26.31
	45.03	1/5/2015	18.67	69.82	0.98	26.36
	45.03	8/10/2015	16.31	70.29	0.51	28.72
	45.03	1/13/2016	16.26	69.93	0.87	28.77
	45.03	7/6/2016	16.47	69.71	1.09	28.56
	45.03	1/12/2017	17.22	70.11	0.69	27.81
	45.03	7/5/2017	17.33	70.34	0.46	27.70
	45.03	9/6/2017	17.36	70.43	-0.87	27.67
45.03	2/8/2018	17.77	70.34	0.46	27.26	
45.03	5/15/2018	NM				
45.13	1/4/2019	18.42	70.41	0.39	26.71	
45.13	7/9/2019	18.27			26.86	
45.13	1/8/2020	16.81			28.32	
MW-45C	44.73	3/15/2004	17.15			27.58
	44.73	9/14/2004	17.82	61.66	9.02	26.91
	44.73	7/18/2005	18.38	60.76	9.89	26.35
	44.73	1/6/2006	19.51	62.87	8.87	25.22
	44.73	7/27/2006	17.92	61.64	8.94	26.81
	44.73	3/7/2007	15.95	60.81	9.79	28.78
	44.73	7/27/2007	14.38			30.35
	44.73	1/29/2008	14.86	61.39	9.46	29.87
	44.73	7/14/2008	17.22	61.25	9.88	27.51
	44.73	2/3/2009	17.00	61.24	9.61	27.73
	44.73	7/23/2009	17.46	61.30	9.55	27.27
	44.73	1/9/2010	14.98	61.56	9.29	29.75
	44.73	5/27/2010	16.31	61.1	9.75	28.42
	44.73	6/28/2010	16.42	63.45	7.40	28.31
	44.73	7/12/2010	16.61	68.8	2.05	28.12
	44.73	8/31/2010	16.46	69.62	1.23	28.27
44.73	1/12/2011	16.31	69.1	1.75	28.42	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-45C	44.73	7/11/2011	18.29	69.3	1.55	26.44
	44.73	3/8/2012	16.31	70.6	0.25	28.42
	44.73	7/10/2012	20.69	70.21	0.64	24.04
	44.73	1/8/2013	21.39	69.91	0.69	23.34
	44.73	7/23/2013	22.72	70.39	0.21	22.01
	44.73	1/8/2014	22.13	70.35	0.25	22.6
	44.73	7/16/2014	21.32	69.91	0.69	23.41
	44.73	1/5/2015	20.19	70.55	0.05	24.54
	44.73	8/10/2015	18.61			26.12
	44.73	1/13/2016	17.49			27.24
	44.73	7/6/2016	17.62			27.11
	44.73	1/12/2017	18.22			26.51
	44.73	7/5/2017	17.96			26.77
	44.73	9/6/2017	18.16			26.57
	44.73	2/8/2018	18.62	70.6	0.00	26.11
	44.73	3/11/2018	18.83			25.9
	44.73	5/15/2018	19.61			25.12
	44.73	1/4/2019	19.02			25.71
	44.73	7/9/2019	18.39			26.34
	44.73	1/7/2020	16.24			28.49
MW-46C	44.94	3/15/2004	16.16	ND	ND	28.78
	44.94	9/14/2004	17.97	ND	ND	26.97
	44.94	7/18/2005	18.50	69.05	3.78	26.44
	44.94	1/13/2006	19.66	70.20	3.22	25.28
	44.94	7/27/2006	17.96	68.89	3.90	26.98
	44.94	3/7/2007	16.01	69.32	3.43	28.93
	44.94	7/27/2007	14.54	69.31	3.59	30.4
	44.94	1/30/2008	15.68	70.81	2.00	29.26
	44.94	7/14/2008	17.38	69.97	2.84	27.56
	44.94	2/3/2009	16.78	69.28	3.53	28.16
	44.94	7/23/2009	17.59	69.35	3.55	27.35
	44.94	1/9/2010	14.53	68.74	4.16	30.41
	44.94	5/27/2010	16.26	69.4	3.50	28.68
	44.94	6/28/2010	16.39	70.85	2.05	28.55
	44.94	7/12/2010	16.29	72.25	0.65	28.65
	44.94	8/31/2010	16.13	72.46	0.44	28.81
	44.94	1/12/2011	15.96	71.75	1.15	28.98
	44.94	7/11/2011	18.07	71.65	1.25	26.87
	44.94	1/26/2012	16.54	ND	ND	28.4
	44.94	7/10/2012	20.34	72.8	0.10	24.6
	44.94	1/8/2013	21.18	71.31	1.59	23.76
	44.94	7/23/2013	21.96	72.16	0.74	22.98
	44.94	1/8/2014	21.81	72.55	0.35	23.13
	44.94	7/16/2014	20.86	71.39	1.51	24.08
	44.94	1/5/2015	20.47	72.06	0.84	24.47
	44.94	8/10/2015	18.39	72.42	0.48	26.55
	44.94	1/13/2016	18.24	72.59	0.31	26.70
	44.94	7/6/2016	18.54	72.49	0.41	26.40
	44.94	1/12/2017	19.27	72.46	0.44	25.67
	44.94	7/5/2017	19.12	72.34	0.56	25.82
44.94	9/6/2017	19.29	72.34	0.56	25.65	
44.94	2/8/2018	19.96	72.46	0.44	24.98	
44.94	3/11/2018	20.04	72.32	0.58	24.90	
44.94	5/15/2018	21.02	72.59	0.31	23.92	
44.94	1/4/2019	20.49	72.46	0.44	24.45	
44.94	7/9/2019	18.72			26.22	
44.94	1/7/2020	16.34			28.6	
MW-47A	43.92	3/20/2020	9.26			34.66
MW-47C	45.61	7/27/2007	16.62			28.99
	45.61	1/29/2008	16.04			29.57
	45.61	7/14/2008	18.15			27.46
	45.61	2/4/2009	18.39			27.22
	45.61	7/23/2009	18.61			27
	45.61	1/9/2010	16.46			29.15
	45.61	7/12/2010	18.33			27.28
	45.61	1/12/2011	17.86			27.75
45.61	7/11/2011	19.94			25.67	
45.61	1/26/2012	18.77			26.84	



**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-47C	45.61	7/9/2012	18.17			27.44
	45.61	1/8/2013	19.47			26.14
	45.61	7/23/2013	20.61			25
	45.61	1/8/2014	19.57			26.04
	45.61	7/16/2014	19.02			26.59
	45.61	1/5/2015	19.07			26.54
	45.61	8/10/2015	17.41			28.20
	45.61	1/13/2016	16.83			28.78
	45.61	7/6/2016	17.01			28.60
	45.61	1/12/2017	17.59			28.02
	45.52	7/5/2017	NM			
	45.52	9/6/2017	NM			
	45.52	1/7/2020	17.28			28.24
	MW-48C	44.68	3/15/2004	17.31		
44.68		9/14/2004	18.60			26.08
44.68		7/18/2005	19.17			25.51
44.68		1/6/2006	20.33			24.35
44.68		7/27/2006	18.73			25.95
44.68		3/7/2007	16.52			28.16
44.68		7/27/2007	15.22			29.46
44.68		1/29/2008	16.32			28.36
44.68		7/14/2008	17.63			27.05
44.68		2/4/2009	17.97			26.71
44.68		7/24/2009	18.39			26.29
44.68		1/9/2010	15.81			28.87
44.68		7/12/2010	17.42			27.26
44.68		1/12/2011	17.52			27.16
44.68		7/11/2011	19.58			25.1
44.68		1/26/2012	18.52			26.16
44.68		7/9/2012	17.12			27.56
44.68		1/8/2013	18.26			26.42
44.68		7/23/2013	20.17			24.51
44.68		1/8/2014	19.19			25.49
44.68		7/16/2014	18.38			26.30
44.68		1/5/2015	18.76			25.92
44.68		8/10/2015	16.34			28.34
44.68		1/13/2016	15.72			28.96
44.68		7/6/2016	16.16			28.52
44.68		1/12/2017	16.71			27.97
44.68		7/5/2017	17.17			27.51
44.68		9/6/2017	17.15			27.53
44.68		2/11/2018	17.36			27.32
44.68		3/11/2018	16.74			27.94
44.68	5/14/2018	17.33			27.35	
44.68	1/4/2019	16.67			28.01	
44.68	7/9/2019	18.23			26.45	
44.68	1/7/2020	17.09			27.59	
MW-49A	46.18	3/7/2007	12.91			33.27
	46.18	7/27/2007	8.86			37.32
	46.18	1/31/2008	12.02			34.16
	46.18	7/15/2008	12.99			33.19
	46.18	2/4/2009	13.29			32.89
	46.18	7/24/2009	13.71			32.47
	46.18	1/9/2010	11.07			35.11
	46.18	7/12/2010	11.62			34.56
	46.18	1/12/2011	10.82			35.36
	46.18	7/11/2011	12.31			33.87
	46.18	1/26/2012	9.48			36.7
	46.18	7/9/2012	9.79			36.39
	46.18	1/8/2013	11.31			34.87
	46.18	7/23/2013	11.92			34.26
	46.18	1/8/2014	11.56			34.62
	46.18	7/16/2014	10.57			35.61
	46.18	1/5/2015	16.12			30.06
	46.18	8/10/2015	9.61			36.57
	46.18	1/13/2016	9.34			36.84
	46.18	7/6/2016	9.57			36.61
	46.18	1/12/2017	10.03			36.15
	46.18	7/5/2017	10.32			35.86
	46.18	9/6/2017	10.24			35.94
	46.18	2/11/2018	10.29			35.89

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-49A	46.18	3/11/2018	10.56			35.62
	46.18	5/14/2018	12.34			33.84
	46.18	1/4/2019	11.81			34.37
	46.18	7/9/2019	11.21			34.97
	46.18	1/7/2020	11.12			35.06
MW-49B	46.22	2/4/2009	11.65			34.57
	46.22	7/24/2009	11.93			34.29
	46.22	1/9/2010	9.73			36.49
	46.22	7/12/2010	11.36			34.86
	46.22	1/12/2011	8.04			38.18
	46.22	7/11/2011	12.29			33.93
	46.22	1/26/2012	10.74			35.48
	46.22	7/9/2012	7.38			38.84
	46.22	1/8/2013	11.27	33.56	1.19	34.95
	46.22	7/23/2013	11.83	33.91	0.84	34.39
	46.22	1/8/2014	11.24			34.98
	46.22	7/16/2014	9.62			36.60
	46.22	1/5/2015	10.74			35.48
	46.22	8/10/2015	8.17			38.05
	46.22	1/13/2016	7.74			38.48
	46.22	7/6/2016	8.02			38.20
	46.22	1/12/2017	8.46			37.76
	46.22	7/5/2017	8.72			37.50
	46.22	9/6/2017	8.67			37.55
	46.22	2/11/2018	10.03			36.19
	46.22	3/11/2018	10.64			35.58
46.22	5/14/2018	13.27			32.95	
46.22	1/4/2019	12.59			33.63	
46.22	7/9/2019	12.02	34.62	0.24	34.2	
46.22	1/7/2020	11.51	20.09	12.96	34.71	
MW-50A	46.96	3/7/2007	8.16			38.8
	46.96	7/27/2007	4.70			42.26
	46.96	1/31/2008	5.68			41.28
	46.96	7/16/2008	7.99			38.97
	46.96	2/4/2009	9.31			37.65
	46.96	7/24/2009	9.49			37.47
	46.96	1/9/2010	7.02			39.94
	46.96	7/12/2010	8.74			38.22
	46.96	1/12/2011	5.61			41.35
	46.96	7/11/2011	9.86			37.1
	46.96	1/26/2012	7.21			39.75
	46.96	7/9/2012	4.63			42.33
	46.96	1/8/2013	5.91			41.05
	46.96	7/23/2013	7.13			39.83
	46.96	1/8/2014	6.71			40.25
	46.96	7/16/2014	6.29			40.67
	46.96	1/5/2015	6.22			40.74
	46.96	8/10/2015	5.01			41.95
	46.96	1/13/2016	4.06			42.90
	46.96	7/6/2016	4.71			42.25
	46.96	1/12/2017	5.21			41.75
	46.96	7/5/2017	5.63			41.33
	46.96	9/6/2017	5.51			41.45
46.96	2/11/2018	4.39			42.57	
46.96	3/11/2018	4.81			42.15	
46.96	5/15/2018	5.27			41.69	
46.96	1/3/2019	4.62			42.34	
46.96	7/9/2019	4.21			42.75	
46.96	1/7/2020	7.09			39.87	
MW-50B	45.89	3/12/2020	8.37			37.52
MW-51A	47.80	3/7/2007	6.96			40.84
	47.80	7/27/2007	5.45			42.35
	47.80	1/31/2008	5.92			41.88
	47.80	7/15/2008	NM			
	47.80	2/4/2009	9.98			37.82
	47.80	7/24/2009	10.34			37.46
	47.80	1/9/2010	7.83			39.97
	47.80	7/12/2010	9.16			38.64
	47.80	1/12/2011	8.56			39.24
	47.80	7/11/2011	12.74			35.06
	47.80	1/26/2012	7.33			40.47

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-51A	47.80	7/9/2012	7.26			40.54
	47.80	1/8/2013	7.62			40.18
	47.80	7/23/2013	10.54			37.26
	47.80	1/8/2014	10.21			37.59
	47.80	7/16/2014	8.51			39.29
	47.80	1/5/2015	9.87			39.29
	47.80	8/10/2015	7.96			39.84
	47.80	1/13/2016	7.13			40.67
	47.80	7/6/2016	7.29			40.51
	47.80	1/12/2017	7.63			40.17
	47.80	7/5/2017	7.74			40.06
	47.80	9/6/2017	7.63			40.17
	47.80	2/11/2018	5.92			41.88
	47.80	3/12/2018	6.41			41.39
	47.80	5/15/2018	7.16			40.64
	47.80	1/4/2019	6.67			41.13
	47.80	7/9/2019	6.29			41.51
47.80	1/7/2020	8.47			39.33	
MW-51C	47.48	7/16/2014	22.21			25.27
	47.48	1/5/2015	NM			
	47.48	8/10/2015	18.79			28.69
	47.48	1/13/2016	18.06			29.42
	47.48	7/6/2016	18.26			29.22
	47.48	1/12/2017	18.68			28.80
	47.48	7/5/2017	19.12			28.36
	47.48	9/6/2017	19.02			28.46
	47.48	2/11/2018	17.63			29.85
	47.48	3/12/2018	18.03			29.45
	47.48	5/15/2018	20.83			26.65
	47.48	1/3/2019	20.17			27.31
	47.48	7/9/2019	20.39			27.09
	47.48	1/7/2020	20.34			27.14
MW-52A	51.91	3/7/2007	13.66			38.25
	51.91	7/27/2007	11.76			40.15
	51.91	1/31/2008	12.60			39.31
	51.91	7/15/2008	14.42			37.49
	51.91	2/5/2009	15.52			36.39
	51.91	7/23/2009	16.39			35.52
	51.91	1/9/2010	12.57			39.34
	51.91	7/12/2010	14.19			37.72
	51.91	1/12/2011	9.06			42.85
	51.91	7/12/2011	16.53			35.38
	51.91	1/26/2012	12.99			38.92
	51.91	7/9/2012	12.43			39.48
	51.91	1/7/2013	14.94			36.97
	51.91	7/22/2013	16.29			35.62
	51.91	1/7/2014	16.01			35.9
	51.91	7/15/2014	15.39			36.52
	51.91	1/5/2015	15.37			36.54
	51.91	8/10/2015	13.61			38.30
51.91	1/13/2016	12.96			38.95	
51.91	7/6/2016	NM			NM	
MW-53C	45.49	3/7/2007	16.12			29.37
	45.49	7/27/2007	14.55			30.94
	45.49	1/29/2008	15.12			30.37
	45.49	7/14/2008	16.86			28.63
	45.49	2/3/2009	16.69			28.8
	45.49	7/23/2009	17.62			27.87
	45.49	1/9/2010	15.19			30.3
	45.49	7/12/2010	15.71			29.78
	45.49	1/12/2011	16.58			28.91
	45.49	7/11/2011	18.61			26.88
	45.49	1/27/2012	17.54			27.95
	45.49	7/10/2012	17.73			27.76
	45.49	1/8/2013	18.14			27.35
	45.49	7/23/2013	19.28			26.21
	45.49	1/8/2014	21.12			24.37
	45.49	7/16/2014	17.37			28.12
	45.49	1/5/2015	20.71			24.78
	45.49	8/10/2015	18.72			26.77
45.49	1/13/2016	18.06			27.43	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-53C	45.49	7/6/2016	18.42			27.07
	45.49	1/12/2017	18.89			26.60
	45.49	7/5/2017	19.16			26.33
	45.49	9/6/2017	19.13			26.36
	45.49	2/11/2018	16.43			29.06
	45.49	3/11/2018	15.54			29.95
	45.49	5/14/2018	16.56			28.93
	45.49	1/4/2019	15.93			29.56
	45.49	7/9/2019	15.86			29.63
	45.49	1/7/2020	16.72			28.77
MW-58B	43.59	3/11/2020	13.42			30.17
MW-54C	44.99	3/7/2007	15.74			29.25
	44.99	7/27/2007	14.63			30.36
	44.99	1/28/2008	15.28			29.71
	44.99	7/14/2008	16.68			28.31
	44.99	2/3/2009	16.87			28.12
	44.99	7/23/2009	17.84			27.15
	44.99	1/9/2010	15.46			29.53
	44.99	7/12/2010	16.49			28.5
	44.99	1/12/2011	16.46			28.53
	44.99	7/11/2011	18.23			26.76
	44.99	1/27/2012	17.42			27.57
	44.99	7/10/2012	17.36			27.63
	44.99	1/8/2013	17.81			27.18
	44.99	7/23/2013	18.89			26.1
	44.99	1/8/2014	18.14			26.85
	44.99	7/16/2014	17.49			27.50
	44.99	1/5/2015	17.86			27.13
	44.99	8/10/2015	16.02			28.97
	44.99	1/13/2016	15.33			29.66
	44.99	7/6/2016	15.66			29.33
	44.99	1/12/2017	16.17			28.82
	44.99	7/5/2017	16.61			28.38
	44.99	9/6/2017	16.59			28.40
	44.99	2/11/2018	15.4			29.59
44.99	3/11/2018	15.68			29.31	
44.99	5/14/2018	16.31			28.68	
44.99	1/4/2019	15.71			29.28	
44.99	7/9/2019	15.27			29.72	
44.99	1/7/2020	15.59			29.4	
MW-55A	52.01	2/4/2009	13.79			38.22
	52.01	7/23/2009	14.06			37.95
	52.01	1/9/2010	10.83			41.18
	52.01	7/12/2010	12.72			39.29
	52.01	1/12/2011	10.13			41.88
	52.01	7/12/2011	15.18			36.83
	52.01	1/26/2012	11.71			40.3
	52.01	7/9/2012	12.29			39.72
	52.01	1/7/2013	13.34			38.67
	52.01	7/22/2013	14.19			37.82
	52.01	1/7/2014	12.73			39.28
	52.01	7/15/2014	11.30			40.71
	52.01	1/5/2015	12.51			39.50
52.01	8/10/2015	10.79			41.22	
MW-55B	52.04	1/26/2012	13.28			38.76
	52.04	7/9/2012	13.93			38.11
	52.04	1/7/2013	13.73			38.31
	52.04	7/22/2013	14.59			37.45
	52.04	1/7/2014	12.89			39.15
	52.04	7/15/2014	12.49			39.55
	52.04	1/5/2015	12.41			39.63
	52.04	8/10/2015	10.19			41.85
MW-57A	47.72	2/5/2009	12.73		0.00	34.99
	47.72	7/23/2009	12.91		0.00	34.81
	47.72	1/9/2010	9.78		0.00	37.94
	47.72	7/12/2010	8.56	24.55	2.55	39.16
	47.72	1/12/2011	9.83	22.76	4.14	37.89
	47.72	7/12/2011	13.88	22.79	4.11	33.84
	47.72	1/26/2012	10.54	22.78	4.12	37.18
	47.72	7/9/2012	9.72	22.65	4.25	38
	47.72	1/7/2013	10.61	22.14	4.76	37.11

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-57A	47.72	7/22/2013	13.21	23.05	3.85	34.51
	47.72	1/7/2014	11.79	26.15	0.75	35.93
	47.72	7/15/2014	10.42	26.09	0.81	37.30
	47.72	1/5/2015	10.13	26.75	0.15	37.59
	47.72	8/10/2015	7.46	26.9	0.00	40.26
	47.72	7/6/2016	7.39			40.33
	47.72	1/12/2017	8.07			39.65
	47.72	7/6/2017	8.41			39.31
	47.72	9/6/2017	8.46			39.26
	47.72	2/7/2018	8.98			38.74
	47.72	3/11/2018	9.24			38.48
	47.72	5/14/2018	9.67			38.05
	47.72	1/4/2019	9.52			38.2
	47.72	7/9/2019	10.11			37.61
	47.72	1/8/2020	14.46			33.26
MW-57B	50.90	1/26/2012	28.83	42.51	0.44	22.07
	50.90	7/9/2012	27.93	42.45	0.50	22.97
	50.90	1/7/2013	28.63	41.36	1.59	22.27
	50.90	7/22/2013	16.34	41.67	1.28	34.56
	50.90	1/7/2014	15.04			35.86
	50.90	7/15/2014	15.71			35.19
	50.90	1/5/2015	14.32			36.58
	50.90	8/10/2015	12.42			38.48
	50.90	7/6/2016	12.44			38.46
	50.90	1/12/2017	13.24			37.66
	50.90	7/6/2017	13.57			37.33
	50.90	9/6/2017	13.79			37.11
	50.90	2/7/2018	12.42			38.48
	50.90	3/11/2018	12.62			38.28
	50.90	5/14/2018	13.29			37.61
50.90	1/4/2019	13.03			37.87	
50.90	7/9/2019	13.71			37.19	
50.90	1/8/2020	13.73			37.17	
MW-58A	47.76	2/5/2009	14.55			33.21
	47.76	7/23/2009	14.04			33.72
	47.76	1/9/2010	12.29			35.47
	47.76	7/12/2010	14.03			33.73
	47.76	1/12/2011	11.88			35.88
	47.76	7/12/2011	16.16			31.6
	47.76	1/26/2012	12.26			35.5
	47.76	7/9/2012	11.62			36.14
	47.76	1/7/2013	11.91			35.85
	47.76	7/22/2013	13.71			34.05
	47.76	1/7/2014	13.26			34.5
	47.76	7/15/2014	13.06			34.70
	47.76	1/5/2015	13.06			34.70
	47.76	8/10/2015	11.29			36.47
	47.76	7/6/2016	7.46			40.30
	47.76	1/12/2017	8.04			39.72
	47.76	7/6/2017	8.39			39.37
	47.76	9/6/2017	8.33			39.43
47.76	2/11/2018	6.47			41.29	
47.76	3/11/2018	12.71			35.05	
47.76	5/14/2018	12.94			34.82	
47.76	1/4/2019	12.29			35.47	
47.76	7/9/2019	12.29			35.47	
47.76	1/7/2020	13.17			34.59	
MW-59A	44.18	2/5/2009	10.71			33.47
	44.18	7/23/2009	9.96			34.22
	44.18	1/9/2010	8.62			35.56
	44.18	7/12/2010	9.97			34.21
	44.18	1/12/2011	8.06			36.12
	44.18	7/11/2011	10.54			33.64
	44.18	1/26/2012	6.36			37.82
	44.18	7/9/2012	7.63			36.55
	44.18	1/8/2013	9.09			35.09
	44.18	7/23/2013	9.76			34.42
	44.18	1/8/2014	9.34			34.84
	44.18	7/16/2014	9.17			35.01
	44.18	1/5/2015	8.71			35.47
44.18	8/10/2015	5.76			38.42	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-59A	44.18	1/13/2016	5.01			39.17
	44.18	7/6/2016	5.26			38.92
	44.18	1/12/2017	5.81			38.37
	44.18	7/5/2017	6.14			38.04
	44.18	9/6/2017	6.09			38.09
	44.18	2/11/2018	6.26			37.92
	44.18	3/12/2018	9.13			35.05
	44.18	5/14/2018	8.81			35.37
	44.18	1/4/2019	8.12			36.06
	44.18	7/9/2019	8.92			35.26
	44.18	1/8/2020	9.21			34.97
MW-59B	44.36	7/12/2010	7.43			36.93
	44.36	1/12/2011	6.89			37.47
	44.36	7/11/2011	11.03			33.33
	44.36	1/26/2012	4.44			39.92
	44.36	7/9/2012	7.48			36.88
	44.36	1/8/2013	9.36			35
	44.36	7/23/2013	9.94			34.42
	44.36	1/8/2014	9.52			34.84
	44.36	7/16/2014	8.67			35.69
	44.36	1/5/2015	8.92			35.44
	44.36	8/10/2015	5.91			38.45
	44.36	1/13/2016	5.22			39.14
	44.36	7/6/2016	5.39			38.97
	44.36	1/12/2017	5.97			38.39
	44.36	7/5/2017	6.27			38.09
	44.36	9/6/2017	6.06			38.30
	44.36	2/11/2018	7.59			36.77
	44.36	3/12/2018	9.61			34.75
	44.36	5/14/2018	9.09			35.27
	44.36	1/4/2019	8.27			36.09
44.36	7/9/2019	9.06			35.3	
44.36	1/8/2020	9.17			35.19	
MW-59D	44.22	2/5/2009	84.17			-39.95
	44.22	7/23/2009	83.53			-39.31
	44.22	1/9/2010	81.73			-37.51
	44.22	7/12/2010	82.16			-37.94
	44.22	1/12/2011	82.83			-38.61
	44.22	7/11/2011	82.89			-38.67
	44.22	1/26/2012	82.93			-38.71
	44.22	7/9/2012	82.36			-38.14
	44.22	1/8/2013	82.81			-38.59
	44.22	7/23/2013	83.04			-38.82
	44.22	1/8/2014	83.14			-38.92
	44.22	7/16/2014	82.67			-38.45
	44.22	1/5/2015	82.07			-37.85
	44.22	8/10/2015	81.77			-37.55
	44.22	1/13/2016	81.03			-36.81
	44.22	7/6/2016	81.62			-37.40
	44.22	1/12/2017	82.09			-37.87
	44.22	7/5/2017	82.17			-37.95
	44.22	9/6/2017	82.16			-37.94
	44.22	2/11/2018	81.09			-36.87
44.22	3/12/2018	81.17			-36.95	
44.22	5/14/2018	81.79			-37.57	
44.22	1/4/2019	81.02			-36.8	
44.22	7/9/2019	81.31			-37.09	
44.22	1/8/2020	83.06			-38.84	
MW-60A	46.79	2/4/2009	9.56			37.23
	46.79	7/23/2009	9.71			37.08
	46.79	1/9/2010	7.72			39.07
	46.79	7/12/2010	8.61			38.18
	46.79	1/12/2011	5.82			40.97
	46.79	7/11/2011	9.86			36.93
	46.79	1/26/2012	4.34			42.45
	46.79	7/9/2012	5.42			41.37
	46.79	1/8/2013	6.91			39.88
	46.79	7/23/2013	10.42			36.37
	46.79	1/8/2014	8.06			38.73
46.79	7/16/2014	7.29			39.50	
46.79	1/5/2015	7.39			39.40	

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-60A	46.79	8/10/2015	6.32			40.47
	46.79	1/13/2016	5.67			41.12
	46.79	7/6/2016	6.13			40.66
	46.79	1/12/2017	--			
	46.79	9/6/2017	NM			
	46.79	2/11/2018	3.49			43.30
	46.79	3/12/2018	3.71			43.08
	46.79	5/14/2018	5.19			41.60
	46.79	1/4/2019	4.33			42.46
	46.79	7/9/2019	5.52			41.27
MW-60AR	45.43	3/20/2020	8.03			37.4
MW-60B	45.38	3/17/2020	24.91			20.47
MW-61A	44.67	2/3/2009	8.35			36.32
	44.67	7/23/2009	8.47			36.2
	44.67	1/9/2010	6.49			38.18
	44.67	7/12/2010	8.09			36.58
	44.67	1/12/2011	6.56			38.11
	44.67	7/11/2011	9.67			35
	44.67	1/26/2012	2.48			42.19
	44.67	7/9/2012	4.55			40.12
	44.67	1/8/2013	6.72			37.95
	44.67	7/23/2013	9.16			35.51
	44.67	1/8/2014	7.04			37.63
	44.67	7/16/2014	6.34			38.33
	44.67	1/5/2015	6.52			38.15
	44.67	8/10/2015	4.02			40.65
	44.67	1/13/2016	3.34			41.33
	44.67	7/6/2016	3.97			40.70
	44.67	1/12/2017	4.34			40.33
	44.67	7/5/2017	4.47			40.20
	44.67	9/6/2017	4.39			40.28
	44.67	2/11/2018	5.52			39.15
44.67	3/12/2018	6.62			38.05	
44.67	5/14/2018	6.27			38.40	
44.67	1/4/2019	5.58			39.09	
44.67	7/9/2019	6.96			37.71	
44.67	1/8/2020	4.87			39.8	
MW-61B	43.43	3/17/2020	8.31			35.12
MW-62B	48.16	2/4/2009	6.99			41.17
	48.16	7/24/2009	7.39			40.77
	48.16	1/8/2010	5.13			43.03
	48.16	7/12/2010	5.79			42.37
	48.16	1/12/2011	4.21			43.95
	48.16	7/12/2011	11.06			37.1
	48.16	1/26/2012	3.18			44.98
	48.16	7/9/2012	4.87			43.29
	48.16	1/8/2013	5.92			42.24
	48.16	7/23/2013	7.01			41.15
	48.16	1/8/2014	6.52			41.64
	48.16	7/15/2014	6.06			42.10
	48.16	1/5/2015	6.02			42.14
	48.16	8/10/2015	4.16			44.00
	48.16	1/13/2016	3.64			44.52
	48.16	7/6/2016	4.09			44.07
	48.16	1/12/2017	4.71			43.45
	48.16	7/6/2017	5.09			43.07
	48.16	9/6/2017	4.71			43.45
	48.16	2/11/2018	4.12			44.04
48.16	3/11/2018	5.37			42.79	
48.16	5/14/2018	6.81			41.35	
48.16	7/2/2018	6.92			41.24	
48.16	1/4/2019	6.03			42.13	
48.16	7/9/2019	4.34			43.82	
48.16	1/8/2020	3.09			45.07	
MW-63B	44.48	2/5/2009	31.54			12.94
	44.48	7/23/2009	9.52			34.96
	44.48	1/9/2010	1.34			43.14
	44.48	7/12/2010	5.71			38.77
	44.48	1/13/2011	7.13			37.35
	44.48	7/11/2011	4.21			40.27
44.48	1/27/2012	2.96			41.52	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-63B	44.48	7/10/2012	1.32			43.16
	44.48	1/8/2013	8.54			35.94
	44.48	7/23/2013	9.43			35.05
	44.48	1/8/2014	7.72			36.76
	44.48	7/16/2014	7.03			37.45
	44.48	1/5/2015	7.09			37.39
	44.48	8/10/2015	5.34			39.14
	44.48	1/13/2016	4.69			39.79
	44.48	7/6/2016	5.01			39.47
	44.48	1/12/2017	5.84			38.64
	44.48	7/5/2017	6.19			38.29
	44.48	9/6/2017	6.12			38.36
	44.48	2/11/2018	5.31			39.17
	44.48	3/11/2018	6.39			38.09
	44.48	5/14/2018	7.19			37.29
	44.48	1/4/2019	6.47			38.01
44.48	7/9/2019	3.96			40.52	
44.48	1/7/2020	4.34			40.14	
MW-64A	48.31	2/4/2009	9.02			39.29
	48.31	7/24/2009	9.13			39.18
	48.31	1/9/2010	6.52			41.79
	48.31	7/12/2010	6.82			41.49
	48.31	1/12/2011	4.77			43.54
	48.31	7/12/2011	8.17			40.14
	48.31	1/26/2012	4.81			43.5
	48.31	7/9/2012	5.93			42.38
	48.31	1/7/2013	7.03			41.28
	48.31	7/22/2013	8.79			39.52
	48.31	1/7/2014	8.39			39.92
	48.31	7/15/2014	7.72			40.59
	48.31	1/5/2015	7.79			40.52
	48.31	8/10/2015	5.71			42.60
	48.31	1/13/2016	5.06			43.25
	48.31	7/6/2016	5.67			42.64
	48.31	1/12/2017	6.07			42.24
	48.31	7/6/2017	6.27			42.04
	48.31	9/6/2017	6.16			42.15
	48.31	2/11/2018	5.46			42.85
48.31	3/12/2018	5.83			42.48	
48.31	5/14/2018	6.39			41.92	
48.31	1/4/2019	5.39			42.92	
44.55	7/9/2019	5.09			39.46	
44.55	1/8/2020	4.91			39.64	
MW-65D	44.55	2/5/2009	86.72			-42.17
	44.55	7/23/2009	86.47			-41.92
	44.55	1/9/2010	84.39			-39.84
	44.55	7/12/2010	84.39			-39.84
	44.55	1/12/2011	83.96			-39.41
	44.55	7/11/2011	85.81			-41.26
	44.55	1/27/2012	85.76			-41.21
	44.55	1/8/2013	85.81			-41.26
	44.55	7/23/2013	85.83			-41.28
	44.55	1/8/2014	85.78			-41.23
	44.55	7/16/2014	84.91			-40.36
	44.55	1/5/2015	85.31			-40.76
	44.55	8/10/2015	85.06			-40.51
	44.55	1/13/2016	84.81			-40.26
	44.55	7/6/2016	85.09			-40.54
	44.55	1/12/2017	85.52			-40.97
	44.55	7/5/2017	85.72			-41.17
	44.55	9/6/2017	85.7			-41.15
	44.55	2/11/2018	83.42			-38.87
	44.55	3/12/2018	83.28			-38.73
44.55	5/14/2018	83.74			-39.19	
44.55	1/4/2019	83.03			-38.48	
44.55	7/9/2019	82.71			-38.16	
44.55	1/8/2020	83.29			-38.74	
MW-66D	46.51	2/5/2009	86.18			-39.67
	46.51	7/23/2009	85.82			-39.31
	46.51	1/9/2010	84.02			-37.51
	46.51	7/12/2010	84.86			-38.35



**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-66D	46.51	1/12/2011	NM			
	46.51	7/11/2011	84.93			-38.42
	46.51	1/26/2012	84.88			-38.37
	46.51	7/9/2012	85.02			-38.51
	46.51	1/8/2013	86.09			-39.58
	46.51	7/23/2013	86.42			-39.91
	46.51	1/8/2014	86.09			-39.58
	46.51	7/16/2014	85.26			-38.75
	46.51	1/5/2015	85.42			-38.91
	46.51	8/10/2015	85.21			-38.70
	46.51	1/13/2016	84.71			-38.20
	46.51	7/6/2016	84.86			-38.35
	46.51	1/12/2017	85.26			-38.75
	46.51	7/5/2017	85.66			-39.15
	46.51	9/6/2017	85.67			-39.16
	46.51	2/11/2018	83.28			-36.77
	46.51	3/12/2018	83.37			-36.86
	MW-67B	43.93	7/12/2010	5.76		
43.93		1/13/2011	10.62			33.31
43.93		7/11/2011	17.64			26.29
43.93		1/27/2012	9.87			34.06
43.93		7/10/2012	11.19			32.74
43.93		1/8/2013	11.72			32.21
43.93		7/23/2013	10.69			33.24
43.93		1/8/2014	10.64			33.29
43.93		7/16/2014	11.22			32.71
43.93		1/5/2015	10.22			33.71
43.93		1/13/2016	6.17			37.76
43.93		7/6/2016	6.39			37.54
43.93		1/12/2017	7.04			36.89
43.93		7/5/2017	7.14			36.79
43.93		9/6/2017	6.97			36.96
43.93		2/11/2018	8.89			35.04
43.93		3/12/2018	9.13			34.80
MW-68A		43.24	05/29/219	5.33		
	43.24	7/9/2019	5.01			38.23
	43.24	1/7/2020	4.28			38.96
MW-68B	44.63	1/27/2012	1.16			43.47
	44.63	7/10/2012	3.82			40.81
	44.63	1/8/2013	6.76			37.87
	44.63	7/23/2013	10.33			34.3
	44.63	1/8/2014	5.82			38.81
	44.63	7/16/2014	7.41			37.22
	44.63	1/5/2015	4.32			40.31
	44.63	8/10/2015	3.56			41.07
	44.63	1/13/2016	2.86			41.77
	44.63	7/6/2016	3.07			41.56
	44.63	1/12/2017	3.86			40.77
	44.63	7/5/2017	3.97			40.66
	44.63	9/6/2017	3.84			40.79
	44.63	2/11/2018	3.07			41.56
MW-68C	44.80	7/12/2010	16.52			28.28
	44.80	1/13/2011	16.92			27.88
	44.80	7/11/2011	19.34			25.46
	44.80	1/27/2012	17.66			27.14
	44.80	7/10/2012	17.96			26.84
	44.80	1/8/2013	19.39			25.41
	44.80	7/23/2013	19.87			24.93
				34.21	3.24	39.81

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-68C	44.80	1/8/2014	19.29			25.51
	44.80	7/16/2014	18.39			26.41
	44.80	1/5/2015	18.71			26.09
	44.80	8/10/2015	16.29			28.51
	44.80	1/13/2016	15.74			29.06
	44.80	7/6/2016	15.94			28.86
	44.80	1/12/2017	16.54			28.26
	44.80	7/5/2017	17.02			27.78
	44.80	9/6/2017	17.01			27.79
	44.80	2/11/2018	16.21			28.59
	44.80	3/12/2018	16.88			27.92
	44.80	5/14/2018	17.35			27.45
	44.80	1/4/2019	16.74			28.06
	44.80	7/9/2019	16.39			28.41
	44.80	1/8/2020	16.54			28.26
MW-69A	45.71	7/12/2010	11.81			33.9
	45.71	1/12/2011	11.16			34.55
	45.71	7/11/2011	NM			
	45.71	1/26/2012	10.44			35.27
	45.71	7/9/2012	4.21			41.5
	45.71	1/8/2013	5.31			40.4
	45.71	7/23/2013	7.34			38.37
	45.71	1/8/2014	7.02			38.69
	45.71	7/16/2014	6.34			39.37
	45.71	1/5/2015	6.71			39.00
	45.71	8/10/2015	3.61			42.10
	45.71	1/13/2016	2.91			42.80
	45.71	7/6/2016	3.79			41.92
	45.71	1/12/2017	4.34			41.37
	45.71	7/5/2017	4.59			41.12
	45.71	9/6/2017	4.43			41.28
	45.71	2/11/2018	11.21			34.50
	45.71	3/11/2018	12.58			33.13
45.71	5/14/2018	11.34			34.37	
45.71	1/4/2019	10.61			35.1	
45.71	7/9/2019	9.71			36	
MW-70B	44.86	1/27/2012	6.51	34.26	1.21	38.35
	44.86	7/10/2012	6.06	34.17	1.30	38.8
	44.86	1/8/2013	6.67	34.02	1.68	38.19
	44.86	7/23/2013	8.22	34.07	1.63	36.64
	44.86	1/8/2014	7.89	35.51	0.14	36.97
	44.86	7/16/2014	6.16	34.71	0.94	38.70
	44.86	1/5/2015	7.07	35.26	0.39	37.79
	44.86	8/10/2015	5.26	35.49	0.16	39.60
	44.86	1/13/2016	4.96	35.39	0.26	39.90
	44.86	7/6/2016	5.34	35.31	0.34	39.52
	44.86	1/12/2017	6.17	35.09	0.56	38.69
	44.86	7/5/2017	6.39	35.14	0.51	38.47
	44.86	9/6/2017	6.56	35.34	0.31	38.30
	44.86	2/8/2018	6.42	35.31	0.34	38.44
	44.86	3/12/2018	6.69	35.21	0.44	38.17
	44.86	5/15/2018	7.52	35.39	0.26	37.34
44.86	1/4/2019	6.96	35.31	0.34	37.9	
44.86	7/9/2019	6.57	35.01	0.64	38.29	
MW-70C	43.41	3/12/2020	16.04			27.37
MW-71B	44.59	1/27/2012	7.08			37.51
	44.59	7/10/2012	8.16			36.43
	44.59	1/8/2013	4.09			40.5
	44.59	7/23/2013	8.61			35.98
	44.59	1/8/2014	16.36			28.23
	44.59	7/16/2014	16.02			28.57
	44.59	1/5/2015	15.83			28.76
	44.59	8/10/2015	13.76			30.83
	44.59	1/13/2016	13.09			31.50
	44.59	7/6/2016	13.31			31.28
	44.59	1/12/2017	13.94			30.65
	44.59	7/5/2017	14.34			30.25
	44.59	9/6/2017	14.21			30.38
	44.59	1/25/2018	0.76			43.83
44.59	3/12/2018	1.61			42.98	
44.59	5/14/2018	2.26			42.33	

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-71B	44.59	1/4/2019	1.58			43.01
	44.59	7/9/2019	0.62			43.97
	44.59	1/8/2020	1.12			43.47
MW-72B	51.97	1/26/2012	38.76			13.21
	51.97	7/9/2012	27.27			24.7
	51.97	1/7/2013	20.08			31.89
	51.97	7/22/2013	18.39			33.58
	51.97	1/7/2014	17.31			34.66
	51.97	7/15/2014	16.91			35.06
	51.97	1/5/2015	16.74			35.23
	51.97	8/10/2015	14.59			37.38
	51.97	1/13/2016	13.93			38.04
	51.97	7/6/2016	NM			
	51.97	2/11/2018	12.26			39.71
	51.97	3/12/2018	19.71			32.26
	51.97	5/14/2018	20.92			31.05
	51.97	1/4/2019	20.13			31.84
	51.97	7/9/2019	15.28			36.69
51.97	1/8/2020	14.81			37.16	
MW-73B	51.42	1/26/2012	25.48			25.94
	51.42	7/9/2012	25.03			26.39
	51.42	1/7/2013	26.11			25.31
	51.42	7/22/2013	26.87			24.55
	51.42	1/7/2014	26.19			25.23
	51.42	7/15/2014	25.14			26.28
	51.42	1/5/2015	25.81			25.61
	51.42	8/10/2015	22.46			28.96
			Plugged and Abandoned			
MW-74B	47.58	1/26/2012	7.63			39.95
	47.58	7/9/2012	7.15			40.43
	47.58	1/8/2013	9.62			37.96
	47.58	7/23/2013	11.72			35.86
	47.58	1/8/2014	9.59			37.99
	47.58	7/16/2014	9.01			38.57
	47.58	1/5/2015	9.07			38.51
	47.58	8/10/2015	7.36			40.22
	47.58	1/13/2016	6.86			40.72
	47.58	7/6/2016	7.39			40.19
	47.58	1/12/2017	7.84			39.74
	47.58	7/5/2017	8.17			39.41
	47.58	9/6/2017	8.02			39.56
	47.58	2/11/2018	6.91			40.67
	47.58	3/12/2018	7.22			40.36
	47.58	5/15/2018	8.33			39.25
	47.58	1/4/2019	7.62			39.96
47.58	7/9/2019	8.59			38.99	
47.58	1/8/2020	8.29			39.29	
MW-75B	46.78	1/26/2012	9.07	35.26	1.84	37.71
	46.78	7/9/2012	9.32	35.2	1.90	37.46
	46.78	1/8/2013	10.16	34.13	2.97	36.62
	46.78	7/23/2013	9.74	35.71	1.39	37.04
	46.78	1/8/2014	10.13	36.72	0.43	36.65
	46.78	7/16/2014	11.41	35.71	1.44	35.37
	46.78	1/5/2015	11.33	36.79	0.36	35.45
	46.78	8/10/2015	8.86	37.07	0.08	37.92
	46.78	1/13/2016	7.81	36.84	0.31	38.97
	46.78	7/6/2016	7.8	36.53	0.62	38.98
	46.78	1/12/2017	8.04	36.36	0.79	38.74
	46.78	7/5/2017	8.04	36.36	0.79	38.74
	46.78	9/6/2017	8.22	36.47	3.15	38.56
	46.78	2/8/2018	8.17	36.91	2.71	38.61
	46.78	3/12/2018	8.37	36.94	2.68	38.41
	46.78	5/15/2018	9.22	37.03	2.59	37.56
	46.78	1/4/2019	9.28	36.96	2.66	37.5
	46.78	7/9/2019	8.89			37.89
	46.78	1/8/2020	8.52	33.52	3.73	38.26
46.32	3/12/2020	5.40			40.92	
MW-76C	47.84	7/16/2014	22.68			25.16
	47.84	1/5/2015	23.41			24.43
	47.84	8/10/2015	21.19			26.65
	47.84	1/13/2016	20.81			27.03

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-76C	47.84	7/6/2016	21.09			26.75
	47.84	1/12/2017	21.67			26.17
	47.84	7/5/2017	21.99			25.85
	47.84	9/6/2017	21.93			25.91
	47.84	2/11/2018	20.74			27.10
	47.84	3/12/2018	21.02			26.82
	47.84	5/15/2018	21.46			26.38
	47.84	1/4/2019	20.67			27.17
	47.84	7/9/2019	20.18			27.66
	47.84	1/7/2020	20.68			27.16
MW-77A	49.05	7/16/2014	6.62			42.43
	49.05	1/5/2015	6.27			42.78
	49.05	8/10/2015	4.34			44.71
	49.05	1/13/2016	3.96			45.09
	49.05	7/6/2016	4.29			44.76
	49.05	1/12/2017	4.73			44.32
	49.05	7/5/2017	4.91			44.14
	49.05	9/6/2017	4.78			44.27
	49.05	2/11/2018	7.62			41.43
	49.05	3/12/2018	8.09			40.96
	49.05	5/15/2018	7.06			41.99
	49.05	1/4/2019	6.34			42.71
	49.05	7/9/2019	6.11			42.94
	49.05	1/7/2020	6.81			42.24
MW-78A	48.68	7/16/2014	8.02	28.72	1.38	40.66
	48.68	1/5/2015	9.17	21.17	8.93	39.51
	48.68	8/10/2015	7.34	23.71	6.39	41.34
	48.68	1/13/2016	6.63	21.77	3.58	42.05
	48.68	7/6/2016	6.71	21.97	3.38	41.97
	48.68	1/12/2017	7.42	22.74	2.61	41.26
	48.68	7/5/2017	7.79	23.59	1.76	40.89
	48.68	9/6/2017	7.81	23.48	6.19	40.87
	48.68	2/11/2018	8.29	23.97	1.38	40.39
	48.68	3/12/2018	8.46	23.91	1.44	40.22
	48.68	5/15/2018	9.28	24.07	1.28	39.4
	48.68	1/4/2019	8.78	24.39	0.96	39.9
	48.68	7/9/2019	9.17	24.67	0.68	39.51
	48.68	1/8/2020	7.96	24.9	0.45	40.72
MW-79A	48.95	7/16/2014	7.26			41.69
	48.95	1/5/2015	5.29			43.66
	48.95	8/10/2015	3.71			45.24
	48.95	1/13/2016	3.06			45.89
	48.95	7/6/2016	3.76			45.19
	48.95	1/12/2017	4.06			44.89
	48.95	7/5/2017	4.31			44.64
	48.95	9/6/2017	4.16			44.79
	48.95	2/11/2018	10.82			38.13
	48.95	3/12/2018	11.26			37.69
	48.95	5/15/2018	9.46			39.49
	48.95	1/4/2019	8.8			40.15
	48.95	7/9/2019	9.68			39.27
	48.95	1/8/2020	9.81			39.14
MW-80B	47.11	7/16/2014	5.29			41.82
	47.11	1/5/2015	6.17			40.94
	47.11	8/10/2015	4.33			42.78
	47.11	1/13/2016	3.96			43.15
	47.11	7/6/2016	4.56			42.55
	47.11	1/12/2017	5.06			42.05
	47.11	7/5/2017	5.34			41.77
	47.11	9/6/2017	5.26			41.85
	47.11	2/11/2018	11.34			35.77
	47.11	3/11/2018	11.77			35.34
	47.11	5/15/2018	11.36			35.75
	47.11	1/4/2019	10.71			36.4
	47.11	7/9/2019	11.02			36.09
	47.11	1/7/2020	10.87			36.24
MW-81B	46.77	7/16/2014	6.47			40.30
	46.77	1/5/2015	7.06			39.71
	46.77	8/10/2015	5.22			41.55
	46.77	1/13/2016	4.77			42.00
MW-81B	46.77	7/6/2016	5.16			41.61

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
	46.77	1/12/2017	5.72			41.05
	46.77	7/5/2017	5.96			40.81
	46.77	9/6/2017	5.71			41.06
	46.77	2/11/2018	7.04			39.73
	46.77	3/11/2018	7.51			39.26
	46.77	5/15/2018	8.23			38.54
	46.77	1/4/2019	7.67			39.1
	46.77	7/9/2019	7.91			38.86
	46.77	1/7/2020	7.12			39.65
MW-82B	44.64	2/11/2018	2.53			42.11
	44.64	3/11/2018	3.44			41.20
	44.64	5/14/2018	5.61			39.03
	44.64	1/4/2019	4.83			39.81
	44.64	7/9/2019	4.03			40.61
	44.64	1/8/2020	2.67			41.97
MW-83B	45.33	2/11/2018	4.06			41.27
	45.33	3/11/2018	4.69			40.64
	45.33	5/14/2018	7.47			37.86
	45.33	7/19/2018	5.87			39.46
	45.33	1/4/2019	6.82			38.51
	45.33	7/9/2019	6.11			39.22
	45.33	1/8/2020	5.33			40
MW-83C	45.42	2/11/2018	17.52			27.90
	45.42	3/11/2018	16.96			28.46
	45.42	5/14/2018	18.11			27.31
	45.42	1/4/2019	17.42			28.00
	45.42	7/9/2019	17.17			28.25
	45.42	1/8/2020	17.74			27.68
MW-84A	43.01	3/12/2020	4.19			38.82
MW-84B	44.50	2/11/2018	4.37			40.13
	44.50	3/11/2018	4.93			39.57
	44.50	5/14/2018	7.36			37.14
	44.50	7/19/2018	6.07			38.43
	44.50	1/4/2019	6.71			37.79
	44.50	7/9/2019	7.09			37.41
	44.50	1/7/2020	4.38			40.12
MW-85C	49.10	2/11/2018	22.51			26.59
	49.10	3/11/2018	22.77			26.33
	49.10	5/15/2018	22.61			26.49
	49.10	1/4/2019	21.92			27.18
	49.10	7/9/2019	21.78			27.32
	49.10	1/7/2020	22.70			26.4
MW-86C	46.61	2/11/2018	20.14			26.47
	46.61	3/11/2018	19.91			26.70
	46.61	5/15/2018	20.26			26.35
	46.61	1/4/2019	19.51			27.10
	46.61	7/9/2019	19.73			26.88
	46.61	1/8/2020	21.09			25.52
MW-87C	44.26	2/11/2018	15.86			28.40
	44.26	3/11/2018	16.29			27.97
	44.26	5/14/2018	16.26			28.00
	44.26	1/4/2019	15.52			28.74
	44.26	7/9/2019	15.83			28.43
	44.26	1/8/2020	15.79			28.47
MW-88A	49.83	3/17/2020	8.00			41.83
MW-88B	49.91	3/17/2020	7.82			42.09
MW-88C	51.17	2/11/2018	24.7			26.47
	51.17	3/11/2018	23.93			27.24
	51.17	5/14/2018	24.67			26.50
	51.17	1/4/2019	24.01			27.16
	51.17	7/9/2019	24.13			27.04
	51.17	1/7/2020	24.64			26.53
MW-89B	44.57	7/19/2018	6.78			37.79
	44.57	1/4/2019	8.21			36.36
	44.57	7/9/2019	7.74			36.83
	44.57	1/7/2020	6.24			38.33
MW-90B	44.39	7/19/2018	5.63			38.76
	44.39	1/4/2019	7.16			37.23
	44.39	7/9/2019	4.13			40.26
	44.39	1/7/2020	0.86			43.53
MW-91A	42.36	3/12/2020	5.21			37.15

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
MW-92B	43.25	3/12/2020	4.19			39.06
MW-93B	43.39	3/12/2020	4.06			39.33
MW-94A	43.55	3/12/2020	3.99			39.56
MW-95A	44.53	3/17/2020	3.88			40.65
MW-96B	45.36	3/17/2020	4.28			41.08
MW-97A	46.21	3/12/2020	5.20			41.01
MW-98A	46.69	3/12/2020	7.19			39.5
MW-98B	46.93	3/12/2020	8.30			38.63
MW-99C	43.67	3/11/2020	14.99			28.68
P-10	47.69	9/2/1993	6.87			40.85
	47.69	12/21/1993	3.32			44.4
	47.69	3/24/1994	3.88			43.84
	47.69	6/22/1994	4.98			42.74
	47.69	9/28/1994	6.38			41.34
	47.69	10/13/1994	7.07			40.65
	47.69	1/24/1995	2.67			45.05
	47.69	4/11/1995	2.59			45.13
	47.69	7/11/1995	4.69			43.03
	47.69	1/23/1996	5.84			41.88
	47.69	7/19/1996	10.04			37.68
	47.69	9/17/1996	8.34			39.38
	47.69	10/31/1996	6.97			40.75
	47.69	11/22/1996	8.84			38.88
	47.69	12/27/1996	6.20			41.52
	47.69	1/22/1997	4.10			43.62
	47.69	2/21/1997	2.86			44.86
	47.69	3/25/1997	3.19			44.53
	47.69	4/23/1997	4.42			43.3
	47.69	4/24/1997	4.57			43.15
	47.69	5/13/1997	3.14			44.58
	47.69	6/20/1997	4.94			42.78
	47.69	6/25/1997	2.74			44.98
	47.69	7/1/1997	4.13			43.59
	47.69	7/24/1997	7.91			39.81
	47.69	8/16/1997	7.86			39.86
	47.69	8/22/1997	8.67			39.05
	47.69	9/25/1997	6.54			41.18
	47.69	10/22/1997	5.36			42.36
	47.69	11/25/1997	5.36			42.36
	47.69	12/19/1997	4.72			43
	47.69	1/20/1998	3.40			44.32
	47.69	1/29/1998	3.11			44.61
	47.69	3/18/1998	2.84			44.88
	47.69	4/24/1998	6.80			40.92
	47.69	5/21/1998	7.35			40.37
	47.69	7/30/1998	8.23			39.49
	47.69	8/25/1998	7.34			40.38
	47.69	9/21/1998	5.25			42.47
	47.69	10/26/1998	6.11			41.61
	47.69	11/23/1998	4.10			43.62
	47.69	2/26/1999	3.21			44.51
	47.69	3/16/1999	4.21			43.51
	47.69	4/29/1999	4.53			43.19
	47.69	6/1/1999	4.53			43.19
	47.69	7/30/1999	6.00			41.72
	47.69	8/27/1999	4.72			43
	47.69	9/27/1999	9.58			38.14
	47.69	10/29/1999	10.61			37.11
	47.69	12/29/1999	11.55			36.17
	47.69	2/4/2000	13.71			34.01
	47.69	2/25/2000	10.44			37.28
	47.69	3/27/2000	7.53			40.19
	47.69	4/7/2000	7.09			40.63
	47.69	5/31/2000	7.14			40.58
	47.69	6/1/2000	7.11			40.61
	47.69	7/28/2000	7.15			40.57
	47.69	8/30/2000	10.15			37.57
	47.69	9/19/2000	11.56			36.16
	47.69	10/27/2000	8.66			39.06
	47.69	11/21/2000	9.64			38.08
P-10	47.69	5/1/2001	6.52			41.2

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
	47.69	10/1/2001	6.85			40.87
	47.69	3/11/2002	3.41			44.31
	47.69	9/23/2002	3.54			44.18
	47.69	3/10/2003	2.43			45.26
	47.69	9/23/2003	1.61			46.08
	47.69	3/15/2004	2.85			44.84
	47.69	9/13/2004	7.99			39.7
	47.69	7/18/2005	4.20			43.49
	47.69	1/4/2006	8.58			39.11
	47.69	7/27/2006	3.46			44.23
	47.69	1/23/2007	2.36			45.33
	47.69	3/7/2007	NM			
	47.69	7/27/2007	3.75			43.94
	47.69	1/29/2008	2.30			45.39
	47.69	7/16/2008	6.91			40.78
	47.69	1/22/2009	6.35			41.34
	47.69	7/23/2009	NM			
	47.69	1/8/2010	4.06			43.63
	47.69	7/12/2010	2.06			45.63
	47.73	1/12/2011	4.13			43.60
	47.73	7/12/2011	9.84			37.89
	47.73	1/27/2012	3.12			44.61
	47.73	7/10/2013	10.79			36.94
	47.73	1/8/2014	5.51			42.22
	47.73	7/2/2014	7.74			39.99
	47.73	1/7/2015	3.96			43.77
	47.73	8/10/2015	5.39			42.34
	47.71	1/12/2016	2.47			45.24
	47.71	7/6/2016	5.18			42.53
	47.71	1/12/2017	4.52			43.19
	47.71	7/12/2017	6.07			41.64
	47.71	1/3/2018	6.71			41.00
	47.71	7/19/2018	5.77			41.94
	47.71	1/3/2019	6.32			41.39
	47.71	7/1/2019	3.12			44.59
	47.71	1/13/2020	3.18			44.53
P-11	48.98	9/2/1993	7.87			41.15
	48.98	12/21/1993	4.57			44.45
	48.98	3/24/1994	5.04			43.98
	48.98	6/22/1994	6.19			42.83
	48.98	9/28/1994	7.40			41.62
	48.98	10/13/1994	8.14			40.88
	48.98	1/24/1995	3.90			45.12
	48.98	4/11/1995	3.77			45.25
	48.98	7/11/1995	5.69			43.33
	48.98	1/23/1996	6.81			42.21
	48.98	7/19/1996	7.81			41.21
	48.98	9/17/1996	9.15			39.87
	48.98	10/31/1996	7.52			41.5
	48.98	11/22/1996	9.46			39.56
	48.98	12/27/1996	6.64			42.38
	48.98	1/22/1997	4.70			44.32
	48.98	2/21/1997	3.88			45.14
	48.98	3/25/1997	4.09			44.93
	48.98	4/23/1997	5.27			43.75
	48.98	4/24/1997	5.41			43.61
	48.98	5/13/1997	4.12			44.9
	48.98	6/20/1997	5.79			43.23
	48.98	6/25/1997	3.83			45.19
	48.98	7/1/1997	5.01			44.01
	48.98	7/24/1997	7.56			41.46
	48.98	8/16/1997	8.74			40.28
	48.98	8/22/1997	9.37			39.65
	48.98	9/25/1997	7.24			41.78
	48.98	10/22/1997	5.98			43.04
	48.98	11/25/1997	6.00			43.02
	48.98	12/19/1997	5.52			43.5
	48.98	1/20/1998	4.30			44.72
	48.98	3/4/1998	4.08			44.94
	48.98	3/18/1998	3.92			45.1
P-11	48.98	4/24/1998	7.61			41.41

**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
	48.98	5/21/1998	8.10			40.92
	48.98	7/30/1998	9.21			39.81
	48.98	8/25/1998	8.44			40.58
	48.98	9/21/1998	5.91			43.11
	48.98	10/26/1998	7.59			41.43
	48.98	11/23/1998	5.41			43.61
	48.98	1/29/1999	4.11			44.91
	48.98	2/26/1999	4.22			44.8
	48.98	3/16/1999	4.96			44.06
	48.98	4/29/1999	5.15			43.87
	48.98	6/1/1999	5.15			43.87
	48.98	7/30/1999	6.66			42.36
	48.98	8/27/1999	5.23			43.79
	48.98	9/27/1999	10.49			38.53
	48.98	10/29/1999	11.91			37.11
	48.98	12/29/1999	11.12			37.9
	48.98	2/4/2000	12.13			36.89
	48.98	2/25/2000	10.46			38.56
	48.98	3/27/2000	8.32			40.7
	48.98	4/7/2000	7.91			41.11
	48.98	5/31/2000	7.96			41.06
	48.98	6/1/2000	7.93			41.09
	48.98	7/28/2000	7.97			41.05
	48.98	8/30/2000	10.88			38.14
	48.98	9/19/2000	12.32			36.7
	48.98	10/27/2000	10.94			38.08
	48.98	11/21/2000	9.77			39.25
	48.98	5/1/2001	7.48			41.54
	48.98	10/1/2001	7.74			41.28
	48.98	3/11/2002	4.51			44.51
	48.98	9/23/2002	4.46			44.56
	48.98	3/10/2003	3.69			45.29
	48.98	9/23/2003	4.54			44.44
	48.98	3/15/2004	4.51			44.47
	48.98	9/13/2004	9.14			39.84
	48.98	7/18/2005	5.27			43.71
	48.98	1/4/2006	9.56			39.42
	48.98	7/27/2006	4.54			44.44
	48.98	3/7/2007	NM			
	48.98	7/27/2007	4.61			44.37
	48.98	1/30/2008	2.71			46.27
	48.98	7/15/2008	7.93			41.05
	48.98	2/4/2009	7.82			41.16
	48.98	7/24/2009	7.74			41.24
	48.98	1/8/2010	5.67			43.31
	48.98	7/12/2010	6.78			42.2
	48.98	1/12/2011	4.21			44.77
	48.98	7/12/2011	11.51			37.47
	48.98	1/26/2012	4.25			44.73
	48.98	1/7/2013	7.96			41.02
	48.98	7/22/2013	10.96			38.02
	48.98	1/7/2014	6.52			42.46
	48.98	7/16/2014	8.87			40.11
	48.98	1/5/2015	5.61			43.37
	48.98	8/10/2015	3.86			45.12
	48.98	1/13/2016	3.26			45.72
	48.98	7/6/2016	3.74			45.24
	48.98	1/12/2017	4.36			44.62
	48.98	7/6/2017	4.62			44.36
	48.98	9/6/2017	4.62			44.36
	48.98	2/11/2018	5.09			43.89
	48.98	3/11/2018	5.54			43.44
	48.98	5/14/2018	7.14			41.84
	48.98	7/2/2018	7.28			41.7
	48.98	1/4/2019	6.43			42.55
	48.98	7/9/2019	5.31			43.67
	48.98	1/7/2020	4.26			44.72
P-12	48.78	9/2/1993	7.02			41.8
	48.78	12/21/1993	4.30			44.52
	48.78	3/24/1994	4.45			44.37
P-12	48.78	6/22/1994	5.06			43.76



**Table 1**  
**GROUNDWATER MEASUREMENTS**  
**Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
	48.78	9/28/1994	6.46			42.36
	48.78	10/13/1994	7.19			41.63
	48.78	1/24/1995	3.63			45.19
	48.78	4/11/1995	3.25			45.57
	48.78	7/11/1995	4.62			44.2
	48.78	1/23/1996	6.62			42.2
	48.78	7/19/1996	8.64			40.18
	48.78	9/17/1996	8.12			40.7
	48.78	10/31/1996	6.81			42.01
	48.78	11/22/1996	8.70			40.12
	48.78	12/27/1996	6.57			42.25
	48.78	1/22/1997	4.93			43.89
	48.78	2/21/1997	3.61			45.21
	48.78	3/25/1997	3.70			45.12
	48.78	4/23/1997	4.58			44.24
	48.78	4/24/1997	4.74			44.08
	48.78	5/13/1997	3.69			45.13
	48.78	6/20/1997	4.86			43.96
	48.78	6/25/1997	3.35			45.47
	48.78	7/1/1997	4.11			44.71
	48.78	7/24/1997	6.58			42.24
	48.78	8/16/1997	7.80			41.02
	48.78	8/22/1997	8.22			40.6
	48.78	9/25/1997	6.54			42.28
	48.78	10/22/1997	5.66			43.16
	48.78	11/25/1997	5.70			43.12
	48.78	12/19/1997	5.13			43.69
	48.78	1/20/1998	4.15			44.67
	48.78	3/4/1998	3.78			45.04
	48.78	3/18/1998	3.61			45.21
	48.78	4/24/1998	6.90			41.92
	48.78	5/21/1998	7.80			41.02
	48.78	7/30/1998	8.15			40.67
	48.78	8/25/1998	8.31			40.51
	48.78	9/21/1998	5.64			43.18
	48.78	10/26/1998	7.66			41.16
	48.78	11/23/1998	5.65			43.17
	48.78	1/29/1999	4.20			44.62
	48.78	2/26/1999	4.31			44.51
	48.78	3/16/1999	4.99			43.83
	48.78	4/29/1999	5.10			43.72
	48.78	6/1/1999	5.10			43.72
	48.78	7/30/1999	6.75			42.07
	48.78	8/27/1999	5.34			43.48
	48.78	9/27/1999	9.36			39.46
	48.78	10/29/1999	10.11			38.71
	48.78	12/29/1999	9.44			39.38
	48.78	2/4/2000	12.10			36.72
	48.78	2/25/2000	8.63			40.19
	48.78	3/27/2000	7.76			41.06
	48.78	4/7/2000	7.35			41.47
	48.78	5/31/2000	7.39			41.43
	48.78	6/1/2000	7.34			41.48
	48.78	7/28/2000	7.37			41.45
	48.78	8/30/2000	10.66			38.16
	48.78	9/19/2000	11.45			37.37
	48.78	10/27/2000	10.94			37.88
	48.78	11/21/2000	8.93			39.89
	48.78	5/1/2001	6.70			42.12
	48.78	10/1/2001	6.93			41.89
	48.78	3/1/2002	4.15			44.67
	48.78	9/23/2002	3.90			44.92
	48.78	3/10/2003	3.13			45.65
	48.78	9/23/2003	3.86			44.92
	48.78	3/15/2004	NM			
	48.78	9/13/2004	7.93			40.85
	48.78	7/18/2005	5.06			43.72
	48.78	1/4/2006	8.98			39.8
	48.78	7/27/2006	4.35			44.43
	48.78	1/22/2007	3.19			45.59
P-12	48.78	3/7/2007	NM			

**Table 1  
GROUNDWATER MEASUREMENTS  
Houston, TX - Wood Preserving Works**

Well ID	TOC Elevation (ft)	Date	Depth to Water (ft)	Depth to DNAPL (ft BTOC)	DNAPL Thickness (ft)	GW Elevation (ft)
	48.78	7/27/2007	4.22			44.56
	48.78	1/29/2008	3.03			45.75
	48.78	7/16/2008	6.78			42
	48.78	1/22/2009	6.99			41.79
	48.78	7/24/2009	NM			
	48.78	1/8/2010	4.13			44.65
	48.78	7/12/2010	3.93			44.85
	48.80	1/12/2011	4.83			43.97
	48.80	7/12/2011	10.02			38.78
	48.80	1/27/2012	4.52			44.28
	48.80	7/9/2012	5.15			43.65
	48.80	7/10/2013	9.73			39.07
	48.80	1/8/2014	6.41			42.39
	48.80	7/2/2014	6.46			42.34
	48.80	1/7/2015	3.19			45.61
	48.80	8/10/2015	4.06			44.74
	48.76	1/12/2016	3.26			45.50
	48.76	7/6/2016	5.09			43.67
	48.76	1/12/2017	5.11			43.65
	48.76	7/12/2017	6.39			42.37
	48.76	1/3/2018	7.14			41.62
	48.76	7/19/2018	6.31			42.45
	48.76	1/3/2019	6.69			42.07
	48.76	7/1/2019	3.06			45.70
	48.76	1/14/2020	3.96			44.80
TW-41B	49.67	2/4/2009	8.44			41.23
	49.67	7/24/2009	8.34			41.33
	49.67	1/8/2010	4.86			44.81
	49.67	7/12/2010	6.12			43.55
	49.67	1/12/2011	5.17			44.5
	49.67	7/12/2011	12.02			37.65
	49.67	1/26/2012	5.27			44.4
	49.67	7/9/2012	6.23			43.44
	49.67	1/7/2013	8.54			41.13
	49.67	7/22/2013	11.53			38.14
	49.67	1/7/2014	7.32			42.35
	49.67	7/16/2014	9.65			40.02
	49.67	1/5/2015	NM			
	49.67	8/10/2015	4.96			44.71
	49.67	1/13/2016	4.13			45.54
	49.67	7/6/2016	4.31			45.36
	49.67	1/12/2017	4.93			44.74
	49.67	7/6/2017	5.32			44.35
	49.67	9/6/2017	5.26			44.41
	49.67	2/11/2018	5.86			43.81
	49.67	3/11/2018	6.69			42.98
	49.67	5/14/2018	8.67			41.00
	49.67	7/2/2018	8.87			40.8
	49.67	1/4/2019	7.97			41.7
	49.67	7/9/2019	8.22			41.45
	49.67	1/7/2020	5.52			44.15
TW-55A	49.67	7/9/2012	13.44			36.23
TW-56A	51.89	2/5/2009	17.48			34.41
	51.89	7/23/2009	17.17			34.72
	51.89	1/8/2010	14.53			37.36
	51.89	7/12/2010	15.78			36.11
	51.89	1/12/2011	14.09			37.8
	51.89	7/12/2011	17.89			34
	51.89	1/26/2012	15.06			36.83
	51.89	1/7/2013	16.92			34.97
	51.89	7/22/2013	18.12			33.77
	51.89	1/7/2014	NM			
	51.89	7/15/2014	16.05			35.84
	51.89	1/5/2015	NM			
	51.89	8/10/2015	6.39			45.50

Notes:

1. The surface completion for MW-23C was repaired and resurveyed in January/February 2009.
2. NM = Not Measured

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-01A 01/28/2008	MW-01A 01/28/2008 Duplicate	MW-01A 07/16/2008	MW-01A 07/16/2008 Duplicate	MW-01A 01/22/2009	MW-01A 01/22/2009 Duplicate	MW-01A 07/22/2009	MW-01A 07/22/2009 Duplicate	MW-01A 01/22/2010	MW-01A 01/22/2010 Duplicate	MW-01A 07/14/2010	MW-01A 07/14/2010 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005			<0.00052	<0.00052								
Benzene	0.005	0.005			<0.00025	<0.00025								
Chlorobenzene	0.1	0.1			<0.00047	<0.00047								
Ethylbenzene	0.7	0.7			<0.00025	<0.00025								
Methylene chloride	0.005	0.005			<0.00054	<0.00054								
Toluene	1	1			<0.00041	<0.00041								
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10			<0.00127	<0.00127								
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.00044	0.000563	0.0109	0.00224	0.0069	0.0016 J	0.0017 J	0.0019 J	0.0019 J	0.0018 J	<0.0009	0.0026 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.0415	0.0409	0.126	0.119	0.054	0.038	0.085	0.091	0.04	0.039	0.068	0.075
Acenaphthylene	1.5	4.4	0.00099	0.000933	0.00143	0.00135	<0.0007	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Anthracene	7.3	22	0.00129	0.00137	0.00267	0.00232	0.0012 J	<0.0007	0.0011 J	0.0014 J	<0.0006	<0.0006	0.0017 J	0.0022 J
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00022	0.0008 J	0.00137 J	0.00126 J	<0.0012	0.0015 J	<0.0033	<0.0033	<0.0033	<0.0033	<0.0033	<0.0033
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.00129	0.00211	0.00774	0.00163	0.0058	0.0018 J	<0.0007	<0.0007	0.0016 J	0.0014 J	0.0044 J	0.0067
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	0.00234	0.00233	0.00923	0.00836	0.0024 J	0.0013 J	0.0037 J	0.0039 J	0.0017 J	0.0015 J	0.004 J	0.0049 J
Fluorene	0.98	2.9	0.0162	0.0167	0.0659	0.0551	0.028	0.018	0.04	0.041	0.022	0.019	0.04	0.047
Naphthalene	0.49	1.5	<0.00044	<0.00038	0.0168	0.00312	<0.0008	<0.0008	0.0029 J	0.0031 J	0.0043 J	0.0036 J	<0.0006	<0.0006
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.00022	0.00035 J	0.00177	0.000783	0.001 J	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	0.0011 J	0.0025 J
Phenol	7.3	22												
Pyrene	0.73	2.2	0.00107	0.00108	0.00417	0.00375	0.001 J	<0.0009	0.0019 J	0.0021 J	<0.0005	<0.0005	0.0021 J	0.0026 J
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-01A 01/11/2011	MW-01A 07/13/2011	MW-01A 07/13/2011 Duplicate	MW-01A 01/31/2012	MW-01A 01/31/2012 Duplicate	MW-01A 07/11/2012	MW-01A 07/11/2012 Duplicate	MW-01A 01/09/2013	MW-01A 01/09/2013 Duplicate	MW-01A 07/11/2013	MW-01A 07/11/2013 Duplicate	MW-01A 01/08/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.0009	0.0068	0.0021 J	<0.0005	<0.0005	0.012	0.011	0.00125	0.00128	0.00193 J	0.0386 J	0.00222 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.07	0.1	0.092	0.029	0.028	0.084	0.083	0.117	0.119	0.098 J	0.132 J	<0.00037
Acenaphthylene	1.5	4.4	0.0011 J	0.0011 J	<0.0005	<0.0005	<0.0005	0.0017 J	0.002 J	0.00222	0.00189	0.00122	0.00137	<0.00093
Anthracene	7.3	22	0.0021 J	0.0029 J	0.0027 J	<0.0005	<0.0005	0.003 J	0.003 J	0.000285 J	0.00373 J	0.0022 J	0.00331 J	<0.003
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0033	0.003 J	0.0012 J	<0.0005	0.0013 J	<0.0005	<0.0005	0.00163	0.00162	<0.000356	<0.000356	0.000838 J
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0007	0.0054	0.0027 J	0.0045 J	0.0044 J	0.025	0.025	0.0141	0.0134	0.00264 J	0.0235 J	<0.00951
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	0.0025 J	0.0062	0.0059	0.0012 J	0.0012 J	0.0047 J	0.0045 J	0.00602	0.00537	0.00399	0.00456	<0.00257
Fluorene	0.98	2.9	0.039	0.056	0.051	0.0013 J	0.0013 J	0.041	0.043	0.0564	0.0556	0.0323 J	0.0545 J	<0.0369
Naphthalene	0.49	1.5	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.00219	0.00245	0.0169 J	0.441 J	<0.0000741
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.0005	0.002 J	0.0011 J	<0.0005	<0.0005	0.0033 J	0.0031 J	0.00388 J	0.0012 J	0.00109 J	0.00928 J	<0.00175
Phenol	7.3	22												
Pyrene	0.73	2.2	0.0011 J	0.0028 J	0.0027 J	<0.0005	<0.0005	0.0021 J	0.0019 J	0.00261	0.00202	0.00165	0.00192	<0.0013
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-01A 01/08/2014	MW-01A 07/02/2014	MW-01A 07/02/2014	MW-01A 01/07/2015	MW-01A 01/07/2015	MW-01A 07/08/2015	MW-01A 07/08/2015	MW-01A 01/12/2016	MW-01A 01/12/2016	MW-01A 07/07/2016	MW-01A 07/07/2016	MW-01A 01/11/2017
			Duplicate		Duplicate		Duplicate		Duplicate		Duplicate		Duplicate	
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	0.0152 J	0.00865 J	0.0000828 J	0.0000693 J	0.000132 J	0.00063 J	0.0041 J	<0.000019	<0.000019	0.000019 J	0.000095 J	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.000741	0.0848 J	0.0228 J	0.0594	0.0607	0.086	0.095	0.048	0.043	0.053	0.053	0.044 J
Acenaphthylene	1.5	4.4	<0.00144	0.00138 J	0.000638 J	0.00104	0.00102	0.0009	0.00099	0.0019	0.0018	0.00063	0.00069	0.00092
Anthracene	7.3	22	<0.00371	0.00326 J	0.00124 J	0.00139	0.00139	0.0018	0.0024	0.00052	0.00059	0.00096	0.001	0.00069
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00067 J	<0.000349	<0.000349	0.00533 J	0.00259 J	<0.0002	<0.000037	0.00042	0.0004	0.000083 J	0.0001 J	0.00031
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0168	0.0132 J	0.00131 J	0.000541	0.000602	0.0032 J	0.009 J	0.00065	0.00054	0.00072	0.00079	0.0011
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	<0.00345	0.0043 J	0.00141 J	0.00246	0.00235	0.0034	0.0038	0.0027	0.0034	0.0027	0.003	0.0021 J
Fluorene	0.98	2.9	<0.0432	0.0369 J	0.0113 J	0.0209	0.0217	0.038	0.045	0.0057	0.0042	0.014	0.014	0.01 J
Naphthalene	0.49	1.5	0.00172 J	0.074 J	0.00469 J	0.000121 J	0.000313 J	0.00083 J	0.0049 J	<0.00002	<0.00002	0.00002 J	0.00022 J	<0.00002
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.00451	0.00537 J	0.000426 J	0.000335 J	0.000326 J	0.0012 J	0.0041 J	0.000067 J	0.000075 J	0.000021 J	0.00027 J	0.000064 J
Phenol	7.3	22		<0.0000377	<0.0000377									
Pyrene	0.73	2.2	<0.00165	0.00204 J	0.000662 J	0.00105	0.00102	0.0015	0.0016	0.0011	0.0014	0.0011	0.0013	0.00095
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-01A 01/11/2017	MW-01A 07/13/2017	MW-01A 07/13/2017	MW-01A 01/04/2018	MW-01A 01/04/2018	MW-01A 07/18/2018	MW-01A 07/18/2018	MW-01A 01/07/2019	MW-01A 01/07/2019	MW-01A 07/02/2019	MW-01A 07/02/2019	MW-01A 01/14/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	Duplicate		Duplicate		Duplicate		Duplicate		Duplicate		Duplicate	
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	0.00018	0.000096 J	0.00024	0.00021	0.00014	0.00021	<0.000019	0.00074	0.00069	0.00019
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.061 J	0.043	0.052	0.041	0.039	0.088	0.12	0.027	0.021	0.063	0.053	0.024
Acenaphthylene	1.5	4.4	0.00098	0.00055	0.00068	0.00086	0.00083	0.00078	0.00086	0.00069	0.00059	0.00071	0.00071	0.00084
Anthracene	7.3	22	0.00093	0.00076	0.00079	0.0012	0.0011	0.002	0.0023	0.00068	0.00046	0.00097	0.00096	<0.000014
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00026	0.00012 J	0.00029	<0.000072	<0.000085	<0.000037	<0.000037	<0.000037	<0.000037	<0.000037	<0.000079	<0.000037
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.001	0.0017	0.0027	0.0016	0.0015	0.01	0.015	0.0019	0.0014	0.0058	0.0054	0.0036
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	0.003 J	0.0025	0.0025	0.0032	0.0029	0.0038	0.0043	0.0018	0.0016	0.0013	0.0012	0.0011
Fluorene	0.98	2.9	0.015 J	0.014	0.014	0.018	0.017	0.04	0.05	0.0037	0.0027	0.019	0.016	0.0064
Naphthalene	0.49	1.5	<0.00002	0.00025	0.00053	<0.00002	<0.00002	0.00027	0.00042	<0.00002	<0.00002	0.00034	0.00034	0.00052
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	0.000021 J	0.00029	0.00045	0.0004	0.00035	0.0021	0.0024	0.00029	<0.000021	0.00076	0.00067	<0.000021
Phenol	7.3	22												
Pyrene	0.73	2.2	0.0012	0.0011	0.001	0.0014	0.0014	0.0019	0.0022	0.00086	0.0007	0.00059	0.00055	0.00052
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
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ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-01A 01/14/2020 Duplicate	MW-02 01/28/2008	MW-02 07/16/2008	MW-02 01/22/2009	MW-02 07/22/2009	MW-02 01/22/2010	MW-02 07/14/2010	MW-02 01/11/2011	MW-02 07/13/2011	MW-02 01/30/2012	MW-02 07/10/2012	MW-02 01/09/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.00038	<0.00039	<0.0008	0.0025 J	<0.0009	<0.0009	<0.0009	0.0021 J	<0.0005	<0.0005	0.00318
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.018	0.017	0.0218	0.014	<0.0009	0.0073	0.018	0.0078	0.026	<0.0005	0.0088	0.0384
Acenaphthylene	1.5	4.4	0.00066	<0.00028	0.0003 J	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.00057
Anthracene	7.3	22	<0.000014	0.000922	0.00042 J	<0.0007	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	0.00129
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000074 J	0.00049 J	<0.00019	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033	0.0021 J	<0.0005	<0.0005	0.000874
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.0021	0.0106	0.00673	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007	0.0038 J	<0.0005	0.0043 J	0.0178
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	0.0012	0.0015	0.000961	<0.0006	0.0011 J	<0.0005	<0.0005	<0.0005	0.0012 J	<0.0005	<0.0005	0.00147
Fluorene	0.98	2.9	0.0038	0.0119	0.0103	0.0039 J	<0.0006	0.0037 J	0.011	0.0049 J	0.015	<0.0005	0.0043 J	0.0201
Naphthalene	0.49	1.5	<0.00002	0.000827	0.00118	<0.0008	0.012	<0.0006	<0.0006	<0.0006	0.0037 J	<0.0005	0.0033 J	0.0211
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.000021	0.000532	<0.00019	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.00241
Phenol	7.3	22												
Pyrene	0.73	2.2	0.00059	0.000816	0.00045 J	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.00087
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-02 07/11/2013	MW-02 01/08/2014	MW-02 07/02/2014	MW-02 01/07/2015	MW-02 07/07/2015	MW-02 01/12/2016	MW-02 07/07/2016	MW-02 01/11/2017	MW-02 07/13/2017	MW-02 01/04/2018	MW-02 07/18/2018	MW-02 01/07/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	0.000897	<0.0000648	0.000509	0.000105 J	0.00013	0.000033 J	<0.000019	<0.000019	0.00093	0.0051	0.00038	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.0179	<0.0000741	0.00452	0.0026	0.001	0.0019	0.00058	0.0035	0.0064	0.015	0.0048	0.0016
Acenaphthylene	1.5	4.4	0.000335 J	<0.000101	0.0000979 J	0.0000892 J	<0.000015	0.000046 J	0.00019	<0.000015	<0.000015	0.00018	<0.000015	<0.000015
Anthracene	7.3	22	0.0013	<0.00131	0.000596	0.000153 J	0.000078 J	0.00011	0.000026 J	0.0001	0.00016	0.00036	0.0002	0.00012
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000356	<0.000343	<0.000349	0.000426 J	<0.00013	<0.00019	<0.000037	0.00011 J	<0.000037	<0.00016	<0.000038	<0.000063
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.00734	<0.000147	0.00301	0.000377 J	0.00075	0.0013	0.00036	0.0019	0.0023	0.0019	0.0016	0.00046
Di-n-butylphthalate (DBP)	2.4	7.3			<0.000104									
Fluoranthene	0.98	2.9	0.00069	<0.000307	0.000368 J	<0.0000693	<0.00001	0.00013	0.000052 J	0.00029	0.00025	0.00071	0.00019	0.00011
Fluorene	0.98	2.9	0.00986	<0.000255	0.00357	0.000681	0.00067	0.001	0.00032	0.0016	0.0041	0.0099	0.0029	0.00081
Naphthalene	0.49	1.5	0.00754	<0.0000741	0.00653	0.000472 J	0.0013	0.0002	0.00013	<0.00002	0.00047	0.00015	0.00025	<0.00002
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	0.000776	<0.000122	0.000594	0.000162 J	0.00011	0.000093 J	0.000066 J	<0.000021	0.0005	0.0011	0.00035	0.000041 J
Phenol	7.3	22			<0.0000377									
Pyrene	0.73	2.2	0.000336 J	<0.000175	0.000201 J	<0.000109	<0.000019	0.000076 J	0.00003 J	0.00017	0.00011	0.00032	0.000082 J	0.000057 J
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray.
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-02 07/02/2019	MW-02 01/14/2020	MW-03 01/30/2008	MW-03 01/24/2018	MW-03 03/27/2018	MW-03 05/25/2018	MW-03 01/09/2019	MW-03 07/12/2019	MW-03 01/21/2020	MW-04 01/29/2008	MW-04 01/24/2018	MW-04 03/23/2018
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005			<0.00052	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.0002	<0.0002
Benzene	0.005	0.005			<0.00025	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.0002	<0.0002
Chlorobenzene	0.1	0.1			<0.00047	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.0003	<0.0003
Ethylbenzene	0.7	0.7			<0.00025	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.0003	<0.0003
Methylene chloride	0.005	0.005			<0.00054	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.001	<0.001
Toluene	1	1			<0.00041	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.0002	<0.0002
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10			<0.00127	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026			<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5			<0.0003	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00029	<0.000041	<0.00004
2,4-Dinitrotoluene	0.0013	0.003			<0.0002	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019	<0.000059	<0.000058
2,6-Dinitrotoluene	0.0013	0.003			<0.0002	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.000043	<0.000042
2-Chloronaphthalene	2	5.8			<0.0004	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00011	<0.000019	<0.0004	<0.000019	<0.000019	<0.000019	<0.000019	<0.000035	0.000045 J	<0.00038	0.00008 J	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073			<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15			<0.00025	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.000048	<0.000047
Acenaphthene	1.5	4.4	0.0019	0.003	0.118	<0.000027	<0.000027	0.00021	<0.000027	<0.000027	0.000061 J	<0.00029	0.00013	<0.000027
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.0003	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029	<0.000015	<0.000015
Anthracene	7.3	22	0.000064 J	0.00011	0.00172	<0.000014	<0.000014	<0.000014	<0.000014	0.000055 J	0.000069 J	0.000739	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02			<0.0002	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00019	<0.000051	<0.00005
Benzo(a)pyrene	0.0002	0.0002			<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000024 J	<0.00019	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019			<0.0004	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.000031	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00015	<0.000037	0.001 J	0.000083 J	<0.000037	<0.000037	<0.000037	<0.000037	0.00011 J	0.00029 J	0.00008 J	<0.000037
Chrysene	0.91	2			<0.0002	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.00041	0.00039	0.00415	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00029	0.000072 J	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3			<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000029 J	<0.00019	0.00003 J	<0.00002
Fluoranthene	0.98	2.9	0.000071 J	0.00024	0.0125	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.000055 J	0.00045 J	<0.00001	<0.00001
Fluorene	0.98	2.9	0.00097	0.0017	0.058	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00019	0.000075 J	<0.00003
Naphthalene	0.49	1.5	0.00011	<0.00002	0.000872	<0.00002	<0.00002	<0.00002	<0.00002	<0.000029	<0.00024	<0.00038	0.0013	0.00055
Nitrobenzene	0.049	0.15			<0.0004	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42			<0.00025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.000026	<0.000025
Pentachlorophenol	0.001	0.001			<0.0002	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019	<0.000081	<0.000079
Phenanthrene	0.73	2.2	0.000054 J	0.00011	0.000592	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00037 J	<0.000021	0.000099 J
Phenol	7.3	22			<0.0002	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	<0.000036	<0.000035
Pyrene	0.73	2.2	0.000037 J	0.00011	0.00538	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000036 J	<0.00019	0.00003 J	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01					0.000895 J	0.00242	0.00363	<b>0.0191</b>	0.000582 J	0.00207		0.00454 0.00092 J

Notes:  
1. All values in milligrams per liter (mg/L).  
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3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-04 05/25/2018	MW-04 01/09/2019	MW-04 07/12/2019	MW-04 01/21/2020	MW-05 01/29/2008	MW-05 07/27/2011	MW-05 02/02/2012	MW-05 07/25/2012	MW-05 02/05/2013	MW-05 08/01/2013	MW-05 01/15/2014	MW-05 07/29/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.0011	<0.0011	<0.0005	<0.000137	<0.00011	<0.00019	<0.00011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015
Vinyl chloride	0.002	0.002									<0.00011			
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	<0.000108
2,4-Dimethylphenol	0.49	1.5	0.00011 J	<0.00004	<0.00004	<0.00004	<0.00029	<0.00005	<0.00005	<0.00005	<0.000292	<0.000295	<0.000292	<0.000304
2,4-Dinitrotoluene	0.0013	0.003	<0.000059	<0.000058	<0.000058	<0.000058	<0.00019	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124	<0.000123	<0.000127
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000762	<0.0000755	<0.0000784
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000784
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000059	0.00026	<0.00038	<0.00005	0.000085 J	<0.00005	0.000468 J	<0.0000667	0.000187 J	<0.0000686
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00008	<0.00008	<0.00008	<0.000783	<0.00079	<0.000783	<0.000814
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00005	<0.00005	<0.00005	<0.000528	<0.000533	<0.000528	<0.000549
Acenaphthene	1.5	4.4	0.00031	<0.000027	0.00052	0.00015	0.000545	0.0053	0.00034	<0.00005	<0.0000755	0.000521	0.000194 J	<0.0000784
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029	<0.00005	<0.00005	<0.00005	<0.0000566	<0.0000571	<0.0000566	<0.0000588
Anthracene	7.3	22	0.000054 J	0.000079 J	<0.000014	0.0001	0.000811	<0.00005	<0.00005	<0.00005	0.000621	0.000427 J	0.000411 J	0.000153 J
Benzo(a)anthracene	0.0091	0.02	<0.000051	<0.00005	<0.00005	<0.00005	<0.00019	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000784
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	0.000033 J	<0.00019	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000784
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124	<0.000123	<0.000127
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	<0.000037	0.0018	0.00034 J	0.00047	<0.0001	0.00019 J	<0.000349	<0.000352	<0.000349	<0.000363
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	0.000034 J	<0.00019	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000784
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.000058	0.000097 J	<0.00029	0.0022	0.00011 J	<0.00005	<0.0000755	0.0000828 J	0.000162 J	<0.0000784
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00005	<0.00005	0.000065 J	<0.000104	<0.000105	<0.000104	<0.000108
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	0.00029	0.00047 J	0.00011 J	<0.00005	<0.00005	<0.000066	0.0000761 J	<0.000066	<0.0000686
Fluorene	0.98	2.9	0.000059 J	<0.00003	0.000053 J	0.000067 J	0.0002 J	0.0012	0.00012 J	<0.00005	<0.000066	0.000166 J	0.000176 J	<0.0000686
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.000053	0.0013	<0.00038	<0.00005	0.00087	<0.00005	0.00133	0.000573 J	0.000969 J	<0.000131
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	<0.000108
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000952	<0.0000943	<0.000098
Pentachlorophenol	0.001	0.001	<0.00008	<0.000079	<0.000079	<0.000079	<0.00019	<0.00005	<0.00005	<0.00005	<0.000575	<0.000581	<0.000575	<0.000598
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	0.00017	0.00039 J	0.00013 J	<0.00005	<0.00005	0.000143 J	<0.0000571	0.000507	<0.0000588
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	<0.00005	<0.00005	<0.00005	0.000193 J	<0.0000381	<0.0000377	<0.0000392
Pyrene	0.73	2.2	0.00011	<0.000019	0.00016	0.0002	0.00045 J	0.00014 J	<0.00005	<0.00005	<0.000104	0.000154 J	<0.000104	<0.000108
<b>Metals</b>														
Arsenic	0.01	0.01	0.00492	0.000963 J	0.0127	0.00223								

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-05 01/24/2018	MW-05 03/20/2018	MW-05 05/24/2018	MW-05 01/09/2019	MW-05 07/11/2019	MW-05 01/13/2020	MW-07 01/29/2008	MW-07 07/16/2008	MW-07 01/22/2009	MW-07 07/22/2009	MW-07 01/22/2010	MW-07 07/14/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002		<0.00052				
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002		<0.00025				
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003		<0.00047				
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003		<0.00025				
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001		<0.00054				
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002		<0.00041				
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003		<0.00127				
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021						
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004		<0.00059				
2,4-Dinitrotoluene	0.0013	0.003	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058		<0.00058				
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042		<0.00042				
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021		<0.00021				
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	0.000087 J	<0.000019	<0.00038	<0.00039	<0.0008	<0.0009	<0.0009	<0.0009
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002		<0.00002				
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047		<0.00047				
Acenaphthene	1.5	4.4	<0.000027	<0.000027	0.00023	<0.000027	0.00043	<0.000027	<0.00028	<0.00029	<0.0008	<0.0009	<0.0009	<0.0009
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00028	0.00044 J	<0.0007	<0.0005	<0.0005	<0.0005
Anthracene	7.3	22	0.000037 J	<0.000014	0.000045 J	<0.000014	0.000097 J	<0.000014	0.000516	0.000982	<0.0007	<0.0006	<0.0006	<0.0006
Benzo(a)anthracene	0.0091	0.02	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005						
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002		<0.00002				
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003		<0.00003				
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00011 J	<0.000037	0.00013 J	<0.000037	<0.000037	<0.000037	<0.00019	<0.00019	<0.0012	<0.0033	<0.0033	0.0049 J
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021		<0.00021				
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	0.000024 J	<0.00002	0.000051 J	<0.00002	<0.00028	<0.00029	<0.0007	<0.0007	<0.0007	<0.0007
Di-n-butylphthalate (DBP)	2.4	7.3	0.000033 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002		<0.00002				
Fluoranthene	0.98	2.9	0.000051 J	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00019	<0.00019	<0.0006	<0.0005	<0.0005	<0.0005
Fluorene	0.98	2.9	<0.00003	<0.00003	0.000039 J	<0.00003	0.000064 J	<0.00003	<0.00019	<0.00019	<0.0008	<0.0006	<0.0006	<0.0006
Naphthalene	0.49	1.5	0.0001	<0.00002	<0.00002	<0.00002	<0.00071	<0.00012	<0.00038	0.000675	<0.0008	<0.0006	<0.0006	<0.0006
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024		<0.00024				
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025		<0.00025				
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079		<0.00079				
Phenanthrene	0.73	2.2	0.000054 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019	0.00036 J	<0.0007	<0.0005	<0.0005	<0.0005
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035		<0.00035				
Pyrene	0.73	2.2	0.000047 J	<0.000019	0.00019	<0.000019	0.00019	<0.000019	<0.00019	<0.00019	<0.0009	<0.0005	<0.0005	<0.0005
<b>Metals</b>														
Arsenic	0.01	0.01	0.00588	0.00255	0.00488	0.00387	0.0171	0.00146 J						

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-07 01/12/2011	MW-07 01/12/2011	MW-07 07/12/2011	MW-07 01/31/2012	MW-07 07/11/2012	MW-07 01/10/2013	MW-07 07/11/2013	MW-07 01/09/2014	MW-07 07/03/2014	MW-07 01/07/2015	MW-07 07/08/2015	MW-07 01/12/2016
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005	<0.000066	<0.0000704	R	<0.000066	<0.0000693	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005	0.00181	<0.0000804	R	<0.0000755	<0.0000792	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.00011 J	<0.0000603	R	<0.0000566	<0.0000594	<0.000015	<0.000015
Anthracene	7.3	22	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	0.000833	0.000749	R	0.000696	0.000271 J	0.00014	0.000085 J
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0033	<0.0005	<0.0005	<0.0005	<0.0005	<0.000349	<0.000372	R	<0.000349	0.000944	<0.00024	0.00024
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000804	R	<0.0000755	<0.0000792	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000066	<0.0000704	R	<0.000066	0.000189 J	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	0.000137 J	<0.0000704	R	<0.000066	0.0000792 J	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000755	0.000111 J	R	<0.0000755	<0.0000792	<0.00002	<0.00002
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000566	<0.0000603	R	<0.0000566	<0.0000594	<0.000021	<0.000021
Phenol	7.3	22												
Pyrene	0.73	2.2	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.000111	R	<0.000104	0.000142 J	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-07 07/07/2016	MW-07 01/12/2017	MW-07 07/13/2017	MW-07 01/04/2018	MW-07 07/19/2018	MW-07 01/08/2019	MW-07 07/01/2019	MW-07 07/01/2019 Duplicate	MW-07 01/13/2020	MW-08 01/29/2008	MW-08 07/16/2008	MW-08 01/22/2009
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												<0.00109
Benzene	0.005	0.005												<0.00112
Chlorobenzene	0.1	0.1												<0.0015
Ethylbenzene	0.7	0.7												<0.00142
Methylene chloride	0.005	0.005												<0.00122
Toluene	1	1												<0.00138
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												<0.00302
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000023 J	0.000066 J	<0.00044	<0.0004	<0.0008
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.00033	<0.0003	<0.0008
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00033	0.00044 J	<0.0007
Anthracene	7.3	22	0.000055 J	0.00025	<0.000014	0.00014	0.00036	<0.000014	0.00014	0.00029	<0.000014	0.00031 J	0.000669	<0.0007
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.00011 J	<0.000037	<0.000088	<0.000037	<0.000037	<0.000037	<0.000037	<0.000037	<0.00022	<0.0002	<0.0012
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000057 J	<0.00033	<0.0003
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.0001	<0.00022	<0.0002	<0.0006
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00022	<0.0002	<0.0008
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00025	0.00017	<0.00044	0.000654	<0.0008
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00014	<0.00022	0.00036 J
Phenol	7.3	22												
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00022	<0.0002	<0.0009
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-08 07/22/2009	MW-08 01/22/2010	MW-08 07/14/2010	MW-08 01/12/2011	MW-08 01/12/2011 Duplicate	MW-08 07/12/2011	MW-08 01/31/2012	MW-08 07/11/2012	MW-08 01/10/2013	MW-08 07/11/2013	MW-08 01/09/2014	MW-08 07/03/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.0009	<0.0009	<0.0009	<0.0009	<0.0009	<0.0005	<0.0005	<0.0005	<0.000066	<0.0000686	<0.0000648	<0.000066
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.0009	<0.0009	<0.0009	<0.0009	<0.0009	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000784	<0.0000741	<0.0000755
Acenaphthylene	1.5	4.4	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000566	<0.0000588	<0.0000556	<0.0000566
Anthracene	7.3	22	<0.0006	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	0.000439 J	0.000101 J	0.000494	<0.0000472
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0033	<0.0033	<0.0033	<0.0033	<0.0033	<0.0005	<0.0005	<0.0005	<0.000349	<0.000363	<0.000343	<0.000349
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000784	<0.0000741	<0.0000755
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000066	<0.0000686	<0.0000648	<0.000066
Fluorene	0.98	2.9	<0.0006	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.000066	<0.0000686	<0.0000648	<0.000066
Naphthalene	0.49	1.5	<0.0006	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000784	<0.0000741	<0.0000755
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000566	<0.0000588	0.0000637 J	<0.0000566
Phenol	7.3	22												<0.0000377
Pyrene	0.73	2.2	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.000108	<0.000102	<0.000104
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-08 01/07/2015	MW-08 07/08/2015	MW-08 01/12/2016	MW-08 07/07/2016	MW-08 01/12/2017	MW-08 07/13/2017	MW-08 01/04/2018	MW-08 07/19/2018	MW-08 01/07/2019	MW-08 07/01/2019	MW-08 01/13/2020	MW-09 01/29/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												<0.00052
Benzene	0.005	0.005												<0.00025
Chlorobenzene	0.1	0.1												<0.00047
Ethylbenzene	0.7	0.7												<0.00025
Methylene chloride	0.005	0.005												<0.00054
Toluene	1	1												<0.00041
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												<0.00127
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												<0.00008
2,4-Dimethylphenol	0.49	1.5												<0.00029
2,4-Dinitrotoluene	0.0013	0.003												<0.00019
2,6-Dinitrotoluene	0.0013	0.003												<0.00019
2-Chloronaphthalene	2	5.8												<0.00038
2-Methylnaphthalene	0.098	0.29	<0.0000693	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00038
4,6-Dinitro-2-methylphenol	0.0024	0.0073												<0.00019
4-Nitrophenol	0.049	0.15												<0.00024
Acenaphthene	1.5	4.4	<0.0000792	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.00067	<0.000027	<0.000027	<0.000027	<0.00029
Acenaphthylene	1.5	4.4	<0.0000594	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029
Anthracene	7.3	22	0.000056 J	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.000045 J	0.000048 J	<0.000014	<0.000014	<0.00019
Benzo(a)anthracene	0.0091	0.02												<0.00019
Benzo(a)pyrene	0.0002	0.0002												<0.00019
bis(2-Chloroethoxy)methane	0.00083	0.0019												<0.00038
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000366	<0.00013	<0.000037	<0.000037	0.00018 J	<0.000037	<0.000097	<0.000037	<0.000037	<0.000037	0.00021	0.00034 J
Chrysene	0.91	2												<0.00019
Dibenzofuran	0.098	0.29	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00011	<0.00002	<0.00002	<0.00002	<0.00029
Di-n-butylphthalate (DBP)	2.4	7.3												<0.00019
Fluoranthene	0.98	2.9	<0.0000693	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.000031 J	<0.00001	<0.00001	<0.00001	0.00045 J
Fluorene	0.98	2.9	<0.0000693	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00017	<0.00003	<0.00003	<0.00003	<0.00019
Naphthalene	0.49	1.5	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00038
Nitrobenzene	0.049	0.15												<0.00038
N-Nitrosodiphenylamine	0.19	0.42												<0.00024
Pentachlorophenol	0.001	0.001												<0.00019
Phenanthrene	0.73	2.2	<0.0000594	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000023 J	<0.000021	<0.000021	0.000046 J	0.000541
Phenol	7.3	22												<0.00019
Pyrene	0.73	2.2	<0.000109	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00019
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-09 07/27/2011	MW-09 02/02/2012	MW-09 07/25/2012	MW-09 04/01/2013	MW-09 01/24/2018	MW-09 03/23/2018	MW-09 05/24/2018	MW-09 01/09/2019	MW-09 07/12/2019	MW-09 01/14/2020	MW-10A 01/28/2008	MW-10A 07/16/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0005	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052
Benzene	0.005	0.005	<0.001	<0.001	<0.0005	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.0005	<0.00012	<0.0003	<0.0003	<0.0003	0.0041	<0.0003	<0.0003	<0.0003	<0.00047
Ethylbenzene	0.7	0.7	<0.0011	<0.0011	<0.0005	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<0.001	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054
Toluene	1	1	<0.001	<0.001	<0.0005	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0031	<0.0031	<0.0015	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<0.000106	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.00005	<0.00005	<0.000298	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00005	<0.000125	<0.000059	<0.000058	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.00006	<0.0000769	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.00005	<0.0000769	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00005	<0.00005	<0.00005	0.000115 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00038
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.0000798	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000538	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.00005	<0.00005	<0.00005	0.000188 J	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.0003
Acenaphthylene	1.5	4.4	<0.00005	<0.00005	<0.00005	<0.0000577	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.0003
Anthracene	7.3	22	0.00036	<0.00005	<0.00005	0.000471 J	<0.000014	0.00013	0.0001	0.000093 J	0.000061 J	<0.000014	<0.000014	<0.0003
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.0000769	<0.000051	<0.00005	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00005	<0.000125	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00018 J	<0.00032	0.00022	<0.000356	0.000082 J	<0.000037	0.00011 J	<0.000037	<0.000037	<0.000037	<0.000037	0.0002 J
Chrysene	0.91	2	<0.00005	<0.00005	<0.00005	<0.0000769	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00005	<0.00005	<0.00005	0.000126 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0003
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	0.000074 J	0.000123 J	<0.00002	<0.00002	0.000031 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00005	<0.00005	<0.00005	<0.0000673	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00019
Fluorene	0.98	2.9	<0.00005	<0.00005	<0.00005	<0.0000673	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00019
Naphthalene	0.49	1.5	<0.00005	<0.00005	<0.00005	0.00431 J	<0.00002	0.00039	0.00008 J	<0.00002	<0.00015	0.00011	<0.00004	<0.00038
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000106	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.0000962	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00005	<0.000587	<0.00008	<0.000079	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.00005	<0.00005	<0.00005	<0.0000577	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019
Phenol	7.3	22	<0.00005	<0.000098	<0.00005	<0.0000385	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.00005	<0.00005	<0.00005	<0.000106	<0.000019	<0.000019	0.00004 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.00019
<b>Metals</b>														
Arsenic	0.01	0.01					0.00104 J	0.0012 J	0.00085 J	0.00202	0.000901 J	0.0043		

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-10A 01/22/2009	MW-10A 07/22/2009	MW-10A 01/21/2010	MW-10A 07/13/2010	MW-10A 01/11/2011	MW-10A 07/13/2011	MW-10A 01/30/2012	MW-10A 07/10/2012	MW-10A 01/09/2013	MW-10A 07/11/2013	MW-10A 01/08/2014	MW-10A 07/15/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.0008	<0.0009	<0.0009	<0.0009	<0.0009	<0.0005	<0.0005	<0.0005	<0.0000667	0.00178	<0.0000648	0.00262 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.0008	<0.0009	<0.0009	<0.0009	0.0017 J	<0.0005	<0.0005	0.0016 J	<0.0000762	0.0306	<0.0000741	0.0306
Acenaphthylene	1.5	4.4	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000571	0.000385 J	<0.0000556	<0.000566
Anthracene	7.3	22	<0.0007	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	0.000468 J	0.00036 J	<0.0000463	<0.000472
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033	<0.0005	<0.0005	<0.0005	0.00171	<0.000356	<0.000343	<0.00349
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007	<0.0005	<0.0005	<0.0005	<0.0000762	0.00866	<0.0000741	0.00862
Di-n-butylphthalate (DBP)	2.4	7.3												
Fluoranthene	0.98	2.9	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000667	0.000186 J	<0.0000648	<0.00066
Fluorene	0.98	2.9	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0000667	0.00631	<0.0000648	0.0111
Naphthalene	0.49	1.5	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0000762	0.199	<0.0000741	0.199
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000571	0.00221	<0.0000556	0.00442 J
Phenol	7.3	22												
Pyrene	0.73	2.2	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000105	<0.000106	<0.000102	<0.00104
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
- All values in milligrams per liter (mg/L).
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  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-10A 01/07/2015	MW-10A 07/07/2015	MW-10A 01/12/2016	MW-10A 07/07/2016	MW-10A 01/11/2017	MW-10A 07/13/2017	MW-10A 01/04/2018	MW-10A 07/18/2018	MW-10A 01/07/2019	MW-10A 07/02/2019	MW-10A 01/14/2020	MW-10B 01/28/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.0000693	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.00272	<0.000027	<0.000027	<0.000027	<0.000028	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.00011	0.0743
Acenaphthylene	1.5	4.4	0.000126 J	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.00122
Anthracene	7.3	22	0.000191 J	0.000069 J	<0.000014	0.000073 J	0.000057 J	<0.000014	<0.000014	<0.000014	0.000065 J	<0.000014	<0.000014	0.00432
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000366	<0.00088	<0.000096	0.000097 J	0.000088 J	<0.000037	<0.000057	<0.0001	<0.000089	<0.000037	<0.000037	<0.00019
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.000349 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.0255
Di-n-butylphthalate (DBP)	2.4	7.3												<0.00019
Fluoranthene	0.98	2.9	<0.0000693	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.00371
Fluorene	0.98	2.9	0.000694	<0.00003	<0.00003	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003	0.000038 J	<0.00003	<0.00003	0.0374
Naphthalene	0.49	1.5	0.000322 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000043 J	<0.00002	0.0185
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	0.000126 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019
Phenol	7.3	22												<0.00019
Pyrene	0.73	2.2	<0.000109	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.00146
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
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  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-10B 07/16/2008	MW-10B 01/22/2009	MW-10B 07/22/2009	MW-10B 01/21/2010	MW-10B 07/13/2010	MW-10B 01/11/2011	MW-10B 07/13/2011	MW-10B 01/30/2012	MW-10B 07/10/2012	MW-10B 01/09/2013	MW-10B 07/11/2013	MW-10B 10/14/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052											
Benzene	0.005	0.005	<0.00025											
Chlorobenzene	0.1	0.1	<0.00047											
Ethylbenzene	0.7	0.7	<0.00025											
Methylene chloride	0.005	0.005	<0.00054											
Toluene	1	1	<0.00041											
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00127											
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29												
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.0975	0.096	<0.0009	0.052	0.069	0.096	0.054	0.1	0.054	0.12	0.977	
Acenaphthylene	1.5	4.4	0.00113	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.0011 J	<0.0005	0.00108	0.00986	
Anthracene	7.3	22	0.00484	0.0043 J	0.0029 J	0.0025 J	0.0038 J	0.0068	0.0033 J	0.0057	0.0032 J	0.00546	0.0391	
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0002 J	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033	0.0013 J	<0.0005	<0.0005	<0.000349	<0.0037	
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.0392	0.035	0.023	0.018	0.025	0.037	0.019	0.038	0.02	0.0401	0.302	0.0334
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.011	
Fluoranthene	0.98	2.9	0.00397	0.0039 J	0.0022 J	0.0017 J	0.0026 J	0.0054	0.0023 J	0.0046 J	0.0028 J	0.00427	0.0274	
Fluorene	0.98	2.9	0.0457	0.051	<0.0006	0.031	0.041	0.059	0.032	0.06	0.031	0.0652	0.468	
Naphthalene	0.49	1.5	0.014	0.0028 J	0.0082	0.0037 J	0.056	0.075	0.0018 J	0.084	0.004 J	0.00399	0.207	
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2												
Phenol	7.3	22	<0.0002	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000377	<0.0004	
Pyrene	0.73	2.2	0.00174	0.002 J	0.0013 J	<0.0005	0.001 J	0.0023 J	0.0011 J	0.002 J	0.0011 J	0.00146	0.0101	
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-10B 01/08/2014	MW-10B 07/15/2014	MW-10B 01/07/2015	MW-10B 01/29/2015	MW-10B 07/07/2015	MW-10B 01/12/2016	MW-10B 07/07/2016	MW-10B 01/11/2017	MW-10B 07/13/2017	MW-10B 01/04/2018	MW-10B 07/18/2018	MW-10B 01/07/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29												
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.0000741	0.0777	0.166	0.0507	0.084	0.11	0.053	0.12	0.051	0.093	0.056	0.07
Acenaphthylene	1.5	4.4	<0.000536	<0.000566	0.00104	0.000597	0.00048	0.00054	0.00032	0.00065	0.00033	0.00068	0.00053	0.00059
Anthracene	7.3	22	<0.00107	0.00352 J	0.00702	0.00179	0.003	0.0057	0.0027	0.0056	0.0023	0.0052	0.0024	0.0041
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000408 J	<0.00349	<0.000366	<0.000366	<0.00014	<0.000037	0.00025	0.00022	0.00028	<0.000059	<0.000064	0.00035
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.00493	0.0258	0.0727	0.0129	0.032	0.049	0.019	0.052	0.021	0.048	0.021	0.028
Di-n-butylphthalate (DBP)	2.4	7.3	0.000275 J	<0.00104	<0.000215	<0.000109	<0.000078	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.000117	0.00211 J	0.00711	0.00117	0.0023	0.0045	0.0023	0.0047	0.0022	0.0053	0.0022	0.0038
Fluorene	0.98	2.9	<0.00429	0.0424	0.0975	0.0202	0.047	0.064	0.029	0.068	0.034	0.06	0.031	0.04
Naphthalene	0.49	1.5	0.0646	0.125	0.556	0.0247	0.077	0.1	0.0054	0.021	0.0033	0.0013	0.00074	0.00056
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2												
Phenol	7.3	22	<0.000037	<0.000377	<0.0000396	<0.0000396	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000102	<0.00104	0.00234	0.000392 J	0.00095	0.0024	0.0009	0.0022	0.00087	0.0019	0.00088	0.0018
<b>Metals</b>														
Arsenic	0.01	0.01												

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  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
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CLASS 2 GROUNDWATER  
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	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-10B 07/02/2019	MW-10B 01/14/2020	MW-11A 01/28/2008	MW-11A 07/16/2008	MW-11A 01/22/2009	MW-11A 07/22/2009	MW-11A 01/21/2010	MW-11A 07/13/2010	MW-11A 01/11/2011	MW-11A 07/12/2011	MW-11A 01/30/2012	MW-11A 07/10/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005						<0.00052						
Benzene	0.005	0.005						<0.00025						
Chlorobenzene	0.1	0.1						<0.00047						
Ethylbenzene	0.7	0.7						<0.00025						
Methylene chloride	0.005	0.005						<0.00054						
Toluene	1	1						<0.00041						
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10						<0.00127						
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29					<0.00038	<0.0004	<0.0008	<0.0009	<0.0009	<0.0009	<0.0005	<0.0005
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.042	0.069	0.0346	0.02	0.0076	<0.0009	<0.0009	0.0028 J	<0.0009	<0.0005	<0.0005	<0.0005
Acenaphthylene	1.5	4.4	0.00031	0.00066	<0.00029	<0.0003	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Anthracene	7.3	22	0.0012	0.0028	0.000798	0.00054	<0.0007	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.0002	0.00028 J	<0.0002	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033	<0.0005	<0.0005	<0.0005
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.013	0.022	0.00276	<0.0003	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007	<0.0005	<0.0005	<0.0005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002										
Fluoranthene	0.98	2.9	0.0012	0.0029	0.00338	0.00387	0.0012 J	0.0011 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Fluorene	0.98	2.9	0.018	0.036	0.0069	0.00089	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005
Naphthalene	0.49	1.5	0.00023	0.0021	<0.00038	<0.0004	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2			0.00036 J	<0.0002	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Phenol	7.3	22	<0.000035	<0.000035										
Pyrene	0.73	2.2	0.00049	0.0013	0.00191	0.00184	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

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<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.000066	<0.0000673	<0.0000648	<0.000066	<0.0000673	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.00175	0.000878	<0.0000741	0.00427	0.000471 J	0.00025	<0.000027	0.0001	<0.000028	<0.000027	<0.000027	0.0017
Acenaphthylene	1.5	4.4	<0.0000566	<0.0000577	<0.0001	0.000185 J	<0.0000577	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.000499	0.00044 J	<0.00125	0.00126	0.000399 J	0.00017	0.0001	0.00026	0.00021	<0.000014	<0.000014	0.00019
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000349	<0.000356	0.00046 J	0.00516	<0.000356	<0.00068	<0.000073	0.00038	0.00025	0.00017 J	<0.000073	0.0007
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0000755	<0.0000769	<0.0000741	0.000618	<0.0000769	0.00012	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	0.00027
Di-n-butylphthalate (DBP)	2.4	7.3				0.000109 J								
Fluoranthene	0.98	2.9	<0.000066	0.000221 J	<0.0000795	0.00215	<0.0000673	0.00028	<0.00001	0.00065	0.000049 J	0.00015	<0.00001	0.00023
Fluorene	0.98	2.9	<0.000066	<0.0000673	<0.0000648	0.00149	<0.0000673	0.00011	<0.00003	0.00013	<0.000031	<0.00003	<0.00003	0.00083
Naphthalene	0.49	1.5	<0.0000755	<0.0000769	<0.0000741	0.000343 J	<0.0000769	<0.00002	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	0.00012
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.0000566	<0.0000577	<0.0000556	0.000384 J	<0.0000577	<0.000021	<0.000021	<0.000021	0.000067 J	<0.000021	<0.000021	0.000086 J
Phenol	7.3	22				<0.0000377								
Pyrene	0.73	2.2	<0.000104	0.000115 J	<0.000102	0.00194	<0.000106	0.00023	<0.000019	0.00073	<0.00002	<0.000019	<0.000019	0.00022
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-11A 01/07/2019	MW-11A 07/02/2019	MW-11A 01/14/2020	MW-11B 01/28/2008	MW-11B 07/16/2008	MW-11B 01/22/2009	MW-11B 07/22/2009	MW-11B 01/21/2010	MW-11B 07/13/2010	MW-11B 01/11/2011	MW-11B 07/12/2011	MW-11B 01/30/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005					<0.00052							
Benzene	0.005	0.005					<0.00025							
Chlorobenzene	0.1	0.1					<0.00047							
Ethylbenzene	0.7	0.7					<0.00025							
Methylene chloride	0.005	0.005					<0.00054							
Toluene	1	1					<0.00041							
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10					<0.00127							
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019									
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.000027	0.00025	<0.000027	0.0649	0.12	0.072	0.12	0.048	0.11	0.039	0.084	0.025
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.00028	0.00126	<0.0007	0.0015 J	0.0013 J	<0.0005	0.0012 J	0.0012 J	0.0011 J
Anthracene	7.3	22	0.00013	0.000097 J	<0.000014	0.00236	0.00472	0.0022 J	0.0043 J	0.0011 J	0.0055	<0.0006	0.0054	<0.0005
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000075	<0.00007	<0.000037	0.00021 J	<0.00021	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033	<0.0005	<0.0005
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	0.0273	0.0649	0.031	0.054	0.012	0.048	0.006	0.038	<0.0005
Di-n-butylphthalate (DBP)	2.4	7.3				<0.00019	<0.00021	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Fluoranthene	0.98	2.9	<0.00001	0.000018 J	<0.00001	0.00175	0.00383	0.0018 J	0.0036 J	0.0014 J	0.0046 J	0.0015 J	0.0046 J	0.0013 J
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	0.0297	0.0578	0.032	0.053	0.013	0.056	0.0038 J	0.046	<0.0005
Naphthalene	0.49	1.5	<0.00002	0.000041 J	<0.00002	0.0354	0.0772	<0.0008	0.048	<0.0006	0.0068	<0.0006	0.06	<0.0005
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021									
Phenol	7.3	22				<0.00019	<0.00021	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Pyrene	0.73	2.2	<0.000019	0.000031 J	<0.000019	0.000848	0.00163	<0.0009	0.002 J	<0.0005	0.0022 J	<0.0005	0.0024 J	<0.0005
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-11B 07/10/2012	MW-11B 01/09/2013	MW-11B 07/11/2013	MW-11B 01/08/2014	MW-11B 07/02/2014	MW-11B 01/07/2015	MW-11B 07/07/2015	MW-11B 01/12/2016	MW-11B 07/07/2016	MW-11B 01/11/2017	MW-11B 07/12/2017	MW-11B 01/03/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29					0.0131							
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	0.1	0.0631	0.108	<0.00037	0.0953	0.0472	0.057	<0.000027	0.039	<0.000027	0.043	<0.000027
Acenaphthylene	1.5	4.4	0.0013 J	0.00136	0.00119	<0.00102	0.00166	0.00113	0.00065	<0.000015	0.00031	<0.000015	0.00032	<0.000015
Anthracene	7.3	22	0.0055	0.000168 J	0.00321	<0.00242	0.00375	0.000945	0.0025	0.00011	0.0018	0.00011	0.0022	0.0001
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0005	0.00195	<0.000356	0.000493 J	<0.000349	<0.000356	<0.00019	<0.00019	0.00018 J	0.00017 J	<0.000037	<0.0001
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	0.04	0.00352	0.0231	<0.0111	0.0199	0.00472	0.014	<0.00002	0.0082	0.000055 J	0.0055	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0005	<0.000104	<0.000106	0.000317 J	0.000109 J	<0.000106	<0.000044	<0.00002	<0.00002	<0.00002	<0.00002	<0.000045
Fluoranthene	0.98	2.9	0.0053	0.00307	0.00383	<0.00267	0.00417	0.00201	0.0034	0.00011	0.0025	0.00021	0.0029	<0.00001
Fluorene	0.98	2.9	0.054	0.00205	0.0388	<0.0195	0.0339	0.00867	0.025	<0.00003	0.019	<0.00003	0.025	<0.00003
Naphthalene	0.49	1.5	0.004 J	<0.0000755	0.00535	0.000382 J	0.0135	<0.0000769	0.0021	<0.00002	0.0082	<0.00002	0.0019	<0.00002
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2					0.012							
Phenol	7.3	22	<0.0005	<0.0000377	<0.0000385	<0.000037	<0.0000377	<0.0000385	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0024 J	0.00154	0.00196	<0.00126	0.00213	0.000935	0.0017	0.00032	0.0012	0.00013	0.0015	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-11B 07/18/2018	MW-11B 01/07/2019	MW-11B 07/02/2019	MW-11B 07/30/2019	MW-11B 10/17/2019	MW-11B 01/14/2020	MW-12A 01/30/2008	MW-12A 07/15/2008	MW-12A 02/04/2009	MW-12A 01/19/2010	MW-12A 06/22/2010	MW-12A 01/18/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005							<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005							<0.00025	<0.00025	0.00073 J	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1							<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7							0.00718	<0.00025	0.0059	0.0029 J	0.00056 J	0.0014 J
Methylene chloride	0.005	0.005							<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005
Toluene	1	1							<0.00041	<0.00041	0.00079 J	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10							0.0105 J	<0.00127	0.012 J	0.0056 J	0.0026 J	0.0025 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026							<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5							<0.0003	<0.00029	<0.00008	<0.00008	<0.00008	0.0001 J
2,4-Dinitrotoluene	0.0013	0.003							<0.0002	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003							<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8							<0.0004	<0.00039	<0.00012	<0.0001	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29							<b>0.174</b>	<b>0.332</b>	<b>0.22</b>	<b>0.15</b>	<b>0.15</b>	0.033
4,6-Dinitro-2-methylphenol	0.0024	0.0073							<0.0002	<0.00049	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15							<0.00025	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	0.074	0.015	0.13		0.033		0.173	0.331	0.25	0.19	0.21	0.19
Acenaphthylene	1.5	4.4	0.00063	0.00054	0.0013		0.016		<0.0003	0.00276	0.0036	0.0026	0.0019	0.0016
Anthracene	7.3	22	0.0037	0.00021	0.0045		<0.000014		0.0103	0.0137	0.0099	0.0093	0.011	0.012
Benzo(a)anthracene	0.0091	0.02							0.00028 J	0.00026 J	<0.00007	<0.00007	<0.00007	0.00082
Benzo(a)pyrene	0.0002	0.0002							<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<b>0.0003</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019							<0.0004	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000085	<0.00026		0.000095 J		0.0013 J	0.00033 J	0.00031	0.0006	<0.0002	0.00036
Chrysene	0.91	2							0.00024 J	0.00021 J	<0.00007	<0.00007	<0.00007	0.00074
Dibenzofuran	0.098	0.29	0.032	<0.00002	0.051		<0.00002		<b>0.125</b>	<b>0.212</b>	<b>0.18</b>	<b>0.14</b>	<b>0.18</b>	<b>0.15</b>
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002		<0.00002		<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.0043	0.0025	0.005		0.0024		0.00693	0.0123	0.0061	0.0059	0.0064	0.0086
Fluorene	0.98	2.9	0.048	0.00015	0.061		0.00035		0.112	0.475	0.15	0.13	0.16	0.14
Naphthalene	0.49	1.5	0.34	<0.00002	<b>0.7</b>	<b>1.1</b>	<b>0.6</b>	<0.00002	<b>2.27</b>	<b>1.47</b>	<b>2.6</b>	<b>1.7</b>	<b>0.6</b>	0.22
Nitrobenzene	0.049	0.15							<0.0004	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42							<0.00025	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001							<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2							0.0833	0.372	0.1	0.087	0.091	0.061
Phenol	7.3	22	<0.000035	<0.000035	<0.000035		<0.000035		<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	0.0023	0.0017	0.0027		0.0023		0.00358	0.00518	0.0025	0.0029	0.0025	0.0044
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

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- Concentrations > C/I AL and non-detects are highlighted dark gray
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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-12A 07/26/2011	MW-12A 02/01/2012	MW-12A 07/19/2012	MW-12A 02/05/2013	MW-12A 07/31/2013	MW-12A 01/14/2014	MW-12A 07/25/2014	MW-12A 01/23/2018	MW-12A 03/19/2018	MW-12A 05/16/2018	MW-12A 01/09/2019	MW-12A 07/11/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0025	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.001	<0.001	<0.0025	0.000237 J	0.0000957 J	0.00048 J	0.000122 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.0025	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.0015 J	0.0042 J	<0.0025	0.000521 J	0.000774 J	0.000257 J	0.000403 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<b>0.0087 J</b>	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.001	<0.001	<0.0025	<0.00015	<0.000166	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.001	<0.001	<0.0025	<0.00011	<0.00011	<0.00011	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0031	0.0048 J	<0.0075	0.00197 J	0.00217 J	0.00145 J	0.00165 J	0.0023	0.00076 J	0.00076 J	<0.0003	0.0006 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<b>&lt;0.00529</b>	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.00005	0.000056 J	<0.000292	<0.000295	<0.0149	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124	<b>&lt;0.00625</b>	<0.000126	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000762	<b>&lt;0.00385</b>	<0.0000777	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.000385	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.014	0.061	<b>0.17</b>	0.0477	<b>0.306</b>	0.0386	<b>0.121</b>	0.067	0.008	0.005	<0.00019	0.013
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.000783	<0.00079	<b>&lt;0.0399</b>	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000528	<0.000533	<0.0269	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.038	0.13	0.2	0.253	0.428	0.342	0.292	0.26	0.23	0.17	<0.00027	0.19
Acenaphthylene	1.5	4.4	<0.00005	0.0015	0.0015	<0.0000566	<0.0000571	<0.00288	0.00225	<0.000015	0.0017	0.0015	<0.000015	0.0013
Anthracene	7.3	22	0.0017	0.028	0.023	0.0179	0.0222	0.0325	0.0179	0.016	0.019	0.014	<0.00014	0.0087
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	0.00011 J	0.000221 J	0.000226 J	<0.00385	0.000268 J	0.00016	0.00026	0.00011	<0.00005	0.00041
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<b>&lt;0.00385</b>	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	0.00012
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124	<b>&lt;0.00625</b>	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00017 J	<0.00027	<0.00012	<0.000349	<0.000352	<b>&lt;0.0178</b>	0.000679	<0.00011	0.0002 J	0.00013 J	<0.000037	0.00031
Chrysene	0.91	2	<0.00005	<0.00005	0.00013 J	0.000186 J	0.000231 J	<0.00385	0.000241 J	0.00015	0.00022	0.00011	<0.000021	0.00034
Dibenzofuran	0.098	0.29	0.025	<b>0.13</b>	<b>0.16</b>	<b>0.17</b>	<b>0.317</b>	<b>0.22</b>	<b>0.193</b>	<b>0.2</b>	<b>0.15</b>	<b>0.11</b>	0.000031 J	<b>0.14</b>
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.00005	<0.000104	<0.00024	<0.00529	0.000797	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0017	0.0031	0.007	0.0111	0.0181	0.018 J	0.0132	0.013	0.0097	0.0084	<0.00001	0.0097
Fluorene	0.98	2.9	0.025	0.067	0.15	0.17	0.316	0.245	0.202	0.21	0.17	0.13	<0.00003	0.16
Naphthalene	0.49	1.5	0.05	<b>1.5</b>	0.36	0.0828 J	<b>0.661</b>	0.0338 J	0.075	0.012	0.0014	<0.0021	<0.00026	<0.0012
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<0.00529	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000952	<0.00481	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00005	<0.0000575	<0.0000581	<b>&lt;0.0293</b>	<0.0000592	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.015	0.078	0.12	0.13	0.234	0.192	0.162	0.13	0.13	0.093	<0.000021	0.063
Phenol	7.3	22	<0.00005	<0.00005	<0.00005	0.000101 J	<0.0000381	<0.00192	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.00068	0.0026	0.0036	0.00515	0.00818	0.00759 J	0.00649	0.0064	0.0049	0.0042	<0.000019	0.0047
<b>Metals</b>														
Arsenic	0.01	0.01								<b>0.017</b>	0.00133 J	0.00093 J	0.00192 J	0.00192 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-12A 01/13/2020	MW-12B 01/31/2008 DNAPL	MW-12B 01/23/2020 DNAPL	MW-12C 01/30/2008	MW-12C 07/15/2008	MW-12C 02/04/2009	MW-12C 01/19/2010	MW-12C 06/22/2010	MW-12C 01/18/2011	MW-12C 07/26/2011	MW-12C 02/01/2012	MW-12C 07/19/2012	
<b>Volatile Organic Compounds</b>															
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.00052	<0.005	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Benzene	0.005	0.005	<0.0002	0.00344 J	<b>0.03</b>	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Chlorobenzene	0.1	0.1	<0.0003	<0.00047	<0.0075	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Ethylbenzene	0.7	0.7	<0.0003	0.0125	0.014 J	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005
Methylene chloride	0.005	0.005	<0.001	<0.00054	<b>&lt;0.025</b>	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001
Toluene	1	1	<0.0002	0.00515	<0.005	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Vinyl chloride	0.002	0.002			<b>&lt;0.005</b>										
Xylenes (total)	10	10	<0.0003	0.0264	<0.0075	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015
<b>Semivolatile Organic Compounds</b>															
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.0008	<b>&lt;0.0021</b>	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.0029	<0.004	<0.00031	<0.00032	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<b>&lt;0.0019</b>	<b>&lt;0.0058</b>	<0.0002	<0.00021	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<b>&lt;0.0019</b>	<b>&lt;0.0042</b>	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.0038	<0.0021	<0.00041	<0.00042	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	0.0023	<b>0.508</b>	<b>320 J</b>	<0.00041	<0.00042	0.00045	0.00024	0.00011 J	0.00012 J	0.00009 J	<0.00005	<0.00005	0.000086 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.0019	<0.002	<0.0002	<0.00053	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.0024	<0.0047	<0.00026	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.28	0.336	<b>350</b>	<0.00031	<0.00032	0.00052	0.00019 J	<0.00009	0.00012 J	<0.00005	<0.00005	<0.00005	0.00011 J
Acenaphthylene	1.5	4.4	0.0012	0.0127	<b>5.7</b>	<0.00031	<0.00032	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Anthracene	7.3	22	0.0097	0.0267	<b>160</b>	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)anthracene	0.0091	0.02	0.00015	0.00746	<b>79</b>	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<b>&lt;0.0019</b>	<b>22</b>	<0.0002	<b>&lt;0.00021</b>	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<b>&lt;0.0038</b>	<b>&lt;0.003</b>	<0.00041	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00011 J	<0.0019	<0.0037	0.00114 J	<0.00021	0.0003	0.00077	<0.00099	<0.0002	0.0004	<0.0001	<0.00011	<0.00011
Chrysene	0.91	2	0.00015	0.00596	<b>64</b>	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	<b>0.23</b>	<b>0.204</b>	<b>320</b>	<0.00031	<0.00032	0.0004	0.00014 J	<0.00008	0.00011 J	<0.00005	<0.00005	<0.00005	0.000054 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.0019	<0.002	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.0067	0.0508	<b>520</b>	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Fluorene	0.98	2.9	0.28	0.196	<b>420</b>	<0.0002	<0.00021	0.00037	0.00014 J	<0.00007	0.000099 J	0.000071 J	0.0001 J	0.000082 J	0.000082 J
Naphthalene	0.49	1.5	<0.00051	<b>5.55</b>	<b>760</b>	0.000734	0.000833	0.003	0.0017	<0.00046	0.00099	0.00048	<0.00054	0.00052	0.00052
Nitrobenzene	0.049	0.15	<0.000024	<0.0038	<0.0024	<0.00041	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.0024	<0.0025	<0.00026	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.000079	<b>&lt;0.0019</b>	<b>&lt;0.0079</b>	<0.0002	<0.00021	<0.00008	<0.00008	<0.00008	0.00015 J	<0.00005	<0.00005	<0.00005	<0.00005
Phenanthrene	0.73	2.2	0.082	0.322	<b>1200</b>	<0.0002	<0.00021	0.00048	0.00015 J	<0.00007	0.00011 J	<0.00005	<0.00005	<0.00005	0.000059 J
Phenol	7.3	22	<0.000035	<0.0019	<0.0035	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Pyrene	0.73	2.2	0.0031	0.033	<b>330</b>	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
<b>Metals</b>															
Arsenic	0.01	0.01	0.00134 J		<b>0.0491 J</b>										

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-12C 02/05/2013	MW-12C 07/31/2013	MW-12C 01/14/2014	MW-12C 07/25/2014	MW-12C 01/23/2018	MW-12C 03/19/2018	MW-12C 05/16/2018	MW-12C 01/09/2019	MW-12C 07/11/2019	MW-12C 01/13/2020	MW-13 01/30/2008	MW-13 07/15/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052
Benzene	0.005	0.005	<0.00008	<0.00008	<0.00008	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025
Chlorobenzene	0.1	0.1	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047
Ethylbenzene	0.7	0.7	<0.000145	<0.00011	<0.00011	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00025
Methylene chloride	0.005	0.005	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054
Toluene	1	1	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041
Vinyl chloride	0.002	0.002	<0.00011	<0.00011	<0.00011	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041
Xylenes (total)	10	10	<0.00026	<0.00026	<0.00026	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00127
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000104	<0.000105	<0.000106	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008
2,4-Dimethylphenol	0.49	1.5	<0.000292	<0.000295	<0.000298	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00029	<0.00029
2,4-Dinitrotoluene	0.0013	0.003	<0.000123	<0.000124	<0.000125	<0.000126	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019	<0.00019
2,6-Dinitrotoluene	0.0013	0.003	<0.000755	<0.000762	<0.000769	0.209	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.00019
2-Chloronaphthalene	2	5.8	<0.000755	<0.000762	<0.000769	<0.000777	0.000079 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.00038
2-Methylnaphthalene	0.098	0.29	0.000146 J	0.000129 J	0.000164 J	<0.000146	0.000092 J	0.000092 J	0.000078 J	0.00039	0.00021	0.00018	<0.00038	<0.00039
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000783	<0.00079	<0.000798	<0.000806	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00019	<0.00049
4-Nitrophenol	0.049	0.15	<0.000528	<0.000533	<0.000538	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00024
Acenaphthene	1.5	4.4	<0.000755	<0.000762	<0.000769	<0.000114	0.0002	0.0001	<0.000027	0.093	0.000071 J	0.00015	<0.00029	<0.00029
Acenaphthylene	1.5	4.4	<0.000566	<0.000571	<0.000577	<0.000583	<0.000015	0.00064	<0.000015	0.00082	<0.000015	<0.000015	<0.00029	<0.00029
Anthracene	7.3	22	0.0000745 J	<0.0000476	<0.0000481	<0.0000485	<0.000014	<0.000014	<0.000014	0.0084	0.00003 J	<0.000014	0.000955	0.000642
Benzo(a)anthracene	0.0091	0.02	<0.000755	<0.000762	<0.000769	<0.000777	<0.00005	<0.00005	<0.00005	0.00014	<0.00005	<0.00005	<0.00019	<0.00019
Benzo(a)pyrene	0.0002	0.0002	<0.000755	<0.000762	<0.000769	<0.000777	<0.00002	<0.00002	<0.00002	0.000041 J	<0.00002	<0.00002	<0.00019	<0.00019
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000123	<0.000124	<0.000125	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.00039
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000349	<0.000352	<0.000356	<0.000359	<0.00009	0.00013 J	<0.000037	0.00011 J	<0.000037	<0.000037	0.00051 J	<0.00019
Chrysene	0.91	2	<0.000755	<0.000762	<0.000769	<0.000777	<0.000021	<0.000021	<0.000021	0.00013	<0.000021	<0.000021	<0.00019	<0.00019
Dibenzofuran	0.098	0.29	0.0000865 J	0.0000857 J	0.0000979 J	<0.000091	0.0001	<0.00002	0.000052 J	0.067	0.00007 J	0.00012	<0.00029	<0.00029
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000104	<0.00011	<0.000106	<0.000107	0.000034 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019
Fluoranthene	0.98	2.9	<0.000066	<0.000067	<0.0000673	<0.000068	<0.00001	<0.00001	<0.00001	0.0071	<0.00001	0.000091 J	0.00046 J	<0.00019
Fluorene	0.98	2.9	0.000149 J	0.000102 J	<0.0000673	<0.000068	<0.00003	<0.00003	0.000087 J	0.085	0.000052 J	0.00019	<0.00019	<0.00019
Naphthalene	0.49	1.5	0.000729	0.000585 J	<0.000853	<0.000598	0.00046	0.0003	<0.00061	<0.00017	<0.0031	0.0017	<0.00038	<0.00039
Nitrobenzene	0.049	0.15	<0.000104	<0.000105	<0.000106	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.00039
N-Nitrosodiphenylamine	0.19	0.42	<0.0000943	<0.0000952	<0.0000962	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00024
Pentachlorophenol	0.001	0.001	<0.000575	<0.000581	<0.000587	<0.000592	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019	<0.00019
Phenanthrene	0.73	2.2	<0.000566	<0.000571	<0.000872	<0.000808	0.000069 J	<0.000021	0.000048 J	0.052	0.000039 J	0.00015	<0.00019	<0.00019
Phenol	7.3	22	<0.0000377	<0.0000381	<0.0000385	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	0.0004	0.00012 J	<0.00019	<0.00019
Pyrene	0.73	2.2	<0.000104	<0.000105	<0.000106	<0.000107	<0.000019	<0.000019	<0.000019	0.0031	<0.000019	0.000068 J	<0.00019	<0.00019
<b>Metals</b>														
Arsenic	0.01	0.01					0.0025	0.00184 J	0.0017 J	0.000796 J	0.002	0.00195 J		

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-13 02/04/2009	MW-13 01/19/2010	MW-13 06/22/2010	MW-13 01/18/2011	MW-13 07/26/2011	MW-13 02/02/2012	MW-13 07/16/2012	MW-13 02/05/2013	MW-13 07/31/2013	MW-13 01/14/2014	MW-13 07/25/2014	MW-13 01/23/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	0.000401 J	<0.00012	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0002
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.000107	<0.000212	<0.000107	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	R	<0.000301	<0.000596	<0.000301	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000126	<0.00025	<0.000126	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	0.00066	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000777	<0.000154	<0.0000777	<0.000042
2-Chloronaphthalene	2	5.8	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.000154	<0.0000777	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00007	0.00076	<0.00007	0.000075 J	0.00026	<0.00005	0.000063 J	<0.000066	<0.000068	0.000141 J	<0.000687	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	R	<0.000806	<0.0016	<0.000806	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	R	<0.000544	<0.00108	<0.000544	<0.000047
Acenaphthene	1.5	4.4	<0.00009	0.00011 J	<0.00009	<0.00009	0.00033	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.000154	<0.000329	<0.000027
Acenaphthylene	1.5	4.4	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000566	0.0000715 J	<0.000115	<0.0000583	0.00003 J
Anthracene	7.3	22	0.0002	0.00043	<0.00007	<0.00007	0.00037	0.00068 J	0.00011 J	0.0011	0.000878	<0.00118	<0.000587	0.00047
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00014 J	<0.00005	<0.0000755	<0.0000777	<0.000154	<0.0000777	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	0.000073 J	<0.00005	<0.0000755	<0.0000777	<0.000154	<0.0000777	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000126	<0.00025	<0.000126	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00035	0.0016	<0.00044	<0.0002	0.00027	0.00043 J	<0.0001	<0.000349	<0.000359	<0.000712	<0.000359	<0.00011
Chrysene	0.91	2	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00017 J	<0.00005	<0.0000755	<0.0000777	<0.000154	<0.0000777	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	0.00019 J	<0.00008	<0.00008	0.00034	0.00063 J	0.00019 J	<0.0000755	<0.0000777	<0.000154	<0.000257	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00007	0.0001 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.000104	<0.000211	<0.000212	0.000122 J	<0.00002
Fluoranthene	0.98	2.9	<0.00007	<0.00007	<0.00007	<0.00007	0.00067 J	0.00015 J	0.00013 J	<0.000066	<0.000068	<0.000135	<0.000068	<0.00001
Fluorene	0.98	2.9	<0.00007	<0.00007	<0.00007	0.000072 J	0.00035	<0.00005	0.00012 J	<0.000066	<0.000068	<0.000135	<0.000188	<0.00003
Naphthalene	0.49	1.5	<0.0001	0.007	<0.0001	0.0005	0.00087	<0.00022	<0.00023	<0.0000755	<0.000226	<0.00126	0.0039	0.00014
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.000107	<0.000212	<0.000107	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000971	<0.000192	<0.0000971	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	R	<0.000592	<0.00117	<0.000592	<0.000079
Phenanthrene	0.73	2.2	<0.00007	0.00014 J	0.0002	<0.00007	0.00029	0.00015 J	0.00049	<0.0000566	<0.0000583	<0.000115	<0.000163	<0.000021
Phenol	7.3	22	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	R	<0.0000388	<0.0000769	<0.0000388	<0.000035
Pyrene	0.73	2.2	<0.00007	<0.00007	<0.00007	<0.00007	0.00011 J	0.0002 J	0.000089 J	<0.000104	<0.000107	<0.000212	<0.000107	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01												0.00303

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-13 03/18/2018	MW-13 05/15/2018	MW-13 01/08/2019	MW-13 07/11/2019	MW-13 01/14/2020	MW-14 01/30/2008	MW-14 07/15/2008	MW-14 02/04/2009	MW-14 02/04/2009 Duplicate	MW-14 01/19/2010	MW-14 01/19/2010 Duplicate	MW-14 06/22/2010
<b><i>Volatile Organic Compounds</i></b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.001
<b><i>Semivolatile Organic Compounds</i></b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.000063	<0.00004	<0.00029	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.00038	<0.00012	<0.00012	<0.0001	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	0.000077 J	0.000054 J	0.00047 J	0.000782	0.00075	0.00078	0.00064	0.0006	0.00049
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00048	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000027	0.000066 J	0.00022	0.00223	0.000515	0.00047	0.0005	0.00043	0.00043	0.00041
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	0.000033 J	<0.000015	<0.00029	<0.00029	<0.00006	<0.00006	<0.00007	<0.00007	<0.00007
Anthracene	7.3	22	0.000039 J	0.000085 J	0.00039	0.00047	<0.000014	0.000678	<0.00019	<0.00007	<0.00007	<0.00007	0.00013 J	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000078 J	0.00011 J	<0.000037	0.000054 J	<0.000037	0.0004 J	<0.00019	0.00081	0.0027	0.0054 J	0.00037 J	<0.00077
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	0.000039 J	0.00016	0.000491	0.000502	0.00045	0.00045	0.0004	0.00044	0.00037
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.000029 J	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Fluoranthene	0.98	2.9	<0.00001	0.000015 J	<0.00001	0.000032 J	<0.00001	0.000506	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	0.000042 J	0.00016	<0.00019	<0.00019	<0.00007	0.00014 J	0.00013 J	0.0001 J	<0.00007
Naphthalene	0.49	1.5	<0.000083	<0.00002	<0.00022	<0.0011	0.00026	0.00222	0.00349	0.0032	0.0033	0.003	0.0026	0.0022
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000599	0.00061	0.00035	0.00039	0.00041	0.00043	0.00044
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.000019	0.00002 J	<0.000019	0.000084 J	<0.000019	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
<b><i>Metals</i></b>														
Arsenic	0.01	0.01	0.00984	0.014	0.0602	0.0715	0.0642							

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-14 01/17/2011	MW-14 07/26/2011	MW-14 07/26/2011 Duplicate	MW-14 02/02/2012	MW-14 02/02/2012 Duplicate	MW-14 07/16/2012	MW-14 07/16/2012 Duplicate	MW-14 02/05/2013	MW-14 07/31/2013	MW-14 01/14/2014	MW-14 07/18/2014	MW-14 01/23/2018
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002
Benzene	0.005	0.005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0011	<0.0011	<0.0011	<0.0011	<0.0005	<0.0005	<0.00011	0.000123 J	<0.00011	<0.00011	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0013	<0.0013	<0.0013	<0.0013	<0.001	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.001
Toluene	1	1	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002
Vinyl chloride	0.002	0.002								<0.00011				<0.00011
Xylenes (total)	10	10	<0.001	<0.0031	<0.0031	<0.0031	<0.0031	<0.0015	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026	<0.0003
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000104	<0.000107	<0.000106	<0.000109	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00005	<0.00005	0.00005 J	0.0013 J	<0.00005	<0.00005	<0.000292	<0.000301	<0.000298	<0.000307	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000123	<0.000126	<0.000125	<0.000129	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000777	<0.0000769	0.0788	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.0000769	<0.0000792	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00039	0.00034	0.00031	0.00064 J	0.00017 J	0.0003	0.00044	0.000402 J	0.000304 J	0.000321 J	0.000336 J	0.00019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000783	<0.000806	<0.000798	<0.000822	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000528	R	<0.000538	<0.000554	<0.000047
Acenaphthene	1.5	4.4	0.00033	0.00032	0.00028	0.00005 J	0.00015 J	0.0003	0.00035	0.0006	0.000549	0.000943	0.000619	0.00019
Acenaphthylene	1.5	4.4	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000566	<0.0000583	<0.0000577	<0.0000594	<0.000015
Anthracene	7.3	22	<0.00007	<0.00005	0.000087 J	<0.00005	<0.00005	0.000069 J	<0.00005	0.000277 J	0.000198 J	<0.000179	0.000139 J	0.000067 J
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.0000769	<0.0000792	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.0000769	<0.0000792	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000123	<0.000126	<0.000125	<0.000129	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00029	0.00047	0.00065	<0.0001	<0.00013	0.00011 J	0.00011 J	<0.000349	<0.000359	<0.000356	0.000615	<0.000091
Chrysene	0.91	2	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.0000769	<0.0000792	<0.000021
Dibenzofuran	0.098	0.29	0.0003	0.00031	0.00027	0.00012 J	0.00015 J	0.00032	0.00034	0.000467 J	0.000372 J	0.000443 J	0.000437 J	0.00024
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000104	<0.000179	<0.000106	<0.000109	<0.00002
Fluoranthene	0.98	2.9	<0.00007	<0.00005	<0.00005	0.00024	0.00031	0.00055 J	<0.00005	0.0000794 J	0.000274 J	0.0000744 J	<0.0000693	0.000028 J
Fluorene	0.98	2.9	0.000079 J	<0.00005	0.000065 J	<0.00005	<0.00005	0.000076 J	<0.00005	<0.000066	<0.000068	0.00008 J	0.0000901 J	<0.00003
Naphthalene	0.49	1.5	0.0024	0.0014	0.0012	0.00034 J	0.0014 J	<0.0015	<0.002	0.00211	0.00216 J	<0.00183	0.00143	0.00067
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	0.00008 J	<0.000104	<0.000107	<0.000106	<0.000109	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000971	<0.0000962	<0.000099	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000575	<0.0000592	<0.0000587	<0.0000604	<0.000079
Phenanthrene	0.73	2.2	0.0003	0.00033	0.00032	0.00011 J	0.00014 J	0.00038	0.00039	0.000484	0.000662	<0.000591	0.000506	0.00035
Phenol	7.3	22	<0.00007	<0.00005	<0.00005	0.00005 J	0.00075 J	<0.00005	<0.00005	<0.0000377	0.000398 J	<0.0000385	<0.0000396	<0.000035
Pyrene	0.73	2.2	<0.00007	<0.00005	<0.00005	0.00029	0.00029	<0.00005	<0.00005	<0.000104	0.000164 J	<0.000106	<0.000109	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01												<0.0004

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-14 03/18/2018	MW-14 05/15/2018	MW-14 01/08/2019	MW-14 07/11/2019	MW-14 01/14/2020	MW-15A 01/30/2008	MW-15A 07/15/2008	MW-15A 02/04/2009	MW-15A 01/18/2010	MW-15A 06/23/2010	MW-15A 01/17/2011	MW-15A 07/13/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.00161 J	<0.00025	0.0018 J	0.0016 J	0.0017 J	0.00074 J	0.0016 J
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.00122 J	<0.00025	0.0019 J	0.0015 J	0.0017 J	<0.0005	0.0019 J
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041	<0.0005	<0.0005	0.00055 J	<0.0005	<0.001
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0056 J	<0.00127	0.0039 J	0.0015 J	0.0047 J	<0.001	0.0038 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00013	<0.00004	<0.0003	0.00042 J	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.0002	<0.0002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0004	<0.00041	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.00014	<0.00019	<0.00019	0.00019	<0.00019	0.0127	<b>0.0995</b>	0.044	0.033	0.042	0.038	<b>0.14</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00051	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00025	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Acenaphthene	1.5	4.4	0.00027	<0.000027	<0.000027	0.0001	0.000091 J	0.134	0.442	0.17	0.17	0.16	0.27	0.2
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.0003	<0.00031	0.0017	0.0015	0.00097	0.0011	0.00097
Anthracene	7.3	22	0.000052 J	<0.000014	0.000052 J	0.000089 J	<0.000014	0.00377	0.00432	0.003	0.0036	0.0049	0.0063	0.0053
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	0.00023 J	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.0004	<0.00041	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.00014 J	<0.000037	<0.000037	<0.000037	<0.0002	<0.0002	0.0026	<0.00073	<0.00084	0.0016	<0.00012
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Dibenzofuran	0.098	0.29	<0.00025	<0.00002	<0.00002	0.000087 J	<0.00002	0.0239	<b>0.156</b>	0.047	0.043	0.048	0.05	0.078
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	0.000022 J	<0.00002	<0.00002	<0.0002	<0.0002	0.00029	0.00011 J	<0.00007	<0.00007	<0.00005
Fluoranthene	0.98	2.9	0.000025 J	<0.00001	<0.00001	0.000067 J	0.000064 J	0.00178	0.00183	0.0011	0.0015	0.002	0.0023	0.0021
Fluorene	0.98	2.9	0.000094 J	<0.00003	<0.00003	0.000069 J	0.000057 J	0.0394	0.18	0.059	0.06	0.062	0.076	0.092
Naphthalene	0.49	1.5	<0.00057	<0.00002	<0.00002	<0.0002	<0.00002	0.00684	0.271	0.048	<0.0018	0.036	0.0023	0.087
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.0004	<0.00041	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00025	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	0.00055
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005
Phenanthrene	0.73	2.2	0.00032	<0.000021	<0.000021	0.000059 J	<0.000021	0.0039	0.0229	0.0095	0.0074	0.012	0.019	0.036
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	0.00019 J	<0.000035	<0.0002	<0.0002	<0.00007	<0.00007	0.0002	<0.00007	<0.00005
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	0.000032 J	0.00006 J	0.00127	0.000664	0.00042	0.00062	0.00076	0.00095	0.00089
<b>Metals</b>														
Arsenic	0.01	0.01	<0.0004	<0.0004	0.000752 J	0.00133 J	0.00185 J							

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-15A 02/02/2012	MW-15A 07/19/2012	MW-15A 01/30/2013	MW-15A 07/30/2013	MW-15A 01/14/2014	MW-15A 07/17/2014	MW-15A 01/23/2018	MW-15A 03/18/2018	MW-15A 05/15/2018	MW-15A 01/08/2019	MW-15A 07/10/2019	MW-15A 01/14/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	0.0012 J	0.0016 J	0.0016	0.0013	0.00106	0.00161	<0.0002	0.00051 J	0.0006 J	<0.0002	0.00074 J	0.00034 J
Chlorobenzene	0.1	0.1	<0.001	<0.0005	<0.00012	0.000121 J	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.0024 J	0.0012 J	0.00066 J	0.000799 J	0.000627 J	0.00101	<0.0003	<0.0003	<0.0003	<0.0003	0.00035 J	<0.0003
Methylene chloride	0.005	0.005	<0.0013	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.001	<0.0005	0.000221 J	0.000199 J	0.00034 J	0.000595 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.0073 J	0.0097 J	0.00417	0.00527	0.00337	0.00854	0.0018	<0.0003	<0.0003	0.0008 J	0.0026	0.00093 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.000104	<0.000107	<0.00529	<0.000104	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00036
2,4-Dimethylphenol	0.49	1.5	0.00059	0.00056	0.002	<0.000301	<0.0149	<0.000292	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.000123	<0.000126	<0.00625	<0.000123	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.0000755	<0.0000777	<0.00385	<0.0000755	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.00385	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.001	0.046	0.00997	0.124	0.075	0.059	0.0076	0.016	0.034	0.098	0.077	0.018
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.000783	<0.000806	<0.0399	<0.000783	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.000528	<0.000544	<0.0269	<0.000528	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.13	0.13	0.141	0.332	0.3	0.205	0.13	0.099	0.1	0.1	0.11	0.14
Acenaphthylene	1.5	4.4	0.00071	0.0012	<0.0000566	<0.0000583	<0.00288	<0.0000566	0.002	0.0007	0.0069	<0.000015	0.00045	0.00066
Anthracene	7.3	22	0.0028	0.0046	0.00313	0.0085	0.0111 J	0.00642	0.0026	0.0024	0.0032	0.0025	0.0026	0.0048
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.00385	<0.0000755	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.00385	<0.0000755	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.000123	<0.000126	<0.00625	<0.000123	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0001	<0.0001	<0.000349	<0.000359	<0.0178	<0.000349	<0.000037	<0.000037	0.0093	<0.000037	<0.000037	<0.000037
Chrysene	0.91	2	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.00385	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.028	0.046	0.0416	0.104	0.0693	0.0572	0.029	0.024	0.018	0.023	0.035	0.036
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.000104	0.000187 J	<0.00529	<0.000104	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.00094	0.0015	0.000885	0.00361	<0.00337	0.00257	0.002	0.0012	0.0016	0.0012	0.0018	0.0031
Fluorene	0.98	2.9	0.043	0.063	0.056	0.139	0.114	0.0822	0.041	0.036	0.029	0.038	0.054	0.065
Naphthalene	0.49	1.5	0.008	0.27	0.0501 J	0.526	0.326	0.248	0.0005	<0.00034	<0.00037	<0.00032	<0.00066	0.00022
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.000104	<0.000107	<0.00529	<0.000104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.0000943	<0.0000971	<0.00481	<0.0000943	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.000575	<0.000592	<0.0293	<0.000575	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.0056	0.014	0.00792	0.052	0.0375	0.0203	0.0046	0.0054	0.0074	0.009	0.0095	0.025
Phenol	7.3	22	<0.00005	<0.00005	<0.0000377	<0.0000388	<0.00192	<0.0000377	<0.000035	<0.000035	<0.000035	<0.000035	0.00039	<0.000035
Pyrene	0.73	2.2	0.00053	0.00084	0.000496	0.00154	<0.00529	0.00101	0.0009	0.00063	0.00078	0.00051	0.00075	0.0013
<b>Metals</b>														
Arsenic	0.01	0.01							0.0264	0.0137	0.019	0.027	0.0251	0.0441

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-15B 02/02/2012	MW-15B 07/19/2012	MW-15B 01/30/2013	MW-15B 07/30/2013	MW-15B 01/14/2014	MW-15B 07/17/2014	MW-15B 01/23/2018	MW-15B 03/18/2018	MW-15B 05/15/2018	MW-15B 01/08/2019	MW-15B 07/10/2019	MW-15B 01/14/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.005	<0.0025	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.005	<b>0.0053 J</b>	0.0022	0.00484	0.00101	0.00292	<0.0002	<0.0002	0.00071 J	<0.0002	0.0023	<0.0002
Chlorobenzene	0.1	0.1	<0.005	<0.0025	<0.00012	0.000124 J	<0.00012	0.000136 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.02 J	0.014 J	0.00159	0.00399	0.00019 J	0.00903	<0.0003	<0.0003	0.001	<0.0003	0.0031	<0.0003
Methylene chloride	0.005	0.005	<b>&lt;0.0065</b>	<0.005	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.005	<0.0025	<0.00015	0.000155 J	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.016	<0.0075	0.000356 J	0.00876	0.000876 J	0.00464	<0.0003	<0.0003	<0.0003	<0.0003	0.0025	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.000104	<0.000107	<0.00106	<0.000104	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00043	<0.00005	<0.000292	<0.000301	<0.00298	<0.000292	<0.00004	<0.00004	<0.00004	<0.00004	<0.00016	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.000123	<0.000126	<0.00125	<0.000123	<0.000059	<0.000058	<0.000059	<0.000059	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.0000755	<0.0000777	<0.000769	<0.0000755	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.0000755	<0.0000777	<0.000769	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<b>0.29</b>	0.023	<0.000744	0.00327	0.00325 J	0.00622	<0.000019	<0.000019	0.00015	<0.000019	0.016	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.000783	<0.000806	<b>&lt;0.00798</b>	<0.000783	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.000528	<0.000544	<0.00538	<0.000528	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.17	0.075	0.0413	0.114	0.134	0.0653	<0.000027	<0.000027	0.012	0.0026	0.029	0.000073 J
Acenaphthylene	1.5	4.4	0.0011	0.0008	0.000987	<0.0000583	0.00148 J	<0.0000566	0.000059 J	<0.000015	0.00027	0.00015	0.00034	<0.000015
Anthracene	7.3	22	0.039	0.0071	0.00179	0.00581	0.00665	0.00517	0.00034	0.00016	0.00058	0.00023	0.0016	0.00011
Benzo(a)anthracene	0.0091	0.02	0.00016 J	0.00017 J	<0.0000755	0.000218 J	0.000868 J	0.00031 J	0.00017	<0.00005	0.00011	<0.000051	0.00016	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.0000755	<0.0000777	<b>&lt;0.000769</b>	<0.0000755	0.00007 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.000123	<0.000126	<b>&lt;0.00125</b>	<0.000123	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0001	<0.00018	<0.000349	<0.000359	<0.00356	0.000548	<0.00007	<0.000037	0.0001 J	<0.000037	<0.000037	0.000055 J
Chrysene	0.91	2	0.00019 J	0.00013 J	<0.0000755	0.000167 J	<0.000769	0.000228 J	0.00014	<0.000021	0.000098 J	<0.000021	0.00013	<0.000021
Dibenzofuran	0.098	0.29	<b>0.15</b>	0.052	0.0127	0.0589	0.0509	0.0272	0.0005	<0.00002	0.0025	0.00014	0.013	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.000104	0.000187 J	<0.00106	<0.000104	<0.00002	<0.00002	0.000042 J	0.000022 J	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.012	0.0062	0.00101	0.00615	0.0131	0.00736	0.0027	<0.00001	0.0031	0.00045	0.0042	0.00011
Fluorene	0.98	2.9	0.084	0.036	0.011	0.0459	0.0443	0.0231	0.00017	<0.00003	0.0028	0.000055 J	0.01	0.000062 J
Naphthalene	0.49	1.5	<b>2.5</b>	<b>0.82</b>	0.0569 J	<b>0.943</b>	0.248	0.452	<0.00002	<0.00002	0.02	<0.00002	0.33	<0.00002
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.000104	<0.000107	<0.00106	<0.000104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.0000943	<0.0000971	<0.000962	<0.0000943	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.000575	<0.000592	<b>&lt;0.00587</b>	<0.000575	<0.00008	<0.000079	<0.00008	<0.00008	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.08	0.052	0.00199	0.0376	0.0257	0.0204	0.00033	<0.000021	0.00016	<0.000021	0.0065	0.00005 J
Phenol	7.3	22	<0.00012	<0.00005	<0.0000377	<0.0000388	0.00141 J	0.00112	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	<0.000035
Pyrene	0.73	2.2	0.005	0.0031	0.000513	0.00291	0.00569	0.00406	0.0016	<0.000019	0.0017	0.00027	0.0018	0.00006 J
<b>Metals</b>														
Arsenic	0.01	0.01							0.00895	0.00329	<b>0.0111</b>	0.00244	<b>0.0133</b>	0.00343

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-15C 01/30/2008	MW-15C 07/15/2008	MW-15C 02/04/2009	MW-15C 01/18/2010	MW-15C 06/23/2010	MW-15C 01/17/2011	MW-15C 07/13/2011	MW-15C 02/02/2012	MW-15C 07/19/2012	MW-15C 01/30/2013	MW-15C 07/30/2013	MW-15C 01/14/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00014
Benzene	0.005	0.005	0.00109 J	<0.00025	0.00096 J	0.0012 J	0.001 J	0.00096 J	<0.001	<0.001	<0.0005	0.000951 J	0.000831 J	0.000863 J
Chlorobenzene	0.1	0.1	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00012
Ethylbenzene	0.7	0.7	0.00135 J	<0.00025	0.00068 J	0.00058 J	<0.0005	<0.0005	<0.0011	0.0017 J	<0.0005	0.000408 J	0.000203 J	0.000275 J
Methylene chloride	0.005	0.005	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00015
Toluene	1	1	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	0.000323 J	0.000263 J	0.000305 J
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	0.000604 J	0.000839 J	0.000581 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0104	<0.000107	<0.00106
2,4-Dimethylphenol	0.49	1.5	<0.00029	<0.00032	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	113	<0.000301	<0.00298
2,4-Dinitrotoluene	0.0013	0.003	<0.0002	<0.00021	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0123	<0.000126	<0.00125
2,6-Dinitrotoluene	0.0013	0.003	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.00755	<0.0000777	<0.000769
2-Chloronaphthalene	2	5.8	<0.00039	<0.00042	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00755	<0.0000777	<0.000769
2-Methylnaphthalene	0.098	0.29	<0.00039	<0.00042	0.000084 J	<0.00007	<0.00007	<0.00007	<0.00005	0.000099 J	0.00022	0.35	<0.000068	<0.000673
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.00053	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.0783	<0.000806	<0.00798
4-Nitrophenol	0.049	0.15	<0.00024	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0528	<0.000544	<0.00538
Acenaphthene	1.5	4.4	0.0293	0.103	0.034	0.0097	0.013	0.032	0.016	0.041	0.042	0.13	0.0574	0.0912
Acenaphthylene	1.5	4.4	<0.00029	0.000651	0.00052	0.00041	0.00062	0.0011	0.0012	0.0013	0.002	<0.00566	0.00268	<0.000577
Anthracene	7.3	22	<0.0002	0.000731	0.00078	0.00031	<0.00007	<0.00007	<0.00005	0.00021	0.00045	0.0191 J	0.0000945 J	<0.000481
Benzo(a)anthracene	0.0091	0.02	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00755	<0.0000777	<0.000769
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00755	<0.0000777	<0.000769
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00039	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0123	<0.000126	<0.00125
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00044 J	<0.00021	<0.0002	<0.00065	<0.00059	0.00044	<0.00057	<0.0001	<0.00012	<0.0349	<0.000359	<0.00356
Chrysene	0.91	2	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00755	<0.0000777	<0.000769
Dibenzofuran	0.098	0.29	0.0336	0.0904	0.034	0.0075	0.005	0.018	0.0046	0.027	0.021	0.116	0.0141	0.0317
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.000059 J	<0.0104	<0.000107	<0.00106
Fluoranthene	0.98	2.9	0.000607	0.00103	0.0006	0.00029	0.0002 J	0.0003	0.00031	0.0016	0.00079	<0.0066	0.000634	0.00158 J
Fluorene	0.98	2.9	0.00328	0.00278	0.0027	0.0011	0.00071	0.0017	0.00074	0.0025	0.0014	0.0769	0.00159	0.00224 J
Naphthalene	0.49	1.5	0.00137	0.00195	0.0016	<0.00057	<0.00094	0.00091	<0.00046	0.0011	0.0018	89.7	<0.00122	<0.00182
Nitrobenzene	0.049	0.15	<0.00039	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0104	<0.000107	<0.00106
N-Nitrosodiphenylamine	0.19	0.42	<0.00024	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00943	<0.0000971	<0.000962
Pentachlorophenol	0.001	0.001	<0.0002	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00755	<0.000592	<0.00587
Phenanthrene	0.73	2.2	<0.0002	0.00028 J	<0.00007	0.00014 J	<0.00007	<0.00007	0.00019 J	0.0015	0.00038	0.0868	0.000354 J	<0.000577
Phenol	7.3	22	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	61.8	<0.000388	<0.000385
Pyrene	0.73	2.2	0.000542	0.00052	0.00027	0.00012 J	0.00011 J	0.00015 J	0.00018 J	0.00093	0.00046	<0.0104	0.00037 J	<0.00106
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-15C 07/17/2014	MW-15C 01/23/2018	MW-15C 03/18/2018	MW-15C 05/15/2018	MW-15C 01/08/2019	MW-15C 07/10/2019	MW-15C 01/14/2020	MW-16 01/31/2008	MW-16 07/15/2008	MW-16 02/05/2009	MW-16 01/18/2010	MW-16 06/23/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0025	<0.0025	<0.0025
Benzene	0.005	0.005	0.000781 J	0.00063 J	0.00053 J	0.00052 J	0.00058 J	0.0005 J	0.00059 J	<b>0.0383</b>	<b>0.11</b>	<b>0.048</b>	<b>0.031</b>	<b>0.058</b>
Chlorobenzene	0.1	0.1	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0025		<0.0025
Ethylbenzene	0.7	0.7	<0.000219 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.037	0.084	0.034	0.021 J	0.032
Methylene chloride	0.005	0.005	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0025	<0.0025	<b>&lt;0.0062</b>
Toluene	1	1	0.00019 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.00619	0.0382	0.0025 J	0.0034 J	0.01 J
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.000392 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.052	0.121	0.036 J	0.027 J	0.04 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000104	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.0004	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.000292	<0.00004	<0.00004	<0.00004	<0.00004	0.0052	<0.00004	<0.00029	<0.0016	0.0039	0.0025	0.0054
2,4-Dinitrotoluene	0.0013	0.003	<0.000123	<0.000058	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.00019	<0.0011	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0000755	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.0011	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.0021	<0.00012	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.000066	<0.000019	<0.000019	<0.000019	<0.000019	<0.000066	<0.000019	0.0747	<b>0.175</b>	<b>0.13</b>	0.079	0.04
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000783	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<b>&lt;0.0026</b>	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000528	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.0013	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	0.0455	0.027	0.021	0.02	0.02	0.017	0.016	0.215	0.939	0.24	0.21	0.18
Acenaphthylene	1.5	4.4	0.00342	0.0027	0.002	0.022	0.014	0.0015	0.00089	<0.00029	0.0067	0.0044	0.0041	0.0031
Anthracene	7.3	22	0.000315 J	0.00047	0.000095 J	0.000067 J	0.00031	0.00064	0.00057	0.0151	0.0321	0.011	0.0084	0.0076
Benzo(a)anthracene	0.0091	0.02	<0.0000755	<0.00005	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00019	<0.0011	0.00014 J	0.00011 J	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.0000755	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<b>&lt;0.0011</b>	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000123	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<b>&lt;0.0021</b>	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000526	<0.000067	<0.000037	0.000072 J	<0.000037	<0.000037	<0.000037	<0.00019	<0.0011	0.0005	<0.0012	<0.0014
Chrysene	0.91	2	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00036 J	<0.0011	0.00014 J	0.000088 J	<0.00007
Dibenzofuran	0.098	0.29	0.0102	0.0081	0.0029	0.0046	0.0053	0.0039	0.0045	<b>0.112</b>	<b>0.253</b>	<b>0.14</b>	<b>0.12</b>	0.091
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000104	<0.00002	<0.00002	0.000037 J	<0.00002	<0.00002	<0.00002	<0.00019	<0.0011	0.00025	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.000763	0.00094	0.00054	0.00074	0.00075	0.00086	0.00066	0.00769	0.0142	0.0064	0.0037	0.0049
Fluorene	0.98	2.9	0.00135	0.001	0.00045	0.00064	0.00056	0.00097	0.00051	0.114	0.222	0.088	0.096	0.086
Naphthalene	0.49	1.5	<0.000748	0.00042	<0.00031	<0.00039	<0.00032	0.019	0.00024	<b>1.9</b>	<b>18.9</b>	<b>4.1</b>	<b>1.9</b>	<b>1.4</b>
Nitrobenzene	0.049	0.15	<0.000104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.0021	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0000943	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.0013	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000575	<0.000079	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.00019	<b>&lt;0.0011</b>	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.0000566	<0.000021	0.00019	0.00041	<0.000021	<0.00039	0.00008 J	0.0421	0.0743	0.04	0.038	0.042
Phenol	7.3	22	<0.0000377	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	0.0047	0.00022	0.00013 J	<0.00007
Pyrene	0.73	2.2	0.00043 J	0.00049	0.00035	0.00043	0.00041	0.00043	0.00038	0.00615	0.0114	0.004	0.0027	0.0025
<b>Metals</b>														
Arsenic	0.01	0.01		0.000738 J	0.000598 J	0.000777 J	0.000629 J	0.000889 J	0.000773 J					

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-16 01/18/2011	MW-16 07/14/2011	MW-16 02/01/2012	MW-16 07/16/2012	MW-16 01/30/2013	MW-17 01/31/2008	MW-17 07/15/2008	MW-17 02/04/2009	MW-17 01/18/2010	MW-17 06/23/2010	MW-17 01/17/2011	MW-17 07/13/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0025	<0.01	<0.01	<0.0005	<0.00014	<0.00052	<0.01	<0.0005	<0.0025	<0.0025	<0.0025	<0.001
Benzene	0.005	0.005	0.025 J	0.068	0.025 J	0.056	0.0376	0.545	0.448	0.65	0.59	0.65	0.31	0.45
Chlorobenzene	0.1	0.1	<0.0025	<0.01	<0.01	<0.0005	0.00051 J	<0.00047	<0.0094	<0.0005	<0.0025	<0.0025	<0.0025	<0.001
Ethylbenzene	0.7	0.7	0.023 J	0.038 J	0.021 J	0.027	0.0211	0.193	0.142	0.26	0.26	0.2	0.21	0.21
Methylene chloride	0.005	0.005	<0.0025	<0.013	<0.013	<0.001	<0.00015	<0.00054	<0.011	<0.0005	<0.0025	<0.0056	<0.0025	<0.0013
Toluene	1	1	<0.0025	<0.01	<0.01	0.007	0.00095 J	0.909	0.728	1.1	1	0.88	0.97	0.85
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.025 J	<0.031	<0.031	0.034	0.0348	0.582	0.44	0.55	0.72	0.61	0.64	0.54
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.002	<0.008	<0.0001	0.001 J	<0.001	<0.0001	<0.00005
2,4-Dimethylphenol	0.49	1.5	0.0022 J	0.012	0.00077	0.0054	<0.000292	11.7	13.4	2.6	3.7	13	3.9 J	2.7
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.0038	<0.021	<0.00009	0.0009 J	<0.0009	<0.00009	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0038	<0.021	<0.00007	0.0007 J	<0.0007	<0.00007	<0.00006
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0076	<0.042	<0.00012	0.001 J	<0.001	<0.0001	<0.00005
2-Methylnaphthalene	0.098	0.29	0.038	0.082	0.034	0.045	0.0467 J	0.42	0.582	0.27	0.56 J	0.39	0.97	0.75
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.000783	<0.0038	<0.053	<0.00008	0.0008 J	<0.0008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.00005	<0.00005	<0.000528	<0.0048	<0.026	<0.00007	0.0007 J	<0.0007	<0.00007	<0.00005
Acenaphthene	1.5	4.4	0.21	0.23	0.21	0.28	0.281	0.137	0.241	0.094	0.17 J	0.071	0.52	0.22
Acenaphthylene	1.5	4.4	0.0035	0.0032	0.0021	0.0028	<0.0000566	<0.0057	<0.032	0.0041	0.0067 J	0.003	0.008	0.0069
Anthracene	7.3	22	0.012	0.017	0.0034	0.017	0.0182 J	0.0115	0.022	0.0099	0.013 J	0.0075	0.12	0.014
Benzo(a)anthracene	0.0091	0.02	0.00017 J	0.00022	0.000059 J	0.00013 J	0.000339 J	<0.0038	<0.021	0.0004	0.0007 J	<0.0007	0.033	0.00047
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0038	<0.021	0.00014 J	0.0008 J	<0.0008	0.0097	0.00027
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.0076	<0.042	<0.00009	0.0009 J	<0.0009	<0.00009	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0002	<0.00025	<0.0002	<0.0001	<0.000349	<0.0038	<0.021	<0.0002	0.002 J	<0.002	0.022	<0.00044
Chrysene	0.91	2	0.00011 J	0.00019 J	0.000053 J	0.0001 J	0.000225 J	<0.0038	<0.021	0.00032	0.0007 J	<0.0007	0.025	0.00047
Dibenzofuran	0.098	0.29	0.13	0.13	0.098	0.17	0.158	0.115	0.195	0.079	0.15 J	0.065	0.47	0.19
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	0.000099 J	<0.00005	<0.00005	<0.000104	<0.0038	<0.021	<0.00007	0.0007 J	<0.0007	<0.00007	<0.00005
Fluoranthene	0.98	2.9	0.0059	0.006	0.0021	0.0059	0.00836	0.0044	<0.021	0.0035	0.0037 J	0.0022	0.17	0.0039
Fluorene	0.98	2.9	0.12	0.14	0.085	0.15	0.147	0.0701	0.109	0.047	0.076 J	0.039	0.42	0.12
Naphthalene	0.49	1.5	1.2	1.8	1.8	2.1	1.81	23.6	25.5	9.7	16	15	16	19
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.0076	<0.042	<0.00009	0.0009 J	<0.0009	<0.00009	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0048	<0.026	<0.00009	0.0009 J	<0.0009	<0.00009	<0.00005
Pentachlorophenol	0.001	0.001	<0.00008	0.000061 J	<0.00005	<0.00005	<0.000575	<0.0038	<0.021	<0.00008	0.0008 J	<0.0008	<0.00008	<0.00005
Phenanthrene	0.73	2.2	0.045	0.058	0.042	0.07	0.0614	0.0502	0.099	0.038	0.06 J	0.033	0.91	0.078
Phenol	7.3	22	0.000074 J	0.000067 J	0.00015 J	0.000088 J	<0.0000377	20.2	16.5	5.5	7.7	19	3.6	3.1
Pyrene	0.73	2.2	0.0034	0.005	0.0017	0.0034	0.0059	0.0076	<0.021	0.002	0.0021 J	0.0012 J	0.12	0.0019
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-17 02/01/2012	MW-17 07/12/2012	MW-17 02/14/2013	MW-17 02/14/2013 Duplicate	MW-17 04/01/2013	MW-17 04/01/2013 Duplicate	MW-17 07/30/2013	MW-17 01/13/2014	MW-17 07/17/2014	MW-17 01/30/2018	MW-17 03/18/2018	MW-17 05/16/2018
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.005	<0.005			<0.007	<0.007	<0.014	<0.007	<0.0028	<0.0002	<0.0002	<0.002
Benzene	0.005	0.005	0.24	0.46			0.435	0.41	0.174	0.324	0.576	0.47	0.55	0.61
Chlorobenzene	0.1	0.1	<0.005	<0.005			<0.006	<0.006	<0.012	<0.006	<0.0024	<0.0003	<0.0003	<0.003
Ethylbenzene	0.7	0.7	0.23	0.21			0.217	0.22	0.279	0.251	0.209	0.26	0.19	0.23
Methylene chloride	0.005	0.005	<0.0065	<0.01			<0.0075	<0.0075	0.115	<0.0075	0.0187 J	<0.001	<0.001	<0.01
Toluene	1	1	0.74	0.81			0.878	0.861	0.68	0.931	0.93	0.97	0.83	0.82
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.63	0.69			0.702	0.672	0.698	0.724	0.641	0.81	0.47	0.77
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0005	<0.00005	<0.038	<0.038	<0.0423	<0.0423	<0.00534	<0.0529	<0.00519	<0.00021	<0.00021	<0.00021
2,4-Dimethylphenol	0.49	1.5	3	4	7.93	7.03	11.3	10.4	3.19	6.75	13.9	4.3	7.7	4.2
2,4-Dinitrotoluene	0.0013	0.003	<0.0005	<0.00005	<0.032	<0.032	<0.05	<0.05	<0.00631	<0.0625	<0.00613	<0.00058	<0.00058	<0.00059
2,6-Dinitrotoluene	0.0013	0.003	<0.0006	<0.00006	<0.029	<0.029	<0.0308	<0.0308	<0.00388	<0.0385	<0.00377	<0.00042	<0.00042	<0.00042
2-Chloronaphthalene	2	5.8	<0.0005	<0.00005	<0.019	<0.019	<0.0308	<0.0308	<0.00388	<0.0385	<0.00377	<0.00021	<0.00021	<0.00021
2-Methylnaphthalene	0.098	0.29	0.29	0.51	0.443	0.421	1.04	1.24	1.04	0.857	0.636	0.3	0.38	0.34
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0008	<0.00008	<0.016	<0.016	<0.0319	<0.0319	<0.0403	<0.0399	<0.0392	<0.0002	<0.0002	<0.0002
4-Nitrophenol	0.049	0.15	<0.0005	<0.00005	<0.033	<0.033	<0.215	<0.215	<0.0272	<0.269	<0.0264	<0.00047	<0.00047	0.0076 J
Acenaphthene	1.5	4.4	0.13	0.14	0.159	0.155	0.36	0.369	0.353	0.315	0.195	0.084	0.13	0.11
Acenaphthylene	1.5	4.4	0.0056	0.005	<0.016	<0.016	<0.0231	<0.0231	0.0147 J	<0.0288	<0.00283	0.0029	<0.00015	0.0041
Anthracene	7.3	22	0.0096	0.014	<0.044	<0.044	0.0192 J	0.0822 J	0.0233 J	0.0278 J	0.0202 J	0.0065	0.011	0.0094
Benzo(a)anthracene	0.0091	0.02	<0.0005	0.00018 J	<0.025	<0.025	<0.0308	<0.0308	<0.00388	<0.0385	<0.00377	<0.0005	<0.0005	<0.00051
Benzo(a)pyrene	0.0002	0.0002	<0.0005	<0.00005	<0.013	<0.013	<0.0308	<0.0308	<0.00388	<0.0385	<0.00377	<0.0002	<0.0002	<0.0002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0005	<0.00005	<0.019	<0.019	<0.05	<0.05	<0.00631	<0.0625	<0.00613	<0.0003	<0.0003	<0.0003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.001	0.00027	<0.059	<0.059	<0.142	<0.142	<0.018	<0.178	<0.0175	<0.00037	<0.00037	<0.00037
Chrysene	0.91	2	<0.0005	0.00011 J	<0.024	<0.024	<0.0308	<0.0308	<0.00388	<0.0385	<0.00377	<0.00021	<0.00021	<0.00021
Dibenzofuran	0.098	0.29	0.093	0.13	0.133 J	0.128 J	0.275	0.234	0.253	0.211 J	0.148	0.071	0.092	0.082
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0005	<0.00005	<0.187	<0.187	<0.0423	<0.0423	<0.00534	<0.0529	<0.00519	<0.0002	<0.0002	<0.0002
Fluoranthene	0.98	2.9	0.0026	0.0034	<0.031	<0.031	<0.0269	<0.0269	0.00667 J	<0.0337	0.00429 J	0.0022	0.003	0.0022
Fluorene	0.98	2.9	0.059	0.082	0.0672 J	0.0489 J	0.146 J	0.0718 J	0.165	0.16 J	0.0943	0.04	0.057	0.051
Naphthalene	0.49	1.5	10	14	15.3	13.8	25.2	25.6	25.8	21.3	15.4	7.4	12	8.1
Nitrobenzene	0.049	0.15	<0.0005	<0.00005	<0.02	<0.02	<0.0423	<0.0423	<0.00534	<0.0529	<0.00519	<0.00024	<0.00024	<0.00024
N-Nitrosodiphenylamine	0.19	0.42	<0.0005	<0.00005	<0.033	<0.033	<0.0385	<0.0385	<0.00485	<0.0481	<0.00472	<0.00025	<0.00025	<0.00025
Pentachlorophenol	0.001	0.001	<0.0005	<0.00005	<0.096	<0.096	<0.235	<0.235	<0.0296	<0.293	<0.0288	<0.00079	<0.00079	<0.0008
Phenanthrene	0.73	2.2	0.042	0.063	0.0564 J	0.0534 J	0.124 J	0.0231 J	0.123	0.0993 J	0.0725	0.034	0.044	0.044
Phenol	7.3	22	3.7	6.1	20.7	16.2	22.2	25.1	1.54	6.46	18.1	7.1	18	6.5
Pyrene	0.73	2.2	0.0018 J	0.0018	<0.033	<0.033	<0.0423	<0.0423	<0.00534	<0.0529	<0.00519	0.0011	0.0015	0.0014
<b>Metals</b>														
Arsenic	0.01	0.01										0.0444	0.0419	0.0415

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-17 01/09/2019	MW-17 07/10/2019	MW-17 01/14/2020	MW-17C 01/30/2008	MW-17C 01/30/2008 Duplicate	MW-17C 07/15/2008	MW-17C 02/04/2009	MW-17C 01/18/2010	MW-17C 06/23/2010	MW-17C 01/17/2011	MW-17C 07/13/2011	MW-17C 02/01/2012
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.002	<0.0002	<0.0002	<0.00052	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.005
Benzene	0.005	0.005	<b>0.35</b>	<b>0.54</b>	<0.0002	<b>0.0565</b>	<b>0.0517</b>	<b>0.0426</b>	<b>0.03</b>	<b>0.0083</b>	<b>0.024</b>	<b>0.023</b>	<b>0.01</b>	<b>0.016 J</b>
Chlorobenzene	0.1	0.1	<0.003	<0.0003	<0.0003	<0.00047	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.005
Ethylbenzene	0.7	0.7	0.21	0.23	<0.0003	0.292	0.312	0.226	0.17	0.053	0.2	0.21	0.021	0.19
Methylene chloride	0.005	0.005	<b>&lt;0.01</b>	<0.001	<0.001	<0.00054	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<b>&lt;0.0065</b>
Toluene	1	1	0.68	0.83	<0.0002	0.0137	0.014	0.0102	0.008	0.0042 J	0.0071	0.0081	0.0046 J	0.0067 J
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.66	0.78	<0.0003	0.485	0.485	0.353	0.25	0.046	0.33	0.42 J	0.029	0.22
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00021	<0.00021	<0.00021	<0.0008	<b>&lt;0.04</b>	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<b>1.9</b>	<b>5.9</b>	<0.00004	<0.0029	<0.14	0.00054 J	0.0028	0.044	0.0018	0.0035 J	<b>1.5</b>	<0.0005
2,4-Dinitrotoluene	0.0013	0.003	<0.00058	<0.00058	<0.000058	<b>&lt;0.0019</b>	<b>&lt;0.096</b>	<0.00021	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00042	<0.00042	<0.000042	<b>&lt;0.0019</b>	<b>&lt;0.096</b>	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.00021	<0.00021	<0.000021	<0.0038	<0.19	<0.00042	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<b>0.23</b>	<b>0.39</b>	<0.000019	<b>1.09</b>	<b>0.418</b>	0.0954	0.085	0.063	<b>0.099</b>	0.075	0.0073	0.062
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.0002	<0.00002	<0.0019	<b>&lt;0.096</b>	<0.00053	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00047	<0.00047	<0.000047	<0.0024	<b>&lt;0.12</b>	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.091	0.16	<0.000027	0.726	0.339	0.227	0.14	0.13	0.14	0.18	0.021	0.097
Acenaphthylene	1.5	4.4	0.0029	0.0042	<0.000015	<0.0029	<0.14	0.00251	0.0012	0.0013	0.0016	0.0017	0.00028	0.0011
Anthracene	7.3	22	0.0057	0.0092	<0.000014	0.178	<0.096	0.00985	0.0084	0.0057	0.0071	0.015	0.0016	0.0048
Benzo(a)anthracene	0.0091	0.02	<0.0005	<0.0005	<0.00005	<b>0.0466</b>	<b>&lt;0.096</b>	<0.00021	0.00018 J	0.00013 J	0.00016 J	0.0012	0.00017 J	0.000091 J
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<0.0002	<0.00002	<b>0.0128</b>	<b>&lt;0.096</b>	<b>&lt;0.00021</b>	<0.00008	<0.00008	<0.00008	<b>0.00027</b>	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0003	<0.0003	<0.00003	<b>&lt;0.0038</b>	<b>&lt;0.19</b>	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00037	<0.00037	<0.000037	<0.0019	<b>&lt;0.096</b>	0.00021 J	<0.0002	<0.00039	<0.0018	0.0015	<b>&lt;0.012</b>	<0.001
Chrysene	0.91	2	<0.00021	<0.00021	<0.000021	0.0428	<0.096	<0.00021	0.00017 J	0.00012 J	0.00017 J	0.001	<0.00005	0.00013 J
Dibenzofuran	0.098	0.29	0.072	<b>0.12</b>	<0.00002	<b>0.61</b>	<b>0.291</b>	<b>0.19</b>	<b>0.13</b>	<b>0.11</b>	<b>0.13</b>	<b>0.19</b>	0.021	0.096
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.0002	<0.00002	<0.0019	<0.096	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.0015	0.0039	0.00011	0.322	<0.096	0.00845	0.007	0.0044	0.005	0.019	0.0018	0.002
Fluorene	0.98	2.9	0.043	0.06	<0.00003	0.422	0.167	0.0799	0.062	0.055	0.069	0.083	0.009	0.054
Naphthalene	0.49	1.5	<b>5.5</b>	<b>8.9</b>	<0.00002	<b>9.8</b>	<b>6.35</b>	<b>5.84</b>	<b>3.4</b>	<b>2.2</b>	<b>3.4</b>	<b>4.1</b>	0.37	<b>3.3</b>
Nitrobenzene	0.049	0.15	<0.00024	<0.00024	<0.000024	<0.0038	<b>&lt;0.19</b>	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.00025	<0.00025	<0.000025	<0.0024	<0.12	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.00079	<0.00079	<0.000079	<b>&lt;0.0019</b>	<b>&lt;0.096</b>	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
Phenanthrene	0.73	2.2	0.028	0.045	<0.000021	<b>1.09</b>	0.316	0.104	0.078	0.058	0.08	0.076	0.014	0.0081
Phenol	7.3	22	2.2	<b>9.9</b>	<0.000035	<0.0019	<0.096	0.0349	0.0013	0.14	<0.00007	0.00078	0.025	0.00014 J
Pyrene	0.73	2.2	0.00081 J	0.0018	0.000066 J	0.19	<0.096	0.00445	0.0033	0.0028	0.0026	0.009	0.00098	0.0025
<b>Metals</b>														
Arsenic	0.01	0.01	<b>0.046</b>	<b>0.046</b>	0.00154 J									

- Notes:
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  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-17C 07/12/2012	MW-17C 02/14/2013	MW-17C 04/01/2013	MW-17C 07/30/2013	MW-17C 01/13/2014	MW-17C 07/17/2014	MW-17C 01/31/2018	MW-17C 03/18/2018	MW-17C 05/16/2018	MW-17C 01/10/2019	MW-17C 07/10/2019	MW-17C 01/15/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005		<0.0014	<0.0014	<0.0007	<0.00014	<0.0002	<0.0002	<0.001	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<b>0.013</b>		<b>0.0114</b>	<b>0.0162</b>	<b>0.00939</b>	<b>0.0132</b>	<b>0.014</b>	<b>0.0067</b>	<b>0.0099</b>	<b>0.012</b>	<b>0.0097</b>	<0.0002
Chlorobenzene	0.1	0.1	<0.0005		<0.0012	<0.0012	<0.0006	<0.00012	<0.0003	<0.0003	<0.0015	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.17		0.161	0.225	0.123	0.0374	0.042	0.16	0.12	0.027	0.027	<0.0003
Methylene chloride	0.005	0.005	<0.001		0.00368 J	<b>0.00786 J</b>	<0.00075	<0.00015	<0.001	<0.001	<0.005	<0.001	<0.001	<0.001
Toluene	1	1	0.0057		0.0049 J	0.00743 J	0.00471 J	0.0073	0.0097	0.0038	0.0069	0.0087	0.0073	<0.0002
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.21		0.217	0.27	0.141	0.0482	0.071	0.1	0.094	0.05	0.039	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<b>&lt;0.038</b>	<0.000106	<b>&lt;0.0107</b>	<b>&lt;0.0106</b>	<b>&lt;0.00519</b>	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.039	0.0418 J	<0.000298	<0.0301	<0.0298	<b>7.09</b>	<b>0.59</b>	0.03	0.08	0.47	0.21	<0.0004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<b>&lt;0.032</b>	<0.000125	<b>&lt;0.0126</b>	<b>&lt;0.0125</b>	<b>&lt;0.00613</b>	<0.00058	<0.00058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<b>&lt;0.029</b>	<0.000769	<b>&lt;0.00777</b>	<b>&lt;0.00769</b>	<b>&lt;0.00377</b>	<0.00042	<0.00042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.019	<0.000769	<0.00777	<0.00769	<0.00377	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<b>0.1</b>	<b>0.173</b>	<b>0.176</b>	<b>0.151</b>	<b>0.144</b>	0.0203 J	0.094	<b>0.12</b>	<b>1.1</b>	0.025	0.026	0.000073 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<b>&lt;0.016</b>	<0.000798	<b>&lt;0.0806</b>	<b>&lt;0.0798</b>	<b>&lt;0.0392</b>	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.033	0.00502	<b>&lt;0.0544</b>	<b>&lt;0.0538</b>	<0.0264	0.0035 J	<0.00047	0.0017	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.14	0.194	0.216	0.239	0.218	0.0299	0.043	0.15	1.4	0.031	0.035	0.00023
Acenaphthylene	1.5	4.4	0.0018	<0.016	<0.000577	<0.00583	<0.00577	<0.00283	0.00076 J	0.002	0.0017	0.00047	0.00037	<0.00015
Anthracene	7.3	22	0.008	<0.044	0.011	0.0144 J	0.0156 J	<0.00236	0.0017	0.01	0.0088	0.0012	0.0013	<0.00014
Benzo(a)anthracene	0.0091	0.02	0.00022	<b>&lt;0.025</b>	0.00016 J	<0.00777	<0.00769	<0.00377	<0.0005	<0.0005	0.00011	0.000062 J	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<b>&lt;0.013</b>	<0.000769	<b>&lt;0.00777</b>	<b>&lt;0.00769</b>	<b>&lt;0.00377</b>	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<b>&lt;0.019</b>	<0.000125	<b>&lt;0.0126</b>	<b>&lt;0.0125</b>	<b>&lt;0.00613</b>	<0.0003	<0.0003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0048	<b>&lt;0.059</b>	0.00148	<b>&lt;0.0359</b>	<b>0.0491 J</b>	<b>&lt;0.0175</b>	0.005	0.0053	0.0018	0.00062	0.00056	<0.00037
Chrysene	0.91	2	0.00016 J	<0.024	0.000167 J	<0.00777	<0.00769	<0.00377	<0.00021	<0.00021	0.00011	0.000059 J	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<b>0.14</b>	<b>0.16</b>	<b>0.185</b>	<b>0.199</b>	<b>0.184</b>	0.0255	0.039	<b>0.13</b>	<b>0.099</b>	0.027	0.027	0.000065 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.187	<0.000106	<0.0107	<0.0106	<0.00519	<0.0002	<0.0002	0.000058 J	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0048	<0.031	0.00784	0.00795 J	0.00707 J	<0.00033	0.00073 J	0.0056	0.0044	0.00078	0.0009	0.00034
Fluorene	0.98	2.9	0.066	0.0785 J	0.0989	0.103	0.0907	0.0118 J	0.019	0.066	0.06	0.012	0.014	0.000061 J
Naphthalene	0.49	1.5	<b>4</b>	<b>0.988</b>	<b>5.9</b>	<b>4.4</b>	<b>6.24</b>	<b>0.772</b>	<b>1.8</b>	<b>4.8</b>	<b>53</b>	<b>1.1</b>	<b>0.97</b>	0.00016
Nitrobenzene	0.049	0.15	<0.00005	<0.02	<0.000106	<0.0107	<0.0106	<0.00519	<0.00024	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.033	<0.0000962	<0.00971	<0.00962	<0.00472	<0.00025	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<b>&lt;0.096</b>	<0.000587	<b>&lt;0.0592</b>	<b>&lt;0.0587</b>	<b>&lt;0.0288</b>	<0.00079	<0.00079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.076	0.106 J	0.12	0.12	0.11	0.0122 J	0.014	0.071	0.07	0.01	0.01	<0.00021
Phenol	7.3	22	<0.0002	0.107 J	<0.000385	<0.00388	<0.00385	<b>8.33</b>	0.0025	0.0036	0.00022	0.033	0.00073	<0.00035
Pyrene	0.73	2.2	0.0028	<0.033	0.00356	<0.0107	<0.0106	<0.00519	0.00031 J	0.0031	0.0025	0.00045	0.00046	0.00021
<b>Metals</b>														
Arsenic	0.01	0.01							0.00112 J	0.00688	0.00479	0.0013 J	0.000854 J	0.00184 J

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-18A 01/30/2008	MW-18A 07/15/2008	MW-18A 02/05/2009	MW-18A 01/18/2010	MW-18A 06/24/2010	MW-18A 01/17/2011	MW-18A 07/13/2011	MW-18A 02/01/2012	MW-18A 07/11/2012	MW-18A 01/31/2013	MW-18A 07/29/2013	MW-18A 01/13/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.00052	<0.0025	<0.005	<0.0025	<0.0025	<0.005	<0.005	<0.005	<0.007	0.00405 J	<0.014
Benzene	0.005	0.005	0.503	0.321	0.48	0.51	0.47	0.55	0.59	0.28	0.68	0.636	0.491	0.239
Chlorobenzene	0.1	0.1	0.0167	0.0074	0.017 J	<0.005	<0.0025	<0.0025	<0.005	<0.005	<0.005	<0.006	<0.0012	<0.012
Ethylbenzene	0.7	0.7	0.555	0.153	0.52	0.48	0.54	0.55	0.39	0.55	0.3	0.316	0.398	0.637
Methylene chloride	0.005	0.005	<0.00054	<0.00054	<0.0025	<0.005	<0.0025	<0.0025	<0.0065	<0.0065	<0.01	<0.0075	0.00976 J	<0.015
Toluene	1	1	0.374	0.0718	0.23	0.32	0.45	0.35	0.23	0.21	0.21	0.154	0.239	0.0731 J
Vinyl chloride	0.002	0.002				0.059		0.07	0.028	0.047	<0.005	0.0181 J	0.029	<0.011
Xylenes (total)	10	10	1.13	0.292	0.98	1.2	1.2	1	0.73	1.1	0.51	0.519	0.991	1.27
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.008	<0.0008	<0.0001	<0.0001	<0.0005	<0.0001	<0.00005	<0.0005	0.00005 J	<0.00108	<0.0214	<0.0106
2,4-Dimethylphenol	0.49	1.5	12.5	2.17	1.9	4.5	7.9	9.6 J	11	5.8	9.4 J	11.8	6.29	2.95
2,4-Dinitrotoluene	0.0013	0.003	<0.02	<0.002	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.0005	0.00005 J	<0.00127	<0.0252	<0.0125
2,6-Dinitrotoluene	0.0013	0.003	<0.02	<0.002	<0.00007	<0.00007	<0.00035	<0.00007	<0.00006	<0.0006	0.00006 J	<0.000784	<0.0155	<0.00769
2-Chloronaphthalene	2	5.8	<0.041	<0.004	<0.00012	<0.0001	<0.0005	<0.0001	<0.00005	<0.0005	0.00005 J	<0.000784	<0.0155	<0.00769
2-Methylnaphthalene	0.098	0.29	0.548	0.594	0.42	0.36	0.4	0.44	0.7	0.23	0.64 J	0.745	0.819	0.996
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.02	<0.005	<0.00008	<0.00008	<0.0004	<0.00008	<0.00008	<0.0008	0.00008 J	<0.00814	<0.161	<0.0798
4-Nitrophenol	0.049	0.15	<0.026	<0.0025	<0.00007	<0.00007	<0.00035	<0.00007	<0.00005	<0.0005	0.00005 J	<0.00549	<0.109	<0.0538
Acenaphthene	1.5	4.4	0.343	0.466	0.19	0.23	0.25	0.24	0.36	0.15	0.3 J	0.464	0.493	0.553
Acenaphthylene	1.5	4.4	<0.031	0.0131	0.0092	0.0062	0.0095	0.0072	0.015	0.007	0.0067 J	0.0151	<0.0117	<0.00577
Anthracene	7.3	22	<0.02	0.0114	0.009	0.0069	0.0075	0.0073	0.013	0.0046	0.009 J	0.0204	0.0204 J	0.0226 J
Benzo(a)anthracene	0.0091	0.02	<0.02	<0.002	<0.00007	<0.00007	<0.00035	<0.00007	<0.00005	<0.0005	0.00005 J	<0.000784	<0.0155	<0.00769
Benzo(a)pyrene	0.0002	0.0002	<0.02	<0.002	<0.00008	<0.00008	<0.0004	<0.00008	<0.00005	<0.0005	0.00005 J	<0.000784	<0.0155	<0.00769
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.041	<0.004	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.0005	0.00005 J	<0.00127	<0.0252	<0.0125
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.02	<0.002	0.00033	<0.0002	<0.001	<0.0002	<0.0001	<0.0011	0.0001 J	<0.00363	<0.0718	<0.0356
Chrysene	0.91	2	<0.02	<0.002	<0.00007	<0.00007	<0.00035	<0.00007	<0.00005	<0.0005	0.00005 J	<0.000784	<0.0155	<0.00769
Dibenzofuran	0.098	0.29	0.233	0.29	0.12	0.15	0.16	0.15	0.23	0.075	0.21 J	0.188	0.279	0.326
Di-n-butylphthalate (DBP)	2.4	7.3	<0.02	<0.002	<0.00007	<0.00007	<0.00035	<0.00007	<0.00005	<0.0005	0.00005 J	<0.00108	<0.0214	<0.0106
Fluoranthene	0.98	2.9	<0.02	0.0021	0.0026	0.0013	0.0013	0.0014	0.0018	0.0013 J	0.0016 J	<0.000686	<0.0136	<0.00673
Fluorene	0.98	2.9	0.155	0.182	0.089	0.096	0.11	0.094	0.18	0.057	0.14 J	0.136	0.214	0.268
Naphthalene	0.49	1.5	7.93	7.43	3.3	4.3	6.1	5.9	7.3	3.6	7.8 J	9.29	11.8	11.4
Nitrobenzene	0.049	0.15	<0.041	<0.004	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.0005	0.00005 J	<0.00108	<0.0214	<0.0106
N-Nitrosodiphenylamine	0.19	0.42	<0.026	<0.0025	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.0005	0.00005 J	<0.00098	<0.0194	<0.00962
Pentachlorophenol	0.001	0.001	<0.02	<0.002	<0.00008	<0.00008	<0.0004	<0.00008	<0.00005	<0.0005	0.00005 J	<0.00598	<0.118	<0.0587
Phenanthrene	0.73	2.2	0.118	0.12	0.078	0.067	0.082	0.063	0.098	0.042	0.083 J	0.101	0.144	0.19
Phenol	7.3	22	0.364	<0.002	0.005	0.043	0.0054	0.02	0.061	0.011	0.12 J	<0.000392	<0.00777	<0.00385
Pyrene	0.73	2.2	<0.02	<0.002	0.0013	0.00075	0.00063 J	0.00085	0.0011	0.00077 J	0.00081 J	<0.00108	<0.0214	<0.0106
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

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- Concentrations > C/I AL and non-detects are highlighted dark gray
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- RAL = Residential Assessment Level, C/I = Commercial/Industrial
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	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-18A 07/16/2014	MW-18A 01/25/2018	MW-18A 03/19/2018	MW-18A 05/16/2018	MW-18A 01/10/2019	MW-18A 07/10/2019	MW-18A 01/08/2020	MW-18C 01/30/2008	MW-18C 07/15/2008	MW-18C 02/05/2009	MW-18C 01/19/2010	MW-18C 06/24/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	0.00482 J	<0.0002	<0.0002	<0.0002	<0.005	<0.0002	<0.005	<0.00052	<0.00052	<0.005	<0.005	<0.0025
Benzene	0.005	0.005	<b>0.483</b>	<b>0.19</b>	<b>0.22</b>	<b>0.17</b>	<b>1.2</b>	<b>0.16</b>	<b>0.41</b>	<b>1.34</b>	<b>0.964</b>	<b>1.4</b>	<b>1.5</b>	<b>1</b>
Chlorobenzene	0.1	0.1	<0.0024	0.0062	0.006	0.0066	<0.0075	0.0017	<0.0075	<0.00047	<0.00047	<0.005	<0.005	<0.0025
Ethylbenzene	0.7	0.7	0.692	0.36	0.36	0.43	0.34	0.28	0.32	0.304	0.178	0.26	0.21	0.13
Methylene chloride	0.005	0.005	<b>0.0138 J</b>	<0.001	<0.001	<0.001	<b>&lt;0.025</b>	<0.001	<b>&lt;0.025</b>	<0.00054	<0.00054	<0.005	<0.005	<0.0025
Toluene	1	1	0.416	0.074	0.068	0.042	0.92	0.057	0.11	<b>1.2</b>	0.691	1	0.96	0.72 J
Vinyl chloride	0.002	0.002	<b>0.02 J</b>	<b>0.0023</b>	0.002	<b>0.0021</b>	<b>&lt;0.005</b>	0.0013	<b>&lt;0.005</b>					
Xylenes (total)	10	10	1.3	0.97	0.79	0.91	1	0.47	0.5	1.1	0.624	1.1	1	1
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<b>&lt;0.00519</b>	<0.00021	<0.00021	<0.00021	<0.00021	<0.000021	<0.000021	<b>&lt;0.04</b>	<0.0008	<0.0001	<0.0001	<0.0005
2,4-Dimethylphenol	0.49	1.5	<b>8.01</b>	0.023	0.12	<b>0.5</b>	0.0054	<b>0.72</b>	<b>1.8</b>	<0.15	<0.0031	0.084	0.0081	0.0078 J
2,4-Dinitrotoluene	0.0013	0.003	<b>&lt;0.00613</b>	<0.00058	<0.00058	<0.00058	<0.00058	<0.000058	<0.000058	<b>&lt;0.098</b>	<b>0.00636</b>	<0.00009	<0.00009	<0.00045
2,6-Dinitrotoluene	0.0013	0.003	<b>&lt;0.00377</b>	<0.00042	<0.00042	<0.00042	<0.00042	<0.000042	<0.000042	<b>&lt;0.098</b>	<b>&lt;0.002</b>	<0.00007	<0.00007	<0.00035
2-Chloronaphthalene	2	5.8	<0.00377	<0.00021	<0.00021	<0.00021	<0.00021	<0.000021	<0.000021	<0.2	<0.0041	<0.00012	<0.0001	<0.0005
2-Methylnaphthalene	0.098	0.29	<b>0.589</b>	<b>0.33</b>	<b>0.34</b>	<b>0.22</b>	<b>0.1</b>	<b>0.47</b>	<b>0.36</b>	<b>0.894</b>	<b>0.674</b>	<b>0.95</b>	<b>0.46</b>	<b>0.2</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<b>&lt;0.0392</b>	<0.0002	<0.0002	<0.0002	<0.0002	<0.00002	<0.00002	<b>&lt;0.098</b>	<b>&lt;0.0051</b>	<0.00008	<0.00008	<0.0004
4-Nitrophenol	0.049	0.15	<0.0264	<0.00047	<0.00047	<0.00047	<0.00047	<0.000047	0.0042	<b>&lt;0.12</b>	<0.0026	<0.00007	<0.00007	<0.00035
Acenaphthene	1.5	4.4	0.352	0.25	0.23	0.24	0.048	0.3	0.33	0.293	0.251	0.18	0.17	0.082
Acenaphthylene	1.5	4.4	0.0155 J	0.0074	0.0092	0.009	0.016	0.0056	0.0081	<0.15	0.00649	0.0036	0.0023	0.0015
Anthracene	7.3	22	0.0192 J	0.0061	0.0079	0.0077	0.0064	0.0063	0.0082	<0.098	0.0321	0.017	0.014	0.0076 J
Benzo(a)anthracene	0.0091	0.02	<0.00377	<0.0005	<0.0005	<0.0005	<0.0005	<0.00005	<0.00005	<b>&lt;0.098</b>	0.0025	0.00039	<0.00007	<0.00035
Benzo(a)pyrene	0.0002	0.0002	<b>&lt;0.00377</b>	<0.0002	<0.0002	<0.0002	<0.0002	<0.00002	<0.00002	<b>&lt;0.098</b>	<b>&lt;0.002</b>	0.00013 J	<0.00008	<b>&lt;0.0004</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019	<b>&lt;0.00613</b>	<0.0003	<0.0003	<0.0003	<0.0003	<0.00003	<0.00003	<b>&lt;0.2</b>	<b>&lt;0.0041</b>	<0.00009	<0.00009	<0.00045
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<b>&lt;0.0175</b>	<0.00037	<0.00037	<0.00037	<0.00037	<0.000037	<0.000043	<b>&lt;0.098</b>	<0.002	0.00023	<0.0002	<0.001
Chrysene	0.91	2	<0.00377	<0.00021	<0.00021	<0.00021	<0.00021	<0.000021	<0.000021	<0.098	0.0021	0.00033	<0.00007	<0.00035
Dibenzofuran	0.098	0.29	<b>0.204</b>	<b>0.16</b>	<b>0.15</b>	<b>0.14</b>	0.047	<b>0.19</b>	<b>0.2</b>	<b>0.263</b>	<b>0.23</b>	<b>0.16</b>	0.091	0.077
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00519	<0.0002	<0.0002	<0.0002	<0.0002	<0.00002	<0.00002	<0.098	<0.002	<0.00007	<0.00007	<0.00035
Fluoranthene	0.98	2.9	<0.0033	0.0021	0.0023	0.0017	0.002	0.0022	0.0023	<0.098	0.0169	0.0047	0.0035	0.0023
Fluorene	0.98	2.9	0.163	0.087	0.12	0.098	0.021	0.091	0.16	0.137	0.153	0.081	0.052	0.034
Naphthalene	0.49	1.5	<b>5.27</b>	<b>4.4</b>	<b>4.9</b>	<b>6.1</b>	<b>3.2</b>	<b>5.6</b>	<b>4.6</b>	<b>16.6</b>	<b>16.7</b>	<b>21</b>	<b>12</b>	<b>6.2</b>
Nitrobenzene	0.049	0.15	<0.00519	<0.00024	<0.00024	<0.00024	<0.00024	<0.000024	<0.000024	<b>&lt;0.2</b>	<0.0041	<0.00009	<0.00009	<0.00045
N-Nitrosodiphenylamine	0.19	0.42	<0.00472	<0.00025	<0.00025	<0.00025	<0.00025	<0.000025	<0.000025	<0.12	<0.0026	<0.00009	<0.00009	<0.00045
Pentachlorophenol	0.001	0.001	<b>&lt;0.0288</b>	<0.00079	<0.00079	<0.00079	<b>0.014</b>	<0.000079	<0.000079	<b>&lt;0.098</b>	<b>0.134</b>	<b>0.026</b>	<b>0.041</b>	<b>0.02</b>
Phenanthrene	0.73	2.2	0.114	0.066	0.08	0.082	0.024	0.075	0.1	0.213	0.177	0.076	0.052	0.032
Phenol	7.3	22	<0.00189	0.0019 J	<0.00035	<0.00035	0.0032	0.00048	0.0014	<0.098	0.0944	0.031	0.059	0.026
Pyrene	0.73	2.2	<0.00519	0.0011	0.0012	0.0008 J	0.0012	0.0011	0.0013	<0.098	0.01	0.0025	0.002	0.0012
<b>Metals</b>														
Arsenic	0.01	0.01		0.0043	<b>0.0239</b>	<b>0.0291</b>	0.0031	<b>0.0248</b>	<b>0.0236</b>					

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-18C 01/17/2011	MW-18C 07/13/2011	MW-18C 02/01/2012	MW-18C 07/11/2012	MW-18C 01/31/2013	MW-18C 07/29/2013	MW-18C 01/13/2014	MW-18C 07/16/2014	MW-18C 01/25/2018	MW-18C 03/19/2018	MW-18C 05/16/2018	MW-18C 01/10/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0025	<0.01	<0.005	<0.005	<0.007	<0.014	<0.007	<0.0028	<0.0002	<0.0002	<0.002	<0.002
Benzene	0.005	0.005	1.3	1.2	1.3	1.2	1.51	1.23	1.51	1.45	1.4	1.3	1.4	0.3
Chlorobenzene	0.1	0.1	<0.0025	<0.01	<0.005	<0.005	<0.006	<0.012	<0.006	<0.0024	<0.0003	0.00052 J	<0.003	<0.003
Ethylbenzene	0.7	0.7	0.18	0.16	0.19	0.15	0.203	0.22	0.245	0.309	0.35	0.29	0.32	0.41
Methylene chloride	0.005	0.005	<0.0025	<0.013	<0.0065	<0.01	<0.0075	0.0688 J	<0.0075	0.0161 J	<0.001	<0.001	<0.01	<0.01
Toluene	1	1	0.83	0.8	0.83	0.72	0.962	0.899	1.07	0.986	1.1	0.96	0.9	0.03
Vinyl chloride	0.002	0.002	<0.0025	<0.01	<0.005	<0.005	<0.0055	<0.011	<0.0055	<0.0022	0.0018	0.0026	<0.002	<0.002
Xylenes (total)	10	10	1	0.9	0.82	0.84	1.01	0.881	1.02	1.36	1	0.93	1.1	0.69
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.0005	<0.00005	<0.0545	<0.00534	<0.0106	<0.00519	<0.00021	<0.00021	<0.00021	<0.00021
2,4-Dimethylphenol	0.49	1.5	0.012 J	0.0031	0.01	<0.0021	<0.153	<0.015	<0.0298	0.0325	<0.0004	<0.0004	0.082	0.29
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.0005	<0.00005	<0.0644	<0.00631	<0.0125	<0.00613	<0.00058	<0.00058	<0.00058	<0.00059
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.0006	<0.00006	<0.0396	<0.00388	<0.00769	<0.00377	<0.00042	<0.00042	<0.00042	<0.00042
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.0005	<0.00005	<0.0396	<0.00388	<0.00769	<0.00377	<0.00021	<0.00021	<0.00021	<0.00021
2-Methylnaphthalene	0.098	0.29	0.31	0.34	0.16	0.46	0.977	0.871	1.06	0.778	0.41	0.44	0.3	0.33
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.0008	<0.00008	<0.411	<0.0403	<0.0798	<0.0392	<0.0002	<0.0002	<0.0002	<0.0002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.0005	<0.00005	<0.277	<0.0272	<0.0538	<0.0264	<0.00047	<0.00047	<0.00047	<0.00047
Acenaphthene	1.5	4.4	0.14	0.12	0.062	0.13	0.32	0.265	0.317	0.246	0.15	0.16	0.15	0.21
Acenaphthylene	1.5	4.4	0.0019	0.0023	0.0018 J	0.0019	<0.0297	<0.00291	<0.00577	<0.00283	0.0025	0.0035	0.003	0.0075
Anthracene	7.3	22	0.015	0.013	0.012	0.008	0.0401 J	0.0284	0.0414 J	0.028	0.017	0.02	0.019	0.007
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00005	<0.0005	0.00014 J	<0.0396	<0.00388	<0.00769	<0.00377	0.0013	<0.0005	<0.0005	<0.00051
Benzo(a)pyrene	0.0002	0.0002	0.00035	0.00015 J	<0.0005	<0.00005	<0.0396	<0.00388	<0.00769	<0.00377	<0.0002	<0.0002	<0.0002	<0.0002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.0005	<0.00005	<0.0644	<0.00631	<0.0125	<0.00613	<0.0003	<0.0003	<0.0003	<0.0003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0002	<0.0001	<0.001	<0.0001	<0.183	<0.018	<0.0356	<0.0175	<0.00037	<0.00037	<0.00037	<0.00037
Chrysene	0.91	2	<0.00007	<0.00005	<0.0005	0.0001 J	<0.0396	<0.00388	<0.00769	<0.00377	0.0008 J	<0.00021	0.00038 J	<0.00021
Dibenzofuran	0.098	0.29	0.13	0.11	0.06	0.14	0.288	0.225	0.276	0.207	0.14	0.15	0.13	0.13
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00005	<0.0005	<0.00005	<0.0545	<0.00534	<0.0106	<0.00519	<0.0002	<0.0002	<0.0002	<0.0002
Fluoranthene	0.98	2.9	0.0059	0.0042	0.0018 J	0.0023	<0.0347	0.00865 J	0.0191 J	0.00957 J	0.0096	0.0071	0.0058	0.0023
Fluorene	0.98	2.9	0.051	0.052	0.028	0.055	0.132 J	0.114	<0.00673	0.116	0.056	0.073	0.062	0.095
Naphthalene	0.49	1.5	13	12	9.7	13	20.2 J	20.9	20.3	14.7	14	12	21	4.4
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.0005	<0.00005	<0.0545	<0.00534	<0.0106	<0.00519	<0.00024	<0.00024	<0.00024	<0.00024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.0005	<0.00005	<0.0495	<0.00485	<0.00962	<0.00472	<0.00025	<0.00025	<0.00025	<0.00025
Pentachlorophenol	0.001	0.001	0.064	0.076 J	0.085	0.075	<0.302	<0.0296	0.188	0.164	0.024	0.041	0.037	<0.0008
Phenanthrene	0.73	2.2	0.055	0.052	0.027	0.055	0.155 J	0.127	0.177	0.122	0.072	0.086	0.078	0.084
Phenol	7.3	22	0.043	0.048	0.027	0.075	0.0601 J	0.0205 J	0.0184 J	0.0285	0.0042	0.0067	<0.00035	<0.00035
Pyrene	0.73	2.2	0.0028	0.0017	0.001 J	0.0011	<0.0545	<0.00534	<0.0106	0.00571 J	0.0055	0.0049	0.0036	0.0012
<b>Metals</b>														
Arsenic	0.01	0.01									0.00467	0.00327	0.00342	0.0257

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-18C 07/10/2019	MW-18C 01/08/2020	MW-19C 01/31/2008	MW-19C 07/15/2008	MW-19C 02/04/2009	MW-19C 01/18/2010	MW-19C 06/23/2010	MW-19C 01/18/2011	MW-19C 01/18/2011 Duplicate	MW-19C 07/14/2011	MW-19C 02/08/2012	MW-19C 07/12/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.002	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Benzene	0.005	0.005	<b>1.2</b>	<b>1.1</b>	0.00398 J	<0.00025	<0.0005	<b>0.0056</b>	<0.0005	<0.0005	<0.0005	<0.001	0.005	<0.0005
Chlorobenzene	0.1	0.1	0.00052 J	<0.003	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Ethylbenzene	0.7	0.7	0.32	0.31	<0.00025	<0.00025	<0.0005	0.0018 J	<0.0005	<0.0005	<0.0005	<0.0011	0.0031 J	<0.0005
Methylene chloride	0.005	0.005	<0.001	<b>&lt;0.01</b>	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001
Toluene	1	1	1	0.97	0.00596	<0.00041	<0.0005	0.0076	<0.0005	<0.0005	<0.0005	<0.001	0.0085	<0.0005
Vinyl chloride	0.002	0.002	<0.0002	<0.002								<0.001	<0.001	<0.0005
Xylenes (total)	10	10	1	0.92	<0.00127	<0.00127	<0.001	0.0043 J	<0.001	<0.001	<0.001	<0.0031	0.0063 J	<0.0015
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.00008	0.00023 J	<0.0001	<0.0001	0.00024	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	0.0051	0.023	<0.00029	<0.00028	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	0.00016 J
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.00019	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.00038	<0.00038	<0.00012	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<b>0.28</b>	<b>0.31</b>	0.00132	<0.00038	0.00025	0.0017	0.000079 J	0.00015 J	0.00022	0.0012	<0.00005	<0.000053
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00019	<0.00047	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.00024	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.058	0.093	0.000562	<0.00028	0.00022	0.001	0.00012 J	0.00015 J	0.0003	0.00067	0.0012	0.00017 J
Acenaphthylene	1.5	4.4	0.005	0.0026	<0.00029	<0.00028	<0.00006	0.00014 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Anthracene	7.3	22	0.0056	0.02	<0.00019	<0.00019	<0.00007	0.0001 J	<0.00007	<0.00007	<0.00007	0.00015 J	<0.00005	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.00005	0.0025	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<b>0.00053</b>	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00038	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	<0.00019	0.00027 J	<0.0002	<0.00028	<0.00036	0.00026	0.00025	<0.00039	<0.00022	0.00014 J
Chrysene	0.91	2	<0.000021	0.0027	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	0.000072 J	0.000084 J	<0.00005	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	0.054	0.085	0.00042 J	<0.00028	0.00017 J	0.00051	<0.00008	0.00013 J	0.0002 J	<0.0006	0.00014 J	<0.00011
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00015 J
Fluoranthene	0.98	2.9	0.0019	0.013	<0.00019	0.00182	0.00015 J	0.00024	0.0021	0.0026 J	0.00062 J	0.0016	<0.00005	0.0018
Fluorene	0.98	2.9	0.024	0.043	0.00044 J	<0.00019	<0.00007	0.00032	0.00028	0.00032	0.00022	<0.00066	<0.00005	0.00033
Naphthalene	0.49	1.5	<b>9.9</b>	<b>14</b>	0.0613	0.000826	0.0077	0.09	<0.0015	0.0061 J	0.0084 J	0.014	0.00077	<0.00048
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.00038	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.00024	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<b>0.034</b>	<b>0.023</b>	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	0.00028 J	<0.00005	<0.00005
Phenanthrene	0.73	2.2	0.026	0.065	<0.00019	<0.00019	<0.00007	0.00016 J	<0.00007	<0.00007	0.00014 J	<0.00053	<0.00005	<0.00005
Phenol	7.3	22	0.005	0.0062	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00024
Pyrene	0.73	2.2	0.0014	0.0066	0.00036 J	0.00117	<0.00007	0.0002	0.0012	0.0016 J	0.0004 J	0.0014	<0.00005	0.0014
<b>Metals</b>														
Arsenic	0.01	0.01	0.00358	0.00283										

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-19C 02/01/2013	MW-19C 07/30/2013	MW-19C 01/15/2014	MW-19C 07/17/2014	MW-19C 02/09/2018	MW-19C 03/18/2018	MW-19C 05/16/2018	MW-19C 01/24/2019	MW-19C 07/10/2019	MW-19C 01/09/2020	MW-20A 01/30/2008	MW-20A 07/14/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.002	<0.0002	<0.0002	<0.0002	<0.00052	<0.005
Benzene	0.005	0.005	0.000558 J	0.00427	0.00028 J	0.0000801 J	0.0013	0.0027	0.0041 J	0.0044	0.0004 J	0.00032 J	0.0609	0.098
Chlorobenzene	0.1	0.1	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.003	<0.0003	<0.0003	<0.0003	<0.00047	<0.005
Ethylbenzene	0.7	0.7	0.000793 J	0.0114	0.000966 J	0.000783 J	0.00091 J	0.0025	<0.003	0.004	0.00096 J	0.00064 J	0.0965	0.077
Methylene chloride	0.005	0.005	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.01	<0.001	<0.001	<0.001	<0.00054	<0.0065
Toluene	1	1	0.00171	0.0155	0.00136	0.000578 J	<0.0002	0.0018	0.0042 J	0.0057	0.00086 J	<0.0002	0.00716	<0.005
Vinyl chloride	0.002	0.002	<0.00011	<0.00011	<0.00011	<0.00011	<0.0002	<0.0002	<0.002	<0.0002	<0.0002	<0.0002		
Xylenes (total)	10	10	0.00151 J	0.0197	0.00207 J	0.00179 J	0.00095 J	0.0044	<0.003	0.0037	0.0022	0.0012	0.113	0.057 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00011	<0.000107	<0.000212	<0.000104	<0.000021	0.000099 J	0.000061 J	<0.000021	0.000082 J	0.000078 J	<0.02	<0.00005
2,4-Dimethylphenol	0.49	1.5	0.00125	<0.000301	<0.000596	<0.000292	0.0028	<0.00038	<0.00004	<0.00032	<0.00056	<0.00004	0.134	0.3
2,4-Dinitrotoluene	0.0013	0.003	<0.00013	<0.000126	<0.00025	<0.000123	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.039	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00008	<0.0000777	<0.000154	<0.0000755	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.039	<0.00006
2-Chloronaphthalene	2	5.8	<0.00008	<0.0000777	<0.000154	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.078	<0.00005
2-Methylnaphthalene	0.098	0.29	0.00084	0.000114 J	0.00142	0.000845	0.000093 J	0.00037	<0.00019	<0.00019	<0.000071	<0.000019	0.204	0.42
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00083	<0.000806	<0.0016	<0.000783	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.039	<0.00008
4-Nitrophenol	0.049	0.15	<0.00056	<0.000544	<0.00108	<0.000528	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	0.177	<0.00005
Acenaphthene	1.5	4.4	<0.000608	0.00279	<0.000154	0.0007	0.0012	0.001	0.00035	0.00078	0.001	0.00099	0.113	0.2
Acenaphthylene	1.5	4.4	<0.00006	<0.0000583	<0.000115	<0.0000566	0.000069 J	<0.000015	<0.000015	<0.000015	0.00003 J	0.00005 J	<0.059	0.0017
Anthracene	7.3	22	0.000115 J	0.000269 J	<0.0000962	<0.0000472	0.000057 J	0.000065 J	<0.000014	0.000057 J	<0.000066	<0.00017	<0.039	0.0058
Benzo(a)anthracene	0.0091	0.02	<0.00008	0.000111 J	<0.000154	<0.0000755	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	0.00026	<0.039	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.0000777	<0.000154	<0.0000755	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.0001	<0.039	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00013	<0.000126	<0.00025	<0.000123	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.078	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00037	0.0012 J	<0.000712	0.000646	0.00019 J	0.0001 J	0.00016 J	0.000096 J	<0.000037	<0.000071	<0.039	<0.00033
Chrysene	0.91	2	<0.00008	<0.0000777	<0.000154	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00024	<0.039	<0.00005
Dibenzofuran	0.098	0.29	0.000367 J	0.000631	0.00116	0.000554	0.00045	0.00081	0.000035 J	<0.00002	<0.00057	0.00069	0.071	0.14
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00011	0.000164 J	<0.000212	<0.000104	<0.00002	<0.00002	0.000031 J	<0.00002	<0.00002	0.000021 J	<0.039	<0.00005
Fluoranthene	0.98	2.9	0.00257	0.000309 J	0.00223	0.00169	0.000081 J	0.00064	0.00023	<0.00001	0.00056	0.0011	<0.039	0.0007
Fluorene	0.98	2.9	0.000605	<0.000068	0.000296 J	0.000485	0.000099 J	0.00037	0.00009 J	<0.00003	<0.00031	0.00047	0.045	0.11
Naphthalene	0.49	1.5	0.0264 J	<0.00196	0.0383	0.0198	0.0022	0.013	<0.00061	<0.00036	<0.0012	<0.00011	4.75	7.7
Nitrobenzene	0.049	0.15	<0.00011	<0.000107	<0.000212	<0.000104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.078	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.0001	<0.0000971	<0.000192	<0.0000943	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.049	<0.00005
Pentachlorophenol	0.001	0.001	<0.00061	<0.000592	<0.00117	<0.000575	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.039	<0.00005
Phenanthrene	0.73	2.2	0.000166 J	0.000201 J	<0.000265	<0.0000566	<0.000021	0.000051 J	<0.000021	<0.000021	<0.00014	<0.00045	<0.039	0.039
Phenol	7.3	22	0.00023 J	0.024	0.000724 J	0.00033 J	0.00071	0.00072	<0.000035	<0.00013	<0.00015	<0.000035	<0.039	<0.00005
Pyrene	0.73	2.2	0.00207	0.000233 J	0.00191	0.00178	0.000053 J	0.00076	0.00037	<0.000019	0.00064	0.0014	<0.039	0.0004
<b>Metals</b>														
Arsenic	0.01	0.01					0.00158 J	0.00107 J	0.00294	0.00149 J	0.00166 J	0.0014 J		

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray.
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-20A 02/01/2012	MW-20A 07/16/2012	MW-20A 01/30/2013	MW-20A 01/23/2018	MW-20A 03/19/2018	MW-20A 05/15/2018	MW-20A 01/08/2019	MW-20A 07/10/2019	MW-20A 01/14/2020	MW-21C 01/29/2008	MW-21C 01/29/2008 Duplicate	MW-21C 07/15/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.005	<0.0005	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.00052
Benzene	0.005	0.005	<b>0.057</b>	<b>0.089</b>	<b>0.0746</b>	<b>0.053</b>	<b>0.05</b>	<b>0.062</b>	<b>0.024</b>	<b>0.013</b>	<b>0.02</b>	<0.00025	<0.00025	<0.00025
Chlorobenzene	0.1	0.1	<0.005	<0.0005	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.00047
Ethylbenzene	0.7	0.7	0.046	0.1	0.0619	0.05	0.027	0.045	0.024	0.0079	0.025	<0.00025	<0.00025	<0.00025
Methylene chloride	0.005	0.005	<b>&lt;0.0065</b>	<0.001	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.00054
Toluene	1	1	<0.005	0.022	0.0028	0.0038	<0.0002	0.0055	0.00077 J	0.0011	<0.0002	<0.00041	<0.00041	<0.00041
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.028 J	0.088	0.0549	0.05	0.033	0.048	0.022	0.022	0.024	<0.00127	<0.00127	<0.00127
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0005	<0.00005	<b>&lt;0.00519</b>	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.00008
2,4-Dimethylphenol	0.49	1.5	0.076	0.1	0.119	0.049	0.072	0.06	0.076	0.0035	0.0018	<0.00029	<0.00029	<0.0003
2,4-Dinitrotoluene	0.0013	0.003	<0.0005	<0.00005	<b>&lt;0.00613</b>	<0.00058	<0.00059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019	<0.00019	<0.0002
2,6-Dinitrotoluene	0.0013	0.003	<0.0006	<0.00006	<b>&lt;0.00377</b>	<0.00042	<0.00042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.00019	<0.0002
2-Chloronaphthalene	2	5.8	<0.0005	<0.00005	<0.00377	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.00038	<0.0004
2-Methylnaphthalene	0.098	0.29	0.064	<b>0.36</b>	<b>0.191</b>	<b>0.15</b>	<b>0.15</b>	<b>0.17</b>	0.069	0.072	0.075	<0.00038	<0.00038	<0.0004
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0008	<0.00008	<b>&lt;0.0392</b>	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	<0.0005
4-Nitrophenol	0.049	0.15	<0.0005	<0.00005	<0.0264	<0.00047	<0.00047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00024	<0.00025
Acenaphthene	1.5	4.4	0.15	0.19	0.12	0.14	0.18	0.15	0.1	0.14	0.11	<0.00029	<0.00029	<0.0003
Acenaphthylene	1.5	4.4	0.0015 J	0.0015	<0.00283	0.0014	<0.00015	0.0013	0.00057	0.00078	0.00092	<0.00029	<0.00029	<0.0003
Anthracene	7.3	22	0.0044	0.0042	0.00589 J	0.021	0.022	0.0081	0.0058	0.0062	0.0075	0.000563	0.000562	<0.0002
Benzo(a)anthracene	0.0091	0.02	<0.0005	<0.00005	<0.00377	<0.0005	<0.00051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00019	<0.00019	<0.0002
Benzo(a)pyrene	0.0002	0.0002	<b>&lt;0.0005</b>	<0.00005	<b>&lt;0.00377</b>	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	<0.0002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0005	<0.00005	<b>&lt;0.00613</b>	<0.0003	<0.0003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.00038	<0.0004
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.001	<0.0001	<b>&lt;0.0175</b>	<0.00037	<0.00037	<0.000037	<0.000037	<0.000037	<0.000037	<b>0.00888</b>	0.00052 J	<0.0002
Chrysene	0.91	2	<0.0005	<0.00005	<0.00377	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019	<0.00019	<0.0002
Dibenzofuran	0.098	0.29	0.073	<b>0.15</b>	0.0799	0.097	<b>0.13</b>	0.081	0.067	0.077	0.079	<0.00029	<0.00029	<0.0003
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0005	<0.00005	<0.00519	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	<0.0002
Fluoranthene	0.98	2.9	0.00082 J	0.00061	<0.0033	0.00078 J	0.00095 J	0.00074	0.00041	0.00058	0.00067	0.00047 J	0.00046 J	<0.0002
Fluorene	0.98	2.9	0.06	0.11	0.0661	0.16	0.16	0.076	0.062	0.068	0.08	<0.00019	<0.00019	<0.0002
Naphthalene	0.49	1.5	<b>0.96</b>	<b>6.1</b>	<b>43.9</b>	<b>2.7</b>	<b>2.9</b>	<b>2.8</b>	<b>1.4</b>	<b>1.9</b>	<b>2</b>	<0.00038	<0.00038	<0.0004
Nitrobenzene	0.049	0.15	<0.0005	<0.00005	<0.00519	<0.00024	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.00038	<0.0004
N-Nitrosodiphenylamine	0.19	0.42	<0.0005	<0.00005	<0.00472	<0.00025	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00024	<0.00025
Pentachlorophenol	0.001	0.001	<0.0005	<0.00005	<b>&lt;0.0288</b>	<0.00079	<0.0008	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019	<0.00019	<0.0002
Phenanthrene	0.73	2.2	0.03	0.043	0.0217 J	0.032	0.037	0.034	0.025	0.026	0.028	0.00039 J	0.00039 J	<0.0002
Phenol	7.3	22	<0.0005	<0.00005	<0.00189	<0.00035	<0.00035	<0.000035	<0.000035	0.00096	<0.000035	<0.00019	<0.00019	<0.0002
Pyrene	0.73	2.2	0.00056 J	0.0003	<0.00519	<0.00019	0.0005 J	0.00048	0.00025	0.00026	0.0003	0.00039 J	0.00039 J	<0.0002
<b>Metals</b>														
Arsenic	0.01	0.01				0.0087	0.00568	0.00895	0.00788	0.00574	0.00808			

- Notes:
1. All values in milligrams per liter (mg/L).
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  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
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  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-21C 07/15/2008	MW-21C 02/04/2009	MW-21C 02/04/2009	MW-21C 01/21/2010	MW-21C 01/21/2010	MW-21C 06/22/2010	MW-21C 06/22/2010	MW-21C 01/19/2011	MW-21C 07/27/2011	MW-21C 07/27/2011	MW-21C 02/02/2012	MW-21C 02/02/2012
			Duplicate		Duplicate		Duplicate		Duplicate			Duplicate		Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001
Benzene	0.005	0.005	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001
Chlorobenzene	0.1	0.1	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001
Ethylbenzene	0.7	0.7	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0011	<0.0011
Methylene chloride	0.005	0.005	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.0013	<0.0013
Toluene	1	1	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0031	<0.0031
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.00039	<0.00012	<0.00012	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.00039	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.000067 J	<0.00005
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00049	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.00269	<0.00009	<0.00009	0.00041	0.00035	<0.00009	<0.00009	0.00034	<0.00005	<0.00005	<0.00005	<0.00005
Acenaphthylene	1.5	4.4	<0.00029	<0.00006	<0.00006	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Anthracene	7.3	22	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	0.00013 J	<0.00005	<0.00005	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00019	<0.0002	0.00038	<0.00072	<0.00054	<0.00023	<0.00041	0.00062	0.00062 J	0.00036 J	<0.0001	<0.00022
Chrysene	0.91	2	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.00019 J	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Fluorene	0.98	2.9	0.00098	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Naphthalene	0.49	1.5	<0.00039	0.00039	0.00033	<0.0001	<0.0001	<0.00021	<0.0001	<0.0001	<0.00005	<0.00005	0.00093 J	0.00015 J
Nitrobenzene	0.049	0.15	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005
Phenanthrene	0.73	2.2	<0.00019	<0.00007	<0.00007	<0.00007	0.0001 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Phenol	7.3	22	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
Pyrene	0.73	2.2	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-21C 07/26/2012	MW-21C 07/26/2012 Duplicate	MW-21C 02/05/2013	MW-21C 02/05/2013 Duplicate	MW-21C 08/01/2013	MW-21C 08/01/2013 Duplicate	MW-21C 01/16/2014	MW-21C 01/16/2014 Duplicate	MW-21C 07/25/2014	MW-21C 07/25/2014 Duplicate	MW-21C 01/24/2018	MW-21C 01/24/2018 Duplicate
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.00014	<0.00014	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008	<0.0002	<0.0002	<0.00008	<0.00008	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012	<0.00018	<0.00018	<0.00012	<0.00012	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.000144	<0.00011	<0.00011	<0.00011	<0.00019	<0.00019	<0.00011	<0.00011	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.00022	<0.00022	<0.00015	<0.00015	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.00017	<0.00017	<0.00015	<0.00015	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.00017	<0.00017	<0.00015	<0.00015	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0015	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026	<0.00058	<0.00058	<0.00026	<0.00026	<0.0003	<0.0003
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.000104	<0.000104	0.000105 J	0.00732 J	<0.000104	<0.000104	<0.000107	<0.000107	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.00005	<0.000292	R	<0.000295	<0.00654	<0.000292	<0.000292	<0.000301	<0.000301	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.000123	<0.000123	0.000124 J	0.00867 J	<0.000123	<0.000123	<0.000126	<0.000126	<0.000059	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.0000755	<0.0000755	0.0000762 J	0.00805 J	<0.0000755	<0.0000755	<0.0000777	<0.0000777	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.0000755	<0.0000755	0.0000762 J	0.00696 J	<0.0000755	<0.0000755	<0.0000777	<0.0000777	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00005	<0.00005	0.000271 J	<0.000066	0.0000667 J	0.00686 J	<0.000157	<0.000187	<0.000068	<0.000068	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.000783	R	0.00079 J	0.00371 J	<0.000783	<0.000783	<0.000806	<0.000806	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.000528	R	0.000533 J	0.00667 J	<0.000528	<0.000528	<0.000544	<0.000544	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.00005	<0.00005	0.000237 J	<0.0000755	0.0000762 J	0.00726 J	0.000254 J	0.000413 J	<0.0000777	<0.0000777	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.00005	<0.00005	<0.0000566	<0.0000566	0.0000571 J	0.00735 J	<0.0000566	<0.0000566	<0.0000583	<0.0000583	<0.000015	<0.000015
Anthracene	7.3	22	<0.00005	<0.00005	0.0000527 J	<0.0000472	0.0000476 J	0.00775 J	0.000243 J	0.000518 J	<0.0000485	<0.0000485	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.0000755	<0.0000755	0.0000762 J	0.00807 J	0.000129 J	0.000137 J	<0.0000777	<0.0000777	<0.000051	<0.000051
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.0000755	<0.0000755	0.0000762 J	0.00775 J	<0.0000755	<0.0000755	<0.0000777	<0.0000777	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.000123	<0.000123	0.000124 J	0.00612 J	<0.000123	<0.000123	<0.000126	<0.000126	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0001	<0.0001	<0.000349	<0.000349	0.000352 J	0.00875 J	<0.000349	<0.000349	<0.000359	<0.000359	0.000058 J	0.000072 J
Chrysene	0.91	2	<0.00005	<0.00005	<0.0000755	<0.0000755	0.0000762 J	0.00794 J	0.0000812 J	0.000132 J	<0.0000777	<0.0000777	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00005	<0.00005	0.000109 J	<0.0000755	0.0000762 J	0.00726 J	<0.000258	<0.000456	<0.0000777	<0.0000777	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000072	<0.000051	<0.000104	<0.000104	0.000105 J	0.00955 J	<0.000104	<0.000104	0.000184 J	0.000107 J	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00005	<0.00005	<0.000066	<0.000066	0.0000667 J	0.00857 J	0.000528 J	0.000866 J	<0.000068	<0.000068	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00005	<0.00005	<0.000066	<0.000066	0.0000667 J	0.00746 J	0.000291 J	0.000705 J	<0.000068	<0.000068	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.00005	<0.00005	0.000429 J	<0.0000755	<0.0021	0.00669 J	0.000523 J	<0.000421	<0.0000777	<0.0000777	<0.00002	<0.00002
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.000104	<0.000104	0.000105 J	0.00628 J	<0.000104	<0.000104	<0.000107	<0.000107	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.0000943	<0.0000943	0.0000952 J	0.00846 J	<0.0000943	<0.0000943	<0.0000971	<0.0000971	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.000575	R	0.000581 J	0.0126 J	<0.000575	<0.000575	<0.000592	<0.000592	<0.00008	<0.000079
Phenanthrene	0.73	2.2	<0.00005	<0.00005	0.000184 J	<0.0000566	0.0000775 J	0.00777 J	0.00128 J	0.00242 J	<0.000087	<0.0000613	<0.000021	<0.000021
Phenol	7.3	22	<0.00005	<0.00005	<0.0000377	R	0.0000381 J	0.00217 J	<0.0000377	<0.0000377	<0.0000388	<0.0000388	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.00005	<0.00005	<0.000104	<0.000104	0.000105 J	0.00813 J	0.000355 J	0.000508 J	<0.000107	<0.000107	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01											0.00128 J	0.00128 J

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-21C 03/20/2018	MW-21C 03/20/2018 Duplicate	MW-21C 05/17/2018	MW-21C 05/17/2018 Duplicate	MW-21C 01/09/2019	MW-21C 01/09/2019 Duplicate	MW-21C 07/16/2019	MW-21C 07/16/2019 Duplicate	MW-21C 01/14/2020	MW-21C 01/14/2020 Duplicate	MW-22A 01/29/2008	MW-22A 07/14/2008
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00109
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00112
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.0015
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00142
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00122
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00138
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00302
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	0.0001 J	<0.00004	<0.00004	<0.00029	<0.00028
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.0002	<0.00019
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	0.023 J	0.000042 J	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	0.003 J	<0.0002	<0.00019
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00039	<0.00038
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	0.000019 J	0.00012 J	<0.000019	<0.000019	0.0004 J	0.00005 J	<0.000019	<0.000019	<0.00039	<0.00038
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00019
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00024
Acenaphthene	1.5	4.4	<0.000027	<0.000027	0.00013	0.00018	<0.000027	<0.000027	0.0057 J	0.00081 J	<0.000027	0.0002	<0.00029	<0.00028
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.000074 J	<0.000015	<0.000015	<0.000015	<0.00029	<0.00028
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.00044 J	0.00005 J	<0.000014	<0.000014	<0.0002	<0.00019
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0002	<0.00019
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00019
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00039	<0.00038
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	0.000048 J	<0.000037	<0.000037	<0.000037	<0.00004	<0.000037	0.00035	0.000065 J	0.0104	<0.00019
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0002	<0.00019
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	0.000052 J	<0.00002	<0.00002	0.0032 J	0.0004 J	<0.00002	0.00017	<0.00029	<0.00028
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00019
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.00081 J	0.000087 J	<0.00001	<0.00001	<0.0002	<0.00019
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.0019 J	0.00025 J	<0.00003	0.00018	<0.0002	<0.00019
Naphthalene	0.49	1.5	<0.00002	<0.00002	0.00002 J	0.0014 J	<0.00002	<0.00002	0.0044 J	0.0008 J	0.00011	0.00013	<0.00039	<0.00038
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00039	<0.00038
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00024
Pentachlorophenol	0.001	0.001	<0.000079	<0.00008	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.0002	<0.00019
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00087 J	0.00011 J	<0.000021	0.000061 J	<0.0002	<0.00019
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	0.00018 J	0.00081 J	<0.000035	<0.000035	<0.0002	<0.00019
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.00039 J	0.000033 J	<0.000019	<0.000019	<0.0002	<0.00019
<b>Metals</b>														
Arsenic	0.01	0.01	0.00109 J	0.00113 J	0.00116 J	0.00108 J	0.00187 J	0.00178 J	0.0013 J	0.00326	0.00109 J	0.00109 J		

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-22A 02/03/2009	MW-22A 01/15/2010	MW-22A 06/29/2010	MW-22A 01/25/2011	MW-22A 07/21/2011	MW-22A 02/15/2012	MW-22A 07/18/2012	MW-22A 01/23/2014	MW-22A 07/30/2014	MW-22AR 02/08/2018	MW-22AR 03/25/2018	MW-22AR 05/31/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.0002	0.00296	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	0.000549	0.0403	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00022	<0.00015	<0.001	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	0.000307 J	0.00925	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	0.000307 J	0.00925	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	0.000834 J	0.0569	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000108	<0.000109	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000304	<0.000307	<0.00004	<0.000041	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000127	<0.000129	<0.000059	<0.000059	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000784	<0.0000792	<0.000042	<0.000043	<0.000042
2-Chloronaphthalene	2	5.8	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000784	<0.0000792	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00007	<0.00007	0.00072	<0.00005	<0.00005	<0.00059	<0.000686	0.0603	<0.00019	<0.00019	<0.00019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000814	<0.000822	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	0.00007 J	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000549	<0.000554	<0.000047	<0.000048	<0.000047
Acenaphthene	1.5	4.4	<0.00009	<0.00009	<0.00009	0.00015 J	<0.00005	0.00009 J	<0.00031	0.00557	0.0783	<0.000027	<0.000028	<0.000027
Acenaphthylene	1.5	4.4	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	0.000742	0.000943	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.0002	<0.00007	<0.00007	0.00011 J	<0.00005	0.000058 J	0.0004	<0.000939	0.00265	<0.000014	<0.000014	0.00017
Benzo(a)anthracene	0.0091	0.02	0.00015 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00011 J	<0.0000784	<0.0000792	<0.000051	<0.000051	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000784	<0.0000792	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000127	<0.000129	<0.00003	<0.000031	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00033	0.0013	<0.0012	<0.0002	0.0015	<0.0011	0.00011 J	<0.000363	0.000703	0.0033	<0.000038	0.0001 J
Chrysene	0.91	2	0.00014 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0001 J	<0.0000784	<0.0000792	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	<0.00008	<0.00008	0.00015 J	<0.00005	0.000074 J	<0.00048	0.001	0.0224	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	0.00017 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000866	<0.000109	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.000086 J	<0.00041	0.000362 J	0.00247	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00007	<0.00007	<0.00007	0.0001 J	<0.00005	0.00008 J	<0.00029	0.000468 J	0.0175	<0.00003	<0.000031	<0.00003
Naphthalene	0.49	1.5	<0.0001	<0.0001	<0.0001	0.0035	0.0001 J	0.00024	<0.0018	<0.0034	<b>0.792</b>	<0.00017	0.00012	<0.00002
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000108	<0.000109	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000098	<0.000099	<0.000025	<0.000026	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	0.00008 J	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000598	<0.000604	<0.00008	<0.000081	<0.000079
Phenanthrene	0.73	2.2	<0.00007	<0.00007	<0.00007	0.00028	<0.00005	0.0002 J	<0.0015	<0.000478	0.000604	<0.000021	0.00012	<0.000021
Phenol	7.3	22	<0.00007	<0.00007	<0.00007	0.00017 J	<0.00005	<0.00005	<0.00005	<0.0000392	<0.0000396	<0.000035	<0.000036	<0.000035
Pyrene	0.73	2.2	0.00013 J	<0.00007	<0.00007	<0.00007	<0.00005	0.000055 J	<0.00033	<0.000108	0.00106	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01										0.000896 J	0.000716 J	0.00293

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-22AR 01/22/2019	MW-22AR 07/31/2019	MW-22AR 02/10/2020	MW-22B 01/29/2008	MW-22B 07/14/2008	MW-22B 02/03/2009	MW-22B 01/15/2010	MW-22B 06/29/2010	MW-22B 01/25/2011	MW-22B 07/21/2011	MW-22B 02/15/2012	MW-22B 07/18/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00052	<0.00109	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00025	0.00313 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	0.0042 J
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.00047	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.00025	0.00262 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	0.0088
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.00054	<0.00122	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.00041	<0.00138	<0.0005	<0.0005	0.00053 J	<0.0005	<0.001	<0.001	0.0033 J
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.00127	0.00339 J	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	0.0057 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.0003	<0.0003	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	0.00014 J	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.0002	<0.0002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.0004	<0.0004	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.00012	<0.0004	<0.0004	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00044	<0.00063
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00025	<0.00025	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.00071	<0.000027	<0.00015	0.0121	0.182	0.022	0.00016 J	0.0093	0.00022	0.003	0.068	0.18
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.0003	0.00192	0.00034	<0.00007	0.00012 J	<0.00007	<0.00005	0.00046	0.0018
Anthracene	7.3	22	0.000028 J	<0.000014	0.0002	0.000948	0.00575	0.00071	<0.00007	0.00031	<0.00007	0.00011 J	0.0017	0.0067
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	0.00029	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	0.0001	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.0004	<0.0004	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000065 J	<0.000037	0.00011 J	0.00172 J	<0.0002	0.00053	0.00022	<0.00061	<0.0002	0.00041	<0.00051	<0.0001
Chrysene	0.91	2	<0.000021	<0.000021	0.00032	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	0.000029 J	<0.00002	<0.00015	0.00363	0.0674	0.0051	0.00026	0.0019	<0.00008	0.00068	0.0079	0.046
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00051	<0.00002	<0.00002	0.00066 J	<0.0002	0.00018 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.000079 J	<0.00001	0.0015	0.00029 J	0.00661	0.0011	0.00011 J	0.00061	<0.00007	0.00019 J	0.0022	0.0065
Fluorene	0.98	2.9	0.00014	<0.00003	0.00019	0.00199	0.0395	0.0018	<0.00007	0.0018	<0.00007	0.00049	0.0035	0.019
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.00044	0.000691	0.00435	0.00017 J	0.00012 J	0.00036	<0.0001	<0.00005	0.0032	0.032
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.0004	<0.0004	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00025	<0.00025	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	0.00029	<0.00005
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
Phenanthrene	0.73	2.2	0.000068 J	<0.000021	0.0015	<0.0002	0.000764	<0.00007	0.00015 J	<0.00007	<0.00007	<0.00005	0.00026	0.0027
Phenol	7.3	22	<0.000035	<0.000035	<0.000041	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0001 J
Pyrene	0.73	2.2	0.000084 J	<0.000019	0.001	<0.0002	0.00239	0.00047	<0.00007	0.00027	<0.00007	0.00012 J	0.001	0.0033
<b>Metals</b>														
Arsenic	0.01	0.01	0.00488	0.000481 J	0.00352									

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-22B 01/23/2014	MW-22B 07/30/2014	MW-22B 08/28/2014	MW-22BR 02/08/2018	MW-22BR 03/25/2018	MW-22BR 05/31/2018	MW-22BR 01/22/2019	MW-22BR 07/31/2019	MW-22BR 02/10/2020	MW-23C 02/04/2009 DNAPL	MW-23C 01/18/2010 DNAPL	MW-23C 06/23/2010 DNAPL
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.005
Benzene	0.005	0.005	0.000304 J	0.00185	0.00238	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<b>0.017</b>	<b>0.012</b>	<b>0.0095 J</b>
Chlorobenzene	0.1	0.1	<0.00018	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.001 J		<0.005
Ethylbenzene	0.7	0.7	0.0022	0.0255	0.0275	0.00034 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.13	0.074	0.12
Methylene chloride	0.005	0.005	<0.00022	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<b>&lt;0.0092</b>
Toluene	1	1	0.00133	0.00584	0.00752	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0023 J	0.0012 J	<0.005
Vinyl chloride	0.002	0.002			<0.00011									
Xylenes (total)	10	10	0.00409	0.0362	0.0383	0.00082 J	<0.0003	0.001	<0.0003	<0.0003	<0.0003	0.073	0.044	0.069 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000108	<0.000109	<0.00011	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<b>&lt;0.005</b>
2,4-Dimethylphenol	0.49	1.5	<0.000304	0.00107	<0.00031	<0.00004	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004	<0.00008	<0.00008	<0.004
2,4-Dinitrotoluene	0.0013	0.003	<0.000127	<0.000129	<0.00013	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<b>&lt;0.0045</b>
2,6-Dinitrotoluene	0.0013	0.003	<0.0000784	<0.0000792	<0.00008	<0.000042	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<b>&lt;0.0035</b>
2-Chloronaphthalene	2	5.8	<0.0000784	<0.0000792	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00011 J	<0.00012	<0.0001	<0.005
2-Methylnaphthalene	0.098	0.29	0.000414 J	0.00721	0.00663	0.00056	0.000067 J	0.000071 J	<0.000019	<0.000019	<0.000043	<b>2.6</b>	<b>0.75</b>	<b>2.7</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000814	<0.000822	<0.00083	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<b>&lt;0.004</b>
4-Nitrophenol	0.049	0.15	<0.000549	<0.000554	<0.00056	<0.000047	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.0035
Acenaphthene	1.5	4.4	0.0244	0.0762	0.123	0.034	0.044	0.047	0.022	<0.000027	0.018	<b>3.4</b>	1.2	<b>3.4</b>
Acenaphthylene	1.5	4.4	0.000886	0.000641	0.00132	<0.000015	0.00096	0.00041	<0.000015	<0.000015	0.00018	0.017	0.01	0.03
Anthracene	7.3	22	<0.00146	0.00292	0.00404	0.00039	0.00062	0.0013	0.00048	<0.000014	0.00042	1.2	0.36	1.2
Benzo(a)anthracene	0.0091	0.02	<0.0000784	<0.0000792	<0.00008	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<b>0.31</b>	<b>0.12</b>	<b>0.3</b>
Benzo(a)pyrene	0.0002	0.0002	<0.0000784	<0.0000792	<0.00008	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<b>0.072</b>	<b>0.029</b>	<b>0.093</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000127	<0.000129	<0.00013	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<b>&lt;0.0045</b>
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000363	0.000672	<0.00037	<0.000037	<0.000038	<0.000037	<0.000037	<0.000037	0.000075 J	0.002	<0.0011	<b>&lt;0.01</b>
Chrysene	0.91	2	<0.0000784	<0.0000792	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.28	0.093	0.27
Dibenzofuran	0.098	0.29	0.00784	0.0238	0.0409	0.00066	0.00049	0.0011	0.00029	<0.00002	<0.00029	<b>3.5</b>	<b>1.2</b>	<b>3.6</b>
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000873	<0.000109	<0.00011	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00007	<0.00007	<0.0035
Fluoranthene	0.98	2.9	0.00187	0.00304	0.00282	0.00098	0.0011	0.0028	0.0011	<0.00001	0.0017	<b>3</b>	0.77	<b>3</b>
Fluorene	0.98	2.9	0.00521	0.0198	0.0355	0.002	0.0017	0.0046	0.0025	<0.00003	0.0026	<b>2.5</b>	0.82	<b>2.6</b>
Naphthalene	0.49	1.5	0.13 J	<b>0.832</b>	<b>0.977</b>	0.0038	0.0008	0.00044	0.00017	<0.00002	<0.00013	<b>9.9</b>	<b>3.9</b>	<b>8.9</b>
Nitrobenzene	0.049	0.15	<0.000108	<0.000109	<0.00011	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.0045
N-Nitrosodiphenylamine	0.19	0.42	<0.000098	<0.000099	<0.0001	<0.000025	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.0045
Pentachlorophenol	0.001	0.001	<0.000598	<0.000604	<0.00061	<0.00008	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<b>&lt;0.004</b>
Phenanthrene	0.73	2.2	<0.000562	0.00053	<0.00006	0.00081	0.00018	0.00039	<0.000021	<0.000021	<0.00013	<b>8.8</b>	<b>2.7</b>	<b>8.2</b>
Phenol	7.3	22	<0.000392	<0.000396	<0.00004	<0.000035	<0.000036	<0.000035	<0.000035	0.00069	<0.000035	<0.00007	<0.00007	<0.0035
Pyrene	0.73	2.2	0.000876	0.00123	0.0023	0.00062	0.00077	0.0015	0.00046	<0.000019	0.00082	<b>1.6</b>	0.59	<b>1.9</b>
<b>Metals</b>														
Arsenic	0.01	0.01				<b>0.0219</b>	<b>0.0159</b>	<b>0.0301</b>	<b>0.0535</b>	0.000559 J	<b>0.016</b>			

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-23C 01/19/2011	MW-23C 07/22/2011	MW-23C 02/02/2012	MW-23C 07/12/2012	MW-23C 02/11/2013	MW-23C 07/31/2013	MW-23C 01/15/2014	MW-23C 08/28/2014	MW-23C 01/09/2020	MW-24AR 02/05/2009	MW-24AR 01/14/2010	MW-24AR 06/29/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.005	<0.01	<0.01	<0.0005	<0.00014	<0.0014	<0.0002	<0.0014	<0.001	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<b>0.0072 J</b>	<0.01	<0.01	<b>0.0071</b>	<b>0.0111</b>	<b>0.0138</b>	<b>0.0126</b>	<b>0.00596 J</b>	0.0027 J	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.005	<0.01	<0.01	<0.0005	0.000279 J	0.00146 J	<0.00018	<0.0012	<0.0015	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	0.13	0.1	0.1	0.17	0.151	0.185	0.165	0.15	0.028	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.005	<b>&lt;0.013</b>	<b>&lt;0.013</b>	<0.001	<0.00015	<0.0015	<0.00022	<0.0015	<0.005	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.005	<0.01	<0.01	0.0025 J	0.00433	0.00819 J	0.00728	0.00378 J	<0.001	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002	<b>&lt;0.005</b>	<b>&lt;0.01</b>	<b>&lt;0.01</b>	<0.0005	<0.00011	<0.0011	<0.00018	<0.0011		<0.0005	<0.0005	<0.0005
Xylenes (total)	10	10	0.059 J	0.048 J	0.039 J	0.11	0.0884	0.0988	0.0959	0.0915	0.025	<0.001	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.0005	<0.001	<0.001	<b>&lt;0.00534</b>	<b>&lt;0.0259</b>	<b>&lt;0.022</b>	<0.00021	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00008	0.0035	0.0014 J	0.028	<0.00282	<0.015	<0.0731	0.202	<0.047	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.0005	<0.001	<0.00118	<b>&lt;0.00631</b>	<b>&lt;0.0307</b>	<b>&lt;0.026</b>	<0.00058	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.0006	<0.0012	<0.000727	<b>&lt;0.00388</b>	<b>&lt;0.0189</b>	<b>&lt;0.016</b>	<0.00042	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.0005	<0.001	<0.000727	<0.00388	<0.0189	<0.016	<0.00021	<0.00012	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<b>1.2</b>	<b>1.3</b>	<b>0.65</b>	<b>28</b>	<b>1.38</b>	<b>1.16</b>	<b>4.52</b>	<b>18.3</b>	<b>0.12</b>	<0.00007	0.00023	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.0008	<0.0016	<b>&lt;0.00755</b>	<b>&lt;0.0403</b>	<b>&lt;0.196</b>	<b>&lt;0.166</b>	<0.0002	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.0005	<0.001	<0.00509	<0.0272	<b>&lt;0.132</b>	<b>&lt;0.112</b>	<0.00047	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<b>1.6</b>	<b>2</b>	0.89	<b>39</b>	<b>1.78</b>	<b>1.58</b>	<b>7.79</b>	<b>25.9</b>	0.21	<0.00009	<0.00009	<0.00009
Acenaphthylene	1.5	4.4	0.012	0.015	0.0068	0.45	<0.000545	<0.00291	<0.0142	0.336	0.0029	<0.00006	<0.00007	<0.00007
Anthracene	7.3	22	0.4	1.7	0.25	<b>16</b>	0.641	0.31	1.49	<b>8.74</b>	0.035	<0.00007	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<b>0.12</b>	<b>0.15</b>	<b>0.046</b>	<b>4.8</b>	<b>0.104</b>	<b>0.0905</b>	<b>0.5</b>	<b>2.63</b>	0.008	<0.00007	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<b>0.04</b>	<b>0.044</b>	<b>0.016</b>	<b>1.2</b>	<b>0.0283</b>	<b>0.0235 J</b>	<b>0.119</b>	<b>0.73</b>	<b>0.0022</b>	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.0005	<b>&lt;0.001</b>	<b>&lt;0.00118</b>	<b>&lt;0.00631</b>	<b>&lt;0.0307</b>	<b>&lt;0.026</b>	<0.0003	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0014	0.0019	<0.001	<b>0.042</b>	<0.00336	<b>&lt;0.018</b>	<b>&lt;0.0873</b>	<b>&lt;0.074</b>	<0.00037	0.00031	<0.00029	<0.00024
Chrysene	0.91	2	0.099	0.21	0.044	<b>4.3</b>	0.103	0.0819	0.476	<b>2.24</b>	0.0072	<0.00007	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<b>1.6</b>	<b>2.7</b>	<b>0.85</b>	<b>46</b>	<b>1.82</b>	<b>1.48</b>	<b>5.45</b>	<b>25.7</b>	<b>0.17</b>	<0.00008	0.000084 J	0.00011 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00005	<0.0005	<0.001	<0.001	<0.00534	<0.0259	<0.022	<0.0002	0.0019	0.0001 J	<0.00007
Fluoranthene	0.98	2.9	<b>0.99</b>	<b>1.8</b>	0.48	<b>34</b>	<b>1.09</b>	0.812	<b>4.42</b>	<b>20.4</b>	0.068	<0.00007	0.00011 J	<0.00007
Fluorene	0.98	2.9	0.88	<b>2</b>	0.57	<b>32</b>	<b>1.19</b>	0.874	<b>3.78</b>	<b>20.5</b>	0.12	<0.00007	<0.00007	<0.00007
Naphthalene	0.49	1.5	<b>8.5</b>	<b>7.5</b>	<b>7.8</b>	<b>83</b>	<b>12.2</b>	<b>13.2</b>	<b>43.8</b>	<b>57.9</b>	<b>0.66</b>	<0.0001	<0.0023	0.00036
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.0005	<0.001	<0.001	<0.00534	<0.0259	<0.022	<0.00024	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.0005	<0.001	<0.000909	<0.00485	<0.0236	<0.02	<0.00025	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.00008	<0.00005	<0.0005	<0.001	<b>&lt;0.00555</b>	<b>&lt;0.0296</b>	<b>&lt;0.144</b>	<b>&lt;0.122</b>	<0.00079	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<b>3.6</b>	<b>3.8</b>	<b>1.9</b>	<b>130</b>	<b>3.48</b>	<b>2.8</b>	<b>18.2</b>	<b>59.4</b>	0.25	<0.00007	0.00018 J	<0.00007
Phenol	7.3	22	<0.00007	<0.00005	0.0011 J	<0.001	<0.000364	<0.00194	<0.00943	<0.008	0.0079	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	0.6	<b>1.1</b>	0.35	<b>21</b>	<b>0.754</b>	0.515	<b>3.04</b>	<b>13.3</b>	0.043	<0.00007	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01									0.00333			

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-24AR 01/25/2011	MW-24AR 07/21/2011	MW-24AR 02/09/2012	MW-24AR 07/25/2012	MW-24AR 02/12/2013	MW-24AR 08/08/2013	MW-24AR 01/23/2014	MW-24B 01/28/2008	MW-24B 07/14/2008	MW-24B 02/03/2009	MW-24B 01/14/2010	MW-24B 06/29/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00245	<0.00109	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00257	<0.00112	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.0005	<0.001	<0.001	<0.0005	0.000201 J	<0.00012	<0.00018	<0.00239	<0.0015	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00203	<0.00142	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00195	<0.00122	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00274	<0.00138	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002							<0.00011					
Xylenes (total)	10	10	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00581	<0.00302	<0.001	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00005	0.00013 J	<0.00005	R	<0.000292	<0.000295	<0.00029	<0.0003	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.00019	<0.0002	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000755	<0.0000762	<0.00019	<0.0002	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.00038	<0.0004	<0.00012	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	0.00018 J	<0.00005	<0.00005	<0.00005	<0.0000667	<0.000066	0.0000924 J	<0.00038	<0.0004	<0.00007	<0.00007	0.000099 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.000783	<0.00079	<0.00019	<0.0002	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.00005	<0.00005	R	<0.000528	<0.000533	<0.00024	<0.00025	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	0.000146 J	<0.00029	<0.0003	<0.00009	<0.00009	<0.00009
Acenaphthylene	1.5	4.4	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000566	<0.0000571	<0.00029	<0.0003	<0.00006	<0.00007	<0.00007
Anthracene	7.3	22	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000476	<0.0000472	<0.000452	0.00066	<0.0002	<0.00007	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.00019	<0.0002	0.00015 J	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.00019	<0.0002	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.00038	<0.0004	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0002	0.00089	0.00048	<0.0001	<0.000352	<0.000349	0.000767 J	<0.00019	<0.0002	0.00046	<0.00021	<0.00074
Chrysene	0.91	2	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.00019	<0.0002	0.00015 J	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	0.000164 J	0.000568	<0.0003	<0.00008	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00005	0.00017 J	<0.000071	<0.000105	0.000168 J	<0.000837	<0.00019	<0.0002	<0.00007	<0.00007	<0.00007
Fluoranthene	0.98	2.9	<0.00007	<0.00005	0.000069 J	<0.00005	<0.0000667	<0.000066	0.000224 J	<0.00019	<0.0002	0.00011 J	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.000066	0.000137 J	0.00026 J	<0.0002	<0.00007	<0.00007	<0.00007
Naphthalene	0.49	1.5	0.00081	<0.00005	<0.00005	<0.00005	<0.000139	<0.0000755	<0.000744	0.00105	<0.0004	<0.0001	<0.0001	0.00083
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.00038	<0.0004	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000943	<0.0000952	<0.00024	<0.00025	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.00008	<0.00005	<0.00005	<0.00005	R	<0.000575	<0.000581	<0.00019	<0.0002	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	0.0001 J	<0.00005	<0.00005	<0.00005	0.000089 J	<0.0000566	<0.000691	0.000676	<0.0002	<0.00007	<0.00007	<0.00007
Phenol	7.3	22	<0.00007	<0.00005	0.00005 J	<0.00005	R	<0.0000377	<0.0000381	<0.00019	<0.0002	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.00007	<0.00005	0.000067 J	<0.00005	<0.000105	<0.000104	0.000172 J	<0.00019	<0.0002	<0.00007	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-24B 01/25/2011	MW-24B 07/21/2011	MW-24B 02/09/2012	MW-24B 07/25/2012	MW-24B 02/12/2013	MW-24B 08/08/2013	MW-24C 01/28/2008	MW-24C 07/14/2008	MW-24C 02/03/2009	MW-24C 01/14/2010	MW-24C 06/29/2010	MW-24C 01/25/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00245	<0.00109	<0.0005	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.00257	<0.00112	<0.0005	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00239	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00203	<0.00142	<0.0005	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00195	<0.00122	<0.0005	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.0005	<0.001	<0.001	<0.0005	<0.00015	0.0209	<0.00274	<0.00138	<0.0005	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002						<0.00011						
Xylenes (total)	10	10	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00581	<0.00302	<0.001	<0.001	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00005	<0.00005	<0.00005	<0.000295	<0.000304	<0.00029	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.00019	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000784	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00038	<0.00038	<0.00012	<0.0001	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	<0.00038	<0.00038	<0.00007	<0.00007	<0.00007	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.000079	<0.0000814	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.000549	<0.00024	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00029	<0.00029	<0.00009	<0.00009	0.00022	<0.00009
Acenaphthylene	1.5	4.4	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000588	<0.00029	<0.00029	<0.00006	<0.00007	<0.00007	<0.00007
Anthracene	7.3	22	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000476	<0.000049	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.00038	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0002	0.00014 J	0.00011 J	<0.00015	<0.000352	<0.000363	<0.00019	<0.00019	0.00055	<0.0002	<0.001	<0.0002
Chrysene	0.91	2	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	0.000571	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00005	<0.00005	<0.00013	<0.000105	<0.000108	0.00063 J	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Fluoranthene	0.98	2.9	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	0.000529	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Naphthalene	0.49	1.5	<0.0001	<0.00005	<0.00005	0.00015 J	<0.0000762	<0.0000784	<0.00038	<0.00038	0.00013 J	<0.00026	<0.0001	<0.0001
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.00038	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.000098	<0.00024	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.000598	<0.00019	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000588	0.000677	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Phenol	7.3	22	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000381	<0.0000392	<0.00019	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.00007	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	0.00041 J	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-24C 07/21/2011	MW-24C 02/09/2012	MW-24C 07/25/2012	MW-24C 02/12/2013	MW-24C 08/08/2013	MW-25A 01/28/2008	MW-25A 07/14/2008	MW-25A 02/03/2009	MW-25A 01/15/2010	MW-25A 06/30/2010	MW-25A 01/26/2011	MW-25A 01/26/2011 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00245	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.00257	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00239	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00203	<0.00025	0.0029 J	<0.0005	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00195	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.001	<0.001	<0.0005	0.000218 J	<0.00015	<0.00274	<0.00041	0.00074 J	<0.0005	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002					<0.00011							
Xylenes (total)	10	10	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00581	<0.00127	0.0047 J	<0.001	<0.001	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.00008	<0.00009	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.00005	<0.00005	<0.000295	<0.000304	<0.00029	<0.00033	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.00019	<0.00022	<0.00009	<0.00009	0.00066	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000784	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00038	<0.00044	<0.00012	<0.0001	<0.0001	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.00005	<0.00005	0.000077 J	<0.0000667	<0.0000686	<0.00038	<0.00044	0.024	<0.00007	<0.00007	<0.00007	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00079	<0.000814	<0.00019	<0.00022	<0.00008	<0.00008	0.00026	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000533	<0.000549	<0.00024	<0.00028	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	0.00038 J	<0.00033	0.034	0.0014	0.0012	0.00054	0.00052
Acenaphthylene	1.5	4.4	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000588	<0.00029	<0.00033	0.0004	<0.00007	0.00034	<0.00007	<0.00007
Anthracene	7.3	22	<0.00005	<0.00005	<0.00005	<0.0000476	<0.000049	0.00067	<0.00022	0.0005	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00019	<b>&lt;0.00022</b>	<0.00008	<0.00008	0.00012 J	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.00038	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00013 J	0.0013	<0.00013	<0.000352	<0.000363	0.00021 J	<0.00022	0.00033	<0.0002	<0.00056	<0.0002	0.00046
Chrysene	0.91	2	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	0.000601	<0.00033	0.018	<0.00008	0.00034	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Fluoranthene	0.98	2.9	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	0.000556	<0.00022	0.00057	0.000084 J	0.000092 J	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	0.00033 J	<0.00022	0.0049	<0.00007	<0.00007	<0.00007	<0.00007
Naphthalene	0.49	1.5	0.0002	<0.00005	0.00019 J	<0.0000762	<0.0000784	0.0011	<0.00044	0.45	<0.0001	0.00024	0.00018 J	0.00027
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.00038	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.0000952	<0.000098	<0.00024	<0.00028	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00005	<0.000581	<0.000598	<0.00019	<0.00022	<0.00008	<0.00008	0.00033	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000588	0.000715	<0.00022	0.0034	<0.00007	<0.00007	<0.00007	<0.00007
Phenol	7.3	22	<0.00005	<0.00005	<0.00005	<0.0000381	<0.0000392	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	0.000484	<0.00022	0.00036	0.00047	0.00015 J	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-25A 07/20/2011	MW-25A 02/08/2012	MW-25A 07/18/2012	MW-25A 02/06/2013	MW-25A 08/06/2013	MW-25A 01/22/2014	MW-25A 07/29/2014	MW-25A 01/31/2018	MW-25A <sup>8</sup> 03/27/2018	MW-25A 05/31/2018	MW-25A 01/15/2019	MW-25A 07/16/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.0005	0.000497 J	0.000121 J	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00018	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<0.000106	<0.00011	<0.000105	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.00005	<0.00005	<0.000298	<0.00031	<0.000295	<0.000301	<0.00004	<0.000041	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00005	<0.000125	<0.00013	<0.000124	<0.000126	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.00006	<0.0000769	<0.00008	<0.0000762	<0.0000777	<0.000042	<0.000043	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000762	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00005	<0.00005	<0.000061	<0.0000673	<0.00007	<0.0000667	<0.000068	0.00014	<0.000019	0.000052 J	<0.000051	0.000057 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.000798	<0.00083	<0.00079	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000538	<0.00056	<0.000533	<0.000544	<0.000047	<0.000048	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.0042	0.00053	<0.00005	0.000171 J	0.000345 J	0.00356	0.0000912 J	0.062	0.054	<0.000027	0.000036 J	<0.000027
Acenaphthylene	1.5	4.4	0.000053 J	<0.00005	<0.00005	<0.0000577	<0.00006	0.000542	<0.0000583	0.00044	0.00053	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	<0.00005	<0.00005	<0.00005	<0.0000481	<0.00005	<0.0000476	<0.0000485	0.00057	0.00068	<0.000014	0.000015 J	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000762	<0.0000777	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000762	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00005	<0.000125	<0.00013	<0.000124	<0.000126	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00023	<0.0001	<0.0001	<0.000356	<0.00037	<0.000352	<0.000359	<0.00008	<0.000038	0.0001 J	<0.000078	<0.000037
Chrysene	0.91	2	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000762	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.0013	0.0005	<0.00005	<0.0000769	<0.00008	<0.0000762	<0.0000777	0.00017	0.00011	0.000033 J	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.00005	<0.000106	<0.00011	<0.000105	<0.000107	<0.00002	<0.00002	<0.00002	0.00002 J	<0.00002
Fluoranthene	0.98	2.9	0.00014 J	<0.00005	<0.00005	<0.0000673	<0.00007	0.00106	<0.000068	0.0062	0.0066	<0.00001	0.000015 J	<0.00001
Fluorene	0.98	2.9	0.00016 J	0.00011 J	<0.00005	<0.0000673	<0.00007	<0.0000667	<0.000068	0.026	0.022	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.00005	0.00017 J	<0.00038	<0.000692	<0.00008	<0.000217	<0.000817	<0.00029	0.0002	0.0006	<0.00029	0.00023
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.000106	<0.00011	<0.000105	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.0000962	<0.0001	<0.0000952	<0.0000971	<0.000025	<0.000026	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00005	<0.000587	<0.00061	<0.000581	<0.000592	<0.000079	<0.000081	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.00005	<0.00005	<0.00005	<0.0000577	<0.00006	<0.0000571	<0.0000583	0.00015	0.00015	<0.000021	0.000029 J	<0.000021
Phenol	7.3	22	<0.00005	0.00005 J	<0.00005	<0.0000385	<0.00004	<0.0000381	<0.0000388	<0.000035	<0.000036	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0009	<0.00005	<0.00005	<0.000106	0.000124 J	0.000585	0.00018 J	0.0033	0.0049	<0.000019	0.000027 J	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01								0.0171	0.00714	0.00171 J	0.00216	0.00285

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-25A 01/15/2020	MW-25C 01/28/2008 DNAPL	MW-25C 01/15/2010 DNAPL	MW-25C 01/26/2011	MW-25C 07/20/2011	MW-25C 02/08/2012	MW-25C 07/18/2012	MW-25C 02/06/2013	MW-25C 08/06/2013	MW-25C 01/22/2014	MW-25C 07/29/2014	MW-25C 01/28/2018
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.001	<0.025	<0.0005	<0.005	<0.01	<0.005	<0.0028	<0.0007	<0.0002	<0.0007	<0.0002
Benzene	0.005	0.005	<0.0002	0.11	0.11 J	0.092	0.076	0.039 J	0.03 J	0.0304	0.0283	0.022	0.0119	0.00089 J
Chlorobenzene	0.1	0.1	<0.0003	<0.00094	<0.025	<0.0005	<0.005	<0.01	<0.005	0.00653 J	<0.0006	0.00034 J	<0.0006	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	0.545	0.47	0.5	0.37	0.34	0.33	0.324	0.173	0.32	0.298	0.042
Methylene chloride	0.005	0.005	<0.001	<0.0011	<0.025	<0.0005	<0.0065	<0.013	<0.01	<0.003	<0.00075	<0.00022	<0.00075	<0.001
Toluene	1	1	<0.0002	0.556	0.52	0.53	0.4	0.31	0.31	0.291	0.204	0.261	0.207	0.015
Vinyl chloride	0.002	0.002	<0.0002				<0.005	<0.01	<0.005	<0.0022	<0.00055	<0.00018	<0.00055	<0.0002
Xylenes (total)	10	10	<0.0003	1.43	1.2	1.2	1	0.98	0.96	1.03	0.575	1.01	1.07	0.29
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.8	0.0001 J	<0.0001	<0.00005	<0.00005	<0.00005	<0.106	<0.011	<0.0105	<0.000107	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	<2.9	0.00008 J	<0.00008	0.0051	<0.00005	<0.00005	<0.298	<0.031	0.372	<0.000301	0.00061
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<1.9	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005	<0.125	<0.013	<0.0124	<0.000126	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<1.9	0.00007 J	<0.00007	<0.00006	<0.00006	<0.00006	<0.0769	<0.008	<0.00762	<0.0000777	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<3.8	0.0001 J	<0.0001	<0.00005	<0.00005	<0.00005	<0.0769	<0.008	<0.00762	<0.0000777	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000019	193	0.76	1.4	1.3	0.92	0.9	0.8	1.32	1.46	0.943	0.48
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<1.9	0.00008 J	<0.00008	<0.00008	<0.00008	<0.00008	<0.798	<0.083	<0.079	<0.000806	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<2.4	0.00007 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.538	<0.056	<0.0533	<0.000544	<0.000047
Acenaphthene	1.5	4.4	<0.000027	97.5	0.21	0.55	0.28	0.26	0.21	0.261 J	0.381	0.416	0.284	0.17
Acenaphthylene	1.5	4.4	<0.000015	<2.9	0.0027 J	0.0041	0.0029	0.0021	0.0021	<0.0577	<0.006	<0.00571	0.00316	0.0017
Anthracene	7.3	22	0.000089 J	42.8	0.035	0.19	0.031	0.021	0.019	<0.0481	0.0377 J	0.0372 J	0.0209	0.015
Benzo(a)anthracene	0.0091	0.02	0.00012	9.37	0.0027 J	0.0047	0.0014	0.00054	0.00086	<0.0769	<0.008	<0.00762	0.000813	0.00079
Benzo(a)pyrene	0.0002	0.0002	0.000038 J	2.3	0.0014 J	0.013	0.00043	0.00017 J	0.0002	<0.0769	<0.008	<0.00762	0.000435 J	0.00028
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<3.8	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005	<0.125	<0.013	<0.0124	0.0016	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<1.9	0.0002 J	<0.0002	<0.002	<0.0001	0.00012 J	<0.356	<0.037	<0.0352	<0.000674	0.00006 J
Chrysene	0.91	2	0.000097 J	8.38	0.0025 J	0.048	0.0012	0.00062	0.00086	<0.0769	<0.008	<0.00762	0.000957	0.00077
Dibenzofuran	0.098	0.29	<0.00002	102	0.22	0.52	0.29	0.26	0.22	0.174 J	0.353	<0.00762	0.276	0.18
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<1.9	0.00007 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.106	<0.011	<0.0105	<0.000107	<0.00002
Fluoranthene	0.98	2.9	0.00045	84.1	0.041	0.32	0.02	0.011	0.0088	<0.0673	0.0149 J	0.018 J	0.0127	0.0092
Fluorene	0.98	2.9	<0.00003	75.5	0.12	0.34	0.14	0.13	0.096	0.102 J	0.163	<0.00667	0.129	0.086
Naphthalene	0.49	1.5	0.0001	750	9.8	18	19	15	13	10.7	19.7	19	10.7	3
Nitrobenzene	0.049	0.15	<0.000024	<3.8	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005	<0.106	<0.011	<0.0105	<0.000107	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<2.4	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005	<0.0962	<0.01	<0.00952	<0.0000971	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<1.9	0.00008 J	<0.00008	<0.00005	<0.00005	<0.00005	<0.587	<0.061	<0.0581	<0.000592	<0.000079
Phenanthrene	0.73	2.2	0.00017	214	0.19	0.7	0.18	0.14	0.12	0.147 J	0.187	0.222	0.14	0.089
Phenol	7.3	22	<0.000035	<1.9	0.00007 J	<0.00007	0.026	0.003	0.0045	<0.0385	0.12	<0.00381	0.00177	<0.000035
Pyrene	0.73	2.2	0.00035	49.5	0.022	0.24	0.0092	0.0047	0.0063	<0.106	<0.011	<0.0105	0.00769	0.0062
<b>Metals</b>														
Arsenic	0.01	0.01	0.00216											0.00283

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray.
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-25C <sup>8</sup> 03/26/2018	MW-25C 05/31/2018	MW-25C 01/15/2019	MW-25C 07/16/2019	MW-25C 01/15/2020	MW-26A 01/29/2008	MW-26A 07/14/2008	MW-26A 02/03/2009	MW-26A 01/13/2010	MW-26A 06/25/2010	MW-26A 01/24/2011	MW-26A 07/19/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Benzene	0.005	0.005	0.00047 J	0.0013	<0.0002	0.0013	0.0017	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<b>0.031</b>
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Ethylbenzene	0.7	0.7	0.043	0.036	0.038	0.035	0.033	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013
Toluene	1	1	0.015	0.013	0.013	0.011	0.01	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002							
Xylenes (total)	10	10	0.32	0.28	0.27	0.25	0.22	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	0.0045 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005
2,4-Dimethylphenol	0.49	1.5	0.0033	0.0075	<0.00004	0.0075	0.0062	<0.0003	<0.0003	0.00054	<0.00008	<0.00008	<0.00008	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.00058	<0.00058	<0.00058	<0.00058	<0.00058	<0.0002	<0.0002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006
2-Chloronaphthalene	2	5.8	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.0004	<0.0004	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005
2-Methylnaphthalene	0.098	0.29	<b>0.69</b>	<b>0.71</b>	<b>0.4</b>	<b>0.54</b>	<b>0.96</b>	<0.0004	<0.0004	0.0024	<0.00007	<0.00007	0.00031	0.00039
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047	<0.00025	<0.00025	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Acenaphthene	1.5	4.4	0.21	0.23	0.13	0.2	0.36	0.0519	0.0173	0.015	0.0097	0.005	0.0039	0.12
Acenaphthylene	1.5	4.4	0.0023	0.0028	0.0012	0.002	0.0025	<0.0003	<0.0003	<0.00006	0.00014 J	<0.00007	<0.00007	0.00047
Anthracene	7.3	22	0.019	0.024	0.011	0.016	0.027	0.0018	0.00047 J	0.00079	<0.00007	0.0002	0.000099 J	0.0026
Benzo(a)anthracene	0.0091	0.02	0.0011	0.0012	0.00065	0.0013	0.00057	<0.0002	<0.0002	0.00016 J	<0.00007	<0.00007	<0.00007	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<b>0.00042 J</b>	<b>0.00021</b>	<b>0.00037 J</b>	0.00019	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0003	<0.0003	<0.00003	<0.0003	<0.00003	<0.0004	<0.0004	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00037	<0.00037	<0.000037	<0.00037	0.00066	<b>0.0214</b>	<0.0002	0.00042	<0.00026	<0.00045	0.00043	<0.00031
Chrysene	0.91	2	0.0013	0.0014	0.00074	0.0014	0.00056	<0.0002	<0.0002	0.00018 J	<0.00007	0.0003	<0.00007	<0.00005
Dibenzofuran	0.098	0.29	<b>0.2</b>	<b>0.21</b>	<b>0.14</b>	<b>0.19</b>	<b>0.34</b>	0.00492	0.00123	0.0026	0.00078	0.00033	0.00038	0.021
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.0002	<0.00002	<0.0002	0.00068	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Fluoranthene	0.98	2.9	0.014	0.015	0.0079	0.014	0.0058	0.00244	0.000523	0.00091	0.0003	0.0004	0.00036	0.0048
Fluorene	0.98	2.9	0.091	0.11	0.062	0.081	0.15	0.00235	0.000742	0.0016	0.00028	0.00034	0.00017 J	0.0057
Naphthalene	0.49	1.5	<b>5.9</b>	<b>5.6</b>	<b>3.5</b>	<b>4.2</b>	<b>7.5</b>	0.00078	<0.0004	0.0074	<0.00051	<0.0001	0.0043	0.0019
Nitrobenzene	0.049	0.15	<0.00024	<0.00024	<0.000024	<0.00024	<0.000024	<0.0004	<0.0004	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.00025	<0.00025	<0.000025	<0.00025	<0.000025	<0.00025	<0.00025	<0.00009	<0.00009	<0.00009	<0.00009	0.00023
Pentachlorophenol	0.001	0.001	<0.00079	<0.00079	<0.000079	<0.00079	<b>0.17</b>	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005
Phenanthrene	0.73	2.2	0.15	0.16	0.077	0.12	0.18	<0.0002	<0.0002	0.003	0.00021	0.00017 J	0.00011 J	0.00029
Phenol	7.3	22	<0.00035	<0.00035	<0.000035	<0.00035	<0.000035	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Pyrene	0.73	2.2	0.011	0.01	0.0052	0.0085	0.0049	0.00126	0.00021 J	0.00069	0.000092 J	0.0002 J	0.00013 J	0.0031
<b>Metals</b>														
Arsenic	0.01	0.01	0.003	0.00305	0.00359	0.00487	0.00391							

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-26A 08/25/2011	MW-26A 10/20/2011	MW-26A 02/15/2012	MW-26A 07/17/2012	MW-26A 02/06/2013	MW-26A 08/07/2013	MW-26A 10/14/2013	MW-26A 01/22/2014	MW-26A 07/24/2014	MW-26A 01/28/2018	MW-26A 03/21/2018	MW-26A 06/05/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005			<0.001	<0.0005	<0.00014	<0.00014		<0.0002	<0.00014	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	0.042	0.004 J	<0.001	<0.0005	0.00118	0.0097	0.00391	0.000434 J	0.000189 J	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1			<0.001	<0.0005	0.000176 J	0.000297 J		<0.00018	0.000205 J	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7			<0.0011	<0.0005	<0.00011	0.000815 J		<0.00019	<0.00011	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005			<0.0013	<0.001	<0.00015	<0.00015		<0.00022	<0.00015	<0.001	<0.001	<0.001
Toluene	1	1			<0.001	<0.0005	<0.00015	<0.000291		<0.00017	<0.00015	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002									<0.00011			
Xylenes (total)	10	10			<0.0031	<0.0015	<0.00026	0.00239 J		<0.00058	<0.00026	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026			<0.00005	<0.00005	<0.00529	<0.000104		<0.00011	<0.000107	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5			<0.00005	<0.00005	<0.0149	<0.000292		<0.00031	<0.000301	0.000056 J	<0.000041	<0.00004
2,4-Dinitrotoluene	0.0013	0.003			<0.00005	0.0001 J	<0.00625	<0.000123		<0.00013	<0.000126	<0.000058	<0.000059	<0.000058
2,6-Dinitrotoluene	0.0013	0.003			<0.00006	<0.00006	<0.00385	<0.0000755		<0.00008	<0.0000777	<0.000042	<0.000043	<0.000042
2-Chloronaphthalene	2	5.8			<0.00005	<0.00005	<0.00385	<0.0000755		<0.00008	<0.0000777	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29			<0.00005	0.000056 J	<0.00337	0.000414 J		<0.00007	<0.000068	0.000053 J	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073			<0.00008	<0.00008	<0.0399	<0.000783		<0.00083	<0.000806	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15			<0.00005	<0.00005	<0.0269	<0.000528		<0.00056	<0.000544	<0.000047	<0.000048	<0.000047
Acenaphthene	1.5	4.4			0.0095	0.0087	0.0481	0.141		0.0699	0.0663	0.0073	<0.000028	0.019
Acenaphthylene	1.5	4.4			0.00013 J	<0.00005	<0.00288	<0.0000566		<0.00006	0.000486	0.00007 J	<0.000015	0.000082 J
Anthracene	7.3	22			0.00025	0.00027	<0.0024	0.00228		0.00136	0.00141	0.00012	0.00015	0.00025
Benzo(a)anthracene	0.0091	0.02			<0.00005	<0.00005	<0.00385	<0.0000755		<0.00008	<0.0000777	<0.00005	<0.000051	<0.00005
Benzo(a)pyrene	0.0002	0.0002			<0.00005	<0.00005	<0.00385	<0.0000755		<0.00008	<0.0000777	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019			<0.00005	<0.00005	<0.00625	<0.000123		<0.00013	<0.000126	<0.00003	<0.000031	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006			<0.0001	0.0002 J	<0.0178	<0.000349		<0.00037	<0.000359	0.000076 J	<0.000038	<0.000037
Chrysene	0.91	2			<0.00005	<0.00005	<0.00385	<0.0000755		<0.00008	<0.0000777	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29			0.0014	0.00084	0.00416 J	0.0151		<0.00008	0.00154	<0.00016	0.000086 J	0.0001
Di-n-butylphthalate (DBP)	2.4	7.3			<0.00005	<0.00005	<0.00529	<0.000104		<0.00011	<0.000107	<0.00002	0.000052 J	<0.00002
Fluoranthene	0.98	2.9			0.00049	0.00092	<0.00337	0.0062		0.00306	0.00465	0.00084	0.0011	0.0014
Fluorene	0.98	2.9			0.0006	0.00041	<0.00337	0.00611		0.0031	0.00245	0.00044	0.00051	0.00084
Naphthalene	0.49	1.5			0.0001 J	0.00027	<0.00385	0.0066		<0.000924	0.000419 J	<0.00014	<0.00002	<0.00002
Nitrobenzene	0.049	0.15			<0.00005	<0.00005	<0.00529	<0.000104		<0.00011	0.00394	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42			<0.00005	<0.00005	<0.00481	<0.0000943		<0.0001	<0.0000971	<0.000025	<0.000026	<0.000025
Pentachlorophenol	0.001	0.001			<0.00005	<0.00005	<0.0293	<0.000575		<0.00061	<0.000592	<0.000079	<0.000081	<0.000079
Phenanthrene	0.73	2.2			<0.00005	<0.00005	<0.00288	<0.0000566		0.000147 J	0.000155 J	<0.000038	<0.000021	<0.000021
Phenol	7.3	22			<0.00005	<0.00005	<0.00192	<0.0000377		<0.00004	<0.0000388	<0.000035	<0.000036	<0.000035
Pyrene	0.73	2.2			0.00024	0.00051	<0.00529	0.00322		0.00159	0.0022	0.00043	0.00042	0.00094
<b>Metals</b>														
Arsenic	0.01	0.01										0.032	0.0427	0.0491

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-26A 01/15/2019	MW-26A 07/17/2019	MW-26A 01/16/2020	MW-27A 01/28/2008	MW-27A 07/14/2008	MW-27A 02/09/2018	MW-27A 03/26/2018	MW-27A 06/01/2018	MW-27A 07/18/2019	MW-27A 01/15/2020	MW-27C 01/28/2008	MW-27C 07/14/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00245	<0.00109	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00245	<0.00109
Benzene	0.005	0.005	<0.0002	0.00036 J	<0.0002	<0.00257	<0.00112	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00257	<0.00112
Chlorobenzene	0.1	0.1	0.00056 J	<0.0003	<0.0003	<0.00239	<0.0015	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00239	<0.0015
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.00203	<0.00142	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00203	<0.00142
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.00195	<0.00122	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00195	<0.00122
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.00274	<0.00138	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00274	<0.00138
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.00581	<0.00302	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00581	<0.00302
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.0001	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00043	<0.0003	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00029	<0.00032
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.00029	<0.0002	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019	<0.00021
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.00029	<0.0002	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.00021
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.00057	<0.0004	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.00042
2-Methylnaphthalene	0.098	0.29	<0.00009	0.000059 J	<0.000042	<0.00057	<0.0004	<0.000019	<0.000019	<0.000019	0.00034	<0.000019	<0.00038	<0.00042
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00029	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00021
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00036	<0.00025	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00026
Acenaphthene	1.5	4.4	0.042	0.028	<0.00017	<0.00043	<0.0003	<0.000027	<0.000027	<0.000027	0.00011	<0.000027	<0.00029	<0.00032
Acenaphthylene	1.5	4.4	0.00027	0.00013	<0.000015	<0.00043	<0.0003	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029	<0.00032
Anthracene	7.3	22	0.00087	0.00047	<0.000087	0.00087	<0.0002	<0.000014	<0.000014	<0.000014	0.000036 J	<0.000014	<0.00019	<0.00021
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00029	<0.0002	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00019	<0.00021
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00029	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	0.000052 J	<0.00019	<0.00021
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00057	<0.0004	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.00042
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000019	<0.000037	0.00044 J	0.00026 J	0.0002	<0.000037	0.0005	0.0002	0.000062 J	0.00039 J	0.00029 J
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.00029	<0.0002	0.00005 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019	<0.00021
Dibenzofuran	0.098	0.29	0.0005	0.00021	<0.000044	0.00044 J	<0.0003	<0.00002	<0.00002	<0.00002	0.00014	<0.00002	<0.00029	<0.00032
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00029	<0.0002	<0.00002	<0.00002	<0.00002	0.000034 J	<0.00002	<0.00019	<0.00021
Fluoranthene	0.98	2.9	0.0044	0.0029	0.00073	0.000756	<0.0002	0.000058 J	<0.00001	<0.00001	0.000024 J	<0.00001	<0.00019	<0.00021
Fluorene	0.98	2.9	0.0039	0.0017	<0.00029	<0.00029	<0.0002	<0.00003	<0.00003	<0.00003	0.000074 J	<0.00003	<0.00019	<0.00021
Naphthalene	0.49	1.5	<0.00049	<0.00032	<0.00017	0.0018	<0.0004	<0.00002	<0.00002	<0.00002	0.0026	0.000053 J	<0.00038	<0.00042
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00057	<0.0004	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.00042
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00036	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00026
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.00029	<0.0002	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019	<0.00021
Phenanthrene	0.73	2.2	0.00012	0.000083 J	<0.00003	0.000746	<0.0002	<0.000021	<0.000021	<0.000021	0.00013	<0.000021	<0.00019	<0.00021
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.00029	<0.0002	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	<0.00021
Pyrene	0.73	2.2	0.0025	0.0015	0.00035	0.00062	<0.0002	0.000078 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.00019	<0.00021
<b>Metals</b>														
Arsenic	0.01	0.01	0.166	0.0933	0.0217			0.000978 J	<0.0004	0.00207	0.000498 J	0.000859 J		

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-27C 02/03/2009	MW-27C 01/14/2010	MW-27C 06/30/2010	MW-27C 01/27/2011	MW-27C 07/20/2011	MW-27C 02/09/2012	MW-27C 07/25/2012	MW-27C 02/12/2013	MW-27C 08/08/2013	MW-27C 01/24/2014	MW-27C 07/25/2014	MW-27C 01/31/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0002
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000108	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000295	<0.000292	<0.000297	<0.000304	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.000127	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000755	<0.0000766	<0.0000784	0.00052
2-Chloronaphthalene	2	5.8	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000766	<0.0000784	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000777	<0.000066	<0.000067	<0.0000686	0.00041
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.000783	<0.000794	<0.000814	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.000528	<0.000536	<0.000549	0.00031 J
Acenaphthene	1.5	4.4	0.00026	0.00015 J	0.00028	0.00019 J	0.00011 J	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000766	<0.0000784	0.0012
Acenaphthylene	1.5	4.4	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000566	<0.0000574	<0.0000588	<0.00015
Anthracene	7.3	22	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000476	<0.0000472	0.000431 J	<0.000049	0.000065 J
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000766	<0.0000784	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000766	<0.0000784	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.000127	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00038	<0.0016	<0.0015	0.00047	<0.00095	0.00014 J	<0.00021	0.000652	<0.000349	<0.000354	<0.000363	<0.000073
Chrysene	0.91	2	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000766	<0.0000784	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000766	<0.0000784	0.00038
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.000055	<0.000105	0.000143 J	<0.000851	<0.000108	<0.00002
Fluoranthene	0.98	2.9	<0.00007	0.00015 J	<0.00007	<0.00007	0.00011 J	<0.00005	<0.00005	<0.0000667	0.000114 J	0.0000881 J	<0.0000686	0.000085 J
Fluorene	0.98	2.9	<0.00007	<0.00007	0.00025	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.000066	<0.000067	<0.0000686	0.00085
Naphthalene	0.49	1.5	0.00037	<0.00013	0.00024	0.00015 J	<0.00005	<0.00005	0.00019 J	<0.0000762	0.000353 J	<0.000953	<0.0000784	<0.00043
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000108	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000943	<0.0000957	<0.000098	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.000575	<0.000584	<0.000598	<0.000079
Phenanthrene	0.73	2.2	<0.00007	0.00014 J	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	0.0000908 J	<0.000406	<0.0000588	0.00027
Phenol	7.3	22	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000381	<0.0000377	<0.0000383	<0.0000392	<0.000035
Pyrene	0.73	2.2	<0.00007	0.0001 J	<0.00007	<0.00007	0.000064 J	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000108	0.000044 J
<b>Metals</b>														
Arsenic	0.01	0.01												0.00261

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-27C 03/26/2018	MW-27C 06/01/2018	MW-27C 01/22/2019	MW-27C 07/18/2019	MW-27C 01/15/2020	MW-28A 01/29/2008	MW-28A 07/14/2008	MW-28A 02/03/2009	MW-28A 01/13/2010	MW-28A 06/30/2010	MW-28A 01/25/2011	MW-28A 07/19/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.0003	<0.00032	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.0002	<0.00021	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	0.000057	<0.000042	<0.000042	<0.000042	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0004	<0.00042	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.0004	<0.00042	<0.00007	<0.00007	<0.00007	0.00064	<0.00005
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00025	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.0003	<0.00032	<0.00009	<0.00009	<0.00009	0.0002	<0.00005
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.0003	<0.00032	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.000611	<0.00021	<0.00007	<0.00007	<0.00007	0.00036	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.0002	<0.00021	0.00013 J	<0.00007	<0.00007	<0.00007	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<b>&lt;0.00021</b>	0.00011 J	<0.00008	<0.00008	<0.00008	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.0004	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000057 J	0.000042	<0.000037	0.000092 J	<0.000037	0.0003 J	<0.00021	0.0037	<0.00022	<0.00019	<0.0002	<0.00032
Chrysene	0.91	2	<0.000021	0.000054 J	<0.000021	<0.000021	<0.000021	<0.0002	<0.00021	0.00013 J	<0.00007	<0.00007	<0.00007	<0.00005
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00032 J	<0.00032	<0.00008	<0.00008	<0.00008	0.0005	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00021	0.00016 J	<0.00007	<0.00007	<0.00007	<0.00005
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.0002	<0.00021	0.00012 J	<0.00007	<0.00007	0.00021	<0.00005
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	0.0003	<0.00005
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.00002	0.000066 J	0.000055 J	0.000699	<0.00042	0.00017 J	<0.0001	<0.0001	0.0023	<0.00005
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.0004	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00025	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005
Pentachlorophenol	0.001	0.001	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.0002	<0.00021	<0.00008	<0.00008	0.00032	<0.00008	<0.00005
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00046 J	<0.00021	<0.00007	<0.00007	<0.00007	0.00097	<0.00005
Phenol	7.3	22	<0.000035	0.000053 J	<0.000035	0.00023	<0.000035	<0.0002	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.00041 J	<0.00021	0.00016 J	<0.00007	<0.00007	0.00013 J	<0.00005
<b>Metals</b>														
Arsenic	0.01	0.01	<0.0004	0.00212	0.000786 J	0.000428 J	0.000623 J							

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-28A 02/16/2012	MW-28A 07/17/2012	MW-28A 02/07/2013	MW-28A 08/07/2013	MW-28A 01/22/2014	MW-28A 07/25/2014	MW-28A 01/25/2018	MW-28A 03/21/2018	MW-28A 05/17/2018	MW-28A 01/14/2019	MW-28A 07/16/2019	MW-28A 01/16/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002						<0.00011						
Xylenes (total)	10	10	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00105	<0.000104	<0.000105	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000034 J
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.00005	<0.00295	<0.000292	<0.000295	<0.000301	<0.00004	<0.00004	0.012	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00124	<0.000123	<0.000124	<0.000126	<0.000059	<0.000059	<0.00058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.000762	<0.0000755	<0.0000762	<0.0000777	<0.000042	<0.000042	<0.00042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.000762	<0.0000755	<0.0000762	<0.0000777	<0.000021	<0.000021	<0.00021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000066	0.00015 J	<0.000667	<0.0000667	<0.0000667	<0.000068	<0.000019	<0.000019	0.001	<0.000055	0.0002	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.0079	<0.000783	<0.00079	<0.000806	<0.00002	<0.00002	<0.0002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00533	<0.000528	<0.000533	<0.000544	<0.000047	<0.000047	<0.00047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.00098	<0.00005	<0.000762	0.000368 J	<0.0000762	<0.0000777	<0.000027	<0.000027	0.00031 J	0.000092 J	0.000082 J	<0.00014
Acenaphthylene	1.5	4.4	<0.00005	<0.00005	<0.000571	<0.0000566	<0.0000571	<0.0000583	<0.000015	<0.000015	<0.00015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	<0.00005	<0.00005	<0.000476	<0.0000472	<0.0000476	<0.0000485	0.000016 J	0.000021 J	<0.00014	<0.000014	0.000036 J	<0.00013
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.000762	<0.0000755	<0.0000762	<0.0000777	<0.000051	<0.000051	<0.0005	<0.00005	0.000055 J	0.000073 J
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.000762	<0.0000755	<0.0000762	<0.0000777	0.000045 J	<0.00002	<0.0002	<0.00002	0.000095 J	0.000057 J
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00124	<0.000123	<0.000124	<0.000126	<0.00003	<0.00003	<0.0003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0001	<0.00043	<0.00352	<0.000349	<0.000352	<0.000359	0.00018 J	0.000062 J	0.00084 J	0.000061 J	<0.00023	0.00009 J
Chrysene	0.91	2	<0.00005	<0.00005	<0.000762	<0.0000755	<0.0000762	<0.0000777	0.000053 J	<0.000021	<0.00021	<0.000021	0.0001	0.000065 J
Dibenzofuran	0.098	0.29	<0.000051	<0.00005	<0.000762	<0.0000755	<0.0000762	<0.0000777	<0.00002	<0.00002	0.00031 J	<0.00011	0.000084 J	<0.0001
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.00105	<0.000104	<0.000105	<0.000107	<0.00003 J	<0.00002	<0.0002	<0.00002	<0.00002	0.000032 J
Fluoranthene	0.98	2.9	<0.00005	<0.00005	<0.000667	<0.000066	<0.0000667	<0.000068	0.000062 J	0.000055 J	<0.0001	<0.00001	0.00012	0.0003
Fluorene	0.98	2.9	<0.00005	<0.00005	<0.000667	<0.000066	<0.0000667	<0.000068	<0.00003	<0.00003	<0.0003	0.000056 J	0.000053 J	<0.00014
Naphthalene	0.49	1.5	<0.00093	0.0013	<0.000762	<0.000173	<0.0004	<0.0000777	<0.00002	<0.00002	0.024	<0.00024	0.0013	<0.0006
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00105	<0.000104	<0.000105	<0.000107	<0.000024	<0.000024	<0.00024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.000952	<0.0000943	<0.0000952	<0.0000971	<0.000025	<0.000025	<0.00025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00581	<0.000575	<0.000581	<0.000592	<0.00008	<0.00008	<0.00079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.00005	<0.00005	<0.000571	<0.0000566	<0.0000571	<0.0000583	0.000032 J	0.00003 J	0.00022 J	<0.000021	0.000049 J	0.00048
Phenol	7.3	22	<0.00005	<0.00005	<0.000381	<0.0000377	<0.0000381	<0.0000388	<0.000035	<0.000035	0.018	<0.000035	<0.000073	<0.000035
Pyrene	0.73	2.2	0.00011 J	<0.00005	<0.00105	0.000246 J	0.000105 J	<0.000107	0.000081 J	0.000079 J	<0.00019	0.000065 J	0.00016	0.00021
<b>Metals</b>														
Arsenic	0.01	0.01							0.0076	0.0053	0.0177	0.0116	0.0092	0.00664

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-28C 01/29/2008	MW-28C 07/14/2008	MW-28C 02/03/2009	MW-28C 01/13/2010	MW-28C 06/30/2010	MW-28C 01/25/2011	MW-28C 07/19/2011	MW-28C 02/16/2012	MW-28C 07/17/2012	MW-28C 02/07/2013	MW-28C 08/07/2013	MW-28C 01/22/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002
Benzene	0.005	0.005	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002
Chlorobenzene	0.1	0.1	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018
Ethylbenzene	0.7	0.7	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019
Methylene chloride	0.005	0.005	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022
Toluene	1	1	<0.00041	<0.00041	0.0026 J	0.0013 J	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.00009	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000101	<0.000104	<0.00011
2,4-Dimethylphenol	0.49	1.5	<0.00029	0.00114 J	<0.00008	0.0002	<0.00008	0.000086 J	<0.00005	<0.00005	<0.00005	<0.000284	<0.000292	<0.00031
2,4-Dinitrotoluene	0.0013	0.003	<0.00019	<0.00022	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000119	<0.000123	<0.00013
2,6-Dinitrotoluene	0.0013	0.003	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000734	<0.0000755	<0.00008
2-Chloronaphthalene	2	5.8	<0.00038	<0.00044	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000734	<0.0000755	<0.00008
2-Methylnaphthalene	0.098	0.29	<0.00038	<0.00044	0.000097 J	0.00024	0.000077 J	0.000079 J	<0.00005	<0.00005	0.00011 J	<0.0000642	<0.000066	0.0000741 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00019	<0.00022	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000761	<0.000783	<0.00083
4-Nitrophenol	0.049	0.15	<0.00024	<0.00028	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000514	<0.000528	<0.00056
Acenaphthene	1.5	4.4	<0.00029	<0.00033	<0.00009	0.00018 J	0.00033	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000734	<0.0000755	<0.00008
Acenaphthylene	1.5	4.4	<0.00029	<0.00033	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000055	<0.0000566	<0.00006
Anthracene	7.3	22	0.00059	<0.00022	<0.00007	<0.00007	0.00014 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000459	<0.0000472	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000734	<0.0000755	<0.00008
Benzo(a)pyrene	0.0002	0.0002	<0.00019	<0.00022	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000734	<0.0000755	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00038	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000119	<0.000123	<0.00013
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00049 J	<0.00022	0.0033	<0.00046	<0.0012	0.00063	<0.00053	<0.00013	<0.0001	<0.000339	<0.000349	<0.00037
Chrysene	0.91	2	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000734	<0.0000755	<0.00008
Dibenzofuran	0.098	0.29	0.00044 J	<0.00033	<0.00008	0.00018 J	<0.00008	<0.00008	0.00019 J	<0.00005	<0.00005	<0.0000734	<0.0000755	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000101	<0.000104	<0.00011
Fluoranthene	0.98	2.9	0.000497	<0.00022	<0.00007	<0.00007	0.00012 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000642	<0.000066	<0.00007
Fluorene	0.98	2.9	0.00022 J	<0.00022	<0.00007	0.00016 J	0.00033	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000642	<0.000066	<0.00007
Naphthalene	0.49	1.5	0.00234	0.00196	0.00057	0.0014	0.00035	0.00029	<0.000091	<0.00031	0.00064	0.000163 J	<0.0000755	<0.000596
Nitrobenzene	0.049	0.15	<0.00038	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000101	<0.000104	<0.00011
N-Nitrosodiphenylamine	0.19	0.42	<0.00024	<0.00028	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000917	<0.0000943	<0.0001
Pentachlorophenol	0.001	0.001	0.00056 J	<0.00022	<0.00008	<0.00008	0.00034	<0.00008	<0.00005	<0.00005	<0.00005	<0.00056	<0.000575	<0.00061
Phenanthrene	0.73	2.2	0.000624	<0.00022	0.00013 J	0.00033	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000055	<0.0000566	0.0000739 J
Phenol	7.3	22	0.00865	0.00648	0.00063	0.0027	0.00075	0.0014	0.00054	<0.00005	<0.00005	<0.0000367	<0.0000377	<0.00004
Pyrene	0.73	2.2	0.00041 J	<0.00022	<0.00007	<0.00007	0.00007 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.000101	<0.000104	<0.00011
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-28C 07/25/2014	MW-28C 01/25/2018	MW-28C 03/21/2018	MW-28C 05/17/2018	MW-28C 01/14/2019	MW-28C 07/16/2019	MW-28C 01/16/2020	MW-29A 01/28/2008	MW-29B 01/28/2008	MW-29C 01/28/2008	MW-30A 01/30/2008	MW-30A 07/14/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00245	<0.00245	<0.00245	<0.00052	<0.01
Benzene	0.005	0.005	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00257	<0.00257	<0.00257	0.147	0.14
Chlorobenzene	0.1	0.1	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00239	<0.00239	<0.00239	<0.00047	<0.01
Ethylbenzene	0.7	0.7	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00203	<0.00203	<0.00203	0.153	0.12
Methylene chloride	0.005	0.005	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00195	<0.00195	<0.00195	<0.00054	<0.013
Toluene	1	1	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00274	<0.00274	<0.00274	0.645	0.51
Vinyl chloride	0.002	0.002	<0.00011											
Xylenes (total)	10	10	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00581	<0.00581	<0.00581	0.386	0.32
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0002	0.00008 J	<0.00008	<0.02	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00057	<0.0003	<0.00029	4.1	2.9
2,4-Dinitrotoluene	0.0013	0.003	<0.000126	<0.000058	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058	<0.00038	<0.0002	<0.00019	<0.038	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.0000777	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00038	0.00374	<0.00019	<0.038	<0.00006
2-Chloronaphthalene	2	5.8	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00076	<0.0004	<0.00038	<0.076	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.000068	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00076	<0.0004	<0.00038	0.573	0.85
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00038	<0.0002	<0.00019	<0.038	<0.00008
4-Nitrophenol	0.049	0.15	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00048	<0.00025	<0.00024	<0.048	<0.00005
Acenaphthene	1.5	4.4	<0.0000777	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.00057	<0.0003	<0.00029	0.215	0.31
Acenaphthylene	1.5	4.4	<0.0000583	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00057	<0.0003	<0.00029	<0.057	0.075
Anthracene	7.3	22	<0.0000485	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.00038	<0.0002	<0.00019	<0.038	0.18
Benzo(a)anthracene	0.0091	0.02	<0.0000777	<0.00005	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005	<0.00038	<0.0002	<0.00019	<0.038	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00038	<0.0002	<0.00019	<0.038	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000126	<0.00003	<0.00003	0.00003 J	<0.00003	<0.00003	<0.00003	<0.00076	<0.0004	<0.00038	<0.076	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000359	0.000064 J	<0.000037	<0.000037	<0.000037	<0.000037	<0.00013	0.000072 J	0.00037 J	0.00042 J	<0.038	<0.0001
Chrysene	0.91	2	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.0002	<0.00019	<0.038	<0.00005
Dibenzofuran	0.098	0.29	<0.0000777	<0.00002	<0.00002	0.00002 J	<0.00002	<0.00002	<0.00002	<0.00057	<0.0003	<0.00029	0.156	0.25
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000107	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00038	<0.0002	<0.00019	<0.038	<0.00005
Fluoranthene	0.98	2.9	<0.000068	<0.00001	<0.00001	<0.00001	<0.00001	0.000015 J	<0.00001	<0.00038	<0.0002	<0.00019	<0.038	0.0041
Fluorene	0.98	2.9	<0.000068	<0.00003	<0.00003	0.00003 J	<0.00003	<0.00003	<0.00003	<0.00038	<0.0002	<0.00019	0.109	0.18
Naphthalene	0.49	1.5	<0.0000777	<0.00021	<0.00002	<0.00002	<0.00002	0.000095 J	<0.00016	<0.00076	<0.0004	0.00047 J	9.77	15
Nitrobenzene	0.049	0.15	<0.000107	<0.000024	<0.000024	0.000024 J	<0.000024	<0.000024	<0.000024	<0.00076	<0.0004	<0.00038	<0.076	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00048	<0.00025	<0.00024	<0.048	<0.00005
Pentachlorophenol	0.001	0.001	<0.000592	<0.000079	<0.00008	<0.00008	<0.000079	<0.000079	<0.000079	<0.00038	<0.0002	<0.00019	<0.038	0.00033 J
Phenanthrene	0.73	2.2	<0.0000583	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.0002	<0.00019	0.075	0.12
Phenol	7.3	22	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	0.0025	0.000086 J	<0.00038	0.00287	<0.00019	0.174	0.14
Pyrene	0.73	2.2	<0.000107	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00038	<0.0002	<0.00019	<0.038	0.022
<b>Metals</b>														
Arsenic	0.01	0.01		0.00206	0.00184 J	0.00184 J	0.000447 J	0.000456 J	0.000937 J					

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-30A 02/03/2012	MW-30A 07/12/2012	MW-30A 02/01/2013	MW-31A 01/31/2008	MW-31A 07/14/2011	MW-31A 02/03/2012	MW-31A 07/12/2012	MW-31A 02/01/2013	MW-32A <sup>7</sup> 01/28/2008 DNAPL	MW-32A <sup>7</sup> 07/14/2008 DNAPL	MW-32A <sup>7</sup> 02/03/2009 DNAPL	MW-32A <sup>7</sup> 01/14/2010 DNAPL
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.01	<0.0025	<0.0028	<0.00052	<0.01	<0.025	<0.005	<0.0014	<0.00052	<0.00109	<0.0005	<0.0005
Benzene	0.005	0.005	0.13	0.14	0.117	0.178	0.14	0.14	0.11	0.135	0.884	0.884	0.69	0.34
Chlorobenzene	0.1	0.1	<0.01	<0.0025	<0.0024	<0.00047	<0.01	<0.025	<0.005	<0.0012	<0.00047	<0.0015	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	0.11	0.13	0.119	0.166	0.19	0.17	0.18	0.171	0.373	0.365	0.34	0.076
Methylene chloride	0.005	0.005	<0.013	<0.005	0.0211	<0.00054	<0.013	<0.032	<0.01	0.00971 J	<0.00054	<0.00122	<0.0005	<0.0005
Toluene	1	1	0.49	0.54	0.443	0.337	0.37	0.36	0.31	0.346	0.95	0.983	0.74	0.36
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.32	0.32	0.302	0.562	0.63	0.71	0.63	0.583	1.02	1.03	0.88	0.35
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0005	<0.00005	<0.0524	<0.002	<0.00005	<0.0005	<0.00005	<0.055	<0.02	<0.008	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	3	2.7	2.94	4.74	5.3	5.1	3.4	4.45	9.57	12.6	2.2	2.1
2,4-Dinitrotoluene	0.0013	0.003	<0.0005	<0.00005	<0.0619	<0.0038	<0.00005	<0.0005	<0.00005	<0.065	<0.044	<0.019	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0006	<0.00006	<0.0381	<0.0038	<0.00006	<0.0006	<0.00006	<0.04	<0.044	<0.019	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0005	<0.00005	<0.0381	<0.0076	<0.00005	<0.0005	<0.00005	<0.04	<0.089	<0.039	<0.00012	<0.0001
2-Methylnaphthalene	0.098	0.29	0.42	0.67	1.01	0.887	1	0.65	0.83	1.17	1.13	0.989	1.2	0.3
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0008	<0.00008	<0.395	<0.0038	<0.00008	<0.0008	<0.00008	<0.415	<0.044	<0.019	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.0005	<0.00005	<0.267	<0.0048	<0.00005	<0.0005	<0.00005	<0.28	<0.056	<0.024	<0.00007	<0.00007
Acenaphthene	1.5	4.4	0.23	0.24	0.436	0.206	0.37	0.3	0.28	0.488	0.341	0.294	0.34	0.13
Acenaphthylene	1.5	4.4	0.0064	0.0074	<0.0286	<0.0057	0.0076	0.0044	0.0032	<0.03	<0.067	<0.029	0.006	0.0019
Anthracene	7.3	22	0.013	0.018	0.0391 J	0.0145	0.032	0.026	0.026	0.056 J	<0.044	0.043	0.077	0.051
Benzo(a)anthracene	0.0091	0.02	<0.0005	<0.00005	<0.0381	<0.0038	0.0038	<0.0005	0.00023	<0.04	<0.044	<0.019	0.0096	0.0067
Benzo(a)pyrene	0.0002	0.0002	<0.0005	<0.00005	<0.0381	<0.0038	0.00089	<0.0005	<0.00005	<0.04	<0.044	<0.019	0.003	0.0023
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0005	<0.00005	<0.0619	<0.0076	<0.00005	<0.0005	<0.00005	<0.065	<0.089	<0.039	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.001	<0.0001	<0.176	<0.0038	<0.0001	<0.001	<0.0001	<0.185	<0.044	<0.019	0.00042	<0.0018
Chrysene	0.91	2	<0.0005	<0.00005	<0.0381	<0.0038	0.0031	<0.0005	0.00017 J	<0.04	<0.044	<0.019	0.0087	0.0064
Dibenzofuran	0.098	0.29	0.21	0.2	0.308	0.169	0.33	0.26	0.26	0.367	0.298	0.26	0.32	0.14
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0005	0.0002	<0.0524	<0.0038	<0.00005	<0.0005	<0.00005	<0.055	<0.044	<0.019	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.0031	0.0038	<0.0333	0.006	0.031	0.0029	0.0052	<0.035	<0.044	0.026	0.098	0.07
Fluorene	0.98	2.9	0.13	0.14	0.247	0.123	0.24	0.17	0.17	0.273	0.163	0.156	0.22	0.087
Naphthalene	0.49	1.5	7.8	12	16.8 J	13.7	21	18	17	19.3 J	25	16.2	16	3.5
Nitrobenzene	0.049	0.15	<0.0005	<0.00005	<0.0524	<0.0076	<0.00005	<0.0005	<0.00005	<0.055	<0.089	<0.039	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0005	<0.00005	<0.0476	<0.0048	<0.00005	<0.0005	<0.00005	<0.05	<0.056	<0.024	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.0005	<0.00005	<0.29	0.0895	0.076 J	0.11	0.094	<0.305	<0.044	<0.019	<0.00008	<0.00008
Phenanthrene	0.73	2.2	0.064	0.1	0.162 J	0.0774	0.24	0.14	0.13	0.268	0.177	0.185	0.45	0.25
Phenol	7.3	22	0.015	0.02	0.0781 J	1.56	0.6	0.76	0.29	0.579	9.01	8.83	1.4	1.3
Pyrene	0.73	2.2	0.0021	0.0018	<0.0524	0.0084	0.018	0.0025	0.002	<0.055	<0.044	<0.019	0.062	0.043
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-32A <sup>7</sup> 07/01/2010 DNAPL	MW-32A <sup>7</sup> 01/25/2011 DNAPL	MW-32A <sup>7</sup> 07/19/2011 DNAPL	MW-32AR 02/09/2012	MW-32AR 07/16/2012	MW-32AR 02/06/2013	MW-32AR 08/07/2013	MW-32AR 01/21/2014	MW-32AR 07/24/2014	MW-32AR 01/28/2018	MW-32AR 03/27/2018	MW-32AR 05/31/2018
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0025	<0.0025	<b>0.03</b>	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<b>1.5</b>	<b>0.61</b>	<b>1.4</b>	<0.001	<0.0005	<b>0.023</b>	0.000475 J	<0.0002	<b>0.0404</b>	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0025	<0.0025	<0.005	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.45	0.41	0.31	<0.0011	<0.0005	0.0082	0.000296 J	<0.00019	0.0208	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0032	<0.0025	<b>&lt;0.0065</b>	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001
Toluene	1	1	<b>1.5</b>	1	<b>1.2</b>	<0.001	<0.0005	0.00338	0.000234 J	<0.00017	0.000849 J	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002							<0.00011		<0.00011			
Xylenes (total)	10	10	1.3	1.1	0.87	<0.0031	<0.0015	0.0176	0.000873 J	<0.00058	0.0336	<0.0003	<0.0003	<0.0003
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.001	<0.001	<0.0005	<0.00005	<0.00005	<b>&lt;0.00529</b>	<0.000104	<0.000104	<0.000104	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<b>15</b>	<b>2.3</b>	<b>31</b>	0.0012	0.000061 J	0.0172 J	<0.000292	<0.000292	0.0722	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.0009	<0.0009	<0.0005	<0.00005	<0.00005	<b>&lt;0.00625</b>	<0.000123	<0.000123	<0.000123	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.0007	<0.0007	<0.0006	<0.00006	<0.00006	<b>&lt;0.00385</b>	<0.0000755	<0.0000755	<0.0000755	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.001	<0.001	<0.0005	<0.00005	<0.00005	<0.00385	<0.0000755	<0.0000755	<0.0000755	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<b>0.48</b>	<b>6.9</b>	<b>0.95</b>	0.000088 J	0.00019 J	0.011 J	<0.000066	0.00021 J	<b>0.206</b>	0.000079 J	0.0068	<0.00019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0008	<0.0008	<0.0008	<0.00008	<0.00008	<b>&lt;0.0399</b>	<0.000783	<0.000783	<0.000783	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.0007	<0.0007	<0.0005	<0.00005	<0.00005	<0.0269	<0.000528	<0.000528	<0.000528	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.19	<b>3.1</b>	0.25	0.0046	0.0029	0.0232 J	0.00625	0.0105	0.111	0.0013	0.0043	0.002
Acenaphthylene	1.5	4.4	0.0079	0.027	0.005	0.00032	<0.00005	<0.00288	0.000258 J	<0.000317	0.00226	0.000054 J	0.000054 J	<0.00015
Anthracene	7.3	22	0.093	0.91	0.09	0.00041	0.00026	<0.0024	0.000223 J	0.000444 J	0.00332	0.000082 J	0.00021	0.000064 J
Benzo(a)anthracene	0.0091	0.02	<b>0.01</b>	<b>0.3</b>	<b>0.038</b>	0.00012 J	0.00015 J	<0.00385	<0.0000755	<0.0000755	0.000218 J	0.00006 J	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<b>0.0067</b>	<b>0.078</b>	<b>0.019</b>	<0.00005	<0.00005	<b>&lt;0.00385</b>	<0.0000755	<0.0000755	<0.0000755	0.000027 J	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<b>&lt;0.0009</b>	<b>&lt;0.0009</b>	<0.0005	<0.00005	<0.00005	<b>&lt;0.00625</b>	<0.000123	<0.000123	0.000452 J	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0041	0.0046	<0.0013	0.00054	0.00014 J	<b>&lt;0.0178</b>	<0.000349	<0.000349	0.000621	0.00011 J	<0.000037	0.00014 J
Chrysene	0.91	2	0.0099	0.28	0.033	0.000093 J	0.00023	<0.00385	<0.0000755	0.0000774 J	0.00016 J	0.000058 J	<0.000021	0.000037 J
Dibenzofuran	0.098	0.29	<b>0.21</b>	<b>3.2</b>	<b>0.27</b>	0.00096	0.0017	0.00936 J	0.000515	0.000664	0.05	<0.000053	0.0024	0.00007 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0007	<0.0007	<0.0005	<0.00005	0.000056 J	<0.00529	<0.000104	<0.000104	<0.000104	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.09	<b>2.5</b>	0.11	0.002	0.003	0.00508 J	0.00102	0.00124	0.00656	0.00048	0.00023	0.00031
Fluorene	0.98	2.9	0.13	<b>2.5</b>	0.18	0.0017	0.0016	0.00932 J	0.00105	0.00317	0.0516	0.00012	0.00093	0.000058 J
Naphthalene	0.49	1.5	<b>11</b>	<b>31</b>	<b>21</b>	0.00044	0.0036	0.406	<0.00297	0.0087	<b>3.83</b>	<0.00058	0.026	0.000043 J
Nitrobenzene	0.049	0.15	<0.0009	<0.0009	<0.0005	<0.00005	<0.00005	<0.00529	<0.000104	<0.000104	<0.000104	0.00035	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	0.014	<0.0009	<0.0005	<0.00005	<0.00005	<0.00481	<0.0000943	<0.0000943	<0.0000943	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.0008	<0.0008	<0.0005	<0.00005	<0.00005	<b>&lt;0.0293</b>	<0.000575	<0.000575	<0.000575	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.19	<b>8.9</b>	0.35	0.00033	0.000055 J	0.00768 J	0.000112 J	0.000101 J	0.0111	0.00022	0.0013	<0.000021
Phenol	7.3	22	<b>14</b>	<b>1.2</b>	<b>21</b>	0.00012 J	0.00029	<0.00192	<0.0000377	<0.0000377	<0.0000377	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.047	<b>1.5</b>	0.072	0.0041	0.0046	<0.00529	0.000617	0.000625	0.00474	0.00031	0.00018	0.00049
<b>Metals</b>														
Arsenic	0.01	0.01										0.00294	<b>0.0228</b>	<b>0.0441</b>

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-32AR 01/23/2019	MW-32AR 07/30/2019	MW-32AR 01/20/2020	MW-33A 01/29/2008	MW-33A 07/14/2008	MW-33A 02/03/2009	MW-33A 02/03/2009 Duplicate	MW-33A 01/13/2010	MW-33A 01/13/2010 Duplicate	MW-33A 06/29/2010	MW-33A 06/29/2010 Duplicate	MW-33A 01/24/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00052	<0.00109	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00025	<b>0.0062</b>	0.00071 J	0.00074 J	0.0025 J	0.0024 J	0.0018 J	0.0034 J	<b>0.0056</b>
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.00047	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.00025	<0.00142	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.0015 J
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.00054	<0.00122	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.00041	<0.00138	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002			<0.0002									
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.00127	<0.00302	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	0.0016 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.0001	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	0.00013 J	<0.00004	<0.00004	<0.0005	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	0.0027
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.00033	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.00033	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.00067	<0.00038	<0.00012	<0.00012	<0.0001	<0.0001	<0.0001	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000028	<0.00067	<0.00038	0.00066	<0.00007	0.0009 J	0.00051 J	0.00092	<0.00007	0.0067
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00033	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00042	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	0.0001	<0.000027	0.0014	0.0133	0.0124	0.013	<0.00009	0.028	0.026	0.012	0.0072	0.014
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.0005	<0.00029	<0.00006	<0.00006	0.00015 J	0.00014 J	<0.00007	<0.00007	0.00014 J
Anthracene	7.3	22	0.00002 J	<0.000014	<0.000014	<0.00033	0.00024 J	0.0002 J	<0.00007	0.00028	0.00024	0.00021	<0.00007	0.00072
Benzo(a)anthracene	0.0091	0.02	<0.00005	0.000054 J	<0.00005	<0.00033	<0.00019	0.0002 J	<0.00007	0.00017 J	0.00017 J	0.00014 J	0.00014 J	0.00025
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	0.000081 J	<b>&lt;0.00033</b>	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00067	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000044	0.000074 J	0.000043 J	0.00137 J	<0.00019	0.00033	0.00035	<0.0003	<0.00038	<0.00035	<0.00035	0.00031
Chrysene	0.91	2	<0.000021	<0.000021	0.000042 J	<0.00033	<0.00019	0.00012 J	<0.00007	0.00012 J	0.000089 J	0.00009 J	<0.00007	0.00016 J
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	0.00074	0.000628	0.00078	<0.00008	0.0019	0.0017	0.0014	0.00035	0.0027
Di-n-butylphthalate (DBP)	2.4	7.3	0.00002 J	<0.00002	<0.00002	0.00055 J	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.000051 J	<0.00001	<0.000029	0.00141	0.00154	0.0022	<0.00007	0.0013	0.0012	0.0012	0.00095	0.003
Fluorene	0.98	2.9	<0.00003	<0.00003	0.000038 J	0.0013	0.000939	0.00067	<0.00007	0.0015	0.0013	0.0012	0.00041	0.0027
Naphthalene	0.49	1.5	0.000067 J	<0.00002	<0.00014	0.00167	0.0047	0.0028	<0.0001	0.02 J	0.009 J	0.0082	0.0013	0.095
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00067	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00042	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.00033	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	0.000034 J	<0.000021	<0.000032	<0.00033	0.00028 J	0.00037	<0.00007	0.00032	0.00024	0.00065	<0.00007	0.0048
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.00033	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	0.000036 J	0.00045	0.00042	0.0019	0.00167	0.0024	<0.00007	0.0019	0.0018	0.0016	0.0015	0.0035
<b>Metals</b>														
Arsenic	0.01	0.01	<b>0.0316</b>	<b>0.0628</b>	0.0077									

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-33A 07/19/2011	MW-33A 07/19/2011 Duplicate	MW-33A 02/15/2012	MW-33A 02/15/2012 Duplicate	MW-33A 07/17/2012	MW-33A 07/17/2012 Duplicate	MW-33A 02/12/2013	MW-33A 02/12/2013 Duplicate	MW-33A 08/07/2013	MW-33A 08/07/2013 Duplicate	MW-33A 01/23/2014	MW-33A 01/23/2014 Duplicate
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.001	<0.005	<0.0005	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002
Benzene	0.005	0.005	<b>0.009</b>	<b>0.0091</b>	<b>0.054</b>	<b>0.051</b>	0.0023 J	0.0022 J	<b>0.00782</b>	<b>0.00782</b>	<b>0.165</b>	<b>0.174</b>	<b>0.223</b>	<b>0.223</b>
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.001	<0.005	<0.0005	<0.0005	<0.00012	<0.00012	0.000185 J	0.000189 J	0.000214 J	<0.00018
Ethylbenzene	0.7	0.7	0.0033 J	0.003 J	0.075	0.061	<0.0005	<0.0005	0.0022	0.00216	0.109	0.117	0.135	0.134
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<0.0013	<b>&lt;0.0065</b>	<0.001	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.00022	<0.00022
Toluene	1	1	<0.001	<0.001	0.019	0.018 J	<0.0005	<0.0005	<0.00015	<0.00015	0.00639	0.00672	0.23	0.229
Vinyl chloride	0.002	0.002					<0.0005	<0.0005			<0.00011	<0.00011		
Xylenes (total)	10	10	<0.0031	<0.0031	0.11	0.092	<0.0015	<0.0015	0.00223 J	0.00229 J	0.172	0.186	0.188	0.19
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000105	<0.000105	<b>&lt;0.00519</b>	<b>&lt;0.00519</b>	<b>&lt;0.104</b>	<b>&lt;0.106</b>
2,4-Dimethylphenol	0.49	1.5	0.0034	0.0046	0.029 J	0.019 J	<0.00005	<0.00005	0.00623	0.0199	<b>0.888</b>	<b>0.801</b>	<b>1.44</b>	<b>1.54</b>
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000124	<0.000124	<b>&lt;0.00613</b>	<b>&lt;0.00613</b>	<b>&lt;0.123</b>	<b>&lt;0.125</b>
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000762	<b>&lt;0.00377</b>	<b>&lt;0.00377</b>	<b>&lt;0.0755</b>	<b>&lt;0.0769</b>
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00377	<0.00377	<0.0755	<0.0769
2-Methylnaphthalene	0.098	0.29	0.022 J	0.031 J	0.03	0.024	0.0015	0.0011	0.00345	0.0149	<b>0.195</b>	<b>0.16</b>	<b>0.263 J</b>	<b>0.27 J</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.00079	<b>&lt;0.0392</b>	<b>&lt;0.0392</b>	<b>&lt;0.783</b>	<b>&lt;0.798</b>
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000533	<0.000533	<0.0264	<0.0264	<b>&lt;0.528</b>	<b>&lt;0.538</b>
Acenaphthene	1.5	4.4	0.037	0.042	0.048	0.064	0.019	0.023	0.0279	0.0374 J	0.157	0.151	0.288 J	0.217 J
Acenaphthylene	1.5	4.4	0.00018 J	0.00021	0.0003	0.0004	<0.00005	<0.00005	<0.0000571	0.000358 J	<0.00283	<0.00283	<0.0566	<0.0577
Anthracene	7.3	22	0.0013	0.0016	0.00082	0.00078	0.0026 J	0.0015 J	0.000748	0.000801	0.0049 J	0.00487 J	<0.0472	<0.0481
Benzo(a)anthracene	0.0091	0.02	0.00019 J	0.00018 J	0.0001 J	0.00011 J	0.00043 J	0.00026 J	0.000174 J	0.000211 J	<0.00377	<0.00377	<b>&lt;0.0755</b>	<b>&lt;0.0769</b>
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.00005	<0.00005	0.00011 J	0.000054 J	<0.0000762	<0.0000762	<b>&lt;0.00377</b>	<b>&lt;0.00377</b>	<b>&lt;0.0755</b>	<b>&lt;0.0769</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000124	<0.000124	<b>&lt;0.00613</b>	<b>&lt;0.00613</b>	<b>&lt;0.123</b>	<b>&lt;0.125</b>
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00056	<0.00068	0.0013 J	0.00084 J	0.00014 J	0.00022	<0.000352	<0.000352	<b>&lt;0.0175</b>	<b>&lt;0.0175</b>	<b>&lt;0.349</b>	<b>&lt;0.356</b>
Chrysene	0.91	2	0.0001 J	0.00005 J	0.000061 J	<0.00005	0.00033 J	0.00016 J	0.000114 J	0.000106 J	<0.00377	<0.00377	<0.0755	<0.0769
Dibenzofuran	0.098	0.29	0.0088 J	0.014 J	0.019 J	0.026 J	0.0049	0.0061	0.00262	0.00699	0.0728	0.0717	<b>0.148 J</b>	<b>0.108 J</b>
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000105	<0.000105	<0.00519	<0.00519	<0.104	<0.106
Fluoranthene	0.98	2.9	0.0021	0.0023	0.0012	0.0011	0.0036 J	0.0026 J	0.00212	0.0025	0.00385 J	0.00435 J	0.0703 J	<0.0673
Fluorene	0.98	2.9	0.0088 J	0.012 J	0.015	0.019	0.0056	0.0073	0.00641	0.00994	0.0668	0.0623	0.145 J	0.0939 J
Naphthalene	0.49	1.5	0.31 J	0.44 J	<b>0.96 J</b>	<b>1.6 J</b>	0.017	0.021	0.112	0.382 J	<b>4.98</b>	<b>4.07</b>	<b>5.82</b>	<b>5.3</b>
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000105	<0.000105	<0.00519	<0.00519	<b>&lt;0.104</b>	<b>&lt;0.106</b>
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000952	<0.00472	<0.00472	<0.0943	<0.0962
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000581	<0.000581	<b>&lt;0.0288</b>	<b>&lt;0.0288</b>	<b>&lt;0.575</b>	<b>&lt;0.587</b>
Phenanthrene	0.73	2.2	0.0046 J	0.0068 J	0.0038	0.0032	0.0058 J	0.0026 J	0.000488	0.000987	0.0232 J	0.0236	0.18 J	0.135 J
Phenol	7.3	22	0.00005 J	0.00091 J	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000381	<0.0000381	<0.00189	<0.00189	0.203 J	0.108 J
Pyrene	0.73	2.2	0.0025	0.0026	0.0021	0.0022	0.0035	0.0032	0.00283	0.00336	<0.00519	<0.00519	<0.104	<0.106
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-33A 08/28/2014	MW-33A 08/28/2014 Duplicate	MW-33A 01/30/2018	MW-33A 01/30/2018	MW-33A 03/27/2018	MW-33A 03/27/2018 Duplicate	MW-33A 06/05/2018	MW-33A 06/05/2018 Duplicate	MW-33A 01/22/2019	MW-33A 01/22/2019 Duplicate	MW-33A 07/17/2019	MW-33A 07/17/2019 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	0.00236 J	0.00389 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.0014 J	0.00224 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002			<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	0.000632 J	0.000872 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000108	<0.000108	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.0212 J	0.0608 J	<0.0002	<0.00021	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000127	<0.000127	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.0000784	<0.0000784	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0000784	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.0216 J	0.0524 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000814	<0.000814	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000549	<0.000549	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.046 J	0.0692 J	0.0005	0.00016	<0.000027	<0.000027	0.00018 J	0.00097 J	<0.000027	<0.000027	0.000098 J	<0.000027
Acenaphthylene	1.5	4.4	0.00041 J	0.000636 J	0.000022 J	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.00132 J	0.00184 J	<0.000014	<0.000014	<0.000014	<0.000014	0.000021 J	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	0.000288 J	0.000318 J	0.000077 J	<0.00005	<0.00005	<0.00005	0.000077 J	0.000091 J	<0.00005	0.000062 J	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.0000784	<0.0000784	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000127	0.000476 J	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000363 J	0.000363 J	<0.000037	<0.000044	<0.000037	<0.000037	0.000078 J	0.00014 J	<0.000037	<0.000037	<0.000022	<0.000084
Chrysene	0.91	2	0.000168 J	0.000178 J	<0.000021	<0.000021	0.000047 J	<0.000021	0.00005 J	0.000061 J	0.000025 J	0.000034 J	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.0164 J	0.0334 J	0.000042 J	<0.00002	<0.00002	0.000036 J	0.00002 J	0.000038 J	<0.00002	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000108	<0.000108	<0.00002	<0.00002	<0.00002	<0.00002	0.000039 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.00395	0.00426	0.0006	0.00034	0.00052	0.00052	0.00045 J	0.0006 J	0.00033	0.00044	0.00016	0.00017
Fluorene	0.98	2.9	0.0146 J	0.0224 J	<0.00003	<0.00003	<0.00003	<0.00003	0.000059 J	0.00017	<0.00003	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	0.489 J	<b>1.02 J</b>	<0.00033	<0.00035	<0.00002	0.00013	<0.00019	<0.00002	<0.00002	<0.00002	<0.000076	<0.000096
Nitrobenzene	0.049	0.15	0.000108 J	0.00936 J	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000098	<0.000098	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000598	<0.000598	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.00427 J	0.00736 J	<0.000021	<0.000021	<0.000021	<0.000021	0.000085 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.0000392	<0.0000392	<0.0002	<0.00025	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	0.00011 J	0.000053 J
Pyrene	0.73	2.2	0.0043	0.00471	0.00028	0.000041 J	0.00057	0.00062	0.00073	0.001	0.00015	0.00034	0.00028	0.00069
<b>Metals</b>														
Arsenic	0.01	0.01			<b>0.0197</b>	<b>0.0202</b>	<b>0.0186</b>	<b>0.0201</b>	0.00573	0.00584	0.01	0.00995	<b>0.0155</b>	<b>0.0147</b>

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-33A 01/20/2020	MW-33A 01/20/2020 Duplicate	MW-34C 01/29/2008 DNAPL	MW-34C 02/08/2012	MW-34CR 07/29/2014	MW-34CR 01/29/2018	MW-34CR 03/27/2018	MW-34CR 06/05/2018	MW-34CR 01/15/2019	MW-35A 01/29/2008	MW-35A 07/14/2008	MW-35A 02/03/2009
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.00052	0.001 J	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005
Benzene	0.005	0.005	<0.0002	<0.0002	<b>0.0287</b>	0.0014 J	0.000154 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025	<0.0005
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.00047	0.001 J	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	0.0903	0.0039 J	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00025	<0.0005
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.00054	0.0013 J	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005
Toluene	1	1	<0.0002	<0.0002	0.0832	0.0041 J	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041	<0.0005
Vinyl chloride	0.002	0.002	<0.0002	<0.0002										
Xylenes (total)	10	10	<0.0003	<0.0003	0.25	0.0077 J	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00127	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<b>&lt;0.002</b>	0.00005 J	<0.000108	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	0.00012 J	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<b>&lt;0.0059</b>	0.00022 J	<0.000304	<0.00004	<0.00004	<0.00004	<0.00004	<0.0003	<0.00028	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<b>&lt;0.004</b>	0.00005 J	<0.000127	<0.000058	<0.000058	<0.000058	<0.000058	<0.0002	<0.00019	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<b>&lt;0.004</b>	0.00006 J	<0.0000784	0.0001 J	<0.000042	<0.000042	<0.000042	<0.0002	<0.00019	<0.00007
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<b>&lt;0.0079</b>	0.00005 J	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021	<0.0004	<0.00038	<0.00012
2-Methylnaphthalene	0.098	0.29	<0.000063	<0.000095	<b>0.437</b>	0.00011 J	0.000255 J	<0.000019	<0.000019	<0.000019	<0.000079	<0.0004	<0.00038	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<b>&lt;0.004</b>	0.00008 J	<0.000814	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00019	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.005	0.00005 J	<0.000549	<0.000047	<0.000047	<0.000047	<0.000047	<0.00025	<0.00024	<0.00007
Acenaphthene	1.5	4.4	0.00073 J	0.0011 J	0.365	0.00005 J	<0.0000784	<0.000027	<0.000027	<0.000027	0.000029 J	0.0176	0.00656	0.0035
Acenaphthylene	1.5	4.4	0.000049 J	0.000038 J	<0.0059	0.00005 J	<0.0000588	<0.000015	<0.000015	<0.000015	<0.000015	<0.0003	<0.00028	<0.00006
Anthracene	7.3	22	0.000057 J	0.000044 J	0.0712	0.00005 J	<0.000049	<0.000014	<0.000014	<0.000014	<0.000014	0.000542	0.00023 J	<0.00007
Benzo(a)anthracene	0.0091	0.02	0.00012	0.00012	<b>0.017</b>	0.00005 J	<0.0000784	<0.00005	<0.00005	<0.00005	<0.00005	<0.0002	<0.00019	<0.00007
Benzo(a)pyrene	0.0002	0.0002	0.000066 J	0.000072 J	<b>0.0065</b>	0.00005 J	<0.0000784	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00019	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<b>&lt;0.0079</b>	0.00005 J	<0.000127	<0.00003	<0.00003	<0.00003	<0.00003	<0.0004	<0.00038	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	<b>0.0076</b>	0.00053 J	<0.000799	<0.000037	<0.000037	0.000081 J	<0.000037	0.00453	<0.00019	0.00024
Chrysene	0.91	2	0.000078 J	0.000093 J	0.0128	0.00005 J	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021	<0.0002	<0.00019	<0.00007
Dibenzofuran	0.098	0.29	0.000096 J	0.000067 J	<b>0.412</b>	0.000071 J	<0.0000784	<0.00002	<0.00002	<0.00002	0.000027 J	0.00819	0.00319	0.0014
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.004	0.00005 J	<0.000108	<0.00002	<0.00002	<0.00002	0.000021 J	0.00046 J	<0.00019	<0.00007
Fluoranthene	0.98	2.9	0.00035 J	0.00057 J	0.14	0.00017 J	<0.0000686	<0.00001	<0.00001	<0.00001	0.000014 J	0.0014	0.00105	0.00034
Fluorene	0.98	2.9	0.000059 J	0.000072 J	0.228	0.00011 J	<0.0000686	<0.00003	<0.00003	<0.00003	<0.00003	0.00328	0.00161	0.00062
Naphthalene	0.49	1.5	<0.0005	<0.00066	<b>5.87</b>	0.00043 J	0.00282	<0.00017	<0.00002	<0.00023	<0.00069	0.0257	0.000704	<0.0001
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.0079	0.00005 J	<0.000108	<0.000024	<0.000024	<0.000024	<0.000024	<0.0004	<0.00038	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.005	0.00005 J	<0.000098	<0.000025	<0.000025	<0.000025	<0.000025	<0.00025	<0.00024	<0.00009
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<b>&lt;0.004</b>	0.00005 J	<0.000598	0.00013 J	<0.000079	<0.000079	<0.000079	<0.0002	<0.00019	<0.00008
Phenanthrene	0.73	2.2	<0.000056	<0.000046	0.431	0.0001 J	<0.0000588	<0.000021	<0.000021	<0.000021	<0.000021	0.00046 J	<0.00019	<0.00007
Phenol	7.3	22	<0.000035	<0.000035	<0.004	0.000072 J	<0.0000392	<0.000035	<0.000035	<0.000035	<0.000035	<0.0002	<0.00019	<0.00007
Pyrene	0.73	2.2	<0.000019	0.00052 J	0.0853	0.00021 J	<0.000108	<0.000019	<0.000019	<0.000019	<0.000019	0.000967	0.000547	0.00025
<b>Metals</b>														
Arsenic	0.01	0.01	0.0081	0.00755				0.00106 J	0.000801 J	0.000689 J	0.00132 J			

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-35A 01/14/2010	MW-35A 06/30/2010	MW-35A 01/27/2011	MW-35A 07/20/2011	MW-35A 02/15/2012	MW-35A 07/18/2012	MW-35A 02/07/2013	MW-35A 08/08/2013	MW-35A 01/24/2014	MW-35A 07/24/2014	MW-35A 01/25/2018	MW-35A 03/22/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	0.000367 J	0.00021 J	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	0.00015 J	<0.0003	0.00042 J
Ethylbenzene	0.7	0.7	0.0015 J	<0.0005	<0.0005	<0.0011	<0.0011	0.0015 J	<0.00011	<0.00011	<0.00019	0.000473 J	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002
Vinyl chloride	0.002	0.002									<0.00011	<0.00011		
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	0.000309 J	<0.00058	<0.00026	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	0.0003	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.000104	<0.000106	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000295	<0.000304	<0.000292	<0.000298	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.000123	<0.000125	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000784	<0.0000755	<0.0000769	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755	<0.0000769	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00061	<0.00007	<0.00007	<0.00005	<0.00005	0.0063	0.000239 J	<0.0000686	0.00035 J	0.000151 J	0.0019	0.000092 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.000814	<0.000783	<0.000798	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.000549	<0.000528	<0.000538	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.017	0.0077	0.00069	0.00091	0.00041	0.0072	0.0196	0.0181 J	0.0551	0.0294	0.0076	0.0064
Acenaphthylene	1.5	4.4	0.00011 J	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000588	0.000754	<0.0000577	<0.000015	0.000088 J
Anthracene	7.3	22	0.00043	0.00035	<0.00007	<0.00005	<0.00005	0.0013	0.000389 J	<0.000049	0.00111	0.000601	0.00038	0.00028
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0003	<0.0000762	<0.0000784	<0.0000755	<0.0000769	0.000054 J	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755	<0.0000769	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.000123	<0.000125	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00045	<0.00048	0.0004	<0.00099	<0.00013	0.00014 J	<0.000352	<0.000363	<0.000349	<0.000356	0.00011 J	<0.000037
Chrysene	0.91	2	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00027	<0.0000762	<0.0000784	<0.0000755	<0.0000769	0.00005 J	<0.000021
Dibenzofuran	0.098	0.29	0.005	0.0026	0.00011 J	0.00013 J	0.00008 J	0.0043	0.000429 J	0.000141 J	0.00177	0.00115	0.0047	0.0011
Di-n-butylphthalate (DBP)	2.4	7.3	0.000092 J	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.000838	<0.000106	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0011	0.00048	0.00021	0.00053	0.00007 J	0.0027	0.000109 J	0.000365 J	0.000484	0.000782	0.00065	0.00039
Fluorene	0.98	2.9	0.0028	0.0014	0.000095 J	0.00012 J	<0.00005	0.0029	<0.0000667	0.002 J	0.0149	0.0071	0.0024	0.00061
Naphthalene	0.49	1.5	0.19	0.0017	0.00028	<0.00005	0.00027	0.05	0.00286 J	0.000557 J	0.00968	0.00293	0.13	0.013
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.000104	<0.000106	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.000098	<0.0000943	<0.0000962	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.000598	<0.000575	<0.000587	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.00039	<0.00007	<0.00007	<0.00005	<0.00005	0.0068	0.000104 J	<0.0000588	<0.000848	0.000449 J	0.0016	0.00028
Phenol	7.3	22	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000381	<0.0000392	<0.0000377	<0.0000385	<0.000035	<0.000035
Pyrene	0.73	2.2	0.00092	0.00031	<0.00007	0.00029	<0.00005	0.0016	0.000305 J	0.000252 J	0.000376 J	0.000548	0.00055	0.00026
<b>Metals</b>														
Arsenic	0.01	0.01											0.0166	0.0714

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-35A 06/05/2018	MW-35A 01/15/2019	MW-35A 07/18/2019	MW-35A 01/10/2020	MW-35B 01/29/2008	MW-35B 07/14/2008	MW-35B 02/03/2009	MW-35B 01/14/2010	MW-35B 07/01/2010	MW-35B 01/27/2011	MW-35B 07/20/2011	MW-35B 02/15/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00109	<0.005	<0.005	<0.0005	<0.0005	<0.005	<0.001
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<b>0.0648</b>	<b>0.0281</b>	<b>0.062</b>	<b>0.064</b>	<b>0.068</b>	<b>0.064</b>	<b>0.056</b>	<b>0.077</b>
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.0015	<0.005	<0.005	<0.0005	<0.0005	<0.005	<0.001
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	0.176	0.113	0.2	0.2	0.21	0.22	0.17	0.19
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00122	<0.005	<0.005	<0.0005	<0.0005	<b>&lt;0.0065</b>	<0.0013
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	0.00494 J	0.00249 J	0.0057 J	<0.005	0.005	0.0045 J	<0.005	0.0042 J
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002								
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	0.135	0.0787	0.15	0.15 J	0.17	0.16	0.12	0.13
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.0001	<0.0001	<b>0.0012</b>	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	0.0021	<0.00004	<0.00004	<0.00029	<0.0003	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019	<0.0002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038	<0.0004	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	0.00032	0.016	0.00012	<0.000019	<b>0.464</b>	0.0561	<b>0.4</b>	<b>0.47</b>	<b>0.36</b>	<b>0.41</b>	<b>0.48</b>	<b>0.18</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024	<0.00025	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.006	0.0039	0.0028	<0.000027	0.217	0.116	0.17	0.22	0.2	0.19	0.2	0.08
Acenaphthylene	1.5	4.4	0.000066 J	0.000068 J	0.000071 J	<0.000015	<0.00029	<0.0003	0.00088	0.0013	0.0011	0.0012	0.00097	0.00063 J
Anthracene	7.3	22	0.00022	0.00044	0.00011	<0.000014	0.0129	0.00842	0.0056	0.008	0.015	0.014	0.016	0.0048 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	0.00044 J	0.0003 J	0.00017 J	0.00032	0.00022	0.00031	0.00021	0.00011 J
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.0002	<0.00008	0.00014 J	0.00012 J	0.00014 J	0.000069 J	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038	<0.0004	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00018 J	<0.000037	0.000057 J	0.000061 J	<b>0.0198</b>	<0.0002	0.00052	<0.00029	<0.00097	0.00041	<0.00056	<0.00088
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	0.00033 J	0.00022 J	0.00015 J	0.00028	0.00017 J	0.00037	0.00025	0.0001 J
Dibenzofuran	0.098	0.29	0.00072	0.0041	0.00039	<0.00002	<b>0.198</b>	<b>0.104</b>	<b>0.16</b>	<b>0.23</b>	<b>0.22</b>	<b>0.2</b>	<b>0.21</b>	0.097
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.00024	0.00015	0.00015	<0.00001	0.00698	0.00624	0.0031	0.0053	0.006	0.0065	0.0056	0.0026 J
Fluorene	0.98	2.9	0.00086	0.0022	0.00036	0.000065 J	0.0912	0.0685	0.063	0.092	0.11	0.09	0.097	0.048
Naphthalene	0.49	1.5	0.0075	0.22	0.00083	<0.0002	<b>9.3</b>	0.365	<b>12</b>	<b>14</b>	<b>11</b>	<b>4.8</b>	<b>12</b>	<b>7.4</b>
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038	<0.0004	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024	<0.00025	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
Phenanthrene	0.73	2.2	<0.000021	0.0025	0.000042 J	<0.000021	0.1	0.0782	0.061	0.086	0.12	0.078	0.12	0.052
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	0.00059	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Pyrene	0.73	2.2	0.00027	0.00007 J	0.000096 J	<0.000019	0.00411	0.0026	0.0017	0.0027	0.0025	0.0032	0.0027	0.0016 J
<b>Metals</b>														
Arsenic	0.01	0.01	<b>0.0189</b>	<b>0.0198</b>	<b>0.0548</b>	<b>0.0219</b>								

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-35B 07/18/2012	MW-35B 02/07/2013	MW-35B 08/08/2013	MW-35B 01/24/2014	MW-35B 07/24/2014	MW-35B 01/25/2018	MW-35B 03/22/2018	MW-35B 06/05/2018	MW-35B 01/15/2019	MW-35B 07/18/2019	MW-35B 01/10/2020	MW-35B 01/10/2020 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.005	<0.0007	<0.0028	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<b>0.064</b>	<b>0.0662</b>	<b>0.0855</b>	<b>0.0664</b>	<b>0.0539</b>	<b>0.078</b>	<b>0.088</b>	<b>0.044</b>	0.0033	0.0045	<b>0.025 J</b>	<b>0.016 J</b>
Chlorobenzene	0.1	0.1	<0.005	<0.0006	<0.0024	0.000241 J	0.000228 J	0.0009 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.19	0.225	0.258	0.187	0.176	0.15	0.15	0.12	0.0094	0.014	0.096	0.083
Methylene chloride	0.005	0.005	<b>0.02 J</b>	<0.00075	<b>0.0234</b>	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.005	0.00437 J	0.00584 J	0.00429	0.00377	0.0057	0.0041	0.0031	<0.0002	0.00054 J	<0.0002	<0.0002
Vinyl chloride	0.002	0.002			<b>&lt;0.0022</b>		<0.00011						<0.0002	<0.0002
Xylenes (total)	10	10	0.13 J	0.153	0.174	0.132	0.114	0.064	0.066	0.056	0.004	0.0062	0.058 J	0.041 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<b>&lt;0.105</b>	<b>&lt;0.00539</b>	<b>&lt;0.106</b>	<0.00107	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.295	<0.0152	<0.298	<0.00301	<0.0004	<0.0004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00021
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<b>&lt;0.124</b>	<b>&lt;0.00637</b>	<b>&lt;0.125</b>	<0.00126	<0.00059	<0.00058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<b>&lt;0.0762</b>	<b>&lt;0.00392</b>	<b>&lt;0.0769</b>	<0.000777	<0.00042	<0.00042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.0762	<0.00392	<0.0769	<0.000777	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<b>0.26</b>	<b>0.295 J</b>	<b>0.431</b>	<b>0.534</b>	<b>0.376</b>	<b>0.13</b>	<b>0.22</b>	<b>0.25</b>	0.011	0.041	<b>0.14 J</b>	0.062 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<b>&lt;0.79</b>	<b>&lt;0.0407</b>	<b>&lt;0.798</b>	<b>&lt;0.00806</b>	<0.0002	0.00078 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<b>&lt;0.533</b>	<0.0275	<b>&lt;0.538</b>	<0.00544	<0.00047	<0.00047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.15	0.173 J	0.258	0.305 J	0.139	0.094	0.18	0.17	0.013	0.029	0.059 J	0.033 J
Acenaphthylene	1.5	4.4	0.00078	<0.0571	<0.00294	<0.0577	0.0015 J	<0.00015	0.0014	0.00076	0.00018	0.00022	0.00038	0.00027
Anthracene	7.3	22	0.0064	<0.0476	0.0202 J	<0.0481	0.0111	0.011	0.014	0.0072	0.0011	0.0022	0.0029 J	0.0018 J
Benzo(a)anthracene	0.0091	0.02	0.0002	<b>&lt;0.0762</b>	<0.00392	<b>&lt;0.0769</b>	<0.000777	<0.00051	<0.0005	0.000075 J	0.000077 J	0.000083 J	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<b>&lt;0.0762</b>	<b>&lt;0.00392</b>	<b>&lt;0.0769</b>	<b>&lt;0.000777</b>	<0.0002	<0.0002	<0.00002	0.000058 J	0.000053 J	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<b>&lt;0.124</b>	<b>&lt;0.00637</b>	<b>&lt;0.125</b>	<b>&lt;0.00126</b>	<0.0003	<0.0003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00018 J	<b>&lt;0.352</b>	<b>&lt;0.0181</b>	<b>&lt;0.356</b>	<0.00359	<0.00037	<0.00037	<0.000037	<0.000014	0.00012 J	<0.000037	<0.000037
Chrysene	0.91	2	0.00023	<0.0762	<0.00392	<0.0769	<0.000777	0.00051 J	<0.00021	0.000086 J	0.000098 J	0.000091 J	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<b>0.14</b>	<b>0.161 J</b>	<b>0.252</b>	<b>0.256 J</b>	<b>0.138</b>	<b>0.13</b>	<b>0.22</b>	<b>0.18</b>	0.015	0.029	0.054 J	0.031 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.105	<0.00539	<0.106	<0.00107	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0039	<0.0667	0.00756 J	0.0698 J	0.00692	0.0081	0.0098	0.0031	0.0013	0.0017	0.0015 J	0.00093 J
Fluorene	0.98	2.9	0.069	<0.0667	0.138	0.167 J	0.076	0.057	0.092	0.095	0.0066	0.014	0.028 J	0.015 J
Naphthalene	0.49	1.5	<b>7.6</b>	<b>8.83</b>	<b>14.1</b>	<b>13.1</b>	<b>9.36</b>	<b>5</b>	<b>13</b>	<b>15</b>	0.079	<b>1.1</b>	<b>2 J</b>	0.13 J
Nitrobenzene	0.049	0.15	<0.00005	<b>&lt;0.105</b>	<0.00539	<b>&lt;0.106</b>	<0.00107	<0.00024	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.0952	<0.0049	<0.0962	<0.000971	<0.00025	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<b>&lt;0.581</b>	<b>&lt;0.0299</b>	<b>&lt;0.587</b>	<b>&lt;0.00592</b>	<0.0008	<0.00079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.066	0.0936 J	0.142	0.27 J	0.0891	0.088	0.18	0.11	0.0086	0.019	0.024 J	0.013 J
Phenol	7.3	22	0.00014 J	<0.0381	<0.00196	0.129 J	<0.000388	<0.00035	<0.00035	<0.000035	<0.000035	0.00017 J	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0019	<0.105	<0.00539	<0.106	0.00327 J	0.0057	0.0053	0.0018	0.00075	0.00079	0.0007 J	0.00045 J
<b>Metals</b>														
Arsenic	0.01	0.01						0.00465	0.00595	<b>0.0116</b>	0.00862	0.0012 J	<b>0.012</b>	<b>0.0159</b>

Notes:

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- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-36A 01/29/2008	MW-36A 07/14/2008	MW-36A 02/03/2009	MW-36A 01/13/2010	MW-36A 06/29/2010	MW-36A 01/20/2011	MW-36A 07/19/2011	MW-36A 02/07/2012	MW-36A 07/17/2012	MW-36A 01/31/2013	MW-36A 08/06/2013	MW-36A 01/16/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002
Benzene	0.005	0.005	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002
Chlorobenzene	0.1	0.1	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018
Ethylbenzene	0.7	0.7	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019
Methylene chloride	0.005	0.005	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022
Toluene	1	1	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017
Vinyl chloride	0.002	0.002	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00018
Xylenes (total)	10	10	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.000104	<0.000104
2,4-Dimethylphenol	0.49	1.5	<0.0003	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000292	<0.000292	<0.000292
2,4-Dinitrotoluene	0.0013	0.003	<0.0002	<0.0002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000123	<0.000123
2,6-Dinitrotoluene	0.0013	0.003	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000755	<0.0000755
2-Chloronaphthalene	2	5.8	<0.0004	<0.00039	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000755	<0.0000755
2-Methylnaphthalene	0.098	0.29	<0.0004	<0.00039	<0.00007	0.0003	0.00023	<0.00007	<0.00005	<0.00005	<0.00005	<0.000066	<0.000066	<0.000066
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000783	<0.000783	<0.000783
4-Nitrophenol	0.049	0.15	<0.00025	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000528	<0.000528	<0.000528
Acenaphthene	1.5	4.4	<0.0003	<0.00029	<0.00009	0.00036	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000755	<0.0000867
Acenaphthylene	1.5	4.4	<0.0003	<0.00029	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000566	<0.0000566	<0.0000566
Anthracene	7.3	22	0.00065	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000472	<0.0000472	<0.0000472
Benzo(a)anthracene	0.0091	0.02	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000755	<0.0000755
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000755	<0.0000755
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0004	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000123	<0.000123
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00061 J	<0.0002	0.00045	<0.00033	<0.00061	<0.00048	<0.0004	0.0025	<0.0001	<0.000349	<0.000349	<0.000349
Chrysene	0.91	2	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000755	<0.0000755
Dibenzofuran	0.098	0.29	0.00049 J	<0.00029	<0.00008	0.0003	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000755	<0.0000755
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000104	<0.000391	<0.000104
Fluoranthene	0.98	2.9	0.000526	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000066	<0.000066	<0.000066
Fluorene	0.98	2.9	0.00028 J	<0.0002	<0.00007	0.00024	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000066	<0.000066	<0.000066
Naphthalene	0.49	1.5	0.00119	<0.00039	0.0006	0.0013	0.0023	<0.0001	<0.00005	<0.00005	0.0003	<0.000211	<0.0000755	<0.0000755
Nitrobenzene	0.049	0.15	<0.0004	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.000104	<0.000104
N-Nitrosodiphenylamine	0.19	0.42	<0.00025	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000943	<0.0000943
Pentachlorophenol	0.001	0.001	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000575	<0.000575	<0.000575
Phenanthrene	0.73	2.2	0.000727	<0.0002	<0.00007	0.00039	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000566	<0.0000861	<0.000201
Phenol	7.3	22	<0.0002	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000377	<0.0000377	<0.0000377
Pyrene	0.73	2.2	0.000531	<0.0002	0.00015 J	0.00021	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000104	<0.000104	0.000155 J
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-36A 07/28/2014	MW-36A 01/25/2018	MW-36A 03/21/2018	MW-36A 05/31/2018	MW-36A 01/14/2019	MW-36A 07/16/2019	MW-36A 01/09/2020	MW-36D 07/15/2010	MW-36D 01/26/2011	MW-36D 07/27/2011	MW-36D 02/14/2012	MW-36D 07/23/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Benzene	0.005	0.005	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	0.00071 J	<0.001	<0.001	<0.0005
Chlorobenzene	0.1	0.1	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Ethylbenzene	0.7	0.7	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005
Methylene chloride	0.005	0.005	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0013	<0.0013	<0.001
Toluene	1	1	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Vinyl chloride	0.002	0.002	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Xylenes (total)	10	10	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.001	<0.001	<0.0031	<0.0031	<0.0015
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.000301	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004	<0.000052	0.00008 J	<0.00008	<0.00005	<0.00005	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000126	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.0000777	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.000177	<0.000019	0.000022 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.00007	0.00013 J	<0.00005	<0.00005	<0.00005
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000544	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Acenaphthene	1.5	4.4	<0.0000777	<0.000028	0.000059	<0.000027	<0.000027	<0.000027	<0.000027	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
Acenaphthylene	1.5	4.4	<0.0000583	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Anthracene	7.3	22	<0.0000485	<0.000014	0.000066 J	<0.000014	<0.000014	<0.000014	<0.000014	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.0000777	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.0000777	<0.00002	0.000064 J	0.00003 J	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000126	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000359	0.000089 J	0.00015 J	0.0001 J	<0.000037	<0.000037	<0.000037	0.005 J	0.00097	0.0012	<0.0012	<0.00035
Chrysene	0.91	2	<0.0000777	<0.000021	<0.000021	0.000027 J	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007	<0.00005	0.000078 J	<0.00005
Dibenzofuran	0.098	0.29	<0.0000777	<0.00002	0.00061	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000107	<0.00002	<0.00002	0.000028 J	<0.00002	<0.00002	<0.00002	<0.00007	<0.00007	0.00032	0.000052 J	<0.00005
Fluoranthene	0.98	2.9	<0.000068	<0.00001	0.00012	0.000033 J	<0.00001	<0.00001	<0.00001	<0.00007	<0.00007	0.000068 J	0.00013 J	0.000054 J
Fluorene	0.98	2.9	<0.000068	<0.000031	0.00034	<0.00003	<0.00003	<0.00003	<0.00003	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Naphthalene	0.49	1.5	<0.00101	<0.00002	<0.00002	<0.00002	<0.000062	0.00015	<0.00034	<0.0001	0.00083	0.000061 J	<0.000072	<0.00014
Nitrobenzene	0.049	0.15	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.0000971	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.000592	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
Phenanthrene	0.73	2.2	<0.0000583	<0.000021	0.00011	0.000026 J	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007	0.000072 J	0.000069 J	<0.00005
Phenol	7.3	22	<0.0000388	<0.000036	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	0.00065	<0.00007	0.000056 J	0.00023	<0.00005
Pyrene	0.73	2.2	<0.000107	<0.000019	0.000076 J	0.000032 J	<0.000019	<0.000019	<0.000019	<0.00007	<0.00007	0.000053 J	0.000087 J	<0.00005
<b>Metals</b>														
Arsenic	0.01	0.01		0.00108 J	0.00753	0.00117 J	0.00107 J	0.00244	0.00354					

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-36D 07/23/2012 Duplicate	MW-36D 02/11/2013	MW-36D 08/05/2013	MW-36D 01/21/2014	MW-36D 08/28/2014	MW-36D 02/07/2018	MW-36D 03/26/2018	MW-36D 05/31/2018	MW-36D 01/24/2019	MW-36D 07/31/2019	MW-36D 01/16/2020	MW-38A 01/29/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052
Benzene	0.005	0.005	<0.0005	<0.00008	<0.00008	<0.0002	0.0000895 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025
Chlorobenzene	0.1	0.1	<0.0005	<0.00012	0.00013 J	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047
Ethylbenzene	0.7	0.7	<0.0005	<0.00011	0.000127 J	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025
Methylene chloride	0.005	0.005	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054
Toluene	1	1	<0.0005	<0.00015	<0.00015	<0.00017	0.00675	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.000105	<0.00011	<0.000104	<0.000109	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.000295	<0.00031	<0.000292	<0.000307	<0.0004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00029
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.000124	<0.00013	<0.000123	<0.000129	<0.00058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.0000762	<0.00008	<0.0000755	<0.0000792	<0.00042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019
2-Chloronaphthalene	2	5.8	<0.00005	<0.0000762	<0.00008	<0.0000755	<0.0000792	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00039
2-Methylnaphthalene	0.098	0.29	<0.00005	<0.0000667	<0.00007	0.000189 J	<0.0000693	<0.00019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00039
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00079	<0.00083	<0.000783	<0.000822	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
4-Nitrophenol	0.049	0.15	<0.00005	<0.000533	<0.00056	<0.000528	<0.000554	<0.00047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024
Acenaphthene	1.5	4.4	<0.00005	<0.0000762	<0.00008	0.00014 J	<0.0000792	0.00056 J	0.002	<0.000027	<0.000027	<0.000027	<0.000027	<0.00029
Acenaphthylene	1.5	4.4	<0.00005	<0.0000571	<0.00006	<0.0000566	<0.0000594	<0.00015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029
Anthracene	7.3	22	<0.00005	<0.0000476	<0.00005	0.000105 J	0.000224 J	<0.00014	0.00016	0.000037 J	<0.000014	0.000017 J	<0.000033	<0.00019
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.0000762	<0.00008	<0.0000755	0.000213 J	<0.0005	0.00024	0.00019	<0.00005	0.000053 J	<0.00005	<0.00019
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.0000762	<0.00008	<0.0000755	0.000192 J	<0.0002	0.0003	0.00024	0.000027 J	0.000087 J	<0.00002	<0.00019
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.000124	<0.00013	<0.000123	<0.000129	<0.0003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00039
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00059	<0.000352	<0.00037	<0.000349	0.00128	<0.00037	0.00054	0.0005	0.000055 J	0.00008 J	<0.000037	0.00078 J
Chrysene	0.91	2	<0.00005	<0.0000762	<0.00008	<0.0000755	0.000347 J	<0.00021	0.00031	0.00026	0.000031 J	0.000058 J	<0.000021	<0.00019
Dibenzofuran	0.098	0.29	<0.00005	<0.0000762	<0.00008	0.00017 J	<0.0000792	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00029
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.000105	<0.00011	<0.000104	<0.000109	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
Fluoranthene	0.98	2.9	0.000093 J	<0.0000667	<0.00007	0.000216 J	0.00111	<0.0001	0.00094	0.00039	0.000048 J	0.00011	<0.00001	<0.00019
Fluorene	0.98	2.9	<0.00005	<0.0000667	<0.00007	0.000147 J	0.0001 J	<0.0003	0.00086	0.00003 J	<0.00003	<0.00003	<0.00003	<0.00019
Naphthalene	0.49	1.5	<0.00015	<0.0000762	<0.00008	0.0011 J	0.0000923 J	<0.0002	0.00015	0.000045 J	<0.00002	<0.00002	<0.000033	<0.00039
Nitrobenzene	0.049	0.15	<0.00005	<0.000105	<0.00011	<0.000104	<0.000109	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00039
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.0000952	<0.0001	<0.0000943	<0.000099	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024
Pentachlorophenol	0.001	0.001	<0.00005	<0.000581	<0.00061	<0.000575	<0.000604	<0.00079	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019
Phenanthrene	0.73	2.2	0.00014 J	<0.0000571	<0.00006	0.000665	0.00102	<0.00021	0.00037	0.00028	<0.000021	0.000066 J	<0.000047	<0.00019
Phenol	7.3	22	<0.00005	<0.0000381	<0.00004	<0.0000377	0.00194	<0.00035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019
Pyrene	0.73	2.2	<0.00005	<0.000105	<0.00011	0.000159 J	0.000881	<0.00019	0.00062	0.00036	0.000042 J	0.000099 J	<0.000019	<0.00019
<b>Metals</b>														
Arsenic	0.01	0.01						0.000773 J	0.00137 J	<0.0004	0.000417 J	<0.0004	0.000533 J	

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-38A 01/29/2008 Duplicate	MW-38A 07/14/2008	MW-38A 07/14/2008 Duplicate	MW-38A 02/03/2009	MW-38A 01/14/2010	MW-38A 06/29/2010	MW-38A 01/25/2011	MW-38A 07/19/2011	MW-38A 08/25/2011	MW-38A 02/15/2012	MW-38A 07/18/2012	MW-38A 02/07/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.00109	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001		<0.001	<0.0005	<0.00014
Benzene	0.005	0.005	<0.00025	<0.00112	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001		<0.001	<0.0005	<0.00008
Chlorobenzene	0.1	0.1	<0.00047	<0.0015	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001		<0.001	<0.0005	<0.00012
Ethylbenzene	0.7	0.7	<0.00025	<0.00142	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011		<0.0011	<0.0005	<0.00011
Methylene chloride	0.005	0.005	<0.00054	<0.00122	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013		<0.0013	<0.001	<0.00015
Toluene	1	1	<0.00041	<0.00138	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001		<0.001	<0.0005	<0.00015
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00127	<0.00302	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031		<0.0031	<0.0015	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005		<0.00005	<0.00005	<0.000105
2,4-Dimethylphenol	0.49	1.5	<0.00029	<0.00029	<0.00028	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005		<0.00005	<0.00005	<0.000295
2,4-Dinitrotoluene	0.0013	0.003	<0.0002	<0.0002	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.000124
2,6-Dinitrotoluene	0.0013	0.003	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006		<0.00006	<0.00006	<0.0000762
2-Chloronaphthalene	2	5.8	<0.00039	<0.00039	<0.00038	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005		<0.00005	<0.00005	<0.0000762
2-Methylnaphthalene	0.098	0.29	<0.00039	<0.00039	<0.00038	0.00044	<0.00007	0.00016 J	0.000085 J	<0.00005		<0.00005	<0.00031	<0.0000667
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008		<0.00008	<0.00008	<0.00079
4-Nitrophenol	0.049	0.15	<0.00024	<0.00024	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00005	<0.000533
Acenaphthene	1.5	4.4	<0.00029	<0.00029	<0.00028	<0.00009	0.00024	<0.00009	<0.00009	0.00043		<0.00005	<0.00025	<0.0000762
Acenaphthylene	1.5	4.4	<0.00029	<0.00029	<0.00028	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00005	<0.0000571
Anthracene	7.3	22	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	0.00023	0.0001 J	<0.00005	<0.00013	0.0000712 J
Benzo(a)anthracene	0.0091	0.02	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	0.00025		<0.00005	<0.00005	<0.0000762
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	0.00052	<0.00005	<0.00005	<0.00005	<0.0000762
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00039	<0.00039	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.000124
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0145	0.00022 J	0.00105 J	<0.00049	<0.00044	<0.00064	<0.00094			0.0016	0.00023	<0.000352
Chrysene	0.91	2	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	0.00022		<0.00005	<0.00005	<0.0000762
Dibenzofuran	0.098	0.29	<0.00029	<0.00029	<0.00028	<0.00008	<0.00008	0.000083 J	<0.00008	<0.00005	0.000055 J	<0.00005	<0.00014	<0.0000762
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	0.00017 J		<0.00005	0.000052 J	<0.000105
Fluoranthene	0.98	2.9	<0.0002	<0.0002	<0.00019	<0.00007	0.00012 J	<0.00007	0.00034	0.00014 J		<0.00005	<0.00017	<0.0000667
Fluorene	0.98	2.9	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00015	<0.0000667
Naphthalene	0.49	1.5	<0.00039	<0.00039	<0.00038	0.006	<0.0001	<0.0001	0.00059	<0.00026		<0.00005	<0.001	<0.0000762
Nitrobenzene	0.049	0.15	<0.00039	<0.00039	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.000105
N-Nitrosodiphenylamine	0.19	0.42	<0.00024	<0.00024	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.0000952
Pentachlorophenol	0.001	0.001	<0.0002	<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	0.000067 J		<0.00005	<0.00005	<0.000581
Phenanthrene	0.73	2.2	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	0.0001 J	<0.00005		0.00011 J	<0.00032	<0.0000571
Phenol	7.3	22	<0.0002	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00005	<0.0000381
Pyrene	0.73	2.2	<0.0002	<0.0002	<0.00019	<0.00007	0.0001 J	<0.00007	0.00021	0.00018 J		<0.00005	<0.00016	<0.000105
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-38A 08/08/2013	MW-38A 01/21/2014	MW-38A 07/25/2014	MW-38A 01/26/2018	MW-38A 03/25/2018	MW-38A 06/05/2018	MW-38A 01/22/2019	MW-38A 07/31/2019	MW-38A 01/21/2020	MW-38B 01/29/2008	MW-38B 07/14/2008	MW-38B 02/03/2009	
<b>Volatile Organic Compounds</b>															
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.00031 J	<0.0002	<0.00052	<0.00109	<0.0005
Benzene	0.005	0.005	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00112	<0.0005
Chlorobenzene	0.1	0.1	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.0015	<0.0005
Ethylbenzene	0.7	0.7	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00142	<0.0005
Methylene chloride	0.005	0.005	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00122	<0.0005
Toluene	1	1	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00138	<0.0005
Vinyl chloride	0.002	0.002	<0.00011	<0.00011	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00138	<0.0005
Xylenes (total)	10	10	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00302	<0.001
<b>Semivolatile Organic Compounds</b>															
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000104	<0.000105	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.000292	<0.000295	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.0003	<0.00029	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000123	<0.000124	<0.000126	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.0002	<0.00019	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0000755	<0.0000762	<0.0000777	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.0002	<0.00019	<0.00007
2-Chloronaphthalene	2	5.8	<0.0000755	<0.0000762	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0004	<0.00038	<0.00012
2-Methylnaphthalene	0.098	0.29	0.000115 J	<0.0000667	<0.000068	<0.000019	0.00014	<0.000019	0.000055 J	<0.000019	0.00017	<0.0004	<0.00038	0.00037	<0.00008
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000783	<0.00079	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<0.00019	<0.00008
4-Nitrophenol	0.049	0.15	<0.000528	<0.000533	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00025	<0.00024	<0.00007	<0.00007
Acenaphthene	1.5	4.4	0.000638	<0.0000762	<0.0000777	<0.000027	0.0039	0.00061	0.014	<0.000027	0.00068	<0.0003	<0.00029	0.0001 J	<0.00006
Acenaphthylene	1.5	4.4	<0.0000566	<0.0000571	<0.0000583	<0.000015	0.000053 J	<0.000015	0.0002	<0.000015	<0.000015	<0.0003	<0.00029	<0.00006	<0.00006
Anthracene	7.3	22	<0.0000472	0.000103 J	<0.0000816	<0.000014	0.00026	0.00015	0.00017	0.000022 J	0.00015	<0.0002	0.00026 J	0.00013 J	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.0000755	<0.0000762	<0.0000777	<0.00005	<0.000051	<0.00005	<0.00005	0.000087 J	<0.00005	<0.0002	<0.00019	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.0000755	<0.0000762	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	0.00012	0.000032 J	<0.0002	<0.00019	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000123	<0.000124	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.0004	<0.00038	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000349	<0.000352	<0.000359	<0.000037	<0.000037	0.0001 J	<0.000037	0.000084 J	0.000063 J	0.00103 J	<0.00019	0.00041	<0.00007
Chrysene	0.91	2	<0.0000755	<0.0000762	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	0.000092 J	0.00006 J	<0.0002	<0.00019	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.0000755	<0.0000762	<0.0000777	<0.00002	<0.00002	0.00006 J	0.00014	<0.00002	0.00036	<0.0003	<0.00029	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	0.000145 J	<0.000105	<0.000107	<0.00002	<0.00002	0.00017 J	0.000068 J	<0.00002	<0.00002	<0.0002	<0.00019	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.000164 J	<0.0000667	<0.000068	<0.00001	0.00023	0.000032 J	0.0013	0.00018	0.00018	<0.0002	<0.00019	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.000066	<0.0000667	<0.000068	<0.00003	0.0003	0.00016	0.0024	<0.00003	0.00029	<0.0002	<0.00019	<0.00007	<0.00007
Naphthalene	0.49	1.5	0.00192 J	<0.000146	<0.0000777	<0.00002	<0.00002	0.00071	0.00015	0.000068 J	0.0011	<0.0004	<0.00038	0.0045	<0.00007
Nitrobenzene	0.049	0.15	<0.000104	<0.000105	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.0004	<0.00038	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0000943	<0.0000952	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	0.00011 J	<0.00025	<0.00024	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000575	<0.000581	<0.000592	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.0002	<0.00019	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.0000566	<0.0000571	<0.0000583	<0.000021	0.00033	<0.000021	0.00058	0.0001	0.00039	<0.0002	<0.00019	<0.00007	<0.00007
Phenol	7.3	22	<0.0000377	<0.0000381	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.0002	<0.00019	<0.00007	<0.00007
Pyrene	0.73	2.2	0.000176 J	<0.000105	<0.000107	<0.000019	0.00021	0.000056 J	0.0011	0.00014	0.00014	<0.0002	<0.00019	<0.00007	<0.00007
<b>Metals</b>															
Arsenic	0.01	0.01				<0.0004	0.0138	0.0124	0.0186	<0.0004	0.0177				

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-38B 01/14/2010	MW-38B 06/29/2010	MW-38B 01/25/2011	MW-38B 07/18/2011	MW-38B 02/15/2012	MW-38B 07/18/2012	MW-38B 02/07/2013	MW-38B 08/08/2013	MW-38B 01/21/2014	MW-38B 07/25/2014	MW-38B 01/26/2018	MW-38B 03/25/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002
Vinyl chloride	0.002	0.002									<0.00011			
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000109	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000295	<0.000292	<0.000295	<0.000307	<0.000041	<0.000041
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.000129	<0.000059	<0.000059
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000755	<0.0000762	<0.0000792	<0.000043	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.0000792	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.0003	<0.0000667	<0.000066	0.000137 J	<0.0000992	0.000096 J	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.000783	<0.00079	<0.000822	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.000528	<0.000533	<0.000554	<0.000048	<0.000047
Acenaphthene	1.5	4.4	<0.00009	0.00047	<0.00009	<0.00005	0.0001 J	0.00096	0.000226 J	<0.0000755	0.0000786 J	<0.000342	<0.000028	0.000095 J
Acenaphthylene	1.5	4.4	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000566	<0.0000571	<0.0000594	<0.000015	<0.000015
Anthracene	7.3	22	<0.00007	<0.00007	0.00011 J	0.00013 J	0.00021	<0.00023	0.000313 J	<0.0000472	0.000141 J	<0.000398	0.000047 J	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.0000792	<0.000051	<0.000051
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.0000792	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	0.000072 J	<0.000124	<0.000123	<0.000124	<0.000129	<0.000031	<0.000031
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00039	<0.00074	<0.0002	<0.001	<0.0001	<0.0001	<0.000352	<0.000349	<0.000352	<0.000366	0.000061 J	0.00012 J
Chrysene	0.91	2	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.0000792	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00028	<0.0000762	<0.0000755	0.0000923 J	<0.00041	0.000053 J	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000109	<0.00002	0.000027 J
Fluoranthene	0.98	2.9	<0.00007	0.00017 J	<0.00007	<0.00005	<0.00005	<0.00032	<0.0000667	<0.000066	0.000101 J	<0.00058	<0.00001	0.000022 J
Fluorene	0.98	2.9	<0.00007	0.00015 J	<0.00007	<0.00005	<0.00005	<0.00027	<0.0000667	<0.000066	0.0000778 J	<0.000219	0.000046 J	<0.00003
Naphthalene	0.49	1.5	<0.00014	<0.0001	0.00031	<0.00005	0.00037	<0.0015	<0.0000762	<0.0000755	0.000466 J	<0.0016	0.00084	0.00016
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000109	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000943	<0.0000952	<0.000099	<0.000026	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.000575	<0.000581	<0.000604	<0.000081	<0.00008
Phenanthrene	0.73	2.2	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00037	<0.0000571	<0.0000566	0.000304 J	<0.000164	0.000061 J	0.000052 J
Phenol	7.3	22	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000381	<0.0000377	<0.0000381	<0.0000396	<0.000036	<0.000035
Pyrene	0.73	2.2	<0.00007	0.00027	<0.00007	<0.00005	<0.00005	<0.00037	<0.000105	<0.000104	<0.000105	0.000472 J	<0.000019	0.000035 J
<b>Metals</b>														
Arsenic	0.01	0.01											0.000636 J	0.000972 J

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-38B 06/05/2018	MW-38B 01/22/2019	MW-38B 07/18/2019	MW-38B 01/21/2020	MW-39B 01/30/2008	MW-39B 07/15/2008	MW-39B 02/04/2009	MW-39B 01/19/2010	MW-39B 06/22/2010	MW-39B 01/18/2011	MW-39B 07/26/2011	MW-39B 02/01/2012
<b><i>Volatile Organic Compounds</i></b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031
<b><i>Semivolatile Organic Compounds</i></b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	0.00025	<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	0.000056 J	<0.00004	<0.0003	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	0.00005 J
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.0002	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	0.00005 J
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	0.00006 J
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.0004	<0.00039	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	0.00005 J
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	0.0003	0.00017	<0.0004	<0.00039	<0.00007	<0.00007	<0.00007	0.000086 J	<0.00005	0.00005 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00049	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.00025	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00005 J
Acenaphthene	1.5	4.4	0.048	<0.000027	0.00012	0.054	0.000664	<0.00029	0.00022	0.00014 J	0.0034	0.00039	0.00028	0.0011 J
Acenaphthylene	1.5	4.4	0.00036	<0.000015	<0.000015	0.00045	<0.0003	<0.00029	<0.00006	<0.00007	<0.00007	<0.00007	0.000053 J	0.00005 J
Anthracene	7.3	22	0.0018	0.0001	0.000037 J	0.0012	0.00106	0.000619	0.00028	<0.00007	<0.00007	<0.00007	0.0004	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	0.000055 J	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00005 J
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.0004	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	0.00005 J
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	0.00031	<0.000037	0.00119 J	<0.00019	0.00046	0.0007	<0.0002	0.00024	0.00092	<0.00015
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	0.000057 J	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	0.008	<0.00002	0.00014	0.0054	<0.0003	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	0.0004 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	0.000036 J	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.0014	<0.00001	<0.00001	0.0025	0.00213	0.000575	0.0014	<0.00007	0.0019	0.00013 J	0.000079 J	0.0011
Fluorene	0.98	2.9	0.019	<0.00003	0.00006 J	0.015	<0.0002	<0.00019	0.00025	0.00021	0.00048	0.00013 J	0.00011 J	0.00032 J
Naphthalene	0.49	1.5	0.00099	<0.00002	0.0026	0.00098	<0.0004	<0.00039	0.00052	0.00018 J	<0.00015	0.00076	<0.00005	0.00034 J
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.0004	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	0.00005 J
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.00025	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	0.000054 J	<0.00005
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.0002	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	0.00005 J
Phenanthrene	0.73	2.2	<0.000021	<0.000021	0.000079 J	0.00095	0.00045 J	<0.00019	<0.00007	0.00025	<0.00007	0.00018 J	<0.00005	<0.00005
Phenol	7.3	22	<0.000035	<0.000035	0.000056 J	<0.000035	<0.0002	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Pyrene	0.73	2.2	0.00086	<0.000019	<0.000019	0.0015	0.00156	0.000956	0.0013	0.00018 J	0.002	<0.00007	0.00017 J	0.0013
<b><i>Metals</i></b>														
Arsenic	0.01	0.01	<b>0.0386</b>	<0.0004	0.000553 J	<b>0.0173</b>								

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-39B 07/19/2012	MW-39B 02/05/2013	MW-39B 07/31/2013	MW-39B 01/14/2014	MW-39B 07/25/2014	MW-39B 01/23/2018	MW-39B 03/19/2018	MW-39B 05/16/2018	MW-39B 01/08/2019	MW-39B 07/11/2019	MW-39B 01/13/2020	MW-40B 01/30/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052
Benzene	0.005	0.005	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0348
Chlorobenzene	0.1	0.1	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047
Ethylbenzene	0.7	0.7	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.162
Methylene chloride	0.005	0.005	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054
Toluene	1	1	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0791
Vinyl chloride	0.002	0.002	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00054
Xylenes (total)	10	10	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.35
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.000104	<0.000105	<0.00106	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.000292	<0.000295	<0.00298	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.000123	<0.000124	<0.00125	<0.000126	<0.000058	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058	<0.00047
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.0000755	<0.0000762	<0.000769	<0.0000777	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00047
2-Chloronaphthalene	2	5.8	<0.00005	<0.0000755	<0.0000762	<0.000769	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00047
2-Methylnaphthalene	0.098	0.29	0.000069 J	<0.000066	<0.0000667	<0.000673	<0.0000708	0.000081 J	<0.000019	0.00008 J	<0.000019	<0.000027	<0.000021	0.522
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.000783	<0.00079	<0.00798	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00047
4-Nitrophenol	0.049	0.15	<0.00005	<0.000528	<0.000533	<0.00538	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00047
Acenaphthene	1.5	4.4	0.0004	0.000756	0.0007	0.00115 J	<0.0012	0.00093	0.001	0.00054	0.00062	0.00065	0.0012	0.365
Acenaphthylene	1.5	4.4	<0.00005	0.00011 J	0.0000676 J	<0.000577	0.0000623 J	<0.000015	<0.000015	<0.000015	<0.000015	0.000021 J	<0.000015	<0.00047
Anthracene	7.3	22	0.0001 J	0.000901	0.000774	<0.00101	<0.000615	0.000053 J	<0.000014	0.0002	0.00016	0.00021	0.00016	<0.00047
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.0000755	<0.0000762	<0.000769	<0.0000777	<0.00005	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005	<0.00047
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.0000755	<0.0000762	<0.000769	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00047
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.000123	<0.000124	<0.00125	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00047
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00015	<0.000349	<0.000352	<0.00356	<0.000359	0.0018	0.00013 J	0.00011 J	<0.000037	0.000042 J	<0.000037	<0.00047
Chrysene	0.91	2	<0.00005	<0.0000755	<0.0000762	<0.000769	<0.0000777	0.000057 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00047
Dibenzofuran	0.098	0.29	0.000067 J	<0.0000755	<0.0000762	<0.000769	<0.0000823	0.0001	<0.00002	0.000054 J	<0.00002	0.000037 J	0.00019	0.239
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.000104	<0.000122	<0.00106	<0.000107	0.000082 J	<0.00002	0.000035 J	<0.00002	<0.00002	<0.00002	<0.00047
Fluoranthene	0.98	2.9	0.00036	0.000112 J	0.000418 J	<0.000673	<0.000605	0.00015	0.000075 J	0.000053 J	0.000067 J	0.000048 J	0.00023	<0.00047
Fluorene	0.98	2.9	0.00019 J	<0.000066	0.000216 J	<0.000673	<0.000311	0.00011	<0.00003	0.000061 J	<0.00003	0.000044 J	0.00032	0.175
Naphthalene	0.49	1.5	0.00018 J	0.000428 J	<0.0000762	<0.000769	<0.0000777	0.00082	<0.00002	<0.0013	<0.000092	<0.00041	0.0023	9.34
Nitrobenzene	0.049	0.15	<0.00005	<0.000104	<0.000105	<0.00106	0.000853	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00047
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.0000943	0.000158 J	<0.000962	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00047
Pentachlorophenol	0.001	0.001	<0.00005	<0.000575	<0.000581	<0.00587	<0.000592	<0.000079	<0.00008	<0.00008	<0.000079	<0.000079	<0.000079	<0.00047
Phenanthrene	0.73	2.2	0.00016 J	<0.0000566	0.0000912 J	<0.000577	<0.0001	<0.000021	<0.000021	0.000056 J	0.000039 J	<0.000021	0.00019	0.173
Phenol	7.3	22	<0.00005	<0.0000377	<0.0000381	<0.000385	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00047
Pyrene	0.73	2.2	0.00052	0.000131 J	0.000655	<0.00106	0.000818	0.00015	0.000074 J	0.000064 J	0.000052 J	0.000072 J	0.00022	<0.00047
<b>Metals</b>														
Arsenic	0.01	0.01						0.0108	0.00188 J	0.00178 J	0.00365	0.00144 J	0.00248	

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray.
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-40B 07/15/2008	MW-40B 02/04/2009	MW-40B 01/19/2010	MW-40B 06/22/2010	MW-40B 01/18/2011	MW-40B 07/14/2011	MW-40B 02/03/2012	MW-40B 07/19/2012	MW-40B 02/05/2013	MW-40B 07/31/2013	MW-40B 01/14/2014	MW-40B 07/18/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.01	<0.01	<0.005	<0.00014	<0.0014	<0.0007	<0.00014
Benzene	0.005	0.005	<b>0.0269</b>	<b>0.026</b>	<b>0.028</b>	<b>0.026</b>	<b>0.019</b>	<b>0.016 J</b>	<b>0.013 J</b>	<b>0.013 J</b>	<b>0.0108</b>	<b>0.0115</b>	<b>0.0109</b>	<b>0.0103</b>
Chlorobenzene	0.1	0.1	<0.00047	0.001 J	<0.0005	<0.0005	<0.0005	<0.01	<0.01	<0.005	<0.00012	<0.0012	<0.0006	<0.00012
Ethylbenzene	0.7	0.7	0.116	0.1	0.12	0.12	0.13	0.081	0.08	0.082	0.0817	0.0798	0.084	0.0825
Methylene chloride	0.005	0.005	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<b>&lt;0.013</b>	<b>&lt;0.013</b>	<b>&lt;0.01</b>	<0.00015	<0.0015	<0.00075	<0.00015
Toluene	1	1	0.059	0.05	0.054	0.05	0.045	0.019 J	0.028 J	0.022 J	0.0118	0.0173	0.0147	0.0154
Vinyl chloride	0.002	0.002									<0.00011			<0.00011
Xylenes (total)	10	10	0.244	0.2	0.22	0.22	0.21	0.12 J	0.13 J	0.14 J	0.116	0.127	0.12	0.126
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0004	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<b>&lt;0.0104</b>	<b>&lt;0.00524</b>	<b>&lt;0.0212</b>	<0.00109
2,4-Dimethylphenol	0.49	1.5	0.0445	0.011	0.014	0.0044	0.00033 J	0.0034	0.004	0.0039	<0.0292	<0.0148	<0.0596	<0.00307
2,4-Dinitrotoluene	0.0013	0.003	<0.001	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<b>&lt;0.0123</b>	<b>&lt;0.00619</b>	<b>&lt;0.025</b>	<0.00129
2,6-Dinitrotoluene	0.0013	0.003	<0.001	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<b>&lt;0.00755</b>	<b>&lt;0.00381</b>	<b>&lt;0.0154</b>	<0.000792
2-Chloronaphthalene	2	5.8	<0.002	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00755	<0.00381	<0.0154	<0.000792
2-Methylnaphthalene	0.098	0.29	<b>4.41</b>	<b>0.58</b>	<b>0.49</b>	<b>0.41</b>	<b>0.27</b>	<b>0.24</b>	<b>0.2</b>	<b>0.28</b>	<b>0.302</b>	<b>0.309</b>	<b>0.35</b>	<b>0.263</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<b>&lt;0.0025</b>	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<b>&lt;0.0783</b>	<b>&lt;0.0395</b>	<b>&lt;0.16</b>	<b>&lt;0.0822</b>
4-Nitrophenol	0.049	0.15	<0.0012	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<b>&lt;0.0528</b>	<0.0267	<b>&lt;0.108</b>	<0.00554
Acenaphthene	1.5	4.4	<b>3.17</b>	0.35	0.33	0.27	0.25	0.17	0.2	0.23	0.315	0.35	0.402	0.236
Acenaphthylene	1.5	4.4	<0.0015	0.0027	0.0025	0.0031	0.0025	0.0019	0.0022	0.0021	<0.00566	<0.00286	<0.0115	0.00335 J
Anthracene	7.3	22	0.0141	0.016	0.0095	0.017	0.017	0.0097	0.019	0.007	0.0183 J	<0.019 J	0.0247 J	0.0142
Benzo(a)anthracene	0.0091	0.02	<0.001	0.00028	0.0001 J	<0.00007	0.00016 J	<0.00005	0.000095 J	<0.00005	<0.00755	<0.00381	<b>&lt;0.0154</b>	<0.000792
Benzo(a)pyrene	0.0002	0.0002	<b>&lt;0.001</b>	0.0002 J	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<b>&lt;0.00755</b>	<b>&lt;0.00381</b>	<b>&lt;0.0154</b>	<b>&lt;0.000792</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019	<b>&lt;0.002</b>	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<b>&lt;0.0123</b>	<b>&lt;0.00619</b>	<b>&lt;0.025</b>	<b>&lt;0.00129</b>
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.001	0.00047	0.0035	<0.0002	0.00075	<0.00053	0.00033	<0.00016	<b>&lt;0.0349</b>	<b>&lt;0.0176</b>	<b>&lt;0.0712</b>	<0.00366
Chrysene	0.91	2	<0.001	0.00023	0.00011 J	<0.00007	0.00013 J	<0.00005	0.00011 J	<0.00005	<0.00755	<0.00381	<0.0154	<0.000792
Dibenzofuran	0.098	0.29	<b>2.13</b>	<b>0.25</b>	<b>0.17</b>	<b>0.22</b>	0.092	<b>0.13</b>	<b>0.15</b>	<b>0.17</b>	<b>0.206</b>	<b>0.242</b>	<b>0.252</b>	<b>0.178</b>
Di-n-butylphthalate (DBP)	2.4	7.3	<0.001	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0104	<0.00524	<0.0212	<0.00109
Fluoranthene	0.98	2.9	0.0067	0.0082	0.0067	0.0064	0.0068	0.0049	0.0042	0.0031	<0.0066	0.0104 J	<0.0135	0.00562
Fluorene	0.98	2.9	0.247	0.2	0.15	0.17	0.093	0.13	0.13	0.15	0.175	0.212	0.217	0.183
Naphthalene	0.49	1.5	<b>94.2</b>	<b>9.7</b>	<b>8</b>	<b>6.8</b>	<b>6.1</b>	<b>4</b>	<b>4.2</b>	<b>6</b>	<b>6.78</b>	<b>7.73</b>	<b>6.07</b>	<b>4.24</b>
Nitrobenzene	0.049	0.15	<0.002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0104	<0.00524	<0.0212	<0.00109
N-Nitrosodiphenylamine	0.19	0.42	<0.0012	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00943	<0.00476	<0.0192	<0.00099
Pentachlorophenol	0.001	0.001	<0.001	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<b>&lt;0.0575</b>	<b>&lt;0.029</b>	<b>&lt;0.117</b>	<b>&lt;0.00604</b>
Phenanthrene	0.73	2.2	0.177	0.16	0.12	0.15	0.083	0.11	0.08	0.1	0.137	0.158	0.197	0.111
Phenol	7.3	22	<0.001	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00377	<0.0019	<0.00769	<0.000396
Pyrene	0.73	2.2	0.0029	0.0043	0.0033	0.0035	0.0039	0.0021	0.0033	0.0019	<0.0104	<0.00524	<0.0212	0.00242 J
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

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	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-40B 01/24/2018	MW-40B 03/19/2018	MW-40B 05/16/2018	MW-40B 01/08/2019	MW-40B 07/11/2019	MW-40B 01/13/2020	MW-41B 01/31/2008 DNAPL	MW-41B 01/23/2020 DNAPL	MW-42B 01/30/2008	MW-42B 01/19/2010	MW-42B 07/14/2011	MW-42B 02/03/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.001	<0.0002	<0.0002	<0.0002	<0.00052	<0.0002	<0.00052	<0.0005	<0.001	<0.001
Benzene	0.005	0.005	<b>0.0091</b>	<b>0.0066</b>	<b>0.014</b>	<b>0.0063</b>	<b>0.0088</b>	<b>0.011</b>	<b>0.0103</b>	<b>0.012</b>	0.00117 J	<0.0005	<0.001	<0.001
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0015	<0.0003	<0.0003	<0.0003	<0.00047	<0.0003	<0.00047	<0.0005	<0.001	<0.001
Ethylbenzene	0.7	0.7	0.049	0.039	0.08	0.041	0.082	0.089	0.0508	0.066	0.00112 J	<0.0005	<0.0011	<0.0011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.005	<0.001	<0.001	<0.001	<0.00054	<0.001	<0.00054	<0.0005	<0.0013	<0.0013
Toluene	1	1	0.0081	0.0049	0.019	0.0048	0.014	0.022	0.0525	0.087	0.00181 J	<0.0005	<0.001	<0.001
Vinyl chloride	0.002	0.002								<0.0002				
Xylenes (total)	10	10	0.066	0.044	0.11	0.052	0.13	0.16	0.127	0.16	0.00377 J	<0.001	<0.0031	<0.0031
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0008	<0.000021	<0.00008	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	0.00034	<0.00004	<0.00004	0.0013	0.013	0.104	0.043	<0.00029	<0.00008	0.00013 J	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<b>&lt;0.0019</b>	<0.000058	<0.00019	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<b>&lt;0.0019</b>	<0.000042	<0.00019	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0038	<0.000021	<0.00039	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<b>0.13</b>	0.056	0.091	0.077	<b>0.18</b>	<b>0.34</b>	<b>0.305</b>	<b>0.67 J</b>	<0.00039	<0.00007	<0.00005	0.000089 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0019	<0.00002	<0.00019	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	0.0015	<0.000047	<0.000047	<0.000047	<0.0024	<0.000047	<0.00024	<0.00007	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.26	0.16	0.14	0.12	0.21	0.39	0.161	0.49	<0.00029	0.00021	<0.00024	0.0017
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	0.0014	0.00083	0.0021	0.0021	<0.0029	0.0089	<0.00029	<0.00007	<0.00005	<0.00005
Anthracene	7.3	22	0.016	0.0082	0.0087	0.007	0.011	0.04	0.0191	0.045	<0.00019	<0.00007	0.00036	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	0.000094 J	0.0036	0.004	<0.00019	<0.00007	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<b>&lt;0.0019</b>	<b>0.0009</b>	<0.00019	<0.00008	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<b>&lt;0.0038</b>	<0.00003	<0.00039	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00038	<0.000037	0.00012 J	0.000079 J	<0.000037	<0.000037	<0.0019	0.00032	0.00135 J	0.00028	<0.0008	<0.0001
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	0.000045 J	<0.000021	0.000051 J	0.003	0.0033	<0.00019	<0.00007	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	<b>0.16</b>	0.085	0.086	0.069	<b>0.13</b>	<b>0.29</b>	<b>0.142</b>	<b>0.39</b>	0.000699	0.00037	<0.00014	0.00016 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0019	<0.00002	<0.00019	<0.00007	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.0081	0.0037	0.0034	0.0041	0.0059	0.017	0.027	0.052	0.000697	0.00059	0.00024	0.00085
Fluorene	0.98	2.9	0.18	0.096	0.1	0.087	0.13	0.3	0.148	0.35	<0.00019	0.00016 J	<0.00026	0.0005
Naphthalene	0.49	1.5	<b>1.5</b>	<b>0.97</b>	<b>1.8</b>	<b>1.3</b>	<b>3.6</b>	<b>7.9</b>	<b>4.57</b>	<b>10</b>	0.000725	0.00035	<0.00048	<0.00063
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.0038	<0.000024	<0.00039	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.0024	<0.000025	<0.00024	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<b>&lt;0.0019</b>	<0.000079	<0.00019	<0.00008	<0.00005	<0.00005
Phenanthrene	0.73	2.2	0.14	0.078	0.1	0.068	0.093	0.22	0.161	0.41	<0.00019	<0.00007	<0.00016	0.00012 J
Phenol	7.3	22	<0.000035	<0.000035	0.00065	<0.000035	0.000053 J	<0.000035	0.00576	<0.000035	<0.00019	<0.00007	<0.00005	<0.00005
Pyrene	0.73	2.2	0.0036	0.0019	0.0016	0.002	0.0024	0.0037	0.017	0.029	0.000624	0.00035	0.00014 J	0.00044
<b>Metals</b>														
Arsenic	0.01	0.01	<b>0.0679</b>	<b>0.0606</b>	<b>0.0494</b>	<b>0.085</b>	<b>0.052</b>	<b>0.0523</b>		<b>0.0842 J</b>				

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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-42B 07/19/2012	MW-42B 02/05/2013	MW-42B 08/01/2013	MW-42B 01/15/2014	MW-42B 07/18/2014	MW-42B 01/24/2018	MW-42B 03/19/2018	MW-42B 05/16/2018	MW-42B 01/08/2019	MW-42B 07/11/2019	MW-42B 01/13/2020	MW-44A 01/30/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0025	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.001	<0.0002	<0.0002	<0.0002	<0.00052
Benzene	0.005	0.005	<0.0025	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.001	<0.0002	<0.0002	<0.0002	<b>0.00751</b>
Chlorobenzene	0.1	0.1	<0.0025	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0015	<0.0003	<0.0003	<0.0003	<0.00047
Ethylbenzene	0.7	0.7	<0.0025	<0.000132	<0.00011	<0.00019	0.000208 J	<0.0003	<0.0003	<0.0015	<0.0003	<0.0003	<0.0003	0.00363 J
Methylene chloride	0.005	0.005	<b>0.0097 J</b>	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.005	<0.001	<0.001	<0.001	<0.00054
Toluene	1	1	<0.0025	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.001	<0.0002	<0.0002	<0.0002	0.002 J
Vinyl chloride	0.002	0.002	<0.00011				<0.00011							
Xylenes (total)	10	10	<0.0075	<0.00026	<0.00026	<0.00058	0.000349 J	<0.0003	<0.0003	<0.0015	<0.0003	<0.0003	<0.0003	0.0186
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.000104	<0.000105	<0.000104	<0.000109	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.000292	<0.000295	0.000916 J	0.000577	<0.00004	<0.00004	<0.00041	<0.00004	<0.00004	<0.00004	<0.0003
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.000123	<0.000124	<0.000123	<0.000129	<0.000059	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.0002
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.0000755	<0.0000762	<0.0000755	<0.0000792	<0.000042	<0.000042	<b>0.012</b>	<0.000042	<0.000042	<0.000042	<0.0002
2-Chloronaphthalene	2	5.8	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000792	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0004
2-Methylnaphthalene	0.098	0.29	0.00015 J	0.000196 J	0.000141 J	0.000317 J	<0.0000693	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000069 J	0.0244
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.0000783	<0.000079	<0.0000783	<0.0000822	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002
4-Nitrophenol	0.049	0.15	<0.00005	<0.000528	<0.000533	<0.000528	<0.000554	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00025
Acenaphthene	1.5	4.4	0.00081	0.00036 J	<0.0000762	0.000355 J	<0.0000792	<0.000027	<0.000027	0.000067 J	<0.000027	0.000068 J	0.00012	0.127
Acenaphthylene	1.5	4.4	<0.00005	<0.0000566	<0.0000571	<0.0000566	<0.0000594	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.0003
Anthracene	7.3	22	<0.00005	0.000194 J	0.000122 J	0.000465 J	<0.0000495	0.000025 J	<0.000014	0.000019 J	<0.000014	0.000064 J	0.000049 J	0.00195
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000792	<0.000051	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.0002
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.000123	<0.000124	<0.000123	<0.000129	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.0004
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00022	<0.000349	<0.000352	<0.000349	0.000513	0.00025	0.00015 J	0.00021	0.000061 J	0.00015 J	0.00013 J	0.00097 J
Chrysene	0.91	2	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000792	0.000038 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0002
Dibenzofuran	0.098	0.29	0.000066 J	0.000217 J	0.000131 J	0.000205 J	<0.0000792	<0.00002	<0.00002	0.000039 J	<0.00002	0.000038 J	0.000065 J	0.0642
Di-n-butylphthalate (DBP)	2.4	7.3	0.000062 J	<0.000104	0.000108 J	<0.000104	<0.000109	0.00006 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002
Fluoranthene	0.98	2.9	0.00041	0.000641	0.000294 J	0.000339 J	<0.0000693	0.00024	0.00014	0.000016 J	0.00011	0.00004 J	0.00018	0.00269
Fluorene	0.98	2.9	0.00016 J	<0.000066	0.000134 J	0.000198 J	<0.0000693	<0.00003	<0.00003	0.000047 J	<0.00003	0.000057 J	0.000067 J	0.045
Naphthalene	0.49	1.5	0.0019	0.000476	0.00288 J	0.00242 J	0.000426 J	<0.00002	0.00049	<0.00061	<0.00002	<0.00043	<0.00076	<b>0.816</b>
Nitrobenzene	0.049	0.15	<0.00005	<0.000104	<0.000105	<0.000104	<0.000109	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.0004
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.0000943	<0.0000952	<0.0000943	<0.000099	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00025
Pentachlorophenol	0.001	0.001	<0.00005	<0.000575	<0.000581	<0.000575	<0.000604	<0.00008	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.0002
Phenanthrene	0.73	2.2	<0.00005	0.000356 J	0.000122 J	0.000501	<0.0000594	0.000077 J	<0.000021	0.000039 J	<0.000021	0.000034 J	0.00012	0.0161
Phenol	7.3	22	<0.00005	<0.0000377	<0.0000381	<0.0000377	0.000801	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.0002
Pyrene	0.73	2.2	0.00023	0.000369 J	0.000127 J	0.000234 J	<0.000109	0.00023	0.00014	<0.000019	0.0001	0.000033 J	0.00013	0.00159
<b>Metals</b>														
Arsenic	0.01	0.01						0.00186 J	0.00108 J	0.00112 J	0.00216	0.0022	0.00133 J	

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray.
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-44A 07/14/2008	MW-44A 02/03/2009	MW-44A 01/13/2010	MW-44A 06/30/2010	MW-44A 01/26/2011	MW-44A 07/20/2011	MW-44A 02/15/2012	MW-44A 07/25/2012	MW-44A 02/12/2013	MW-44A 08/05/2013	MW-44A 01/17/2014	MW-44A 08/28/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014
Benzene	0.005	0.005	<b>0.00635</b>	0.003 J	<0.0005	0.0026 J	<0.0005	0.002 J	0.0042 J	0.0044 J	0.00206	<b>0.00849</b>	<b>0.00727</b>	0.0042
Chlorobenzene	0.1	0.1	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012
Ethylbenzene	0.7	0.7	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	0.000624 J	0.00172	0.00067	0.000344 J
Methylene chloride	0.005	0.005	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015
Toluene	1	1	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	0.000252 J	<0.000705	0.000418 J	0.000329 J
Vinyl chloride	0.002	0.002	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00018	<0.00011
Xylenes (total)	10	10	0.006 J	0.0013 J	<0.001	0.0026 J	<0.001	<0.0031	0.0052 J	0.0033 J	0.00469	0.0207	0.00805	0.00561
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00105	<0.0011	<0.000106	<0.000108
2,4-Dimethylphenol	0.49	1.5	<0.00032	<0.00008	<0.00008	<0.00008	0.00081	<0.00005	<0.00005	<0.00005	<0.00295	<0.0031	<0.000298	<0.000304
2,4-Dinitrotoluene	0.0013	0.003	<0.00021	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00124	<0.0013	<0.000125	<0.000127
2,6-Dinitrotoluene	0.0013	0.003	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.000762	<0.0008	<0.0000769	<0.0000784
2-Chloronaphthalene	2	5.8	<0.00042	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000762	<0.0008	<0.0000769	<0.0000784
2-Methylnaphthalene	0.098	0.29	0.00779	0.00097	0.00012 J	0.004 J	<0.00007	0.0023	0.0048	0.0095 J	<0.000667	<b>0.109</b>	0.0106	0.00902
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<b>&lt;0.0079</b>	<b>&lt;0.0083</b>	<0.000798	<0.000814
4-Nitrophenol	0.049	0.15	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00533	<0.0056	<0.000538	<0.000549
Acenaphthene	1.5	4.4	0.202	0.12	0.13	0.2	0.023	0.23	0.21	0.22	0.07	0.546	0.394	0.197
Acenaphthylene	1.5	4.4	<0.00032	0.0012	0.00079	0.00096 J	0.0014	0.0013	0.001	0.0013	0.00276 J	<0.0006	<0.0000577	0.0014
Anthracene	7.3	22	0.00393	0.0046	0.0077	0.0067 J	0.00055	0.0058	0.0068	0.004	<0.000476	0.017	0.017	0.00868
Benzo(a)anthracene	0.0091	0.02	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000762	<0.0008	<0.0000769	<0.0000784
Benzo(a)pyrene	0.0002	0.0002	<b>&lt;0.00021</b>	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<b>&lt;0.000762</b>	<b>&lt;0.0008</b>	<0.0000769	<0.0000784
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00042	<0.00009	<0.00009	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005	<b>&lt;0.00124</b>	<b>&lt;0.0013</b>	<0.000125	0.00014 J
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00021	0.00043	<0.00031	<0.00028	0.00048	<0.00075	<0.00011	<0.0001	<0.00352	<0.0037	<0.000356	<0.000363
Chrysene	0.91	2	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000762	<0.0008	<0.0000769	<0.0000784
Dibenzofuran	0.098	0.29	<b>0.125</b>	0.054	0.0087	0.0043 J	0.00072	0.0014	0.001	0.0031	<0.000762	<b>0.135</b>	<b>0.107</b>	0.0159
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00105	<0.0011	<0.000106	<0.000108
Fluoranthene	0.98	2.9	0.00367	0.0032	0.0056	0.006 J	0.0014	0.0095	0.0094	0.0065	0.00257 J	0.0137	0.0154	0.00749
Fluorene	0.98	2.9	0.0865	0.056	0.069	0.097	0.00027	0.094	0.11	0.091 J	0.00495	0.172	0.178	0.0987
Naphthalene	0.49	1.5	0.287	0.021	<0.0011	0.16	0.00035	0.042	0.32	0.39	<0.000941	<b>1.72</b>	0.235 J	0.0163
Nitrobenzene	0.049	0.15	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00105	<0.0011	<0.000106	0.0012
N-Nitrosodiphenylamine	0.19	0.42	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	0.00065	<0.00005	<0.000952	<0.001	<0.0000962	<0.000098
Pentachlorophenol	0.001	0.001	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<b>&lt;0.00581</b>	<b>&lt;0.0061</b>	<0.000587	<0.000598
Phenanthrene	0.73	2.2	0.0184	0.02	0.0055	0.0025 J	<0.00007	0.0047	0.0073	0.0064	<0.000571	0.0416	0.0438	0.0217
Phenol	7.3	22	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.000062 J	<0.000381	<0.0004	<0.0000385	<0.0000392
Pyrene	0.73	2.2	0.00156	0.0016	0.0032	0.003 J	0.001	0.0046	0.0054	0.0038	0.00139 J	0.00732	0.0083	0.0041
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-44A 01/31/2018	MW-44A 03/26/2018	MW-44A 06/01/2018	MW-44A 01/22/2019	MW-44A 07/17/2019	MW-44A 01/09/2020	MW-44C 01/29/2008 DNAPL	MW-44C 07/20/2011 DNAPL	MW-44C 07/18/2012 DNAPL	MW-44C 02/06/2013 DNAPL	MW-44C 01/15/2020	MW-45C 01/29/2008 DNAPL
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.001	<0.005	<0.0014	<0.0002	<0.00052
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.001	<0.005	0.000964 J	<0.0002	0.0448
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.001	<0.005	0.00293 J	<0.0003	<0.00047
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0204	<0.0011	0.32	0.233	<0.0003	0.111
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.0013	<0.01	<0.0015	<0.001	<0.00054
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0117	<0.001	0.16	0.0895	<0.0002	0.112
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002						
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0487	<0.0031	0.84	0.688	<0.0003	0.298
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0002	<0.00005	<0.00075	<0.529	<0.000021	<0.002
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004	0.00219 J	<0.00005	<0.00075	<1.49	0.00011 J	0.0104
2,4-Dinitrotoluene	0.0013	0.003	0.00017 J	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00039	<0.00005	<0.00075	<0.625	<0.000058	<0.004
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.00039	<0.00006	<0.0009	<0.385	<0.000042	<0.004
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00078	<0.00005	<0.00075	<0.385	<0.000021	<0.008
2-Methylnaphthalene	0.098	0.29	<0.000093	<0.000019	<0.000019	<0.000019	<0.000019	<0.00035	0.00551	<0.00005	62	1.15 J	0.00012	1.01
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00039	<0.00008	<0.0012	<3.99	<0.00002	<0.004
4-Nitrophenol	0.049	0.15	<0.000047	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047	<0.00049	<0.00005	<0.00075	<2.69	<0.000047	<0.005
Acenaphthene	1.5	4.4	0.062	0.064	0.09	0.037	0.056	0.064	0.0531	0.00012 J	31	0.632 J	0.00045	0.472
Acenaphthylene	1.5	4.4	0.00078	0.0006	<0.000015	0.00036	0.00028	0.00097	0.00083	0.000097 J	0.29	<0.288	0.00065	0.0085
Anthracene	7.3	22	0.00066	0.0024	0.0022	0.00044	0.00089	0.00027	0.0131	0.00014 J	19	<0.24	0.0086	0.18
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.000051	<0.00005	<0.00005	0.000085 J	0.000054 J	0.00327	0.00017 J	3.5	<0.385	0.0018	0.0416
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00039	0.00022	0.87	<0.385	0.00065	<0.004
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.00078	<0.00005	<0.00075	<0.625	<0.00003	<0.008
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0001	<0.000038	0.0001 J	<0.000037	<0.00025	<0.000037	0.00235 J	<0.00087	0.013	<1.78	0.0012	<0.004
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	0.000026 J	0.000052 J	<0.000021	0.00278	0.00032	3.3	<0.385	0.0022	0.0347
Dibenzofuran	0.098	0.29	0.00019	0.00029	<0.00002	<0.00002	<0.00002	<0.00033	0.046	<0.00005	38	0.453 J	0.00034	0.488
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00039	<0.00005	<0.00075	<0.529	0.0016	<0.004
Fluoranthene	0.98	2.9	0.0067	0.0095	0.012	0.0058	0.0091	0.0059	0.0264	0.00016 J	28	<0.337	0.0095	0.384
Fluorene	0.98	2.9	0.024	0.03	0.039	0.0097	0.014	0.0078	0.0346	<0.00005	26	<0.337	0.00067	0.357
Naphthalene	0.49	1.5	<0.00028	0.00036	0.00038	0.00011	<0.00019	<0.003	<0.00078	<0.00016	230	18 J	0.000097 J	6.05
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00078	<0.00005	<0.00075	<0.529	<0.000024	<0.008
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.00049	<0.00005	<0.00075	<0.481	<0.000025	<0.005
Pentachlorophenol	0.001	0.001	<0.000079	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.00039	<0.00005	<0.00075	<2.93	<0.000079	<0.004
Phenanthrene	0.73	2.2	0.00012	0.00018	0.0002	<0.000021	0.00012	<0.00022	0.0668	0.000081 J	88	0.498 J	0.0077	0.903
Phenol	7.3	22	<0.000035	<0.000036	<0.000035	<0.000035	<0.000035	<0.00013	<0.00039	<0.00005	<0.00075	<0.192	<0.000035	<0.004
Pyrene	0.73	2.2	0.0037	0.0069	0.0073	0.0033	0.0057	0.0045	0.0159	0.00013 J	19	<0.529	0.006	0.234
<b>Metals</b>														
Arsenic	0.01	0.01	0.0275	0.0169	0.0165	0.0101	0.0303	0.00966					0.00314	

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-45C 01/20/2020	MW-46C 01/30/2008 DNAPL	MW-46C 01/15/2020	MW-47A 03/20/2020	MW-47C 07/14/2008	MW-47C 02/04/2009	MW-47C 01/20/2010	MW-47C 06/24/2010	MW-47C 01/19/2011	MW-47C 07/21/2011	MW-47C 02/07/2012	MW-47C 07/27/2012
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.00052	<0.0002	<0.0002	<0.00109	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Benzene	0.005	0.005	0.0005 J	<b>0.0222</b>	<b>0.028</b>	<0.0002	<0.00112	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Chlorobenzene	0.1	0.1	<0.0003	<0.00047	<0.0003	<0.0003	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Ethylbenzene	0.7	0.7	0.00044 J	0.0249	0.051	<0.0003	<0.00142	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005
Methylene chloride	0.005	0.005	<0.001	<0.00054	<0.001	<0.001	<0.00122	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001
Toluene	1	1	<0.0002	0.0167	0.0047	<0.0002	<0.00138	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005
Vinyl chloride	0.002	0.002	<0.0002											
Xylenes (total)	10	10	0.0051	0.0377	0.11	<0.0003	<0.00302	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000085	<0.0004	<0.000021	<0.000021	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.0014	0.00016 J	<0.00004	<0.00032	<0.00008	<0.00008	0.00011 J	<0.00008	<0.00005	<0.00005	0.00042
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.00096	<0.000058	<0.000058	<0.00021	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<b>0.0022 J</b>	<0.00096	<0.000042	<0.000042	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.0019	<0.000021	<0.000021	<0.00042	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.00013	0.0825	<b>0.2</b>	<0.000019	<0.00042	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.0044	<0.00005
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00096	<0.00002	<0.00002	<0.00021	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	0.000067 J	<0.0012	<0.000047	<0.000047	<0.00026	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.000084 J	0.0635	0.16	<0.000027	<0.00032	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	0.00017 J	0.000058 J
Acenaphthylene	1.5	4.4	0.000037 J	<0.0014	0.0016	<0.000015	<0.00032	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Anthracene	7.3	22	0.0009	0.0222	0.073	<0.000014	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.000074 J	<0.00005
Benzo(a)anthracene	0.0091	0.02	0.00039	0.00597	0.0071	<0.00005	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<b>0.00042</b>	<b>0.00215</b>	<b>0.0021</b>	<0.00002	<b>&lt;0.00021</b>	<0.00008	0.000099 J	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<b>&lt;0.0019</b>	<0.00003	<0.00003	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000067 J	<0.00096	0.0011	0.00011 J	<0.00021	0.0036	<0.00065	<0.00021	<0.0002	0.0001 J	<b>0.011</b>	<0.0001
Chrysene	0.91	2	0.00062	0.00521	0.0068	<0.000021	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	0.000075 J	0.0636	<b>0.16</b>	<0.00002	<0.00032	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00096	<0.00002	0.000076 J	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00015 J	<0.00005
Fluoranthene	0.98	2.9	0.00062	0.0426	0.084	<0.00001	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005
Fluorene	0.98	2.9	0.00023	0.048	0.12	<0.00003	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00025	<0.00005
Naphthalene	0.49	1.5	<0.00061	<b>1.1</b>	<b>1.8</b>	<0.00002	<0.00042	0.00019 J	<0.0001	0.00046	0.00021	<0.00005	0.0041	0.00046
Nitrobenzene	0.049	0.15	<0.000024	<0.0019	<0.000024	<0.000024	<0.00042	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.0012	<0.000025	<0.000025	<0.00026	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.000079	<0.00096	<0.000079	<0.000079	<0.00021	<0.00008	0.0004	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005
Phenanthrene	0.73	2.2	<0.00027	0.119	0.24	<0.000021	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.0003	<0.00005
Phenol	7.3	22	<0.000035	0.0133	<0.000035	<0.000035	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00044	0.00056
Pyrene	0.73	2.2	0.00068	0.0252	0.053	<0.000019	<0.00021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00015 J	<0.00005
<b>Metals</b>														
Arsenic	0.01	0.01	0.00073 J		0.00272	0.000566 J								

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-47C 02/07/2013	MW-47C 08/06/2013	MW-47C 01/17/2014	MW-47C 07/30/2014	MW-47C 01/23/2019	MW-47C 07/17/2019	MW-47C 01/16/2020	MW-48C 01/29/2008	MW-48C 01/29/2008	MW-48C 07/14/2008	MW-48C 02/04/2009	MW-48C 01/21/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.00109	<0.0005	<0.0005
Benzene	0.005	0.005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025	<0.00112	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0015	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.00025	<0.00025	<0.00142	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.00122	<0.0005	<0.0005
Toluene	1	1	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.00041	<0.00041	<0.00138	<0.0005	<0.0005
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.00127	<0.00127	<0.00302	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000105	<0.00011	R	<0.000104	<0.000021	<0.000021	<0.000021	<0.00008	<0.00008	<0.00009	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	R	<0.00031	R	<0.000292	0.000095 J	<0.00004	0.0002	<0.00029	<0.00029	<0.00033	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000124	<0.00013	R	<0.000123	<0.000058	<0.000058	<0.000058	<0.00019	<0.00019	<0.00022	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0000762	<0.00008	R	<0.0000755	<0.000042	<0.000042	<0.000042	<0.00019	<0.00019	<0.00022	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0000762	<0.00008	R	<0.0000755	<0.000021	<0.000021	<0.000021	<0.00038	<0.00038	<0.00044	<0.00012	<0.0001
2-Methylnaphthalene	0.098	0.29	0.000098 J	<0.00007	R	<0.000066	<0.000019	<0.000019	<0.0001	<0.00038	<0.00038	<0.00044	<0.00007	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	R	<0.00083	R	<0.000783	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	<0.00022	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000533	<0.00056	R	<0.000528	<0.000047	<0.000047	<0.000047	<0.00024	<0.00024	<0.00028	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.0000762	<0.00008	R	<0.0000755	<0.000027	<0.000027	<0.000057	<0.00029	<0.00029	<0.00033	<0.00009	<0.00009
Acenaphthylene	1.5	4.4	<0.0000571	<0.00006	R	<0.0000566	<0.000015	<0.000015	<0.000015	<0.00029	<0.00029	<0.00033	<0.00006	<0.00007
Anthracene	7.3	22	0.000107 J	<0.00005	R	<0.0000472	<0.000014	<0.000014	<0.000014	0.000584	0.000589	<0.00022	0.00012 J	<0.00007
Benzo(a)anthracene	0.0091	0.02	0.000137 J	<0.00008	R	<0.0000755	<0.000005	<0.000005	<0.000005	<0.00019	<0.00019	<0.00022	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	0.000306 J	<0.00008	R	<0.0000755	<0.000002	<0.000002	0.00004 J	<0.00019	<0.00019	<0.00022	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000124	<0.00013	R	<0.000123	<0.000003	<0.000003	<0.00003	<0.00038	<0.00038	<0.00044	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000594	<0.00037	R	<0.000349	<0.000056	<0.000027	0.00079	0.00045 J	<0.00019	0.00028 J	0.00034	<0.0018
Chrysene	0.91	2	0.000127 J	<0.00008	R	<0.0000755	<0.000021	<0.000021	0.00006 J	<0.00019	<0.00019	<0.00022	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	0.000104 J	<0.00008	R	<0.0000755	0.000034 J	<0.00002	<0.000049	0.00031 J	<0.00029	<0.00033	0.00025	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000105	<0.00011	R	<0.000104	<0.00002	<0.00002	0.0014	<0.00019	<0.00019	<0.00022	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.000289 J	0.000186 J	0.0000718 J	<0.000066	0.000028 J	<0.00001	<0.00011	0.00047 J	0.000687	0.00033 J	<0.00007	0.00013 J
Fluorene	0.98	2.9	0.000116 J	<0.00007	R	<0.000066	<0.00003	<0.00003	<0.000042	<0.00019	<0.00019	<0.00022	<0.00007	<0.00007
Naphthalene	0.49	1.5	<0.000401	<0.00008	0.000297 J	<0.0000755	0.00083	<0.00002	<0.00056	0.00119	0.00062	<0.00044	0.00052	0.0002 J
Nitrobenzene	0.049	0.15	<0.000105	<0.00011	R	<0.000104	<0.000024	<0.000024	<0.000024	<0.00038	<0.00038	<0.00044	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0000952	<0.0001	R	<0.0000943	<0.000025	<0.000025	<0.000025	<0.00024	<0.00024	<0.00028	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	R	<0.00061	R	<0.000575	<0.000079	<0.000079	<0.000079	<0.00019	<0.00019	<0.00022	<0.00008	<0.00008
Phenanthrene	0.73	2.2	0.000406 J	<0.00006	0.000185 J	<0.0000566	0.000052 J	<0.000021	<0.000088	0.00046 J	0.00043 J	<0.00022	0.00032	<0.00007
Phenol	7.3	22	R	<0.00004	R	<0.0000377	<0.000035	<0.000035	<0.000035	<0.00019	<0.00019	<0.00022	<0.00007	<0.00007
Pyrene	0.73	2.2	0.000388 J	0.000131 J	R	<0.000104	0.000021 J	<0.000019	<0.0001	0.00039 J	0.000528	<0.00022	<0.00007	0.0001 J
<b>Metals</b>														
Arsenic	0.01	0.01					<0.0004	0.00044 J	0.00234					

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-48C 06/24/2010	MW-48C 07/15/2010	MW-48C 01/19/2011	MW-48C 07/18/2011	MW-48C 02/06/2012	MW-48C 07/24/2012	MW-48C 01/31/2013	MW-48C 08/01/2013	MW-48C 01/16/2014	MW-48C 07/16/2014	MW-48C 01/28/2018	MW-48C 03/20/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002
Vinyl chloride	0.002	0.002									<0.00018			
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0005	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	<0.000104	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.0073	<0.00008	<0.00008	<0.00005	<0.00005	0.00014 J	<0.000292	<0.000295	<0.000292	<0.000292	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124	<0.000123	<0.000123	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00035	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0005	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.18	<0.00007	<0.00007	<0.00005	<0.00005	0.0013	<0.0000667	<0.0000667	<0.000066	<0.000066	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0004	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000783	<0.000079	<0.000783	<0.000783	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00035	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000528	<0.000533	<0.000528	<0.000528	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.073	<0.00009	<0.00009	<0.00005	<0.00005	0.0011	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	0.014	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000566	<0.0000571	<0.0000566	<0.0000566	<0.000015	<0.000015
Anthracene	7.3	22	0.007	<0.00007	<0.00007	<0.00005	<0.00005	0.00077	<0.0000472	<0.0000476	<0.0000472	<0.0000472	0.000025 J	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00035	<0.00007	<0.00007	<0.00005	<0.00005	0.000066 J	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.00005	0.000057 J
Benzo(a)pyrene	0.0002	0.0002	<0.0004	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124	<0.000123	<0.000123	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.001	0.0013	0.001	<0.00043	<0.0001	<0.00024	<0.000349	<0.000352	<0.000349	<0.000349	0.000079 J	0.00011 J
Chrysene	0.91	2	<0.00035	<0.00007	<0.00007	<0.00005	<0.00005	0.000073 J	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.000021	0.000044 J
Dibenzofuran	0.098	0.29	0.065	<0.00008	<0.00008	<0.00005	<0.00005	0.00096	<0.0000755	<0.0000762	<0.0000755	<0.0000755	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00035	<0.00007	<0.00007	<0.00005	<0.00005	0.000053 J	<0.000104	<0.000105	<0.000104	<0.000104	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0021	0.00019 J	0.00019 J	<0.00013	<0.00005	0.00095	<0.000066	0.000134 J	0.000153 J	<0.000066	0.000049 J	0.000099 J
Fluorene	0.98	2.9	0.032	<0.00007	<0.00007	<0.00005	<0.00005	0.0011	<0.000066	<0.0000667	<0.000066	<0.000066	<0.00003	<0.00003
Naphthalene	0.49	1.5	5	<0.0001	<0.0001	<0.00005	<0.00005	0.0071	<0.000495	<0.000158	<0.0000755	<0.0000755	<0.00002	<0.00002
Nitrobenzene	0.049	0.15	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	<0.000104	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000952	<0.0000943	<0.0000943	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	0.019	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000575	<0.000581	<0.000575	<0.000575	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.03	<0.00007	<0.00007	<0.00005	<0.00005	0.0034	<0.0000566	<0.0000571	<0.0000566	<0.0000566	<0.000022	<0.000021
Phenol	7.3	22	0.024	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000377	<0.0000381	<0.0000377	<0.0000377	<0.000035	<0.000035
Pyrene	0.73	2.2	0.001	0.00015 J	0.00012 J	<0.0001	<0.00005	0.00052	<0.000104	<0.000105	<0.000104	<0.000104	0.000052 J	0.000087 J
<b>Metals</b>														
Arsenic	0.01	0.01											0.000831 J	0.000581 J

Notes:  
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2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
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ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-48C 05/24/2018	MW-48C 01/10/2019	MW-48C 07/17/2019	MW-48C 01/16/2020	MW-49A 01/31/2008	MW-49A 07/15/2008	MW-49A 02/04/2009	MW-49A 01/21/2010	MW-49A 06/25/2010	MW-49A 01/20/2011	MW-49A 07/22/2011	MW-49A 02/07/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005	<0.005	<0.005	<0.01	<0.05
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	0.0108	0.165	0.24	0.2	0.29	0.057	0.2	<0.05
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	0.00865	0.00702	0.0053	0.0024 J	<0.005	0.0084 J	<0.01	<0.05
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	0.0238	0.0837	0.084	0.085	0.14	0.04 J	0.094	<0.055
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005	<0.005	<0.005	<0.013	<0.065
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	0.00805	0.0415	0.077	0.083	0.13	0.021 J	0.11	<0.05
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	0.0352	0.187	0.2	0.21	0.34	0.079 J	0.2	<0.16
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.004	<0.002	<0.0001	<0.0001	<0.0005	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00004	0.001	<0.00004	<0.00004	0.025	6.08	6.8	0.86	3.7	0.18	3	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.0095	<0.004	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.0095	<0.004	<0.00007	<0.00007	<0.00035	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.019	<0.008	<0.00012	<0.0001	<0.0005	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.000019	0.00038	<0.000019	<0.000076	0.0693	0.492	0.6	0.35	0.44	0.13	0.27	<0.00005
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.0095	<0.01	<0.00008	<0.00008	<0.0004	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.012	<0.005	<0.00007	<0.00007	<0.00035	<0.00007	<0.00005	<0.00005
Acenaphthene	1.5	4.4	<0.000027	0.000098 J	<0.000027	<0.000073	0.215	0.468	0.32	0.2	0.21	0.13	0.13	<0.00005
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.014	<0.006	0.0039	0.0032	0.0052	0.0018	0.0029	<0.00005
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.0095	0.0164	0.01	0.0071	0.0099	0.0096	0.011	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.0095	<0.004	0.00066	<0.00007	<0.00035	<0.00007	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	0.000021 J	<0.0095	<0.004	0.00024	<0.00008	<0.0004	<0.00008	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.019	<0.008	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000063 J	<0.000037	<0.0003	0.000039 J	<0.0095	<0.004	0.0009	<0.0015	<0.001	<0.00029	<0.0001	<0.0001
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	0.000029 J	<0.0095	<0.004	0.0006	<0.00007	<0.00035	<0.00007	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.000054	0.148	0.293	0.21	0.14	0.16	0.075	0.09	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.0095	<0.004	<0.00007	<0.00007	<0.00035	<0.00007	<0.00005	<0.00005
Fluoranthene	0.98	2.9	<0.00001	<0.00001	0.000035 J	<0.000061	<0.0095	0.0063	0.0058	0.0025	0.0034	0.0038	0.0033	<0.00005
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00005	0.102	0.205	0.15	0.11	0.13	0.073	0.092	<0.00005
Naphthalene	0.49	1.5	<0.00002	0.0085	<0.00002	<0.000052	2.13	11	9	5.1	10	1.8	7.4	<0.00005
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.019	<0.008	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.012	<0.005	<0.00009	<0.00009	<0.00045	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.0095	<0.004	<0.00008	<0.00008	<0.0004	<0.00008	<0.00005	<0.00005
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000052	0.0939	0.147	0.096	0.072	0.086	0.062	0.07	<0.00005
Phenol	7.3	22	<0.000035	0.002	0.000074 J	<0.000035	<0.0095	0.0111	<0.00007	0.00077	0.0011	0.0058	0.0095	<0.00005
Pyrene	0.73	2.2	<0.000019	<0.000019	0.000036 J	<0.000065	<0.0095	<0.004	0.0046	0.0017	0.0018	0.002	0.0016	<0.00005
<b>Metals</b>														
Arsenic	0.01	0.01	0.000562 J	0.000924 J	0.00167 J	0.00126 J								

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-49A 07/26/2012	MW-49A 02/07/2013	MW-49A 08/01/2013	MW-49A 01/16/2014	MW-49A 07/16/2014	MW-49A 01/29/2018	MW-49A 04/01/2018	MW-49A 05/31/2018	MW-49A 01/23/2019	MW-49A 07/31/2019	MW-49A 01/07/2020	MW-50A 01/31/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.00014	<0.0014	<0.0002	<0.0028	<0.0002	<0.001	<0.0002	<0.0002	<0.0002	<0.002	<0.00052
Benzene	0.005	0.005	<b>0.042</b>	<b>0.114</b>	<b>0.094</b>	<b>0.0565</b>	<b>0.108</b>	<b>0.013</b>	<b>0.016</b>	<b>0.01</b>	0.004	<0.0002	<b>0.3</b>	<0.00025
Chlorobenzene	0.1	0.1	0.0037 J	<b>0.299</b>	<b>0.476</b>	<b>0.304</b>	<b>0.211</b>	<0.0003	<0.0015	<0.0003	<0.0003	<0.0003	0.0063 J	<0.00047
Ethylbenzene	0.7	0.7	0.037	0.0321	0.0499	0.0331	0.0701	0.01	0.01	0.0067	0.0031	<0.0003	0.1	<0.00025
Methylene chloride	0.005	0.005	<0.001	<0.00015	<0.0015	<0.00022	<b>0.0212</b>	<0.001	<0.005	<0.001	<0.001	<0.001	<b>&lt;0.01</b>	<0.00054
Toluene	1	1	0.031	0.0343	0.0347	0.0296	0.0593	0.003	0.0075	0.0065	0.0023	<0.0002	0.19	<0.00041
Vinyl chloride	0.002	0.002	<0.0005	<0.00011	<0.0011	<0.00018	<b>&lt;0.0022</b>	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.002	<0.00041
Xylenes (total)	10	10	0.082	0.0777	0.106	0.0699	0.157	0.023	0.023	0.015	0.0087	<0.0003	0.24	<0.00127
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<b>&lt;0.00524</b>	<b>&lt;0.0105</b>	<b>&lt;0.00519</b>	<b>&lt;0.0104</b>	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008
2,4-Dimethylphenol	0.49	1.5	0.037	<b>1.42</b>	<b>0.903</b>	<b>2.1 J</b>	<b>1.23</b>	0.097	0.033	0.097	<0.00004	<0.00004	<b>5.7</b>	<0.00029
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<b>&lt;0.00619</b>	<b>&lt;0.0124</b>	<b>&lt;0.00613</b>	<b>&lt;0.0123</b>	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<b>&lt;0.00381</b>	<b>&lt;0.00762</b>	<b>&lt;0.00377</b>	<b>&lt;0.00755</b>	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00019
2-Chloronaphthalene	2	5.8	<0.00005	<0.00381	<0.00762	<0.00377	<0.00755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038
2-Methylnaphthalene	0.098	0.29	<0.00005	<b>0.218</b>	<b>0.216</b>	<b>0.267</b>	<b>0.293</b>	0.000078 J	0.000067 J	0.0079	<0.00019	<0.00019	<b>0.13</b>	<0.00038
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<b>&lt;0.0395</b>	<b>&lt;0.079</b>	<b>&lt;0.0392</b>	<b>&lt;0.0783</b>	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
4-Nitrophenol	0.049	0.15	<0.00005	<0.0267	<b>&lt;0.0533</b>	<0.0264	<b>&lt;0.0528</b>	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024
Acenaphthene	1.5	4.4	0.033	0.134	0.126	0.18	0.126	0.0036	0.0049	0.007	<0.000027	<0.000027	0.057	<0.00029
Acenaphthylene	1.5	4.4	0.00062	<0.00286	<0.00571	0.00528 J	<0.00566	0.00012	0.00012	0.00015	<0.000015	<0.000015	0.00097	<0.00029
Anthracene	7.3	22	0.00076	0.00824 J	0.0119 J	0.0132 J	<0.00472	0.00028	0.00035	0.00062	<0.000014	<0.000014	0.0028	<0.00019
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00381	<0.00762	<0.00377	<0.00755	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00019
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<b>&lt;0.00381</b>	<b>&lt;0.00762</b>	<b>&lt;0.00377</b>	<b>&lt;0.00755</b>	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<b>&lt;0.00619</b>	<b>&lt;0.0124</b>	<b>&lt;0.00613</b>	<b>&lt;0.0123</b>	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0001 J	<b>&lt;0.0176</b>	<b>&lt;0.0352</b>	<b>&lt;0.0175</b>	<b>&lt;0.0349</b>	0.00013 J	0.00014 J	0.0002 J	<0.000055	<0.000037	<0.00011	<0.00019
Chrysene	0.91	2	<0.00005	<0.00381	<0.00762	<0.00377	<0.00755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00019
Dibenzofuran	0.098	0.29	0.0099	0.0851	0.0812	0.0902	0.0941	0.0021	0.0029	0.0049	<0.00002	<0.00002	0.037	<0.00029
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00524	<0.0105	<0.00519	<0.0104	0.000069 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
Fluoranthene	0.98	2.9	0.0018	<0.00333	<0.00667	0.00521 J	<0.0066	0.00035	0.00039	0.00049	<0.00001	<0.00001	0.00056	<0.00019
Fluorene	0.98	2.9	0.015 J	0.0717	0.0662	0.0864	0.0651	0.0025	0.0027	0.0039	<0.00003	<0.00003	0.027	<0.00019
Naphthalene	0.49	1.5	<0.00005	<b>2.88</b>	<b>3.3</b>	<b>5.86</b>	<b>5.13</b>	<0.00016	0.0002	0.046	0.000089 J	<0.00002	<b>4.5</b>	<0.00038
Nitrobenzene	0.049	0.15	<0.00005	<0.00524	<0.0105	<0.00519	<0.0104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00024	<0.00038
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00476	<0.00952	<0.00472	<0.00943	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00024
Pentachlorophenol	0.001	0.001	<0.00005	<b>&lt;0.029</b>	<b>&lt;0.0581</b>	<b>&lt;0.0288</b>	<b>&lt;0.0575</b>	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.00019
Phenanthrene	0.73	2.2	0.002	0.0455	0.0684	0.0564	0.0519	0.00066	0.00055	0.0034	<0.000021	<0.000021	0.016	0.00026 J
Phenol	7.3	22	<0.00023	<0.0019	<0.00381	<0.00189	<0.00377	<0.000035	<0.000035	0.0001 J	<0.000035	0.00064	0.17	<0.00019
Pyrene	0.73	2.2	0.00095	<0.00524	<0.0105	<0.00519	<0.0104	0.00018	0.00029	0.00034	<0.000019	<0.000019	0.00052	<0.00019
<b>Metals</b>														
Arsenic	0.01	0.01						0.00163 J	0.00233	0.000922 J	0.0012 J	0.000658 J	0.00463	

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-50A 07/16/2008	MW-50A 02/04/2009	MW-50A 01/20/2010	MW-50A 06/25/2010	MW-50A 01/27/2011	MW-50A 07/28/2011	MW-50A 02/09/2012	MW-50A 07/24/2012	MW-50A 04/02/2013	MW-50A 08/09/2013	MW-50A 01/29/2014	MW-50A 08/28/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00109	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014
Benzene	0.005	0.005	<0.00112	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008
Chlorobenzene	0.1	0.1	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012
Ethylbenzene	0.7	0.7	<0.00142	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011
Methylene chloride	0.005	0.005	<0.00122	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00015	<0.00015
Toluene	1	1	<0.00138	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015
Vinyl chloride	0.002	0.002										<0.00011		
Xylenes (total)	10	10	<0.00302	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00009	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000106	<0.000108	<0.000104	<0.000109
2,4-Dimethylphenol	0.49	1.5	<0.00033	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	0.000083 J	<0.000298	<0.000304	<0.000292	<0.000307
2,4-Dinitrotoluene	0.0013	0.003	<0.00022	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.000127	<0.000123	<0.000129
2,6-Dinitrotoluene	0.0013	0.003	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000769	<0.0000784	<0.0000755	<0.0000792
2-Chloronaphthalene	2	5.8	<0.00044	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	<0.0000755	<0.0000792
2-Methylnaphthalene	0.098	0.29	<0.00044	<0.00007	<0.00007	<0.00007	0.00019 J	<0.00005	<0.00005	0.0039	0.000107 J	<0.0000686	0.000264 J	<0.0000693
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00056	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.0000798	<0.0000814	<0.0000783	<0.0000822
4-Nitrophenol	0.049	0.15	<0.00028	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000538	<0.000549	<0.000528	<0.000554
Acenaphthene	1.5	4.4	<0.00033	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	0.0029	<0.0000769	<0.0000784	<0.0000755	<0.0000792
Acenaphthylene	1.5	4.4	<0.00033	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000577	<0.0000588	<0.0000566	<0.0000594
Anthracene	7.3	22	<0.00022	0.00011 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0006	<0.0000481	<0.000049	<0.0000472	<0.0000495
Benzo(a)anthracene	0.0091	0.02	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	0.0000972 J	<0.0000784	<0.0000755	<0.0000792
Benzo(a)pyrene	0.0002	0.0002	<0.00022	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	0.00015 J	<0.0000784	<0.0000755	<0.0000792
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.000127	<0.000123	<0.000129
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00022	0.0035	<0.0002	<0.0003	0.00029	<0.00077	0.00032	<0.00012	0.000512	0.000409 J	<0.000349	<0.000366
Chrysene	0.91	2	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	0.000157 J	<0.0000784	<0.0000755	<0.0000792
Dibenzofuran	0.098	0.29	<0.00033	0.00025	<0.00008	<0.00008	0.00011 J	<0.00005	<0.00005	0.0024	<0.0000769	<0.0000784	0.000134 J	<0.0000792
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.000086 J	0.000194 J	0.000147 J	<0.000104	<0.000109
Fluoranthene	0.98	2.9	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00061	<0.0000673	<0.0000686	<0.000066	<0.0000693
Fluorene	0.98	2.9	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0022	<0.0000673	<0.0000686	<0.000066	<0.0000693
Naphthalene	0.49	1.5	<0.00044	0.0003	<0.0001	<0.0004	0.0026 J	<0.00005	<0.00005	0.02	<0.0000769	0.000265 J	0.00129	0.00071
Nitrobenzene	0.049	0.15	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000106	<0.000108	<0.000104	<0.000109
N-Nitrosodiphenylamine	0.19	0.42	<0.00028	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000962	<0.000098	<0.0000943	<0.000099
Pentachlorophenol	0.001	0.001	<0.00022	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	0.000921 J	<0.000598	<0.000575	<0.000604
Phenanthrene	0.73	2.2	<0.00022	0.00031	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0045	0.000164 J	<0.0000588	0.0000703 J	<0.0000594
Phenol	7.3	22	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00038	<0.00005	<0.0000385	<0.0000392	<0.0000377	<0.0000396
Pyrene	0.73	2.2	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00031	0.000138 J	<0.000108	<0.000104	<0.000109
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray.
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-50A 01/30/2018	MW-50A 03/28/2018	MW-50A 04/01/2018	MW-50A 05/25/2018	MW-50A 01/10/2019	MW-50A 07/19/2019	MW-50A 01/09/2020	MW-50A 03/12/2020	MW-51A 01/31/2008	MW-51A 01/31/2008 Duplicate	MW-51A 02/04/2009	MW-51A 01/20/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002		<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.00052	<0.0005	<0.0005
Benzene	0.005	0.005	<0.0002		<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00025	<0.00025	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.0003		<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	<0.00047	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.0003		<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0068	<0.00025	<0.00025	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.001		<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.00054	<0.0005	<0.0005
Toluene	1	1	<0.0002		<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0033	<0.00041	<0.00041	<0.0005	<0.0005
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003		<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0056	<0.00127	<0.00127	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.00008	<0.0002	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.000042	<0.00029	<0.00057	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00006	<0.00019	<0.00038	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000044	<0.00019	<0.00038	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.00038	<0.00076	<0.00012	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00019	0.016	<0.00038	<0.00076	<0.00007	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.000021	<0.00019	<0.00038	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000049	<0.00024	<0.00048	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.00037	0.02	<0.00029	<0.00057	<0.00009	<0.00009
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.000022 J	0.00028	<0.00029	<0.00057	<0.00006	<0.00007
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.0016	<0.00019	<0.00038	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000052	<0.00019	<0.00038	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.000021	<0.00019	<b>&lt;0.00038</b>	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.000031	<0.00038	<0.00076	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	0.0001 J	<0.000037	<0.000037	<0.000037	<0.000085	<0.000039	<0.00019	<0.00038	0.00034	<0.00019
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.00019	<0.00038	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00013	0.015	0.000566	<0.00057	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000038 J	<0.00019	<0.00038	<0.00007	<0.00007
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.000052 J	0.0011	<0.00019	<0.00038	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00011	0.011	0.000602	0.00042 J	<0.00007	<0.00007	<0.00007
Naphthalene	0.49	1.5	<0.000046	<0.00002	0.00018	<0.00002	0.00027	<0.002	0.31	0.00182	0.00131	0.00029	<0.0001	<0.0001
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000025	<0.00038	<0.00076	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000026	<0.00024	<0.00048	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000082	<0.00019	<0.00038	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00007	0.016	0.00097	0.00075	<0.00007	<0.00007	<0.00007
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	0.00016 J	<0.000066	<0.000036	0.00044 J	<0.00038	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000034	0.00059	<0.00019	<0.00038	<0.00007	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01	0.00205		<0.0004	0.00857	0.00134 J	0.000642 J	0.000718 J	0.00372				

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-51A 06/24/2010	MW-51A 01/20/2011	MW-51A 07/28/2011	MW-51A 02/15/2012	MW-51A 07/24/2012	MW-51A 04/02/2013	MW-51A 08/09/2013	MW-51A 01/29/2014	MW-51A 07/24/2014	MW-51A 01/29/2018	MW-51A 03/28/2018	MW-51A 05/24/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000106	<0.000108	<0.000104	<0.000104	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000298	<0.000304	<0.000292	<0.000292	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.000127	<0.000123	<0.000123	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000769	<0.0000784	0.00292	<0.0000755	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	<0.0000755	<0.0000755	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00013 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000673	<0.0000686	<0.000066	<0.000066	0.000069 J	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000798	<0.000814	<0.000783	<0.000783	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000538	<0.000549	<0.000528	<0.000528	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.00013 J	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	<0.0000755	<0.0000755	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000577	<0.0000588	<0.0000566	<0.0000566	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.00017 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.000481	<0.00049	<0.000472	<0.000472	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	0.00014 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	<0.0000755	<0.0000755	<0.00005	<0.000051	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	<0.0000755	<0.0000755	<0.00002	<0.00002	0.000022 J
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.000127	<0.000123	<0.000123	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00035	<0.00029	<0.0018	<0.0001	<0.00033	<0.000356	<0.000363	0.00121	0.000804	<0.000037	0.000069 J	0.000069 J
Chrysene	0.91	2	0.00013 J	<0.00007	<0.00005	<0.00005	0.00011 J	<0.0000769	<0.0000784	<0.0000755	<0.0000755	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.00012 J	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	<0.0000755	<0.0000755	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00007	<0.00005	<0.00005	0.000051 J	<0.000106	0.00011 J	<0.000104	<0.000104	<0.00002	<0.00002	0.000025 J
Fluoranthene	0.98	2.9	0.00072	<0.00007	<0.00005	<0.00005	0.00012 J	<0.0000673	<0.0000686	<0.000066	<0.000066	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	0.00011 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000673	<0.0000686	<0.000066	<0.000066	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	0.00087	<0.00011	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000784	0.000118 J	0.000162 J	<0.00018	<0.00023	0.000087 J
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000106	<0.000108	<0.000104	<0.000104	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000962	<0.000098	<0.0000943	<0.0000943	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000587	<0.000598	<0.000575	<0.000575	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.00068	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000577	0.0000699 J	<0.0000566	<0.0000566	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000385	<0.0000392	<0.0000377	<0.0000377	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.00037	<0.00007	<0.00005	<0.00005	0.000088 J	<0.000106	<0.000108	<0.000104	<0.000104	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01										<0.0004	<0.0004	<0.0004

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	Residential Assessment Level	C/I Assessment Level	MW-51A 01/10/2019	MW-51A 07/19/2019	MW-51A 01/09/2020	MW-51C 07/24/2014	MW-51C 01/29/2018	MW-51C 03/28/2018	MW-51C 05/24/2018	MW-51C 01/10/2019	MW-51C 07/19/2019	MW-51C 01/09/2020	MW-52A 01/31/2008	MW-52A 01/18/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00052	<0.0025
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	0.000104 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0576	0.0047 J
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00047	
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0892	0.014 J
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00054	<0.0025
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.103	0.012 J
Vinyl chloride	0.002	0.002				<0.00011								
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.24	0.044 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000104	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0008	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.000292	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	1.54	0.0046
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000123	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.019	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.0000755	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.019	<0.00007
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.038	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.000019	0.000067 J	<0.000044	<0.000066	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.929	0.54
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.0000783	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.019	<0.00008
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000528	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.024	<0.00007
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.00004	<0.0000755	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.494	0.36
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.0000566	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.029	0.0045
Anthracene	7.3	22	<0.000014	<0.000014	<0.000063	<0.0000472	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.0001	0.046	0.022
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.0000755	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.019	0.00047
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	0.000027 J	<0.0000755	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.019	0.00013 J
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.000123	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.038	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000078 J	0.00013 J	<0.00014	0.00111	<0.000037	<0.000037	0.000073 J	0.00013 J	0.000079 J	<0.000085	<0.019	<0.00032
Chrysene	0.91	2	<0.000021	<0.000021	0.00005 J	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.019	0.00041
Dibenzofuran	0.098	0.29	<0.00002	0.00003 J	<0.000049	<0.0000755	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00011	0.373	0.28
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	0.000034 J	<0.000104	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.019	<0.00007
Fluoranthene	0.98	2.9	<0.00001	<0.00001	0.00013	<0.000066	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.00018	0.027	0.015
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.000043	<0.000066	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00011	0.263	0.23
Naphthalene	0.49	1.5	0.00012	0.00055	<0.00032	0.000553	<0.0002	<0.00021	0.00029	0.00017	0.00023	<0.0003	10.3	3.9
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.038	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.0000943	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.024	<0.00009
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000575	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.019	<0.00008
Phenanthrene	0.73	2.2	<0.000021	0.000026 J	<0.00021	<0.0000566	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00056	0.24	0.24
Phenol	7.3	22	<0.000035	0.00019 J	<0.000035	0.000628	<0.000035	<0.000035	<0.000035	<0.000035	0.00027	<0.000035	0.038	<0.00007
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000083	<0.000104	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00011	0.039	0.0066
<b>Metals</b>														
Arsenic	0.01	0.01	<0.0004	<0.0004	0.00188 J		0.000614 J	0.0004 J	<0.0004	<0.0004	<0.0004	0.000452 J		

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray.
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-52A 07/14/2011	MW-52A 02/03/2012	MW-52A 07/12/2012	MW-52A 02/01/2013	MW-53C 01/29/2008	MW-53C 07/14/2008	MW-53C 02/03/2009	MW-53C 01/13/2010	MW-53C 06/30/2010	MW-53C 01/26/2011	MW-53C 07/20/2011	MW-53C 02/09/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0005	<0.00014	<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Benzene	0.005	0.005	0.0025 J	0.0017 J	<b>0.0053</b>	0.00461	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.0005	<0.00012	<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Ethylbenzene	0.7	0.7	0.011	0.0053	0.0099	0.00677	<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<0.001	<0.00015	<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013
Toluene	1	1	0.0089	0.0034 J	0.0084	0.00679	<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Vinyl chloride	0.002	0.002	<0.001		<0.0005	0.000661 J								
Xylenes (total)	10	10	0.025	0.011 J	0.021	0.0147	<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<b>&lt;0.00524</b>	<0.00008	<0.00009	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	0.0045	0.0034	0.029	0.0479	<0.00029	<0.00033	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.00005	<0.00005	<b>&lt;0.00619</b>	<0.00019	<0.00022	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.00006	<0.00006	<b>&lt;0.00381</b>	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.00005	<0.00005	<0.00005	<0.00381	<0.00038	<0.00044	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<b>0.33</b>	0.096	<b>0.165</b>		<0.00038	<0.00044	<0.00007	0.000071 J	<0.00007	<0.00007	<0.00005	0.00008 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<b>&lt;0.0395</b>	<0.00019	<0.00022	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.0267	<0.00024	<0.00028	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Acenaphthene	1.5	4.4	0.26	0.19	0.15	0.271	<0.00029	<0.00033	<0.00009	0.0002	0.00032	<0.00009	0.00032	0.0002 J
Acenaphthylene	1.5	4.4	0.004	0.0024	0.0025	<0.00286	<0.00029	<0.00033	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Anthracene	7.3	22	0.041	0.036	0.021	0.0231 J	0.000569	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Benzo(a)anthracene	0.0091	0.02	0.00063	0.00031	0.00022	<0.00381	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	0.00017 J	0.000066 J	<0.00005	<b>&lt;0.00381</b>	<0.00019	<b>&lt;0.00022</b>	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.00005	<0.00005	<b>&lt;0.00619</b>	<0.00038	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00042	0.00043	<0.0001	<b>&lt;0.0176</b>	<0.00019	0.00026 J	0.00072	<0.00024	<0.00032	0.00037	<0.00014	<0.0001
Chrysene	0.91	2	0.0006	0.00033	0.00028	<0.00381	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Dibenzofuran	0.098	0.29	<b>0.2</b>	<b>0.14</b>	<b>0.13</b>	<b>0.178</b>	<0.00029	<0.00033	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.00005	<0.00524	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Fluoranthene	0.98	2.9	0.024	0.013	0.017	0.0245	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Fluorene	0.98	2.9	0.18	0.12	0.11	0.167	<0.00019	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Naphthalene	0.49	1.5	<b>1.9</b>	<b>0.77</b>	<b>0.83</b>	<b>0.878</b>	0.00075	0.00161	0.0012	<0.00027	<0.0001	0.00015 J	<0.00005	0.00066
Nitrobenzene	0.049	0.15	<0.00005	<0.00005	<0.00005	<0.00524	<0.00038	<0.00044	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.00476	<0.00024	<0.00028	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.00005	<0.00005	<0.00005	<b>&lt;0.029</b>	<0.00019	<0.00022	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
Phenanthrene	0.73	2.2	0.22	0.081	0.12	0.226	0.00043 J	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Phenol	7.3	22	0.000066 J	0.000052 J	<0.00015	<0.0019	<0.00019	0.00027 J	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Pyrene	0.73	2.2	0.011	0.0054	0.0071	0.0124 J	0.00039 J	<0.00022	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-53C 07/18/2012	MW-53C 02/06/2013	MW-53C 08/06/2013	MW-53C 01/22/2014	MW-53C 07/25/2014	MW-53C 01/28/2018	MW-53C 03/21/2018	MW-53C 05/31/2018	MW-53C 01/14/2019	MW-53C 07/16/2019	MW-53C 01/09/2020	MW-54C 01/28/2008
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	0.000644 J	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00245
Benzene	0.005	0.005	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00257
Chlorobenzene	0.1	0.1	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00239
Ethylbenzene	0.7	0.7	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.00584
Methylene chloride	0.005	0.005	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00195
Toluene	1	1	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00274
Vinyl chloride	0.002	0.002					<0.00011							
Xylenes (total)	10	10	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00581
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.000106	<0.00011	<0.000524	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00008
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.000298	<0.00031	<0.00148	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.0011	<0.00029
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.000125	<0.00013	<0.000619	<0.000126	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00019
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.0000769	<0.00008	<0.000381	<0.0000777	<0.000042	<0.000042	0.0016	<0.000042	0.0028	<0.000042	<0.00019
2-Chloronaphthalene	2	5.8	<0.00005	<0.0000769	<0.00008	<0.000381	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00038
2-Methylnaphthalene	0.098	0.29	<0.000091	<0.0000673	<0.00007	0.000358 J	<0.0000826	<0.000019	<0.000019	<0.000019	<0.000019	0.000071 J	<0.00025	0.109
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.000798	<0.00083	<0.00395	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
4-Nitrophenol	0.049	0.15	<0.00005	<0.000538	<0.00056	<0.00267	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00024
Acenaphthene	1.5	4.4	<0.00005	<0.0000769	<0.00008	0.000856 J	<0.0000777	<0.000027	<0.000027	0.000044 J	<0.000027	<0.000027	<0.00032	0.074
Acenaphthylene	1.5	4.4	<0.00005	<0.0000577	<0.00006	<0.000286	<0.0000583	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029
Anthracene	7.3	22	<0.00005	<0.0000481	<0.00005	<0.000238	<0.0000485	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.00268
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.0000769	<0.00008	<0.000381	<0.0000777	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	0.00024 J
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.0000769	<0.00008	<0.000381	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.000125	<0.00013	<0.000619	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00038
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0001	<0.000356	<0.00037	<0.00176	<0.000359	0.00012 J	<0.000037	0.0001 J	<0.000037	<0.000037	<0.000037	0.00105 J
Chrysene	0.91	2	<0.00005	<0.0000769	<0.00008	<0.000381	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00021 J
Dibenzofuran	0.098	0.29	<0.00005	<0.0000769	<0.00008	<0.000381	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.0611
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.000106	<0.00011	<0.000524	<0.000107	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00069 J
Fluoranthene	0.98	2.9	<0.00005	<0.0000673	<0.00007	<0.000333	<0.000068	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.000071 J	0.00426
Fluorene	0.98	2.9	<0.00005	<0.0000673	<0.00007	0.000355 J	<0.000068	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.0002	0.0323
Naphthalene	0.49	1.5	<0.00048	<0.000183	<0.00008	<0.00212	<0.00194	<0.00002	<0.000034	0.00023	<0.00025	0.00013	<0.0012	0.892
Nitrobenzene	0.049	0.15	<0.00005	<0.000106	<0.00011	<0.000524	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00038
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.0000962	<0.0001	<0.000476	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	0.000056 J	<0.00024
Pentachlorophenol	0.001	0.001	<0.00005	<0.000587	<0.00061	<0.0029	<0.000592	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	0.00025 J
Phenanthrene	0.73	2.2	<0.00005	<0.0000577	<0.00006	0.000939 J	<0.0000665	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00025	0.0389
Phenol	7.3	22	<0.00005	<0.0000385	<0.00004	<0.00019	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	<0.00016	0.0005	<0.00019
Pyrene	0.73	2.2	<0.00005	<0.000106	<0.00011	<0.000524	<0.000107	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000061	0.00227
<b>Metals</b>														
Arsenic	0.01	0.01						0.000502 J	0.000443 J	0.000694 J	<0.0004	0.000569 J	0.000728 J	

**Notes:**

- All values in milligrams per liter (mg/L).
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- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-54C 07/14/2008	MW-54C 02/03/2009	MW-54C 01/21/2010	MW-54C 06/30/2010	MW-54C 01/26/2011	MW-54C 07/20/2011	MW-54C 02/08/2012	MW-54C 07/25/2012	MW-54C 02/12/2013	MW-54C 08/06/2013	MW-54C 01/23/2014	MW-54C 07/25/2014
<b><i>Volatile Organic Compounds</i></b>														
1,2-Dichloroethane	0.005	0.005	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014
Benzene	0.005	0.005	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008
Chlorobenzene	0.1	0.1	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.000128 J	0.000128 J	<0.00018	<0.00012
Ethylbenzene	0.7	0.7	0.00391 J	0.0029 J	<0.0005	0.0024 J	<0.0005	0.0018 J	0.0011 J	0.0011 J	0.000187 J	0.00062 J	0.000527	0.000282 J
Methylene chloride	0.005	0.005	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015
Toluene	1	1	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015
Vinyl chloride	0.002	0.002												<0.00011
Xylenes (total)	10	10	<0.00127	0.0027 J	<0.001	0.0011 J	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	0.00076 J	0.00062 J	<0.00026
<b><i>Semivolatile Organic Compounds</i></b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.00011	<0.000104	<0.000109
2,4-Dimethylphenol	0.49	1.5	<0.00029	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	0.000098 J	<0.00005	<0.000295	<0.00031	<0.000292	<0.000307
2,4-Dinitrotoluene	0.0013	0.003	<0.0002	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.00013	<0.000123	<0.000129
2,6-Dinitrotoluene	0.0013	0.003	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.00008	<0.0000755	<0.0000792
2-Chloronaphthalene	2	5.8	<0.00039	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	<0.0000755	<0.0000792
2-Methylnaphthalene	0.098	0.29	<b>0.14</b>	<b>0.13</b>	<0.00007	0.0096	0.00025	0.022	0.0065	0.0054 J	0.00392	0.0173	0.0176	0.00834
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000079	<0.00083	<0.000783	<0.000822
4-Nitrophenol	0.049	0.15	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.00056	<0.000528	<0.000554
Acenaphthene	1.5	4.4	0.0738	0.067	0.00016 J	0.024	0.0023	0.039	0.035	0.022	0.0219	0.0749	0.062	0.0367
Acenaphthylene	1.5	4.4	<0.00029	0.00072	<0.00007	0.00042	<0.00007	0.00045	0.00051	0.00039	<0.0000571	<0.00006	0.00105	0.000526
Anthracene	7.3	22	0.00293	0.003	<0.00007	0.005	0.00027	0.0029	0.0024	0.0019	0.00183	0.00389	0.00445	0.00261
Benzo(a)anthracene	0.0091	0.02	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	0.0000993 J	<0.0000792
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	<0.0000755	<0.0000792
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.00013	<0.000123	<0.000129
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00029 J	0.00072	<0.00077	<0.00037	0.0016	<0.00015	<0.0001	0.00017 J	<0.000352	<0.00037	<0.000349	<0.000366
Chrysene	0.91	2	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	0.0000758 J	<0.0000792
Dibenzofuran	0.098	0.29	0.0739	0.064	<0.00008	0.028	0.0018	0.046	0.047	0.029	0.0223	0.0878	0.0695	0.0471
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.000064 J	<0.000105	<0.00011	<0.000824	<0.000109
Fluoranthene	0.98	2.9	0.00349	0.0032	<0.00007	0.0032	0.00016 J	0.0034	0.0026	0.002	0.00246	0.00474	0.00575	0.00302
Fluorene	0.98	2.9	0.0351	0.03	<0.00007	0.015	0.001	0.022	0.021	0.011	0.0092	0.0409	0.0321	<0.0208
Naphthalene	0.49	1.5	<b>0.912</b>	<b>1.1</b>	<0.0001	0.21	0.0055	0.47	0.35	0.15	0.0681	0.383 J	0.315 J	0.18
Nitrobenzene	0.049	0.15	<0.00039	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.00011	<0.000104	<0.000109
N-Nitrosodiphenylamine	0.19	0.42	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	0.00014 J	<0.0000952	<0.0001	<0.0000943	<0.000099
Pentachlorophenol	0.001	0.001	<0.0002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.00061	<0.000575	<0.000604
Phenanthrene	0.73	2.2	0.0495	0.042	<0.00007	0.024	0.0011	0.04	0.034	0.019	0.0128	0.04	0.042	0.0148
Phenol	7.3	22	<0.0002	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00011 J	<0.00005	<0.0000381	<0.00004	<0.0000377	<0.0000396
Pyrene	0.73	2.2	0.00163	0.0018	<0.00007	0.0016	<0.00007	0.0017	0.0015	0.0013	0.00138	0.00248	0.00373	0.00169
<b><i>Metals</i></b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-54C 01/28/2018	MW-54C 03/20/2018	MW-54C 05/31/2018	MW-54C 01/14/2019	MW-54C 07/16/2019	MW-54C 01/09/2020	MW-55A 02/04/2009	MW-55A 01/18/2010	MW-55A 07/14/2011	MW-55A 02/03/2012	MW-55A 07/12/2012	MW-55A 01/30/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0025	<0.005	<0.01	<0.005	<0.0028
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.19	0.072	0.07	0.15	0.17	0.133
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.005	<0.01	<0.005	<0.0024
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.15	0.2	0.17	0.2	0.24	0.228
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0025	<0.0065	<0.013	<0.01	<0.003
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.44	0.29	0.24	0.41	0.39	0.385
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.35	0.47	0.42	0.48	0.62	0.575
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.00005	<0.0005	<0.00005	<0.000104
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	1.2	0.28	0.48	1.8	0.96	<0.000292
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00005	<0.0005	<0.00005	<0.000123
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00006	<0.0006	<0.00006	<0.0000755
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001	<0.00005	<0.0005	<0.00005	<0.0000755
2-Methylnaphthalene	0.098	0.29	0.00048	0.00075	0.0059	0.0014	0.0021	0.002	0.63	0.39	0.33	0.25	0.31	<0.000066
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.0008	<0.00008	<0.000783
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00005	<0.0005	<0.00005	<0.000528
Acenaphthene	1.5	4.4	0.023	0.03	0.04	0.014	0.041	0.021	0.28	0.19	0.16	0.14	0.11	0.0573
Acenaphthylene	1.5	4.4	0.00034	0.00036	0.00054	<0.0002	0.00037	0.00023	0.0037	0.0028	0.003	0.0019 J	0.0017	0.0021
Anthracene	7.3	22	0.0015	0.0027	0.0041	0.0013	0.0023	0.0015	0.047	0.021	0.016	0.016	0.0075	0.00062
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	0.000065 J	<0.00005	<0.00005	<0.00005	0.01	0.0018	0.0014	<0.0005	0.00034	<0.0000755
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	0.000039 J	<0.00002	<0.00002	0.0069	0.00081	0.00062	<0.0005	0.000081 J	<0.0000755
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	0.000056 J	<0.00003	<0.00003	<0.00009	0.00009 J	<0.00005	<0.0005	<0.00005	<0.000123
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000055 J	<0.000037	0.00011 J	0.000096 J	<0.000037	<0.00018	0.00073	<0.00031	<0.0001	<0.001	<0.0001	<0.000349
Chrysene	0.91	2	<0.000021	0.000059 J	0.000053 J	0.000038 J	<0.000021	0.000054 J	0.0099	0.0017	0.0014	<0.0005	0.00025	<0.0000755
Dibenzofuran	0.098	0.29	0.025	0.029	0.045	0.015	0.032	0.02	0.2	0.13	0.12	0.084	0.078	0.0265
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	0.000029 J	<0.00002	<0.00002	<0.00002	<0.00007	<0.00007	<0.00005	<0.0005	<0.00005	<0.000104
Fluoranthene	0.98	2.9	0.0023	0.004	0.0049	0.002	0.0033	0.0026	0.052	0.0081	0.009	0.0044	0.0042	0.000459 J
Fluorene	0.98	2.9	0.011	0.014	0.022	0.0085	0.018	0.0095	0.16	0.083	0.08	0.057	0.048	0.00213
Naphthalene	0.49	1.5	0.022	0.029	0.068	0.019	0.032	<0.035	17	11	8.6	9.9	9.7	<0.00227
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	0.00009 J	<0.00005	<0.0005	<0.00005	<0.000104
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00005	<0.0005	<0.00005	<0.0000943
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	0.00053	<0.00008	<0.00005	<0.0005	<0.00005	<0.0000575
Phenanthrene	0.73	2.2	0.0023	0.0041	0.023	0.0052	0.0065	0.0056	0.2	0.084	0.083	0.047	0.045	<0.0000566
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	0.00044	<0.000064	0.15	0.025 J	0.0038	0.079	0.046	<0.0000377
Pyrene	0.73	2.2	0.0012	0.0018	0.0028	0.001	0.0016	0.0014	0.032	0.0052	0.0061	0.0041	0.0021	0.000223 J
<b>Metals</b>														
Arsenic	0.01	0.01	0.00128 J	0.00133 J	0.0012 J	0.00123 J	0.00103 J	0.00122 J						

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-55A 07/30/2013	MW-55A 01/14/2014	MW-55A 07/17/2014	MW-55B 02/02/2012	MW-55B 07/12/2012	MW-55B 01/30/2013	MW-55B 07/30/2013	MW-55B 01/14/2014	MW-55B 07/17/2014	MW-57A 02/05/2009	MW-57A 01/20/2010	MW-57A 06/23/2010 DNAPL
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.014	<0.007	<0.0028	<0.01	<0.005	<0.007	<0.014	<0.007	<0.0028	<0.0005	<0.0025	<0.005
Benzene	0.005	0.005	0.145	0.0715	0.0881	0.78	0.89	0.881	0.809	0.648	0.846	0.26	0.17	0.47
Chlorobenzene	0.1	0.1	<0.012	<0.006	<0.0024	<0.01	<0.005	<0.006	<0.012	<0.006	<0.0024	<0.0005	<0.0025	<0.005
Ethylbenzene	0.7	0.7	0.26	0.2	0.368	0.13	0.21	0.162	0.173	0.134	0.126	0.34	0.32	0.45
Methylene chloride	0.005	0.005	0.0894 J	<0.0075	0.0179 J	<0.013	<0.01	0.0213 J	0.0517 J	<0.0075	0.0155 J	<0.0005	<0.0025	0.014 J
Toluene	1	1	0.431	0.311	0.409	0.65	0.9	0.76	0.782	0.597	0.591	0.63	0.13	0.86
Vinyl chloride	0.002	0.002	<0.011						<0.011				<0.0025	
Xylenes (total)	10	10	0.584	0.486	0.869	0.39	0.68	0.623	0.624	0.481	0.443	0.92	0.6	1.2
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00534	<0.0529	<0.00519	<0.0005	<0.0005	<0.0208	<0.0534	<0.529	<0.013	<0.0001	<0.0001	0.0005 J
2,4-Dimethylphenol	0.49	1.5	0.956	0.519	0.463	35	30	2.06	25.2	44.2	35.6	1.8	3	2.7 J
2,4-Dinitrotoluene	0.0013	0.003	<0.00631	<0.0625	<0.00613	<0.0005	<0.0005	<0.0245	<0.0631	<0.625	<0.0153	<0.00009	<0.00009	0.00045 J
2,6-Dinitrotoluene	0.0013	0.003	<0.00388	<0.0385	<0.00377	<0.0006	<0.0006	<0.0151	<0.0388	<0.385	<0.00943	<0.00007	<0.00007	0.00035 J
2-Chloronaphthalene	2	5.8	<0.00388	<0.0385	<0.00377	<0.0005	<0.0005	<0.0151	<0.0388	<0.385	<0.00943	<0.00012	<0.0001	0.0005 J
2-Methylnaphthalene	0.098	0.29	0.468	0.463	0.486	0.28	0.64	0.757	0.868	0.901 J	0.512	0.73	0.89	3.5 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0403	<0.399	<0.0392	<0.0008	<0.0008	<0.157	<0.403	<3.99	<0.0979	<0.00008	<0.00008	0.0004 J
4-Nitrophenol	0.049	0.15	<0.0272	<0.269	<0.0264	<0.0005	<0.0005	<0.106	<0.272	<2.69	<0.066	<0.00007	<0.00007	0.00035 J
Acenaphthene	1.5	4.4	0.207	0.251	0.219	0.19	0.26	0.347	<0.0388	<0.385	0.19	0.24	0.31	2 J
Acenaphthylene	1.5	4.4	<0.00291	<0.0288	<0.00283	0.0057	0.01	<0.0113	<0.0291	<0.288	<0.00708	0.0056	0.0061	0.02 J
Anthracene	7.3	22	0.0336	0.083 J	0.032	0.016	0.03	0.0492 J	0.0437 J	<0.24	0.027 J	0.044	0.022	0.9 J
Benzo(a)anthracene	0.0091	0.02	<0.00388	<0.0385	<0.00377	0.0011 J	0.0012 J	<0.0151	<0.0388	<0.385	<0.00943	0.011	0.00051	0.15 J
Benzo(a)pyrene	0.0002	0.0002	<0.00388	<0.0385	<0.00377	<0.0005	<0.0005	<0.0151	<0.0388	<0.385	<0.00943	0.0045	0.00012 J	0.037 J
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00631	<0.0625	<0.00613	<0.0005	<0.0005	<0.0245	<0.0631	<0.625	<0.0153	<0.00009	<0.00009	0.00045 J
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.018	<0.178	<0.0175	<0.0001	<0.0001	<0.0698	<0.18	<1.78	<0.0436	0.002	<0.0004	0.001 J
Chrysene	0.91	2	<0.00388	<0.0385	<0.00377	<0.0005	0.0009 J	<0.0151	<0.0388	<0.385	<0.00943	0.0094	0.00034	0.14 J
Dibenzofuran	0.098	0.29	<0.00388	0.15 J	0.14	0.15	0.23	<0.0151	0.309	<0.385	0.138	0.21	0.17	1.9 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00534	<0.0529	<0.00519	<0.0005	<0.0005	<0.0208	<0.0534	<0.529	<0.013	<0.00007	<0.00007	0.00035 J
Fluoranthene	0.98	2.9	0.0148 J	0.0595 J	0.0175 J	0.0047	0.014	0.0153 J	<0.034	<0.337	0.0119 J	0.054	0.0063	1.4 J
Fluorene	0.98	2.9	0.0996	0.172 J	0.1	0.09	0.15	0.166	0.195 J	<0.337	0.0816	0.083	0.11	1.6 J
Naphthalene	0.49	1.5	13.8	11.7	11.6	21	24	2.3	21.9	24.3	13.5	16	7.4	20 J
Nitrobenzene	0.049	0.15	<0.00534	<0.0529	<0.00519	<0.0005	<0.0005	<0.0208	<0.0534	<0.529	<0.013	<0.00009	<0.00009	0.00045 J
N-Nitrosodiphenylamine	0.19	0.42	<0.00485	<0.0481	<0.00472	<0.0005	<0.0005	<0.0189	<0.0485	<0.481	<0.0118	<0.00009	<0.00009	0.00045 J
Pentachlorophenol	0.001	0.001	<0.0296	<0.293	<0.0288	<0.0005	0.00069 J	<0.115	<0.296	<2.93	<0.0719	<0.00008	<0.00008	0.0004 J
Phenanthrene	0.73	2.2	0.078	0.174 J	0.0893	0.057	0.13	0.13	0.228 J	<0.288	0.1	0.22	0.088	4 J
Phenol	7.3	22	<0.00194	<0.0192	<0.00189	150	130	0.0999	103	454	127	0.052	0.0099	0.042 J
Pyrene	0.73	2.2	0.00729 J	<0.0529	0.0101 J	0.0042	0.0076	<0.0208	<0.0534	<0.529	<0.013	0.038	0.0037	0.84 J
<b>Metals</b>														
Arsenic	0.01	0.01												

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-57A 01/18/2011	MW-57A 07/22/2011	MW-57A 02/02/2012	MW-57A 07/24/2012	MW-57A 02/11/2013	MW-57A 07/31/2013	MW-57A 01/15/2014	MW-57A 07/29/2014	MW-57A 07/10/2019	MW-57A 01/08/2020	MW-58A 02/05/2009	MW-58A 01/20/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.005	<0.01	<0.01	<0.0005	<0.0014	<0.007	<0.0002	<0.00014	<0.0002	<0.0002	<0.0025	<0.0025
Benzene	0.005	0.005	0.23	0.084	0.14	0.064	0.138	0.137	0.109	0.0412	0.00057 J	0.038	0.052	0.038
Chlorobenzene	0.1	0.1	<0.005	<0.01	<0.01	<0.0005	<0.0012	<0.006	0.000465 J	0.000625 J	<0.0003	0.00046 J	<0.0025	0.0093 J
Ethylbenzene	0.7	0.7	0.29	0.13	0.22	0.17	0.24	0.283	0.198	0.274	0.00085 J	0.015	0.079	0.063
Methylene chloride	0.005	0.005	<0.005	<0.013	<0.013	<0.001	0.00367 J	<0.0075	<0.00022	<0.00015	<0.001	<0.001	<0.0025	<0.0025
Toluene	1	1	0.38	0.055	0.23	0.1	0.244	0.308	0.198	0.0355	<0.0002	0.01	0.022 J	0.02 J
Vinyl chloride	0.002	0.002	<0.005	<0.01	<0.01	0.0016 J	<0.0011	<0.0055	0.00154		<0.0002	<0.0002		<0.0025
Xylenes (total)	10	10	0.68	0.19	0.4	0.33	0.591	0.572	0.454	0.455	<0.0003	0.026	0.1	0.04 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0005	<0.0005	<0.0005	<0.075	<0.0267	<0.0259	<0.000109	<0.000021	<0.000021	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	2	1	1.7	0.2	1.62	0.994	7.91 J	0.0443 J	<0.00059	0.41	0.047	0.097
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.0005	<0.0005	<0.0005	<0.0886	<0.0316	<0.0307	<0.000129	<0.000058	<0.000058	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.0006	<0.0006	<0.0006	<0.0545	<0.0194	<0.0189	<0.0000792	<0.000042	<0.000042	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0001	<0.0005	<0.0005	<0.0005	<0.0545	<0.0194	<0.0189	<0.0000792	<0.000021	<0.000021	<0.00012	<0.0001
2-Methylnaphthalene	0.098	0.29	3.5	13	1.9	3.1	13.9	1.5	8.24	0.616	0.19	0.18	0.22	0.1
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.0008	<0.0008	<0.0008	<0.566	<0.201	<0.196	<0.000822	<0.00002	<0.00002	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00007	<0.0005	<0.0005	<0.0005	<0.382	<0.136	<0.132	<0.000554	<0.000047	<0.000047	<0.00007	<0.00007
Acenaphthene	1.5	4.4	1.9	8.6	1.2	1.8	8.56	0.997	5.69	0.335	0.099	0.19	0.31	0.18
Acenaphthylene	1.5	4.4	0.022	0.091	0.014	0.024	<0.0409	<0.0146	<0.0142	0.00779	0.0019	0.0025	0.0012	0.0013
Anthracene	7.3	22	0.62	8.4	0.34	0.55	3.09	0.337	2.02	0.0557	0.0084	0.048	0.0045	0.0098
Benzo(a)anthracene	0.0091	0.02	0.12	0.45	0.047	0.074	0.605	0.0521 J	0.361	0.0072	0.001	0.0036	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	0.028	0.16	0.014	0.024	0.165 J	<0.0194	0.0962 J	0.00385	0.00054	0.0011	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.0005	<0.0005	<0.0005	<0.0886	<0.0316	<0.0307	0.0021	<0.00003	<0.00003	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0002	<0.001	0.0016 J	<0.001	<0.252	<0.0898	<0.0873	<0.000808	0.00021	<0.00037	0.0003	<0.0002
Chrysene	0.91	2	0.11	0.53	0.046	0.089	0.602	0.0482 J	0.36	0.00625	0.001	0.003	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	1.7	8.6	0.86	1.7	7.28	0.799	4.69	0.257	0.076	0.15	0.23	0.14
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.0005	<0.0005	<0.0005	<0.075	<0.0267	<0.0259	<0.000109	<0.00002	<0.00002	0.0012	<0.00007
Fluoranthene	0.98	2.9	0.99	6	0.48	0.74	4.98	0.412	3.19	0.0561	0.0091	0.038	0.0025	0.0058
Fluorene	0.98	2.9	1.4	7.9	0.72	1.4	6.54	0.713	4.16	0.21	0.059	0.13	0.15	0.12
Naphthalene	0.49	1.5	18	71	9.2	22	60.7	13.5	56.9	7.27	0.61	1	2.4	0.67
Nitrobenzene	0.049	0.15	<0.00009	<0.0005	<0.0005	<0.0005	<0.075	<0.0267	<0.0259	<0.000109	<0.000024	<0.000024	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.0005	<0.0005	<0.0005	<0.0682	<0.0243	<0.0236	<0.000099	<0.000025	<0.000025	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.00008	<0.0005	<0.0005	<0.0005	<0.416	<0.148	<0.144	<0.000604	0.00059	<0.000079	<0.00008	<0.00008
Phenanthrene	0.73	2.2	3.5	13	2	3	17	1.61	13.1	0.271	0.059	0.17	0.041	0.049
Phenol	7.3	22	0.02	<0.0005	0.0089	<0.0005	<0.0273	<0.00971	<0.00943	<0.0000396	<0.00016	0.0049	0.00029	0.0074
Pyrene	0.73	2.2	0.67	3.3	0.34	0.42	3.12	0.264	2.29	0.0308 J	0.0055	0.025	0.0012	0.0034
<b>Metals</b>														
Arsenic	0.01	0.01									0.00447	0.0545		

Notes:

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- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-58A 06/23/2010	MW-58A 01/19/2011	MW-58A 07/27/2011	MW-58A 02/03/2012	MW-58A 07/24/2012	MW-58A 02/11/2013	MW-58A 08/06/2013	MW-58A 01/29/2014	MW-58A 08/28/2014	MW-58A 01/31/2018	MW-58A 03/19/2018	MW-58A 05/16/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0025	<0.0025	<0.005	<0.01	<0.0005	<0.00014	<0.00014	<0.00014	<0.0014	<0.0002	<0.0002	<0.001
Benzene	0.005	0.005	0.075	0.034	<0.005	0.12	0.16	0.0943	0.0000807 J	<0.00008	0.259	0.0048	0.012	0.012
Chlorobenzene	0.1	0.1	0.01 J	0.0029 J	<0.005	<0.01	0.0018 J	0.00295	<0.00012	<0.00012	<0.0012	<0.0003	0.00054 J	<0.0015
Ethylbenzene	0.7	0.7	0.11	0.03	<0.0055	0.085	0.099	0.0648	<0.00011	<0.00011	0.167	0.0066	0.038	0.035
Methylene chloride	0.005	0.005	<0.0048	<0.0025	<0.0065	<0.013	<0.001	<0.00015	<0.00015	<0.00015	<0.0015	<0.001	<0.001	<0.005
Toluene	1	1	0.045	0.0059 J	<0.005	0.043 J	0.041	0.0176	<0.00015	<0.00015	0.135	0.00091 J	0.00063 J	0.0027 J
Vinyl chloride	0.002	0.002		<0.0025	<0.005		0.011	0.00281	<0.00011	<0.00011	0.0101 J	<0.0002	<0.0002	<0.001
Xylenes (total)	10	10	0.15	0.029 J	<0.016	0.23	0.31	0.122	<0.00026	<0.00026	0.352	0.012	0.068	0.015
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	0.0001 J	<0.0001	<0.00005	<0.00005	<0.00005	<0.01	<0.00011	<0.000104	<0.00539	<0.00021	<0.00021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.61	0.68	<0.00005	1.1	2.4	0.95	<0.00031	<0.000292	9.19	0.0015 J	0.00053	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0118	<0.00013	<0.000123	<0.00637	<0.00058	<0.00058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.00727	<0.00008	<0.0000755	<0.00392	<0.00042	<0.00042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00727	<0.00008	<0.0000755	<0.00392	<0.00021	<0.00021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.21	0.057	<0.00005	0.082	0.0076	0.243	<0.00007	<0.000066	0.373	0.038	0.045	0.14
4,6-Dinitro-2-methylphenol	0.0024	0.0073	0.00008 J	<0.00008	<0.00008	<0.00008	<0.00008	<0.0755	<0.00083	<0.000783	<0.0407	<0.0002	<0.0002	<0.00002
4-Nitrophenol	0.049	0.15	0.00007 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0509	<0.00056	<0.000528	<0.0275	<0.00047	<0.00047	<0.000047
Acenaphthene	1.5	4.4	0.28	0.12	<0.00005	0.16	0.057	0.205	<0.00008	<0.0000755	0.221	0.1	0.17	0.19
Acenaphthylene	1.5	4.4	0.0015	0.00072	<0.00005	0.0011	0.0011	<0.00545	<0.00006	<0.0000566	0.00996 J	0.0012	0.001	0.0017
Anthracene	7.3	22	0.017 J	0.0051	0.00039	0.0055	0.0069	0.0245 J	<0.00005	<0.0000472	0.0126 J	0.0055	0.008	0.011
Benzo(a)anthracene	0.0091	0.02	0.00007 J	<0.00007	<0.00005	<0.00005	0.00072	<0.00727	<0.00008	<0.0000755	<0.00392	<0.0005	0.000083 J	0.000083 J
Benzo(a)pyrene	0.0002	0.0002	0.00008 J	<0.00008	<0.00005	<0.00005	0.00027	<0.00727	<0.00008	<0.0000755	<0.00392	<0.0002	<0.0002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	0.0321 J	<0.00013	<0.000123	<0.00637	<0.0003	<0.0003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00046 J	0.00035	0.00071	<0.0001	<0.00094	<0.0336	<0.00037	<0.000349	<0.0181	<0.00037	<0.00037	0.000098 J
Chrysene	0.91	2	0.00007 J	<0.00007	<0.00005	<0.00005	0.0011	<0.00727	<0.00008	<0.0000755	<0.00392	<0.00021	0.000083 J	0.000082 J
Dibenzofuran	0.098	0.29	0.23	0.079	0.0017	0.13	0.0088	0.128	<0.00008	<0.0000755	0.136	0.036	0.08	0.091
Di-n-butylphthalate (DBP)	2.4	7.3	0.00007 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.01	<0.00011	<0.000104	<0.00539	<0.0002	<0.0002	<0.00002
Fluoranthene	0.98	2.9	0.009 J	0.0049	0.001	0.0036	0.0099	0.0102 J	<0.00007	<0.000066	<0.00343	0.0065	0.0074	0.0067
Fluorene	0.98	2.9	0.16	0.065	<0.00005	0.08	0.027	0.12	<0.00007	<0.000066	0.109	0.084	0.096	0.12
Naphthalene	0.49	1.5	1.5	0.45	<0.00005	2.2	0.068	2.96 J	0.00036 J	<0.0000755	4.05	0.0037	0.95	0.32
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.01	<0.00011	<0.000104	<0.00539	<0.00024	<0.00024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	0.00009 J	<0.00009	<0.00005	<0.00005	<0.00005	<0.00909	<0.0001	<0.0000943	<0.0049	<0.00025	<0.00025	<0.000025
Pentachlorophenol	0.001	0.001	0.00008 J	<0.00008	<0.00005	<0.00005	0.00017 J	<0.0555	<0.00061	<0.000575	<0.0299	<0.00079	<0.00079	<0.000079
Phenanthrene	0.73	2.2	0.061 J	0.037	<0.00005	0.039	0.036	0.0563	<0.00006	<0.0000566	0.0702	0.024	0.042	0.038
Phenol	7.3	22	0.0065 J	0.00037	0.000077 J	0.0038	0.00074	<0.00364	<0.00004	<0.0000377	<0.00196	0.00054 J	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0042 J	0.0022	0.00073	0.0022	0.0069	<0.01	<0.00011	<0.000104	<0.00539	0.0031	0.0039	0.0036
<b>Metals</b>														
Arsenic	0.01	0.01										0.000713 J	0.00106 J	0.00143 J

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-58A 01/23/2019	MW-58A 07/10/2019	MW-58A 01/08/2020	MW-59A 02/05/2009	MW-59A 01/20/2010	MW-59A 06/24/2010	MW-59A 01/20/2011	MW-59A 07/18/2011	MW-59A 02/06/2012	MW-59A 07/27/2012	MW-59A 01/31/2013	MW-59A 08/01/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014
Benzene	0.005	0.005	0.0011	0.0049	0.0053	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008
Chlorobenzene	0.1	0.1	0.00046 J	0.0008 J	0.00072 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012
Ethylbenzene	0.7	0.7	0.0032	0.017	0.0098	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015
Toluene	1	1	0.0014	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011
Xylenes (total)	10	10	0.005	0.028	0.028	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.000108
2,4-Dimethylphenol	0.49	1.5	0.0001 J	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	<0.00008	0.000066 J	<0.00005	<0.00005	<0.000292	<0.000304
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000127
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000784
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000784
2-Methylnaphthalene	0.098	0.29	0.000072 J	0.015	0.08	<0.00007	<0.00007	0.0002	0.0018	<0.00005	<0.00005	<0.00005	<0.000066	<0.0000686
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.0000783	<0.0000814
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000528	<0.000549
Acenaphthene	1.5	4.4	0.023	0.11	0.27	<0.00009	<0.00009	0.0003	0.00079	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000784
Acenaphthylene	1.5	4.4	0.00038	0.00087	0.0015	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000566	<0.0000588
Anthracene	7.3	22	0.002	0.062	0.018	<0.00007	<0.00007	0.00026	0.0004	<0.00005	<0.00005	<0.00005	<0.0000472	0.0000519 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	0.00082	0.000091 J	<0.00007	<0.00007	<0.00007	0.00015 J	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000784
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000784
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000127
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00007	<0.000037	<0.00011	0.0006	<0.00065	<0.00023	<0.00031	<0.00054	0.00015 J	<0.0001	<0.000349	<0.000363
Chrysene	0.91	2	0.00003 J	0.00084	0.000071 J	<0.00007	<0.00007	<0.00007	0.00014 J	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000784
Dibenzofuran	0.098	0.29	0.013	0.059	0.13	<0.00008	<0.00008	0.0007	0.00099	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000784
Di-n-butylphthalate (DBP)	2.4	7.3	0.000032 J	<0.00002	<0.00002	0.00077	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.000075	<0.000104	0.000169 J
Fluoranthene	0.98	2.9	0.002	0.02	0.0099	<0.00007	<0.00007	0.0005	0.0012	<0.00012	<0.00005	<0.00005	<0.000066	<0.0000686
Fluorene	0.98	2.9	0.015	0.15	0.19	<0.00007	<0.00007	0.00045	0.00084	<0.00005	<0.00005	<0.00005	<0.000066	<0.0000686
Naphthalene	0.49	1.5	0.00042	0.46	1.2	<0.0001	<0.0001	0.00047	0.0066	<0.00005	<0.00005	0.000051 J	<0.0000755	<0.0000784
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.000108
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.000098
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000575	<0.000598
Phenanthrene	0.73	2.2	0.0038	0.036	0.062	<0.00007	<0.00007	0.0017	0.0024	<0.00018	<0.00005	<0.00005	<0.0000566	0.000075 J
Phenol	7.3	22	0.000074 J	<0.000035	<0.000035	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.000065 J	<0.0000377	<0.0000392
Pyrene	0.73	2.2	0.00088	0.0099	0.005	<0.00007	<0.00007	0.00029	0.00059	<0.00005	<0.00005	<0.00005	<0.000104	<0.000108
<b>Metals</b>														
Arsenic	0.01	0.01	0.00232	0.00748	0.000906 J									

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-59A 01/16/2014	MW-59A 07/30/2014	MW-59A 01/29/2018	MW-59A 03/20/2018	MW-59A 05/24/2018	MW-59A 01/23/2019	MW-59A 07/17/2019	MW-59A 01/16/2020	MW-59D 02/05/2009	MW-59D 02/05/2009 Duplicate	MW-59D 01/20/2010	MW-59D 01/20/2010 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	0.0011 J	0.0014 J	<0.0005	<0.0005
Toluene	1	1	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.00064 J	0.0007 J	<0.0005	<0.0005
Vinyl chloride	0.002	0.002	<0.00018	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005
Xylenes (total)	10	10	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.001	<0.001	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000104	<0.000109	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.000292	<0.000307	<0.00004	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000123	<0.000129	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0000755	<0.0000792	<0.000042	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0000755	<0.0000792	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	<0.00012	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.0000902	<0.0000693	0.00011	<0.000019	<0.000019	<0.000019	<0.000019	0.000036 J	0.00015 J	0.0003	<0.00007	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0000783	<0.0000822	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000528	<0.000554	<0.000047	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.0000831	<0.0000792	0.0001	<0.000028	<0.000027	<0.000027	<0.000027	<0.000029	0.00015 J	0.00014 J	<0.00009	<0.00009
Acenaphthylene	1.5	4.4	<0.0000566	<0.0000594	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00006	<0.00006	<0.00007	<0.00007
Anthracene	7.3	22	0.000119 J	<0.0000495	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.0000755	<0.0000792	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.0000755	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000123	<0.000129	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000349	<0.000366	0.000094 J	<0.000038	<0.000037	<0.000037	<0.000037	<0.00005	0.006	0.0055	<0.00023	<0.00087
Chrysene	0.91	2	<0.0000755	<0.0000792	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.000136	<0.0000792	0.00017	<0.00002	<0.00002	<0.00002	<0.00002	<0.00016	0.00014 J	<0.00008	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	0.000178 J	<0.000109	<0.00002	<0.00002	<0.00002	<0.00002	0.000064 J	<0.00002	0.0029	0.0023	<0.00007	<0.00007
Fluoranthene	0.98	2.9	0.000199 J	<0.0000693	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.000032	<0.00007	<0.00007	<0.00007	<0.00007
Fluorene	0.98	2.9	0.000176 J	<0.0000693	0.00011	<0.000031	<0.00003	<0.00003	<0.00003	<0.00011	0.00013 J	0.00012 J	<0.00007	<0.00007
Naphthalene	0.49	1.5	<0.000381	0.000219 J	0.00068	<0.00002	<0.00002	<0.00002	<0.00012	<0.0012	0.0019	0.0029	<0.0001	<0.0001
Nitrobenzene	0.049	0.15	<0.000104	<0.000109	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0000943	<0.000099	<0.000025	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000575	<0.000604	<0.000079	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.000587	<0.0000594	0.00018	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	0.0002	0.00025	<0.00007	<0.00007
Phenol	7.3	22	<0.0000377	<0.0000396	<0.000035	<0.000036	0.000089 J	<0.000035	0.000058 J	<0.000035	<0.00007	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	0.00012 J	<0.000109	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000023	<0.00007	<0.00007	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01				0.00181 J	0.00131 J	<b>0.0101</b>	0.00243	0.00455	0.00368			

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray.
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-59D 07/01/2010	MW-59D 07/01/2010 Duplicate	MW-59D 01/20/2011	MW-59D 07/27/2011	MW-59D 07/27/2011 Duplicate	MW-59D 02/14/2012	MW-59D 02/14/2012 Duplicate	MW-59D 07/23/2012	MW-59D 02/11/2013	MW-59D 02/11/2013 Duplicate	MW-59D 08/05/2013	MW-59D 08/05/2013 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014
Benzene	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00015	<0.00015
Toluene	1	1	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015
Vinyl chloride	0.002	0.002								<0.001				
Xylenes (total)	10	10	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000105	<0.000105	<0.00011	<0.00011
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000295	<0.000295	<0.00031	<0.00031
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000124	<0.000124	<0.00013	<0.00013
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000762	<0.00008	<0.00008
2-Chloronaphthalene	2	5.8	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00008	<0.00008
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00007	0.00046	<0.00005	<0.00005	<0.00005	<0.00005	0.000071 J	<0.0000667	<0.0000667	0.00016 J	0.00007 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.00079	<0.00083	<0.00083
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000533	<0.000533	<0.00056	<0.00056
Acenaphthene	1.5	4.4	<0.00009	<0.00009	0.00095	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00008	<0.00008
Acenaphthylene	1.5	4.4	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000571	<0.00006	<0.00006
Anthracene	7.3	22	<0.00007	<0.00007	0.00069	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000476	<0.0000476	<0.00005	<0.00005
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00007	0.00027	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00008	<0.00008
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000124	<0.000124	<0.00013	<0.00013
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00031	<0.00049	<0.0015	0.0011	0.0018	<0.00094	<0.00068	<0.00014	<0.000352	0.000538	0.000805 J	0.000812 J
Chrysene	0.91	2	<0.00007	<0.00007	0.00024	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00008	<0.00008
Dibenzofuran	0.098	0.29	<0.00008	<0.00008	0.0011	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000762	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000105	<0.000105	<0.00011	<0.00011
Fluoranthene	0.98	2.9	<0.00007	<0.00007	0.0018	<0.00005	<0.00005	<0.00005	<0.00005	0.000075 J	<0.00005	<0.0000667	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.00007	<0.00007	0.00079	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000667	<0.00007	<0.00007
Naphthalene	0.49	1.5	<0.00022	<0.0001	0.0034	<0.00005	<0.00005	<0.00005	<0.000064	<0.00036	<0.0000762	<0.0000762	0.00226 J	0.00008 J
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000105	<0.000105	<0.00011	<0.00011
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000952	<0.0001	<0.0001
Pentachlorophenol	0.001	0.001	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000581	<0.000581	<0.00061	<0.00061
Phenanthrene	0.73	2.2	<0.00007	<0.00007	0.0037	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000571	<0.00006	<0.00006
Phenol	7.3	22	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00014 J	0.000087 J	<0.00005	<0.0000381	<0.0000381	<0.00004	<0.00004
Pyrene	0.73	2.2	<0.00007	<0.00007	0.0011	<0.00005	<0.00005	<0.00005	0.000054 J	<0.00005	<0.000105	<0.000105	<0.00011	<0.00011
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-59D 01/23/2014	MW-59D 01/23/2014 Duplicate	MW-59D 08/28/2014	MW-59D 08/28/2014 Duplicate	MW-59D 02/07/2018	MW-59D 02/07/2018 Duplicate	MW-59D 03/26/2018	MW-59D 03/26/2018 Duplicate	MW-59D 06/01/2018	MW-59D 06/01/2018 Duplicate	MW-59D 01/24/2019	MW-59D 01/24/2019 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	0.000135 J	0.000114 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.00018	<0.00018	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.00019	<0.00019	<0.00011	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.00022	<0.00022	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.00017	<0.00017	0.000258 J	0.000249 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00058	<0.00058	<0.00026	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000104	<0.000104	<0.00011	<0.00011	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.000292	<0.000292	<0.00031	<0.00031	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000123	<0.000123	<0.00013	<0.00013	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000066	<0.000066	0.000334 J	0.00007 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000083 J	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000783	<0.000783	<0.00083	<0.00083	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000528	<0.000528	<0.00056	<0.00056	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.00012	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.0000566	<0.0000566	<0.00006	<0.00006	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	<0.000425	<0.000472	<0.00005	<0.00005	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.000024 J	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000123	<0.000123	<0.00013	<0.00013	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000425 J	0.000364 J	0.00306 J	0.00598 J	0.000063 J	0.000054 J	<0.000037	0.000095 J	0.00024	0.00026	<0.000037	<0.000037
Chrysene	0.91	2	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.0000755	<0.0000755	<0.00008	<0.00008	<0.00002	<0.00002	<0.00002	<0.00002	0.000092 J	0.00024 J	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000825	<0.000827	<0.00011	<0.00011	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0000789 J	0.0000746 J	0.00018 J	0.00007 J	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.000066	0.0000688 J	<0.00007	<0.00007	<0.00003	<0.00003	<0.00003	<0.00003	0.000064 J	0.00017 J	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.000727	<0.000796	0.00576 J	0.00008 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.0002	<0.00002	<0.00002
Nitrobenzene	0.049	0.15	<0.000104	<0.000104	<0.00011	<0.00011	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.0000943	<0.0000943	<0.0001	<0.0001	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000575	<0.000575	<0.00061	<0.00061	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.000455	<0.000464	0.00018 J	0.0000608 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00018	<0.000021	<0.000021
Phenol	7.3	22	<0.0000377	<0.0000377	<0.00004	<0.00004	<0.000035	<0.000035	<0.000035	<0.000035	0.00063 J	0.00097 J	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000104	<0.000104	0.000131 J	<0.00011	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01					<0.0004	<0.0004	<0.0004	<0.0004	0.00111 J	0.00101 J	0.000765 J	0.000637 J

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-59D 07/31/2019	MW-59D 07/31/2019 Duplicate	MW-59D 01/16/2020	MW-59D 01/16/2020 Duplicate	MW-60A 02/04/2009	MW-60A 01/20/2010	MW-60A 06/24/2010	MW-60A 01/19/2011	MW-60A 07/18/2011	MW-60A 02/07/2012	MW-60A 07/23/2012	MW-60A 02/14/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	
Vinyl chloride	0.002	0.002									<0.0005	<0.001	<0.001	<0.0005
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00038
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	0.001	<0.00018
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00032
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.00029
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00019
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	0.00028	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0021	0.000146 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00016
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00033
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000036	<0.000027	0.00045	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	0.0012	<0.00016
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00016
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	0.00034	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00027	<0.00044
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00025
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00013
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00019
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	<0.000037	<0.000037	0.002	<0.0025	<0.0002	0.0031	<0.00017	0.00023	<0.0001	<0.00059
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00024
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.000028	<0.00002	0.00035	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	0.00099	<0.00016
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	0.0023	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.000076	<0.00187
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	0.00039	<0.00007	0.0003	0.00029	<0.00005	0.00028	0.0003	<0.00031
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.000031	<0.00003	0.00044	<0.00007	<0.00007	<0.00007	<0.00005	0.00016 J	0.00089	<0.00012
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.000013	<0.00002	0.0095	<0.0001	0.0015	<0.0001	<0.00005	<0.00005	0.025	0.00043 J
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0002
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00033
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00096
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000043	<0.000021	0.0011	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.0015	<0.00029
Phenol	7.3	22	0.00011 J	<0.000035	<0.000035	<0.000035	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	0.000275 J
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	0.00029	<0.00007	0.0002 J	0.00079	<0.00005	0.0013	0.00033	<0.00033
<b>Metals</b>														
Arsenic	0.01	0.01	<0.0004	0.000502 J	0.000508 J	0.000478 J								

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-60A 04/02/2013	MW-60A 08/02/2013	MW-60A 01/15/2014	MW-60A 07/16/2014	MW-60A 02/08/2018	MW-60A 03/20/2018	MW-60A 05/25/2018	MW-60A 01/11/2019	MW-60A 07/17/2019	MW-60AR 03/20/2020	MW-61A 02/03/2009	MW-61A 01/20/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005
Benzene	0.005	0.005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005
Toluene	1	1	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005
Vinyl chloride	0.002	0.002	<0.00011	<0.00011	<0.00018	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005
Xylenes (total)	10	10	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000106	<0.000105	<0.000104	<0.000104	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.000298	<0.000295	<0.000292	<0.000292	<0.00004	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000125	<0.000124	<0.000123	<0.000123	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0000769	<0.0000762	<0.0000755	<0.0000755	<0.000042	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0000769	<0.0000762	<0.0000755	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.0000667	<0.0000667	0.000143 J	0.000516	<0.000019	<0.000019	<0.000019	<0.000019	0.000072 J	<0.000019	0.00041	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000798	<0.00079	<0.000783	<0.000783	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000538	<0.000533	<0.000528	<0.000528	<0.000047	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.0000769	<0.0000762	0.000157 J	0.000167 J	<0.000027	<0.000028	<0.000027	<0.000027	<0.000027	<0.000027	0.00017 J	<0.00009
Acenaphthylene	1.5	4.4	<0.0000577	<0.0000571	<0.0000566	<0.0000566	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00006	<0.00007
Anthracene	7.3	22	<0.0000481	0.0000883 J	0.000158 J	<0.0000472	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.0000769	<0.0000762	<0.0000755	<0.0000755	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.0000769	<0.0000762	<0.0000755	<0.0000755	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000125	<0.000124	<0.000123	<0.000123	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000356	<0.000352	<0.000349	<0.000349	<0.000037	<0.000038	0.000076 J	<0.000092	<0.000055	0.000063 J	0.0017	<0.002
Chrysene	0.91	2	<0.0000769	<0.0000762	<0.0000755	<0.0000755	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.0000769	<0.0000762	0.000145 J	0.000116 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000106	<0.000105	<0.000104	<0.000104	<0.00002	<0.00002	<0.00002	0.000064 J	0.000075 J	<0.00002	0.011	<0.00007
Fluoranthene	0.98	2.9	<0.0000673	<0.0000667	0.0000894 J	<0.000066	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.0000673	<0.0000667	0.000162 J	<0.000066	<0.00003	0.000054 J	<0.00003	<0.00003	<0.00003	<0.00003	0.00011 J	<0.00007
Naphthalene	0.49	1.5	<0.0000769	<0.0000762	0.000668 J	0.00653	<0.00002	<0.00002	<0.00002	<0.00002	<0.00013	<0.00002	0.0066	<0.0001
Nitrobenzene	0.049	0.15	<0.000106	<0.000105	<0.000104	<0.000104	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0000962	<0.0000952	<0.0000943	<0.0000943	<0.000025	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000587	<0.000581	<0.000575	<0.000575	<0.000079	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.0000577	<0.0000571	0.000345 J	<0.0000566	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00021	<0.00007
Phenol	7.3	22	<0.0000385	<0.0000381	<0.0000377	<0.0000377	<0.000035	<0.000036	<0.000035	<0.000035	<0.000035	<0.000035	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.000106	<0.000105	<0.000104	<0.000104	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01					0.000649 J	0.000706 J	0.000636 J	0.00453	0.00044 J	0.00189 J		

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-61A 07/01/2010	MW-61A 01/27/2011	MW-61A 07/21/2011	MW-61A 02/07/2012	MW-61A 07/27/2012	MW-61A 04/02/2013	MW-61A 08/01/2013	MW-61A 01/23/2014	MW-61A 08/28/2014	MW-61A 02/08/2018	MW-61A 03/20/2018	MW-61A 05/25/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001
Toluene	1	1	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00018	<0.00011	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000106	<0.000105	<0.000104	<0.00011	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000298	<0.000295	<0.000292	<0.00031	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.000124	<0.000123	<0.00013	<0.000058	<0.000059	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.000021	<0.000021	0.000021 J
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000667	<0.0000666	<0.00007	<0.000019	<0.000019	0.000019 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000798	<0.00079	<0.000783	<0.00083	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000538	<0.000533	<0.000528	<0.00056	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.000027	<0.000027	0.000027 J
Acenaphthylene	1.5	4.4	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000577	<0.0000571	<0.0000566	<0.00006	<0.000015	<0.000015	0.000015 J
Anthracene	7.3	22	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000481	<0.0000476	<0.0000434	<0.00005	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.00005	<0.000051	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.000124	<0.000123	<0.00013	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00021	<0.0002	0.00023	0.00038	0.00027	<0.000356	<0.000352	0.00163 J	0.000536 J	<0.000037	<0.000037	0.000054 J
Chrysene	0.91	2	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000769	<0.0000762	<0.0000755	<0.00008	<0.00002	<0.00002	0.00002 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000106	<0.000105	<0.0000827	<0.00011	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000673	<0.0000667	0.0000806 J	<0.00007	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000673	<0.0000667	<0.0000666	<0.00007	<0.00003	<0.00003	0.00003 J
Naphthalene	0.49	1.5	<0.00018	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00012	<0.0000668	<0.00008	<0.00002	<0.00002	<0.00002
Nitrobenzene	0.049	0.15	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000106	<0.000105	<0.000104	<0.00011	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000962	<0.0000952	<0.0000943	<0.0001	<0.000025	<0.000025	0.00015 J
Pentachlorophenol	0.001	0.001	<0.00032	<0.00008	<0.00005	<0.00005	<0.00005	<0.000587	<0.000581	<0.000575	<0.00061	<0.000079	<0.00008	<0.000079
Phenanthrene	0.73	2.2	<0.00007	<0.00007	<0.00005	<0.00005	0.00016 J	<0.0000577	0.0000586 J	<0.000046	<0.00006	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000385	<0.0000381	<0.0000377	<0.00004	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000106	<0.000105	<0.000104	<0.00011	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01										0.000743 J	0.00116 J	0.00172 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-61A 01/23/2019	MW-61A 07/17/2019	MW-61A 01/16/2020	MW-62B 02/04/2009	MW-62B 01/21/2010	MW-62B 07/14/2010	MW-62B 01/27/2011	MW-62B 07/27/2011	MW-62B 08/25/2011	MW-62B 02/08/2012	MW-62B 07/26/2012	MW-62B 02/11/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001		<0.001	<0.0005	<0.00014
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	0.0043 J		<0.001	0.002 J	<0.00008
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005	<0.001		<0.001	<0.0005	<0.00012
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	0.00071 J	<0.0005	<0.0005	<0.0005	0.041		<0.0011	0.0021 J	<0.00011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013		<0.0013	<0.001	<0.00015
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	0.0095		<0.001	0.0012 J	<0.00015
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002									
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.001	<0.001	<0.001	<0.001	0.025		<0.0031	0.0053 J	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005		<0.00005	<0.00005	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005		<0.00005	<0.00005	<0.000282
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.000118
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006		<0.00006	<0.00006	<0.0000727
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005		<0.00005	<0.00005	<0.0000727
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	0.00012 J	0.0016	0.00064	<0.00007	<0.00005		<0.00005	<0.00006	<0.000174
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008		<0.00008	<0.00008	<0.000755
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00005	<0.000509
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000052	0.0078	0.039	0.00041	<0.00009	0.21		0.026	0.085	0.000242 J
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.00006	0.00066	<0.00007	<0.00007	0.0026		0.0013	0.00084	0.000112 J
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	0.00024	0.0011	<0.00007	<0.00007	0.013		<0.00005	0.0032	0.000723
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00005	<0.0000727
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005	<0.0000727
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.000118
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	0.000081 J	0.00041	<0.00098	0.0016	0.00022	0.00042		<0.00029	0.00013 J	<0.000336
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.00005	<0.0000727
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	0.0024	0.013	0.00034	<0.00008	0.15	0.23	0.00012 J	0.038	0.000174 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.000056 J	<0.00002	0.00065	<0.00007	<0.00007	<0.00007	<0.00005		<0.00005	<0.000078	<0.0001
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00006	0.00012 J	0.0011	<0.00007	0.00014 J	0.0079		0.00053	0.004	0.00033 J
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	0.0012	0.015	0.00016 J	<0.00007	0.058		0.0002	0.0087 J	<0.0000636
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.00021	0.0027	0.00028	0.0096	<0.0001	0.035		<0.00021	0.0056	<0.00129
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.0001
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005		<0.00005	<0.00005	<0.0000909
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005		<0.00005	<0.00005	<0.000555
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.0001	0.00087	0.0025	0.00025	<0.00007	0.035		0.00014 J	0.0026	<0.000472
Phenol	7.3	22	<0.000035	0.00011 J	<0.000035	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005		0.000053 J	<0.00005	<0.0000364
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000059	<0.00007	0.00047	<0.00007	0.000077 J	0.0033		0.00037	0.0021	0.000387 J
<b>Metals</b>														
Arsenic	0.01	0.01	0.00069 J	0.00117 J	0.00107 J									

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-62B 08/02/2013	MW-62B 01/29/2014	MW-62B 07/29/2014	MW-62B 01/24/2018	MW-62B 03/20/2018	MW-62B 05/24/2018	MW-62B 01/23/2019	MW-62B 07/16/2019	MW-62B 01/27/2020	MW-64A 02/04/2009	MW-64A 01/21/2010	MW-64A 07/14/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005
Benzene	0.005	0.005	<0.00008	<0.00008	<0.00008	<0.0002	<0.0002	0.0021	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005
Chlorobenzene	0.1	0.1	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005
Ethylbenzene	0.7	0.7	<0.00011	<0.00011	<0.00011	<0.0003	<0.0003	0.013	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005
Methylene chloride	0.005	0.005	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0005
Toluene	1	1	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	0.003	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00026	<0.00026	<0.00026	<0.0003	<0.0003	0.014	<0.0003	<0.0003	<0.0003	<0.001	<0.001	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000105	<0.000104	<0.000109	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	0.49	1.5	<0.000295	<0.000292	<0.000307	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008
2,4-Dinitrotoluene	0.0013	0.003	<0.000124	<0.000123	<0.000129	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.0000762	<0.0000755	<0.0000792	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	2	5.8	<0.0000762	<0.0000755	<0.0000792	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00016 J	<0.00012	<0.0001	<0.0001
2-Methylnaphthalene	0.098	0.29	<0.0000667	<0.000066	<0.0000693	<0.000019	<0.000019	0.00011	<0.000019	0.000096 J	0.00011	0.00014 J	<0.00007	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000079	<0.0000783	<0.0000822	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.000533	<0.000528	<0.000554	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007
Acenaphthene	1.5	4.4	<0.0000762	<0.0000755	0.000235 J	0.00006 J	0.023	0.11	<0.000027	0.000099 J	0.04	0.00029	<0.00009	<0.00009
Acenaphthylene	1.5	4.4	<0.0000571	<0.0000566	<0.0000594	0.000061 J	0.00063	0.0013	<0.000015	<0.000015	0.00029	<0.00006	<0.00007	<0.00007
Anthracene	7.3	22	<0.0000476	<0.0000472	0.0000699 J	0.00036	0.00051	0.0044	<0.000014	<0.000014	0.00096	0.00016 J	<0.00007	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.0000762	<0.0000755	<0.0000792	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.0000762	<0.0000755	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000124	<0.000123	<0.000129	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000352	<0.000349	<0.000366	0.00026	0.00018 J	0.00013 J	<0.000037	<0.000037	0.000095 J	0.0004	<0.00016	0.002
Chrysene	0.91	2	<0.0000762	<0.0000755	<0.0000792	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000027 J	<0.00007	<0.00007	<0.00007
Dibenzofuran	0.098	0.29	<0.0000762	<0.0000755	0.0000916 J	<0.00002	0.0031	0.048	<0.00002	0.000048 J	0.0042	0.00012 J	<0.00008	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	0.000107 J	<0.000104	<0.000109	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00014 J	0.02	<0.00007	<0.00007
Fluoranthene	0.98	2.9	<0.0000667	<0.000066	<0.0000693	0.00052	0.0015	0.0041	<0.00001	0.000017 J	0.002	0.00076	<0.00007	<0.00007
Fluorene	0.98	2.9	<0.0000667	<0.000066	0.000126 J	0.00015	0.0047	0.034	<0.00003	0.000036 J	0.011	0.00018 J	<0.00007	<0.00007
Naphthalene	0.49	1.5	<0.0000762	<0.0000755	<0.0000792	<0.00002	<0.00002	0.038	<0.00002	0.00028	0.001	0.00092	<0.0001	<0.0001
Nitrobenzene	0.049	0.15	<0.000105	<0.000104	<0.000109	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.0000952	<0.0000943	<0.000099	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.001	0.001	<0.000581	<0.000575	<0.000604	<0.000079	0.00033	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008
Phenanthrene	0.73	2.2	<0.0000571	<0.0000566	<0.000144	<0.000021	0.00023	0.0083	<0.000021	0.000021 J	0.00087	0.00055	<0.00007	<0.00007
Phenol	7.3	22	<0.0000381	<0.0000377	<0.0000396	<0.000035	<0.000035	<0.000035	<0.000035	0.00056	<0.000035	<0.00007	<0.00007	<0.00007
Pyrene	0.73	2.2	<0.000105	<0.000104	<0.000109	0.00039	0.00077	0.002	<0.000019	<0.000019	0.0013	0.00063	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01				0.00842	0.0173	0.028	<0.0004	0.00194 J	0.0161			

Notes:  
1. All values in milligrams per liter (mg/L).  
2. Concentrations > RAL and non-detects are highlighted light gray.  
3. Concentrations > C/I AL and non-detects are highlighted dark gray  
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.  
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial  
6. J = Estimated value, < = not detected at the specified detection limit.  
7. MW-32A was screened in the B-CZ & replaced with MW-32AR  
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-64A 01/27/2011	MW-64A 07/27/2011	MW-64A 02/08/2012	MW-64A 07/25/2012	MW-64A 04/01/2013	MW-64A 08/06/2013	MW-64A 01/29/2014	MW-64A 07/29/2014	MW-64A 01/31/2018	MW-64A 03/25/2018	MW-64A 05/31/2018	MW-64A 01/23/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	0.000154 J	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00015	<0.00015	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.00005	<0.00005	<0.000106	<0.00011	<0.000104	<0.000108	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00008	<0.00005	<0.00005	<0.00005	<0.000298	<0.00031	<0.000292	<0.000304	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.00013	<0.000123	<0.000127	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000769	<0.00008	<0.0000755	<0.0000784	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.00005	0.000053 J	<0.00005	<0.0000673	<0.00007	<0.000066	<0.0000686	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.0000798	<0.00083	<0.0000783	<0.000814	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.00005	<0.00005	<0.000538	<0.00056	<0.000528	<0.000549	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.00009	<0.00005	0.0096	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000784	<0.000027	0.035	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.00007	<0.00005	0.0005	<0.00005	<0.0000577	<0.00006	<0.0000566	<0.0000588	<0.000015	0.00086	<0.000015	<0.000015
Anthracene	7.3	22	<0.00007	0.00036	<0.00005	<0.00005	0.000158 J	<0.00005	<0.0000472	0.000127 J	<0.000014	0.00056	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000784	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000784	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.00005	<0.00005	<0.000125	<0.00013	<0.000123	<0.000127	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00049	0.00076	<0.00013	0.00021	<0.000356	<0.00037	<0.000349	<0.000363	<0.0001	<0.000037	0.00023	<0.000037
Chrysene	0.91	2	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	0.0013	<0.00005	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000784	0.00003 J	0.00038	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	0.000079 J	<0.00005	0.000084 J	<0.000106	<0.00011	0.000117 J	<0.000108	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00007	0.00057	0.00021	<0.00005	<0.0000673	<0.00007	<0.000066	<0.0000686	<0.00001	0.00095	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00007	<0.00005	0.00012 J	<0.00005	<0.0000673	<0.00007	<0.000066	<0.0000686	<0.00003	0.0013	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.0001	<0.00005	<0.00063	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.000317	<0.000094	0.00058	<0.00002	<0.00002
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.00005	<0.00005	<0.000106	<0.00011	<0.000104	<0.000108	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000962	<0.0001	<0.0000943	<0.000098	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00005	<0.00005	<0.00005	<0.000587	<0.00061	<0.000575	<0.000598	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000577	<0.00006	<0.0000566	<0.0000588	0.000032 J	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.00007	0.000077 J	<0.00005	<0.00005	<0.0000385	<0.00004	<0.0000377	<0.0000392	<0.000035	<0.000035	0.00051	<0.000035
Pyrene	0.73	2.2	<0.00007	0.00042	0.00013 J	<0.00005	<0.000106	<0.00011	<0.000104	<0.000108	<0.000019	0.00059	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01									0.000419 J	0.0117	0.00111 J	<0.0004

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-64A 07/11/2019	MW-64A 01/27/2020	MW-65D 02/05/2009	MW-65D 01/21/2010	MW-65D 07/01/2010	MW-65D 01/26/2011	MW-65D 07/27/2011	MW-65D 02/14/2012	MW-65D 07/23/2012	MW-65D 02/11/2013	MW-65D 08/05/2013	MW-65D 01/21/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	0.0013 J	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019
Methylene chloride	0.005	0.005	<0.001	<0.001	0.00095 J	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022
Toluene	1	1	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.000104
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000295	<0.000304	<0.000292
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.000123
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000784	<0.0000755
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755
2-Methylnaphthalene	0.098	0.29	<0.000019	0.00037	0.00012 J	<0.00007	0.00014 J	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	0.0000808 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.000814	<0.000783
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.000549	<0.000528
Acenaphthene	1.5	4.4	0.0003	0.00016	0.00019 J	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000588	<0.0000566
Anthracene	7.3	22	<0.000014	0.000096 J	0.000078 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000476	<0.000049	0.0000574 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000127	<0.000123
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.00031	0.0019	<0.0027	<0.001	0.001	<0.00013	<0.00025	0.000593	<0.000363	<0.000349	
Chrysene	0.91	2	<0.000021	<0.000021	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755
Dibenzofuran	0.098	0.29	<0.00002	0.0001 J	0.00016 J	0.00012 J	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000784	<0.0000755
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.000079 J	0.00029	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	0.000135 J	0.000148 J	<0.000104
Fluoranthene	0.98	2.9	<0.00001	0.00028	0.000097 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	0.000117 J
Fluorene	0.98	2.9	0.000031 J	0.000082 J	0.00016 J	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000667	<0.0000686	<0.000066
Naphthalene	0.49	1.5	<0.00033	0.002	0.00051	0.00026	<0.00059	0.00019 J	<0.00005	<0.00005	<0.000094	<0.0000762	<0.0000784	0.000529 J
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.000104
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.000098	<0.0000943
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.000598	<0.000575
Phenanthrene	0.73	2.2	<0.000021	<0.000021	0.00014 J	<0.00007	<0.00007	<0.00007	0.000065 J	<0.00005	<0.00005	<0.0000571	0.000093 J	0.000294 J
Phenol	7.3	22	<0.000035	<0.000035	<0.00007	0.0015	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.0000381	<0.0000392	<0.0000377
Pyrene	0.73	2.2	0.00015	0.00015	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000105	<0.000108	<0.000104
<b>Metals</b>														
Arsenic	0.01	0.01	0.00939	0.00126 J										

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-65D 08/28/2014	MW-65D 02/07/2018	MW-65D 03/26/2018	MW-65D 06/01/2018	MW-65D 01/24/2019	MW-65D 07/31/2019	MW-65D 01/16/2020	MW-66D 02/05/2009	MW-66D 01/20/2010	MW-66D 07/01/2010	MW-66D 07/27/2011	MW-66D 02/14/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Benzene	0.005	0.005	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Chlorobenzene	0.1	0.1	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Ethylbenzene	0.7	0.7	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011
Methylene chloride	0.005	0.005	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013
Toluene	1	1	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Vinyl chloride	0.002	0.002	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Xylenes (total)	10	10	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.001	<0.001	<0.001	<0.0031	<0.0031
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00011	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	0.49	1.5	<0.00031	<0.00004	<0.00004	<0.00004	<0.00009	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005
2,4-Dinitrotoluene	0.0013	0.003	<0.00013	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.0013	0.003	<0.00008	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	2	5.8	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	0.098	0.29	<0.00007	<0.000019	<0.000019	0.000089 J	0.00016	<0.000019	<0.000019	0.00062	<0.00007	<0.00007	<0.00005	<0.00005
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00083	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	0.049	0.15	<0.00056	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Acenaphthene	1.5	4.4	<0.00008	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.0004	<0.00009	<0.00009	<0.00005	<0.00005
Acenaphthylene	1.5	4.4	<0.00006	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00006	<0.00007	<0.00007	<0.00005	<0.00005
Anthracene	7.3	22	<0.00005	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.00015 J	<0.00007	<0.00007	0.00022	0.00027
Benzo(a)anthracene	0.0091	0.02	<0.00008	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00007	<0.00007	<0.00007	0.00011 J	0.00012 J
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	0.00016 J	0.00013 J
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00013	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00244	<0.000037	<0.000037	0.0002 J	0.00006 J	0.00016 J	0.00026	<b>0.0064</b>	<0.0028	<0.00096	0.0019	<0.0002
Chrysene	0.91	2	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00007	<0.00007	<0.00007	0.00046	0.00052
Dibenzofuran	0.098	0.29	<0.00008	<0.00002	<0.00002	0.000061 J	0.000039 J	<0.00002	<0.00002	0.00036	<0.00008	0.000083 J	<0.00005	<0.00005
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00011	<0.00002	<0.00002	0.000022 J	<0.00002	<0.00002	<0.00002	0.00044	0.000086 J	<0.00007	0.000056 J	<0.00005
Fluoranthene	0.98	2.9	<0.00007	<0.00001	<0.00001	0.000013 J	0.000027 J	<0.00001	<0.00001	0.00026	<0.00007	<0.00007	0.00035	0.00057
Fluorene	0.98	2.9	<0.00007	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00033	<0.00007	<0.00007	<0.00005	<0.00005
Naphthalene	0.49	1.5	0.00071	<0.00002	<0.00002	0.00029	<0.0026	<0.00002	<0.00002	0.0058	<0.0001	<0.0002	<0.00005	<0.00005
Nitrobenzene	0.049	0.15	<0.00011	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	0.19	0.42	<0.0001	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.001	0.001	<0.00061	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	0.000084 J	<0.00005
Phenanthrene	0.73	2.2	<0.00006	<0.000021	<0.000021	0.000035 J	<0.000021	<0.000021	<0.000021	0.00073	0.00012 J	<0.00007	0.00011 J	0.00011 J
Phenol	7.3	22	<0.00004	<0.000035	<0.000035	<0.000035	<0.00019	<0.000035	<0.000035	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005
Pyrene	0.73	2.2	<0.00011	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.00017 J	<0.00007	<0.00007	0.00036	0.00051
<b>Metals</b>														
Arsenic	0.01	0.01		<0.0004	0.00761	0.00292	0.00202	0.00135 J	0.000507 J					

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-66D 07/23/2012	MW-66D 04/02/2013	MW-66D 08/05/2013	MW-66D 01/29/2014	MW-66D 08/28/2014	MW-66D 10/03/2014	MW-66D 02/07/2018	MW-66D 03/26/2018	MW-66D 06/01/2018	MW-66D 01/24/2019	MW-66D 07/31/2019	MW-66D 01/16/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.00014	<0.00014	<0.00014	<0.00014		<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0005	<0.00008	<0.00008	<0.00008	<0.00008		<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.00012	<0.00012	<0.00012	<0.00012		<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.00011	<0.00011	<0.00011	<0.00011		<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.00015	<0.00015	<0.00015	<0.00015		<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0005	<0.00015	<0.00015	<0.00015	<0.00015		<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002							<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0015	<0.00026	<0.00026	<0.00026	<0.00026		<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.000106	<0.00011	<0.000104	<0.000109	<0.000107	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	<0.000298	<0.00031	<0.000292	<0.000307	<0.000301	<0.0004	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	<0.000125	<0.00013	<0.000123	<0.000129	<0.000126	<0.00058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	<0.0000769	<0.00008	<0.0000755	<0.0000792	<0.0000777	<0.00042	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	<0.0000769	<0.00008	<0.0000755	<0.0000792	<0.0000777	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.000085 J	<0.0000673	<0.00007	<0.000066	0.000211 J	<0.000068	<0.00019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.0000798	<0.00083	<0.0000783	<0.0000822	<0.0000806	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	<0.000538	<0.00056	<0.000528	<0.000554	<0.000544	<0.00047	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.000054 J	<0.0000769	<0.00008	0.000145 J	0.000141 J	<0.0000777	<0.00027	<0.000028	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	0.000081 J	<0.0000577	<0.00006	<0.0000566	0.000411 J	0.000206 J	<0.00015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.00059	<0.0000481	<0.00005	<0.0000472	0.000304	0.00256	<0.00014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	0.00036	<0.0000769	<0.00008	<0.0000755	0.00041 J	0.000245 J	<0.0005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<b>0.00067</b>	<0.0000769	<0.00008	<0.0000755	<b>0.000436 J</b>	<b>0.000439 J</b>	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	<0.000125	<0.00013	<0.000123	<0.000129	<0.000126	<0.0003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0032	<0.000356	<0.00037	<0.000349	0.000585	0.000415 J	<0.00037	<0.000038	0.00019 J	0.00017 J	0.00022	0.00046 J
Chrysene	0.91	2	0.0018	<0.0000769	<0.00008	<0.0000755	0.00104	0.000582	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.000066 J	<0.0000769	<0.00008	<0.0000755	0.000133 J	<0.0000777	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000078	<0.000106	<0.00011	<0.000104	0.000121 J	<0.000107	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0019	<0.0000673	<0.00007	<0.000066	0.00116	0.000346 J	<0.0001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00005	<0.0000673	<0.00007	<0.000066	0.000143 J	<0.000068	<0.0003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.0004	<0.0000769	0.0000999 J	0.000367 J	0.00118	<0.0000777	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00013
Nitrobenzene	0.049	0.15	<0.00005	<0.000106	<0.00011	<0.000104	<0.000109	<0.000107	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.0000962	<0.0001	<0.0000943	<0.000099	<0.0000971	<0.00025	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	<0.000587	<0.00061	<0.000575	<0.000604	<0.000592	<0.00079	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.00058	<0.0000577	<0.00006	0.000132 J	0.000295 J	0.0000729 J	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.00005	<0.0000385	<0.00004	<0.0000377	<0.0000396	<0.0000388	<0.00035	<0.000036	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0019	<0.000106	<0.00011	<0.000104	0.00118	0.000388 J	<0.00019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01							0.000711 J	0.00663	0.00223	0.00204	0.00124 J	0.00138 J

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-68A 05/29/2019	MW-68A 05/29/2019 Duplicate	MW-68A 07/18/2019	MW-68A 01/17/2020	MW-68B 02/16/2012	MW-68B 07/16/2012	MW-68B 02/06/2013	MW-68B 08/08/2013	MW-68B 01/22/2014	MW-68B 01/22/2014 Duplicate	MW-68B 07/24/2014	MW-68B 07/24/2014 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.005	<0.005	<0.014	<0.014	<0.0002	<0.0002	<0.0028	<0.0028
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	2.7	2.4	2.35	2.88	1.5	1.51	2.18	2.1
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.005	<0.005	0.0273 J	<0.012	0.000458 J	0.000468 J	<0.0024	<0.0024
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	0.45	0.49	0.449	0.55	0.364	0.363	0.403	0.453
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.0065	<0.01	<0.015	0.101	<0.00022	<0.00022	<0.003	<0.003
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	0.91	0.93	0.701	0.625	0.329	0.329	0.538	0.57
Vinyl chloride	0.002	0.002								<0.011			0.007 J	0.0022 J
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	1.2	1.3	1.04	1.28	0.857	0.862	1.08	1.22
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.00005	<0.00025	<0.00524	<0.00519	<0.0105	<0.0105	<0.0055	<0.00529
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	0.19	0.27	0.273	<0.0146	0.536	0.457 J	0.445	0.451
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000059	<0.000058	<0.000058	<0.00005	<0.00025	<0.00619	<0.00613	<0.0124	<0.0124	<0.0065	<0.00625
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.00006	<0.0003	<0.00381	<0.00377	<0.00762	<0.00762	<0.004	<0.00385
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.00005	<0.00025	<0.00381	<0.00377	<0.00762	<0.00762	<0.004	<0.00385
2-Methylnaphthalene	0.098	0.29	<0.000019	0.000026 J	0.00012	0.000048 J	0.66	1.3	0.952	1.41	1.1	1.19	0.852	0.906
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.0004	<0.00395	<0.00392	<0.0079	<0.0079	<0.00415	<0.00399
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.00005	<0.00025	<0.0267	<0.0264	<0.0533	<0.0533	<0.028	<0.0269
Acenaphthene	1.5	4.4	0.0012	0.0011	0.0019	0.0003	0.15	0.23	0.261	0.304	0.263	0.26	0.178	0.181
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	0.000054 J	0.0023	0.003	<0.00286	<0.00283	<0.00571	<0.00571	<0.003	<0.00288
Anthracene	7.3	22	<0.000014	<0.000014	0.000028 J	0.000059 J	0.046	0.034	0.0194 J	0.023 J	0.0428 J	0.0292 J	0.0169 J	0.0162 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.000051	<0.00005	<0.00005	0.006	0.0054	<0.00381	<0.00377	0.0123 J	0.00762 J	<0.004	<0.00385
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	0.0017	0.0016	<0.00381	<0.00377	<0.00762	<0.00762	<0.004	<0.00385
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00005	<0.00025	<0.00619	<0.00613	<0.0124	<0.0124	<0.0065	<0.00625
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00015 J	0.000074 J	0.00024	<0.000061	<0.0001	<0.0005	<0.0176	<0.0175	<0.0352	<0.0352	<0.0185	<0.0178
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	0.0052	0.005	<0.00381	<0.00377	0.00806 J	0.00976 J	<0.004	<0.00385
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	0.000082 J	<0.00002	0.19	0.3	0.26	0.325	0.284	0.276	0.198	0.196
Di-n-butylphthalate (DBP)	2.4	7.3	0.00087	0.00064	0.00015 J	0.000034 J	<0.00005	<0.00025	<0.00524	<0.00519	<0.0105	<0.0105	<0.0055	<0.00529
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	0.00013	0.05	0.044	<0.00333	0.00764 J	0.052	0.0357 J	0.00825 J	0.00751 J
Fluorene	0.98	2.9	0.000044 J	0.000042 J	0.00012	0.000069 J	0.096	0.13	0.118	0.154	0.149	0.143	0.0966	0.0953
Naphthalene	0.49	1.5	0.00019	0.00025	0.00035	<0.00022	14	26	11.8	31.2	17	17.6	10.5	12.6
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.00005	<0.00025	<0.00524	<0.00519	<0.0105	<0.0105	<0.0055	<0.00529
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	0.0011	<0.00025	<0.00476	<0.00472	<0.00952	<0.00952	<0.005	<0.00481
Pentachlorophenol	0.001	0.001	<0.000079	<0.00008	<0.000079	<0.000079	<0.00005	<0.00025	<0.029	<0.0288	<0.0581	<0.0581	<0.0305	<0.0293
Phenanthrene	0.73	2.2	<0.000021	<0.000021	0.000062 J	0.00016	0.19	0.24	0.12	0.136	0.263	0.196	0.106	0.103
Phenol	7.3	22	0.000041 J	<0.000035	<0.000035	<0.000035	0.035	0.058	0.0421	0.0795	0.0862	0.0929	0.00988 J	0.00192 J
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	0.000068 J	0.031	0.024	<0.00524	<0.00519	0.0341 J	0.0227 J	<0.0055	<0.00529
<b>Metals</b>														
Arsenic	0.01	0.01	0.00966	0.00894	0.0353	0.0423								

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-68B 01/29/2018	MW-68B 01/29/2018 Duplicate	MW-68B 03/21/2018	MW-68B 03/21/2018 Duplicate	MW-68B 06/06/2018	MW-68B 06/06/2018 Duplicate	MW-68B 01/15/2019	MW-68B 01/15/2019 Duplicate	MW-68B 07/18/2019	MW-68B 07/18/2019 Duplicate	MW-68B 01/23/2020 DNAPL	MW-68C 07/15/2010
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.002	<0.002	<0.005	<0.0005
Benzene	0.005	0.005	2.1	2.2	1.4	1.4	1.9	2	2	1.9	1.4	1.4	1.4	0.00081 J
Chlorobenzene	0.1	0.1	<0.0015	<0.0015	<0.0003	<0.0003	<0.0003	<0.0003	0.00056 J	0.00056 J	<0.003	<0.003	<0.0075	<0.0005
Ethylbenzene	0.7	0.7	0.61	0.56	0.29	0.28	0.5	0.5	0.5	0.49	0.52	0.51	0.39	<0.0005
Methylene chloride	0.005	0.005	<0.005	<0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.01	<0.01	<0.025	<0.0005
Toluene	1	1	0.45	0.42	0.2	0.19	0.45	0.44	0.086	0.084	0.37	0.37	0.23	<0.0005
Vinyl chloride	0.002	0.002											<0.005	
Xylenes (total)	10	10	1.6	1.5	0.83	0.84	1.4	1.4	1.2	1.2	1.5	1.4	1.1	<0.001
<b>Semivolatiles Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.0001
2,4-Dimethylphenol	0.49	1.5	0.051 J	0.071 J	0.07	0.065	0.075 J	0.23 J	0.05	0.058	0.076	0.076	0.075	<0.0008
2,4-Dinitrotoluene	0.0013	0.003	<0.00058	<0.00058	<0.00059	<0.00059	<0.00058	<0.00058	<0.00058	0.0013 J	<0.00058	<0.00058	<0.00058	<0.00009
2,6-Dinitrotoluene	0.0013	0.003	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00007
2-Chloronaphthalene	2	5.8	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.0001
2-Methylnaphthalene	0.098	0.29	0.6	0.6	0.67	0.67	1.4	1.3	0.33	0.31	0.72	0.66	1.1 J	<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0011 J	<0.0002	<0.0002	<0.00002	<0.00008
4-Nitrophenol	0.049	0.15	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047	0.0074 J	<0.00047	<0.00047	<0.00047	<0.00007
Acenaphthene	1.5	4.4	0.13	0.15	0.21	0.22	0.34	0.36	0.1	0.13	0.2	0.17	0.44	<0.00009
Acenaphthylene	1.5	4.4	0.002	0.0024	0.0021	0.0016	0.0022	0.0029	0.0012	0.0014	0.0027	0.0022	0.0037	<0.00007
Anthracene	7.3	22	0.014	0.014	0.014	0.012	0.015	0.021	0.008	0.0098	0.03 J	0.022 J	0.15	<0.00007
Benzo(a)anthracene	0.0091	0.02	<0.0005	<0.0005	<0.00051	<0.00051	<0.0005	<0.0005	<0.0005	<0.0005	0.0051 J	0.003 J	0.047	<0.00007
Benzo(a)pyrene	0.0002	0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0018	0.0012	0.0089	<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0036	<0.0003	<0.0003	<0.00003	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00037	<0.00037	<0.00037	<0.00037	<0.00037	<0.00037	<0.00037	<0.00037	<0.00037	<0.00037	0.0011	0.00098
Chrysene	0.91	2	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	0.005 J	0.0029 J	0.04	<0.00007
Dibenzofuran	0.098	0.29	0.16	0.16	0.24	0.25	0.38	0.38	0.1	0.12	0.21	0.18	0.5	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.00018 J	<0.00007
Fluoranthene	0.98	2.9	0.0052	0.0049	0.006	0.005	0.0063	0.0088	0.0031	0.0039	0.043 J	0.027 J	0.31	<0.00007
Fluorene	0.98	2.9	0.082	0.086	0.092	0.078	0.082 J	0.18 J	0.057	0.065	0.11	0.083	0.32	<0.00007
Naphthalene	0.49	1.5	9.2	10	12	13	23	23	5	3.8	9.6	9.9	12	0.00083
Nitrobenzene	0.049	0.15	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00024	<0.00009
N-Nitrosodiphenylamine	0.19	0.42	<0.00025	<0.00025	<0.00025	<0.00025	<0.00025	<0.00025	<0.00025	0.0051	<0.00025	<0.00025	<0.00025	<0.00009
Pentachlorophenol	0.001	0.001	<0.00079	<0.00079	<0.0008	<0.0008	<0.00079	<0.00079	<0.00079	<0.00079	<0.00079	<0.00079	<0.00079	<0.00008
Phenanthrene	0.73	2.2	0.099	0.098	0.15	0.088	0.096 J	0.22 J	0.065	0.073	0.21	0.16	0.95	<0.00007
Phenol	7.3	22	<0.00035	<0.00035	<0.00035	<0.00035	<0.00035	<0.00035	0.0019 J	0.0015 J	<0.00035	<0.00035	<0.00035	0.0005
Pyrene	0.73	2.2	0.0025	0.0024	0.0032	0.0029	0.0045 J	0.0058 J	0.0015	0.0018	0.025 J	0.015 J	0.19	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01	0.0114	0.0117	0.014	0.0151	0.0112	0.0107	0.0123	0.0125	0.0134	0.0129	0.00944 J	

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-68C 01/25/2011	MW-68C 07/21/2011	MW-68C 02/16/2012	MW-68C 07/17/2012	MW-68C 02/06/2013	MW-68C 08/07/2013	MW-68C 01/22/2014	MW-68C 07/24/2014	MW-68C 08/28/2014	MW-68C 01/29/2018	MW-68C 03/21/2018	MW-68C 06/06/2018
<b>Volatiles Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.00014	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	0.0021 J	0.0032 J	<b>0.0069</b>	<b>0.0079</b>	0.00134	0.00364	0.00225	<b>0.0073</b>	0.00118	0.0028	0.0049	<0.0002
Chlorobenzene	0.1	0.1	<0.0005	<0.001	<0.001	<0.0005	<0.000352	<0.00012	<0.00018	<0.00012	<0.00012	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0005	<0.0011	<0.0011	<0.0005	0.000363 J	0.000517 J	0.00024 J	0.000419 J	0.00014 J	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.00015	<0.001	<0.001	<0.001
Toluene	1	1	0.00067 J	0.0011 J	0.0019 J	0.0023 J	0.000632 J	<0.0016	0.00059	0.00138	0.000442 J	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002								<0.00011	<0.00011			
Xylenes (total)	10	10	<0.001	<0.0031	<0.0031	<0.0015	0.000873 J	0.000879 J	<0.00058	0.000649 J	<0.00026	<0.0003	0.00046 J	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000104		<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.00012 J	0.00031	0.00095	0.0014	<b>R</b>	<0.000292	0.000454 J	<0.000292		<0.00004	0.0015	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.000123		<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.0000755	<0.0000762	<0.0000755		<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000762	<0.0000755		<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00016 J	0.00024	<0.00011	0.0025	0.00132	0.000301 J	0.00331	0.00188 J		<0.000019	0.00014	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.00008	<0.00079	<0.000783	<0.00079	<0.000783		<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00007	<0.00005	<0.00005	<0.00005	<b>R</b>	<0.000528	<0.000533	<0.000528		<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.00009	0.00013 J	<0.00005	0.0013	0.000647	<0.0000755	0.00183	0.000235 J		<0.000027	0.00017	<0.000027
Acenaphthylene	1.5	4.4	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000566	<0.0000571	<0.0000566		<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	<0.00007	<0.00005	<0.00005	0.00089	<0.0000476	<0.0000472	0.00106	<0.0000472		<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00007	<0.00005	<0.00005	0.00018 J	<0.0000762	<0.0000755	0.000276 J	<0.0000755		<0.000005	<0.000005	<0.000005
Benzo(a)pyrene	0.0002	0.0002	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	0.000171 J	<0.0000755		<0.000002	<0.000002	<0.000002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000124	<0.000123		<0.000003	<0.000003	<0.000003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.006	0.001	<0.0015	0.0018	0.000637	0.00157 J	<0.000352	<0.000349		0.00015 J	<0.000037	<0.000056
Chrysene	0.91	2	<0.00007	<0.00005	<0.00005	0.00016 J	<0.0000762	<0.0000755	0.000301 J	<0.0000755		<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00008	0.0002 J	<0.000078	0.0018	0.000168 J	<0.0000755	0.00192	0.0000942 J		<0.000002	<0.000002	<0.000002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00007	<0.00005	<0.00005	0.00011 J	<0.000105	<0.000104	<0.000105	<0.000104		<0.000002	<0.000002	0.0001 J
Fluoranthene	0.98	2.9	<0.00007	<0.00005	<0.00005	0.0016	<0.0000667	<0.000066	0.00233	<0.000066		<0.000001	<0.000001	0.00021
Fluorene	0.98	2.9	<0.00007	0.0001 J	<0.00005	0.0012	0.00034 J	0.000135 J	0.00167	0.000155 J		<0.000003	0.00012	<0.000003
Naphthalene	0.49	1.5	0.0014	0.0027	<0.0015	0.015	0.0129	0.00643 J	0.0112	0.00274		0.00088	0.0032	0.00035
Nitrobenzene	0.049	0.15	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000105	<0.000104		<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000943	<0.0000952	<0.0000943		<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00008	<0.00005	<0.00005	<0.00005	<b>R</b>	<0.000575	<0.000581	<0.000575		<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.00007	0.00016 J	<0.00005	0.005	0.000499	<0.0000566	0.00585	<0.0000566		<0.000021	0.0001	<0.000021
Phenol	7.3	22	0.0039	0.0049	0.0074	0.000062 J	<b>R</b>	<0.0000377	<0.0000381	<0.0000377		<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.00007	<0.00005	<0.00005	0.00086	<0.000105	<0.000104	0.0014	<0.000104		<0.000019	<0.000019	0.00015
<b>Metals</b>														
Arsenic	0.01	0.01										<0.0004	0.000618 J	<0.0004

Notes:

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- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
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- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
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**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-68C 01/15/2019	MW-68C 07/18/2019	MW-68C 01/17/2020	MW-69A 07/15/2010	MW-69A 01/19/2011	MW-69A 07/21/2011	MW-69A 02/08/2012	MW-69A 07/24/2012	MW-69A 02/07/2013	MW-69A 08/06/2013	MW-69A 01/24/2014	MW-69A 07/16/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014
Benzene	0.005	0.005	<0.0002	<0.0002	0.00056 J	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	0.00047 J	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00018	<0.00011
Xylenes (total)	10	10	0.0011	<0.0003	0.00062 J	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000105	<0.00011	<0.000106	<0.000104
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	0.0036	<0.00008	<0.00005	<0.00005	0.000078 J	<0.000295	<0.00031	<0.000298	<0.000292
2,4-Dinitrotoluene	0.0013	0.003	<0.000059	<0.000058	<0.000058	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.00013	<0.000128	<0.000123
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000762	<0.00008	<0.0000769	<0.0000755
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	<0.0000769	<0.0000755
2-Methylnaphthalene	0.098	0.29	<0.000077	0.000072 J	0.00033	0.0038	0.000074 J	<0.00005	<0.00005	0.0009	<0.0000667	<0.00007	<0.0000673	<0.000066
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000079	<0.000083	<0.0000798	<0.0000783
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000533	<0.00056	<0.000538	<0.000528
Acenaphthene	1.5	4.4	<0.000027	0.00012	0.000036 J	0.0037	0.00025	<0.00005	<0.00005	0.00082	<0.0000762	<0.00008	<0.0000769	<0.0000755
Acenaphthylene	1.5	4.4	<0.000015	0.000053 J	0.000023 J	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000571	<0.00006	<0.0000577	<0.0000566
Anthracene	7.3	22	<0.000014	<0.000014	0.000036 J	0.00039	0.00024	<0.00005	<0.00005	0.00047	<0.0000476	<0.00005	0.000497	<0.0000472
Benzo(a)anthracene	0.0091	0.02	<0.000051	<0.00005	0.000055 J	0.00049	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	<0.0000769	<0.0000755
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	0.000063 J	0.00013 J	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	<0.0000769	<0.0000755
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000124	<0.00013	<0.000125	<0.000123
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000044	0.00032	0.00029	0.0059	0.00081	0.00086	<0.00018	<0.0003	<0.000352	<0.00037	<0.000356	<0.000349
Chrysene	0.91	2	<0.000021	<0.000021	0.000054 J	0.00032	0.00011 J	<0.00005	<0.00005	<0.00005	<0.0000762	<0.00008	<0.0000769	<0.0000755
Dibenzofuran	0.098	0.29	0.000066 J	0.000064 J	0.00003 J	0.003	0.00022	<0.00005	<0.00005	0.00071	<0.0000762	<0.00008	<0.0000769	<0.0000755
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.000067 J	<0.00002	<0.00007	<0.00007	0.00069 J	<0.00005	<0.00005	<0.000105	<0.00011	<0.0000855	<0.000104
Fluoranthene	0.98	2.9	<0.00001	<0.00001	0.000074 J	0.0025	0.00057	0.000059 J	<0.00005	0.00045	<0.0000667	<0.00007	<0.0000673	<0.000066
Fluorene	0.98	2.9	0.000057 J	0.000051 J	0.000036 J	0.0033	0.00036	<0.00005	<0.00005	0.00085	<0.0000667	<0.00007	<0.0000673	<0.000066
Naphthalene	0.49	1.5	<0.00079	0.0011	<0.00052	0.026	0.00011 J	<0.00005	<0.00029	<0.004	<0.000142	<0.00008	<0.000713	<0.000155
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000105	<0.00011	<0.000106	<0.000104
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0001	<0.0000962	<0.0000943
Pentachlorophenol	0.001	0.001	<0.00008	<0.000079	<0.000079	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000581	<0.00061	<0.000587	<0.000575
Phenanthrene	0.73	2.2	0.000062 J	<0.000021	0.000042 J	0.0083	0.0012	<0.00005	<0.00005	0.0022	<0.0000571	<0.00006	<0.000039	<0.0000566
Phenol	7.3	22	<0.000035	0.00026	<0.000035	0.0069	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000381	<0.00004	<0.0000385	<0.0000377
Pyrene	0.73	2.2	<0.000019	<0.000019	0.000092 J	0.0022	0.00037	<0.00005	<0.00005	0.00033	<0.000105	<0.00011	<0.000106	<0.000104
<b>Metals</b>														
Arsenic	0.01	0.01	<0.0004	<0.0004	<0.0004									

**Notes:**

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
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HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-69A 01/28/2018	MW-69A 03/20/2018	MW-69A 05/24/2018	MW-69A 01/10/2019	MW-69A 07/17/2019	MW-70C 03/12/2020	MW-70C 03/12/2020 Duplicate	MW-73B 02/02/2012	MW-73B 07/16/2012	MW-73B 01/30/2013	MW-73B 07/30/2013	MW-73B 01/15/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.001	<0.0005	<0.00014	0.000678 J	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<b>0.01</b>	<b>0.0089</b>	<b>0.0097</b>	<0.0005	0.000218 J	0.000156 J	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.001	<0.0005	<0.00012	<0.00012	<0.00018
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.057	0.052	0.0059	<0.0005	<0.00011	<0.00011	0.000437 J
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0013	<0.001	<0.00015	<0.00015	<0.00022
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.034	0.032	0.015	<0.0005	0.000336 J	<0.00015	0.000575
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.064	0.061	0.0059 J	<0.0015	<0.00026	<0.00026	0.00133 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.000022	<0.0005	<0.00005	<0.000104	<0.000107	<0.000104
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	0.00041	<0.000041	0.007	0.028	<0.000292	<0.000301	0.000946 J
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00006	<0.00006	<0.0005	<0.00005	<0.000123	<0.000126	<0.000123
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000044	<0.000043	<0.0006	<0.00006	<0.0000755	<0.0000777	<0.0000755
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.000022	<0.0005	<0.00005	<0.0000755	<0.0000777	<0.0000755
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.067 J	0.098 J	0.00055 J	0.00011 J	<0.000066	0.0000878 J	0.0161
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.000021	<0.000021	<0.0008	<0.00008	<0.000783	<0.000806	<0.000783
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000049	<0.000048	<0.0005	<0.00005	<0.000528	<0.000544	<0.000528
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.059 J	0.087 J	0.012	0.00016 J	<0.000185	0.000118 J	0.0112
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000016	0.00074	0.0013 J	<0.00005	0.0000696 J	<0.0000583	<0.0000566
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.0049 J	0.0082 J	<0.0005	0.00012 J	0.000186 J	0.000245 J	0.00462
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000052	0.000079 J	<0.0005	0.000057 J	<0.0000755	<0.0000777	0.00131
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.000021	<0.000021	<b>&lt;0.0005</b>	<0.00005	<0.0000755	<0.0000777	<b>0.00039 J</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.000031	<0.000031	<0.0005	<0.00005	<0.000123	<0.000126	<0.000123
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00028	0.00038	0.00025	<0.000037	<0.00034	<0.000039	0.000052 J	<0.001	0.00012 J	<0.000349	<0.000359	0.0015 J
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	0.000061 J	<0.0005	0.000096 J	<0.0000755	<0.0000777	0.00119
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.055 J	0.078 J	0.00078 J	0.000067 J	<0.0000755	<0.0000777	0.0102
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000061 J	0.0001 J	<0.0005	<0.00005	<0.000104	0.000133 J	0.000169 J
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.003 J	0.0053 J	0.01	0.000059 J	0.000138 J	<0.000068	0.00937
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.028 J	0.038 J	0.0041	0.00021	<0.000066	0.0000805 J	0.00951
Naphthalene	0.49	1.5	<0.00019	<0.00002	0.00008 J	<0.00002	<0.000078	<b>1.6 J</b>	<b>2.5 J</b>	0.0014 J	<0.00064	<0.000436	<0.000674	0.0906
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000025	<0.000025	<0.0005	<0.00005	<0.000104	<0.000107	<0.000104
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000026	<0.000026	<0.0005	<0.00005	<0.0000943	<0.0000971	<0.0000943
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000082	<0.000081	<0.0005	<0.00005	<0.0000575	<0.0000592	<0.0000575
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.04	0.054	0.00087 J	0.000089 J	<0.0000566	0.000228 J	0.0348
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000036	<0.000036	0.0053	0.00015 J	<0.0000377	<0.0000388	0.000522
Pyrene	0.73	2.2	<0.000019	0.000055 J	<0.000019	<0.000019	<0.000019	0.0017 J	0.0029 J	0.0077	<0.00005	<0.000104	<0.000107	0.00725
<b>Metals</b>														
Arsenic	0.01	0.01	0.00916	0.0017 J	<b>0.0142</b>	0.000717 J	0.000642 J	0.00598	0.0047					

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	Residential Assessment Level	C/I Assessment Level	MW-73B 07/18/2014	MW-74B 02/09/2012	MW-74B 07/26/2012	MW-74B 04/02/2013	MW-74B 01/29/2014	MW-74B 08/28/2014	MW-74B 01/30/2018	MW-74B 03/28/2018	MW-74B 06/07/2018	MW-74B 01/23/2019	MW-74B 07/30/2019	MW-74B 01/17/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.00014	<0.01	<0.005	<0.0028	<0.0007	<0.0028	<0.0002	<0.0002	<0.0002	<0.002	<0.001	<0.0002
Benzene	0.005	0.005	0.00309	0.35	0.71	0.552	0.795	0.652	0.47	0.58	0.71	0.83	0.59	0.12
Chlorobenzene	0.1	0.1	<0.00012	<0.01	<0.005	<0.0024	<0.0006	<0.0024	<0.0003	<0.0003	<0.0003	<0.003	<0.0015	<0.0003
Ethylbenzene	0.7	0.7	<0.00011	0.086	0.14	0.147	0.203	0.2	0.25	0.12	0.17 J	0.22	0.15	0.076
Methylene chloride	0.005	0.005	<0.00015	<0.013	<0.01	<0.003	<0.00075	<0.003	<0.001	<0.001	<0.001	<0.01	<0.005	<0.001
Toluene	1	1	<0.00015	0.32	0.56	0.533	0.774	0.741	0.75	0.56	0.74	0.69	0.52	0.29
Vinyl chloride	0.002	0.002	<0.00011											
Xylenes (total)	10	10	<0.00026	0.25	0.38	0.427	0.553	0.558	0.53	0.33	0.51 J	0.63	0.42	0.2
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000109	<0.0005	<0.0005	<0.106	<0.208	<0.216	<0.0021	<0.00021	<0.0021	<0.00021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.000307	55	41	56.9	525	70.6	59	30	57	9	37	0.21
2,4-Dinitrotoluene	0.0013	0.003	<0.000129	<0.0005	<0.0005	<0.125	<0.245	<0.255	<0.0058	<0.00058	<0.0058	<0.00058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000792	<0.0006	<0.0006	<0.0769	<0.151	<0.157	<0.0042	<0.00042	<0.0042	<0.00042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000792	<0.0005	<0.0005	<0.0769	<0.151	<0.157	<0.0021	<0.00021	<0.0021	<0.00021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000693	0.39	0.43 J	0.673	5.52	0.95 J	0.3	2.4	0.99	0.22	0.48	1.3
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000822	<0.0008	<0.0008	<0.798	<1.57	<1.63	<0.002	<0.0002	<0.002	<0.0002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000554	<0.0005	<0.0005	<0.538	<1.06	<1.1	0.033 J	<0.00047	<0.0047	<0.00047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.000792	0.29	0.21	0.31 J	2.4	0.413 J	0.31	1.4	0.33	0.098	0.24	0.98
Acenaphthylene	1.5	4.4	<0.000594	0.0058	0.0062	<0.0577	<0.113	<0.118	0.012	0.019	0.0098 J	0.0032	0.0037	0.018
Anthracene	7.3	22	0.00015 J	0.037	0.024	<0.0481	0.282 J	<0.098	0.027	0.58	0.034	0.0074	0.0071	0.42
Benzo(a)anthracene	0.0091	0.02	<0.000792	<0.0005	0.0022	<0.0769	<0.151	<0.157	<0.005	0.22	<0.005	<0.0005	0.00022	0.13
Benzo(a)pyrene	0.0002	0.0002	<0.000792	<0.0005	0.00085 J	<0.0769	<0.151	<0.157	<0.002	0.064	<0.002	<0.0002	<0.00002	0.034
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000129	<0.0005	<0.0005	<0.125	<0.245	<0.255	<0.003	<0.0003	<0.003	<0.0003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000603	<0.001	<0.001	<0.356	<0.698	<0.725	<0.0037	<0.00037	<0.0037	<0.00037	<0.000037	0.0014
Chrysene	0.91	2	<0.000792	<0.0005	0.0018 J	<0.0769	<0.151	<0.157	<0.0021	0.23	<0.0021	<0.00021	0.00016	0.082
Dibenzofuran	0.098	0.29	<0.000792	0.25	0.19	0.252 J	1.84	<0.157	0.24	1.4	0.24	0.079	0.19	0.72
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000109	<0.0005	<0.0005	<0.106	<0.208	<0.216	<0.002	<0.0002	<0.002	<0.0002	<0.00002	0.022
Fluoranthene	0.98	2.9	<0.000693	0.0044	0.018	<0.0673	<0.132	<0.137	0.015	1.4	0.017	0.0038	0.0035	0.67
Fluorene	0.98	2.9	<0.000693	0.17	0.14 J	0.196 J	1.34	0.263 J	0.19	1.4	0.19	0.056	0.14	0.88
Naphthalene	0.49	1.5	<0.000792	16	10	13.9	139	17.9	18	21	19	4	13 J	8.5
Nitrobenzene	0.049	0.15	<0.000109	<0.0005	<0.0005	<0.106	<0.208	<0.216	<0.0024	<0.00024	<0.0024	<0.0024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000099	<0.0005	<0.0005	<0.0962	<0.189	<0.196	<0.0025	<0.00025	<0.0025	<0.00025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000604	<0.0005	<0.0005	<0.587	<1.15	<1.2	<0.0079	<0.00079	<0.0079	<0.00079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.000594	0.15	0.15	0.169 J	1.28	0.307 J	0.16	3.7	0.17	0.046	0.087	2.3
Phenol	7.3	22	<0.000396	43	38	63.2	420	53.3	56	25	39	5	33	0.48
Pyrene	0.73	2.2	<0.000109	0.005	0.01	<0.106	<0.208	<0.216	0.0079 J	0.83	0.0077 J	0.002	0.0023	0.44
<b>Metals</b>														
Arsenic	0.01	0.01							0.00162 J	0.00142 J	0.00131 J	0.0014 J	0.00128 J	0.000443 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-75B 02/08/2012	MW-75B 07/26/2012	MW-75B 04/02/2013	MW-75B 01/29/2014	MW-75B 07/24/2014	MW-75B 01/17/2020	MW-76B 03/12/2020	MW-76C 07/24/2014	MW-76C 10/03/2014	MW-76C 01/30/2018	MW-76C 03/28/2018	MW-76C 05/25/2018
			DNAPL	DNAPL	DNAPL	DNAPL	DNAPL							
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.01	<0.0025	<0.0028	<0.0007	<0.00014	<0.0002	<0.0002	<0.00014		<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	0.61	0.85	0.369	0.502	0.298	0.12	<0.0002	0.000149 J		<0.0002	<0.0002	0.00021 J
Chlorobenzene	0.1	0.1	<0.01	<0.0025	<0.0024	<0.0006	<0.00012	<0.0003	<0.0003	<0.00012		<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.13	0.1	0.069	0.0773	0.0737	0.037	<0.0003	<0.00011		<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.013	<0.005	<0.003	<0.00075	<0.00015	<0.001	<0.001	<0.00015		<0.001	<0.001	<0.001
Toluene	1	1	0.51	0.5	0.282	0.328	0.273	0.13	<0.0002	0.000156 J		<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002					<0.00011			<0.00011				
Xylenes (total)	10	10	0.41	0.33	0.247	0.276	0.255	0.11	<0.0003	<0.00026		<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.0005	<0.0005	<0.00212	<0.0519	<0.00214	<0.000021	<0.000022	<0.000104	<0.000107	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.18	0.64	0.0695	6.35	<0.00602	0.86	<0.000042	<0.000292	<0.000301	0.0018	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.0005	<0.0005	<0.0025	<0.0613	<0.00252	<0.000058	<0.00006	<0.000123	<0.000126	<0.000058	<0.000058	<0.000059
2,6-Dinitrotoluene	0.0013	0.003	<0.0006	<0.0006	<0.00154	<0.0377	<0.00155	<0.000042	<0.000044	<0.0000755	<0.0000777	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.0005	<0.0005	<0.00154	<0.0377	<0.00155	<0.000021	<0.000022	<0.0000755	<0.0000777	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.62	0.6 J	0.101	3.18	0.546	0.81	<0.00002	0.000392 J	0.0000976 J	0.0001	0.00012	0.000032 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.0008	<0.0008	<0.016	<0.392	<0.0161	<0.00002	<0.000021	<0.000783	<0.000806	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.0005	<0.0005	<0.0108	<0.264	<0.0109	<0.000047	<0.000049	<0.000528	<0.000544	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.34	0.26	0.0697	2.57	0.429	0.67	<0.000028	0.000696	0.00024 J	0.00015	0.00023	0.000071 J
Acenaphthylene	1.5	4.4	0.013	0.0058	<0.00115	0.0672 J	0.0121	0.0086	<0.000016	<0.0000566	<0.0000583	0.0002	<0.000015	<0.000015
Anthracene	7.3	22	0.035	0.045	0.00948 J	0.605	0.0626	0.39	<0.000015	0.000234 J	0.00011 J	0.00006 J	0.000055 J	0.000048 J
Benzo(a)anthracene	0.0091	0.02	0.00064 J	0.0047	<0.00154	0.0667 J	0.00748 J	0.12	<0.000052	<0.0000755	<0.0000777	<0.00005	<0.00005	<0.000051
Benzo(a)pyrene	0.0002	0.0002	<0.0005	0.0013 J	<0.00154	<0.0377	<0.00155	0.035	<0.000021	<0.0000755	0.000276 J	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.0005	<0.0005	<0.0025	<0.0613	<0.00252	<0.00003	<0.000031	<0.000123	<0.000126	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.001	<0.001	<0.00712	<0.175	<0.00718	0.0021	<0.000039	0.000803	0.000714	<0.00024	0.00015 J	0.000096 J
Chrysene	0.91	2	0.00062 J	0.0042	<0.00154	0.0704 J	0.00677 J	0.095	<0.000022	<0.0000755	<0.0000777	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.29	0.23	0.0533	1.56	0.214	0.54	<0.000021	0.000507	0.000159 J	0.00012	0.00012	0.000056 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0005	<0.0005	<0.00212	<0.0519	<0.00214	0.0016	0.000025 J	<0.000104	0.000124 J	<0.00002	<0.00002	0.00004 J
Fluoranthene	0.98	2.9	0.016	0.04	0.0103	0.708	0.0914	0.75	<0.00001	0.000322 J	0.000188 J	0.00019	0.00018	0.00007 J
Fluorene	0.98	2.9	0.19	0.17 J	0.0425	1.59	0.218	0.61	<0.000031	0.000778	0.000264 J	0.00016	0.00014	0.000076 J
Naphthalene	0.49	1.5	8.9	9.3	0.211	27.1	5.7	7.5	0.000096 J	0.00176	0.000506	<0.0028	0.0019	0.00036
Nitrobenzene	0.049	0.15	<0.0005	<0.0005	<0.00212	<0.0519	<0.00214	<0.000024	<0.000025	<0.000104	0.000124 J	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.0005	<0.0005	<0.00192	<0.0472	<0.00194	<0.000025	<0.000026	<0.0000943	<0.0000971	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.0005	<0.0005	<0.0117	<0.288	<0.0118	<0.000079	<0.000082	0.000272	<0.000592	<0.000079	<0.000079	<0.00008
Phenanthrene	0.73	2.2	0.24	0.27	0.0606	2.13	0.238	1.9	<0.000022	0.00183	0.000611	0.00051	0.00044	0.00023
Phenol	7.3	22	0.0066	0.0027	0.0069 J	0.108 J	<0.000777	0.0029	<0.000036	0.00284	0.00163	0.0032	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0098	0.026	0.00617 J	0.416	0.0537	0.51	<0.00002	0.000194 J	0.000161 J	0.00016	0.00012	0.000048 J
<b>Metals</b>														
Arsenic	0.01	0.01							0.00167 J	0.00115 J		0.00157 J	0.000631 J	0.000527 J

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	Residential Assessment Level	C/I Assessment Level	MW-76C 01/23/2019	MW-76C 07/30/2019	MW-76C 01/09/2020	MW-77A 07/24/2014	MW-77A 01/30/2018	MW-77A 03/28/2018	MW-77A 05/24/2018	MW-77A 02/01/2019	MW-77A 07/30/2019	MW-77A 01/09/2020	MW-78A 07/24/2014 DNAPL	MW-78A 01/17/2020 DNAPL
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0007	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.00008	<b>0.054</b>	<b>0.063</b>	<b>0.053</b>	<0.0002	<b>0.063</b>	<b>0.058</b>	<b>0.0571</b>	<b>0.018</b>
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0006	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.00011	0.059	0.063	0.044	<0.0003	0.072	0.063	0.0637	0.011
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.00075	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.00015	0.011	0.011	0.006	<0.0002	0.012	0.0069	0.1	0.033
Vinyl chloride	0.002	0.002				<0.00011							<0.00055	
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.00026	0.1	0.11	0.058	<0.0003	0.097	0.065	0.158	0.028
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<b>&lt;0.00519</b>	<0.00021	<0.00021	<0.00021	<0.000022	<0.000021	<0.000021	<b>&lt;0.0259</b>	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.0041	<0.00004	<0.00038	<0.0146	0.014	0.015	0.007	<0.000042	0.062	0.055	<b>6.66</b>	<b>2.5</b>
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<b>&lt;0.00613</b>	<0.00058	<0.00058	<0.00058	<0.00006	<0.000058	<0.000058	<b>&lt;0.0307</b>	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<b>&lt;0.00377</b>	<0.00042	<0.00042	<0.00042	<0.000044	<0.000042	<0.000042	<b>&lt;0.0189</b>	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.00377	<0.00021	<0.00021	<0.00021	<0.000022	<0.000021	<0.000021	<0.0189	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00031	<0.000019	<0.00045	0.0571	<b>0.2</b>	<b>0.28</b>	0.085	<0.00002	<b>0.2</b>	<b>0.22</b>	<b>0.879</b>	<b>0.13</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<b>&lt;0.0392</b>	<0.0002	<0.0002	<0.0002	<0.000021	<0.00002	<0.00002	<b>&lt;0.196</b>	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.0264	0.0044 J	<0.00047	<0.00047	<0.000049	<0.000047	<0.000047	<b>&lt;0.132</b>	<0.000047
Acenaphthene	1.5	4.4	0.00011	<0.000027	<0.00041	0.0456	0.2	0.23	0.079	<0.000028	0.16	0.13	0.497	0.071
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	0.000035 J	<0.00283	0.0032	0.0035	0.0012	<0.000016	0.0013	0.0013	<0.0142	0.0015
Anthracene	7.3	22	0.000041 J	<0.000014	<0.00017	<0.00236	0.0034	0.0052	0.0025	<0.000015	0.0028	0.0036	0.105 J	0.004
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	0.00014	<0.00377	<0.0005	<0.0005	<0.0005	<0.000052	<0.00005	<0.00005	<b>0.0336 J</b>	0.00043
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	0.000037 J	<b>&lt;0.00377</b>	<0.0002	<0.0002	<0.0002	<0.000021	<0.00002	<0.00002	<b>&lt;0.0189</b>	0.00014
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<b>&lt;0.00613</b>	<0.0003	<0.0003	<0.0003	<0.000031	<0.00003	<0.00003	<b>&lt;0.0307</b>	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000091	<0.000037	<0.00016	<b>&lt;0.0175</b>	<0.00037	<0.00037	<0.00037	0.0001 J	<0.000037	<0.000077	<b>&lt;0.0873</b>	0.00092
Chrysene	0.91	2	<0.000021	<0.000021	0.00013	<0.00377	<0.00021	<0.00021	<0.00021	<0.000022	<0.000021	0.000041 J	0.0248 J	0.00039
Dibenzofuran	0.098	0.29	0.00011	<0.00002	<0.00029	0.0229 J	0.086	0.09	0.04	<0.000021	0.096	0.079	<b>0.411</b>	0.055
Di-n-butylphthalate (DBP)	2.4	7.3	0.000027 J	0.000054 J	<0.00002	<0.00519	<0.0002	<0.0002	<0.0002	0.000081 J	<0.00002	<0.00002	<0.0259	0.0016
Fluoranthene	0.98	2.9	<0.00001	<0.00001	0.001	<0.0033	0.0014	0.0013	0.00067 J	<0.00001	0.00052	0.00071	0.165	0.0029
Fluorene	0.98	2.9	0.000076 J	<0.00003	<0.00032	0.024	0.076	0.083	0.037	<0.000031	0.078	0.067	0.382	0.04
Naphthalene	0.49	1.5	0.007	0.00018 J	<0.0038	<b>0.884</b>	<b>7.8</b>	<b>6</b>	<b>1.5</b>	<0.000021	<b>4.9 J</b>	<b>4.7</b>	<b>7.18</b>	<b>3.1</b>
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00519	<0.00024	<0.00024	<0.00024	<0.000025	<0.000024	<0.000024	<0.0259	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00472	<0.00025	<0.00025	<0.00025	<0.000026	<0.000025	<0.000025	<0.0236	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<b>&lt;0.0288</b>	<0.00079	<0.00079	<0.00079	<0.000082	<0.000079	<0.000079	<b>&lt;0.144</b>	<0.000079
Phenanthrene	0.73	2.2	0.000086 J	<0.000021	<0.00069	0.0262	0.026	0.035	0.019	<0.000022	0.032	0.031	0.604	0.031
Phenol	7.3	22	0.0012	<0.000035	<0.000035	<0.00189	<0.00035	<0.00035	<0.00035	<0.000036	<0.000035	<0.000035	0.192	0.13
Pyrene	0.73	2.2	<0.000019	<0.000019	0.0007	<0.00519	0.00068 J	0.001	0.00037 J	<0.00002	0.00041	0.00046	0.0967 J	0.0018
<b>Metals</b>														
Arsenic	0.01	0.01	0.000579 J	0.00216	0.00299		<b>0.0263</b>	<b>0.0187</b>	<b>0.019</b>	0.00207	<b>0.0231</b>	<b>0.0237</b>		0.00991

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-79A 08/28/2014	MW-79A 01/30/2018	MW-79A 03/28/2018	MW-79A 05/25/2018	MW-79A 01/23/2019	MW-79A 07/30/2019	MW-79A 01/17/2020	MW-80B 08/28/2014	MW-80B 01/30/2018	MW-80B 03/28/2018	MW-80B 05/24/2018	MW-80B 01/10/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0007	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<b>0.0485</b>	<b>1</b>	<b>0.3</b>	<b>0.36</b>	<b>0.45</b>	<b>0.013</b>	<b>0.12</b>	0.0000898 J	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0006	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	0.0215	0.18	0.12	0.14	0.19	0.0067	0.075	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.00075	<0.001	<0.001	<0.001	<b>&lt;0.01</b>	<0.001	<0.001	<0.00015	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	0.076	0.99	0.44	0.48	0.55	0.018	0.26	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	0.0763	0.48	0.31	0.41	0.54	0.023	0.2	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<b>&lt;0.00539</b>	<0.00021	<0.00021	<0.00021	<0.00021	<0.000021	<0.000021	<0.00011	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<b>6.11</b>	<b>11</b>	<b>11</b>	<b>20</b>	<b>2.5</b>	0.33	<b>6.7</b>	<0.00031	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<b>&lt;0.00637</b>	<0.00058	<0.00058	<0.00058	<0.00058	<0.000058	<0.000058	<0.00013	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<b>&lt;0.00392</b>	<0.00042	<0.00042	<0.00042	<0.00042	<0.000042	<0.000042	<0.00008	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00392	<0.00021	<0.00021	<0.00021	<0.00021	<0.000021	<0.000021	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<b>0.654</b>	<b>0.17</b>	<b>0.42</b>	<b>0.44</b>	<b>0.1</b>	0.051	<b>0.43</b>	0.000158 J	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<b>&lt;0.0407</b>	<0.0002	<0.0002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00083	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.0275	<0.00047	<0.00047	0.016 J	<0.00047	<0.000047	<0.000047	<0.00056	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.427	0.13	0.17	0.16	0.039	0.034	0.19	0.0000835 J	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	0.0112 J	0.0045	<0.00015	0.0056	0.0015	0.00076	0.0036	<0.00006	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.0673	0.0057	0.0092	0.0084	0.0021	0.002	0.0094	<0.00005	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<b>0.00985 J</b>	<0.0005	<0.0005	<0.0005	<0.0005	0.00023	0.00081	<0.00008	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<b>&lt;0.00392</b>	<0.0002	<0.0002	<0.0002	<0.0002	<0.00002	<b>0.00029</b>	<0.00008	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<b>&lt;0.00637</b>	<0.0003	<0.0003	<0.0003	<0.0003	<0.00003	<0.00003	<0.00013	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<b>&lt;0.0181</b>	<0.00037	<0.00037	<0.00037	<0.00037	0.00055	<0.000037	0.00106	<0.000069	<0.000037	0.00012 J	<0.000037
Chrysene	0.91	2	0.00948 J	<0.00021	<0.00021	<0.00021	<0.00021	0.00027	0.00072	<0.00008	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<b>0.342</b>	0.097	<b>0.14</b>	0.092	0.037	0.026	<b>0.14</b>	<0.00008	<0.00002	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00539	<0.0002	<0.0002	<0.0002	<0.0002	<0.00002	0.00098	<0.00011	<0.00002	<0.00002	0.000047 J	<0.00002
Fluoranthene	0.98	2.9	0.0713	0.0036	0.0051	0.0023	0.001	0.0012	0.0076	<0.00007	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	0.291	0.063	0.081	0.056	0.022	0.017	0.094	<0.00007	<0.00003	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	<b>6.89</b>	<b>7.9</b>	<b>8.5</b>	<b>15</b>	<b>1.9</b>	<b>0.74 J</b>	<b>9</b>	0.00157	<0.00002	<0.00002	<0.00002	0.000068 J
Nitrobenzene	0.049	0.15	<0.00539	<0.00024	<0.00024	<0.00024	<0.00024	<0.000024	<0.000024	<0.00011	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.0049	<0.00025	<0.00025	<0.00025	<0.00025	<0.000025	<0.000025	<0.0001	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<b>&lt;0.0299</b>	<0.00079	<0.00079	<0.00079	<0.00079	<0.000079	<0.000079	<0.00061	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.355	0.038	0.049	0.039	0.012	0.0082	0.081	0.0000792 J	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	1.13	2.6	4	4.2	0.51	0.063	0.35	0.00018 J	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.0434	0.0022	0.0038	0.0022	0.00063 J	0.0011	0.0035	<0.00011	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01		<b>0.0184</b>	<b>0.0149</b>	<b>0.0134</b>	<b>0.0133</b>	0.00991	0.00893		0.00286	0.00187 J	0.00202	0.0018 J

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-80B 07/30/2019	MW-80B 01/07/2020	MW-81B 07/24/2014	MW-81B 01/29/2018	MW-81B 03/28/2018	MW-81B 05/25/2018	MW-81B 01/10/2019	MW-81B 07/19/2019	MW-81B 01/09/2020	MW-82B 02/01/2018	MW-82B 03/22/2018	MW-82B 06/06/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.00008	<0.0002	<0.0002	0.00021 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.00012	<0.0003	<0.0003	0.00036 J	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002			<0.00011									
Xylenes (total)	10	10	<0.0003	<0.0003	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.00065 J	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	0.00028	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000126	<0.000058	<0.000059	0.00079	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.0000777	<0.000042	<0.000042	0.001	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.0000777	<0.000021	<0.000021	0.0018	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000021 J
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000036	<0.000068	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000019 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.000027	0.000046 J	<0.0000777	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.000027 J
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.0000583	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015 J
Anthracene	7.3	22	<0.000014	0.000026 J	<0.0000485	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.000068 J	0.000091 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.0000777	<0.00005	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00007 J	<0.000069	<0.000359	<0.000037	<0.000037	0.00006 J	0.000058 J	<0.000037	<0.000066	0.00011 J	<0.000037	<0.000056
Chrysene	0.91	2	<0.000021	<0.000021	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00002	0.000048 J	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00002 J
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.000107	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.000068	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00003	0.000032 J	<0.000068	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00003 J
Naphthalene	0.49	1.5	<0.00002	<0.000034	<0.0000777	<0.00002	<0.00002	<0.00002	0.00016	0.000099 J	<0.00023	0.00019	<0.00015	0.00002 J
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000592	<0.000079	<0.00008	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.000021	0.000027 J	0.0000944 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.000035	<0.000035	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000107	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01	0.00162 J	0.00212		0.00207	0.00134 J	0.00203	0.00116 J	0.000563 J	0.0013 J	0.00271	0.00175 J	<b>0.0103</b>

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 2 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-82B 01/22/2019	MW-82B 07/30/2019	MW-82B 01/21/2020	MW-83B 02/08/2018	MW-83B 03/22/2018	MW-83B 06/07/2018	MW-83B 07/19/2018	MW-83B 01/15/2019	MW-83B 07/18/2019	MW-83B 01/10/2020	MW-83C 02/08/2018	MW-83C 03/22/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	0.018	0.019	0.02	0.03	0.032	0.021	0.021	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	0.08	0.1	0.085	0.068	0.091	0.055	0.078	0.00066 J	0.0005 J
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	0.0055	0.0046	0.0049	0.007	0.0082	0.0042	0.0052	<0.0002	<0.0002
Vinyl chloride	0.002	0.002										<0.0002		
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	0.1	0.11	0.091	0.066	0.1	0.061	0.098	0.0014	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.0004	<0.0004	<0.00004	<0.00004	<0.00004	<0.00004	0.000072 J	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.00058	<0.00059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.00042	<0.00042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	0.0001	0.15	0.75	0.23	0.076	0.086	0.05	0.078	0.0015	0.0015
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.00047	<0.00047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	0.0003 J	<0.000047
Acenaphthene	1.5	4.4	<0.000027	<0.000027	0.000043 J	0.098	0.33	0.11	0.027	0.026	0.021	0.066	0.00083	0.001
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	0.00086 J	0.0016	0.00064	<0.000015	0.00034	0.00025	0.00047	<0.000015	<0.000015
Anthracene	7.3	22	0.000042 J	<0.000014	0.000031 J	0.01	0.011	0.007	0.014	0.012	0.00099	0.0039	0.000034 J	0.000068 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.0005	<0.00051	0.000058 J	<0.00005	<0.00005	<0.00005	0.00007 J	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.0003	<0.0003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000037	<0.000037	<0.00037	<0.00037	0.00012 J	<0.000037	<0.000064	0.00036	<0.000037	0.00019 J	<0.000037
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.00021	<0.00021	0.00009 J	<0.000021	<0.000021	<0.000021	0.000065 J	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	0.00004 J	0.043	0.17	0.061	0.023	0.02	0.014	0.023	0.00061	0.00044
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	0.000094 J	0.00021	0.00015 J	0.00012 J	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	0.0035	0.0043	0.0046	0.00057	0.00051	0.00038	0.0033	0.000044 J	<0.00001
Fluorene	0.98	2.9	<0.00003	<0.00003	0.000037 J	0.046	0.072	0.039	0.0083	0.0099	0.0076	0.024	0.00035	0.00034
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.00041	2.6	14	2	1.5	1.6	0.77	1.2	0.012	0.016
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.00024	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.00025	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.00079	<0.0008	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.000021	<0.000021	0.00012	0.04	0.071	0.045	0.0086	0.0074	0.0059	0.021	0.00044	0.00053
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.00035	<0.00035	<0.000035	<0.000035	<0.000035	0.00051	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	0.0029	0.0023	0.0026	0.00037	0.0003	0.0002	0.0022	0.000027 J	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01	0.00838	0.00873	0.00484	0.0353	0.0185	0.0673	0.0731	0.0916	0.0648	0.0709	0.000609 J	<0.0004

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.



**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-83C 06/07/2018	MW-83C 01/15/2019	MW-83C 07/18/2019	MW-83C 01/10/2020	MW-84A 03/12/2020	MW-84B 02/08/2018	MW-84B 03/27/2018	MW-84B 06/07/2018	MW-84B 07/19/2018	MW-84B 07/19/2018 Duplicate	MW-84B 01/24/2019	MW-84B 07/30/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<b>0.0097</b>	<b>0.0086</b>	0.0017	0.002	0.0019	0.0024	<b>0.01</b>
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.039	0.037	0.0036	0.0029	0.0026	0.0051	0.013
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0025	0.00099 J	<0.0002	<0.0002	<0.0002	0.00056 J	0.0015
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.035	0.031	0.003	0.0019	0.0017	0.0033	0.0041
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	0.000087 J	<0.000021	<0.000021	<0.000021	<b>0.0014 J</b>	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.000041	<0.0004	<0.0004	0.0017 J	0.00048	<0.00004	<0.00016	0.00081
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000059	<0.00058	<0.00058	<0.00058	<0.00058	<0.00058	<0.00058	<0.00058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000043	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042	<0.00042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021	<0.00021
2-Methylnaphthalene	0.098	0.29	0.000089 J	<0.000019	0.00012	0.00021	<0.000019	<b>0.55</b>	<b>0.58</b>	0.025	0.074 J	<b>0.2 J</b>	<0.000019	0.0027
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<b>0.0038</b>	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000048	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047	<0.00047
Acenaphthene	1.5	4.4	0.00017	<0.000027	0.000094 J	0.00018	<0.000028	0.22	0.27	0.031	0.048 J	0.14 J	0.000032 J	0.008
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.003	0.0032	0.00061 J	0.00062	<0.000015	0.000043 J	0.0003
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.02	0.0092	0.0022	0.0027 J	0.0042 J	<0.000014	0.00042
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.000051	<0.0005	<0.0005	<0.0005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.000031	<0.0003	<0.0003	<0.0003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00012 J	<0.000037	0.00023	<0.000037	<0.000038	0.00056 J	<0.00037	<0.00037	<0.000037	<0.000037	<0.000037	0.00026
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00021	<0.00021	<0.00021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.000046 J	<0.00002	0.00007 J	0.00015	<0.00002	<b>0.13</b>	<b>0.22</b>	0.019	0.037 J	0.077 J	<0.00002	0.00054
Di-n-butylphthalate (DBP)	2.4	7.3	0.000092 J	0.00006 J	0.00004 J	<0.00002	<0.00002	<0.0002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00002	0.000055 J
Fluoranthene	0.98	2.9	0.000018 J	<0.00001	<0.00001	<0.00001	<0.00001	0.0039	0.0029	0.00069 J	0.0011 J	0.0016 J	<0.00001	0.00015
Fluorene	0.98	2.9	0.000067 J	<0.00003	<0.00003	0.00012	<0.000031	0.074	0.076	0.011	0.018 J	0.039 J	<0.00003	0.0026
Naphthalene	0.49	1.5	0.00039	<0.00036	0.0014	0.0018	0.000023 J	<b>2.4</b>	<b>2.2</b>	0.066	0.49 J	<b>0.95 J</b>	<0.00002	0.06 J
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00024	<0.00024	<0.00024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000026	<0.00025	<0.00025	<0.00025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000081	<0.00079	<0.00079	<0.00079	<b>0.005</b>	0.000094 J	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.000038 J	<0.000021	<0.000021	0.00017	<0.000021	0.088	0.072	0.013	0.024 J	0.043 J	<0.000021	0.0021
Phenol	7.3	22	<0.000035	0.000038 J	0.000068 J	<0.000035	<0.000036	<0.00035	<0.00035	0.00062 J	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.0025	0.0018	0.00036 J	0.00062 J	0.00097 J	<0.000019	0.00011
<b>Metals</b>														
Arsenic	0.01	0.01	0.00139 J	0.00616	0.00617	0.00564	0.00464	0.00269	0.00277	<0.0004	<0.0004	<0.0004	0.00219	0.00838

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-84B 01/16/2020	MW-85C 02/01/2018	MW-85C 03/28/2018	MW-85C 05/24/2018	MW-85C 02/01/2019	MW-85C 07/30/2019	MW-85C 01/09/2020	MW-86C 02/01/2018	MW-86C 02/01/2018	MW-86C 03/28/2018	MW-86C 03/28/2018	MW-86C 05/25/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	0.0064	<0.0002	0.013	0.0026	<0.0002	0.00093 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0094	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	0.0032	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000069 J	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	0.00048	<0.00004	<0.00004	<0.00004	<0.000041	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.0065	0.0001 J	0.000049 J	0.000067 J	<0.000019	<0.000019	<0.000041	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	0.00035	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000048	<0.000047	0.00077 J	<0.000047	0.0003 J	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.025	0.00013	<0.000027	<0.000027	<0.000028	<0.000027	<0.000077	0.00091	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	0.00031	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.00095	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000031 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.000091 J	0.0002	0.00013 J	<0.000038	<0.000037	<0.00016	0.000053 J	<0.000037	<0.000037	<0.000037	0.000052 J
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000053 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.015	0.000087 J	<0.00002	0.000036 J	<0.00002	<0.00002	<0.000059	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.00023	<0.00002	0.00003 J	<0.00002	<0.00002	<0.00002	0.000093 J	0.000051 J	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.00037	0.000015 J	<0.00001	<0.00001	<0.00001	<0.00001	0.00025	0.00011	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	0.0063	0.00011	<0.00003	0.000034 J	<0.000031	<0.00003	<0.000055	0.00038	<0.00003	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	0.13	0.00069	0.017	0.0026	<0.00002	0.0012 J	<0.00046	0.000054 J	<0.00002	<0.00002	<0.00002	0.000079 J
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	0.00015 J	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.0065	0.0001	<0.000021	0.0001	<0.000021	<0.000021	<0.000067	<0.000021	<0.000021	<0.000021	<0.000021	0.000032 J
Phenol	7.3	22	<0.000035	0.00011 J	<0.000035	<0.000035	<0.000036	<0.000035	<0.000042	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.00025	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.00021	0.00006 J	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01	0.00363	0.00152 J	0.00287	0.00588	0.00136 J	0.000633 J	0.00272	0.00146 J	0.00156 J	0.00612	0.00608	0.00768

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	Residential Assessment Level	C/I Assessment Level	MW-86C	MW-86C	MW-86C	MW-86C	MW-86C	MW-86C	MW-86C	MW-86C	MW-87C	MW-87C	MW-87C	MW-87C	MW-87C
			05/25/2018	01/11/2019	01/11/2019	07/30/2019	07/30/2019	01/17/2020	01/17/2020	02/08/2018	03/27/2018	06/07/2018	01/22/2019	07/17/2019	
			Duplicate		Duplicate		Duplicate		Duplicate						
<b>Volatiles Organic Compounds</b>															
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0039	0.0045	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>															
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.0028	0.00041 J	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	0.00018 J	<0.00004	<0.00004	<0.00004	0.00004 J	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	0.000058 J	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	0.000042 J	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000021 J	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000061 J	0.000066 J	0.00043	0.00014	0.000019 J	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00002 J	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	0.00033 J	<0.000047	0.000047 J	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	0.00042	0.00011	0.000027 J	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.000015 J	0.000015 J	<0.000015	<0.000015
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.00012	0.000081 J	<0.000014	0.000014 J	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	0.00005 J	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00002 J	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00003 J	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000061 J	<0.000057	<0.0001	0.00006 J	0.00005 J	<0.000037	<0.000053	0.0003	<0.000037	0.00012 J	<0.000037	<0.000037	<0.000037
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.000021 J	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00032	0.000063 J	0.00002 J	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.000072 J	0.00011 J	0.00011 J	0.000073 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00002 J	<0.00002	0.00017 J
Fluoranthene	0.98	2.9	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	0.00001 J	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00027	<0.00003	0.00003 J	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.00002	0.000079 J	<0.00002	0.00012 J	0.00031 J	<0.00075	<0.00022	0.0014	0.00038	0.00002 J	<0.00002	<0.00017	<0.00017
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	0.000024 J	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	0.000025 J	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	0.000079 J	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00036	0.00048	0.00015	0.000021 J	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00011	<0.000035	0.000035 J	<0.000035	0.00013 J
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000019 J	<0.000019	<0.000019	<0.000019
<b>Metals</b>															
Arsenic	0.01	0.01	0.00824	0.00402	0.00405	0.00236	0.0019 J	0.000645 J	0.000699 J	<0.0004	<0.0004	<0.0004	0.000587 J	<0.0004	<0.0004

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray.
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	MW-87C 01/20/2020	MW-88A 03/17/2020	MW-88B 03/17/2020	MW-88C 02/01/2018	MW-88C 03/19/2018	MW-88C 05/24/2018	MW-88C 01/08/2019	MW-88C 07/31/2019	MW-88C 01/14/2020	MW-89B 07/19/2018	MW-89B 01/22/2019	MW-89B 07/18/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000059	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.00023	<0.000019	<0.000019	<0.000019	0.000052 J	<0.000019	<0.000019	<0.000019	<0.000019	0.000054 J	<0.000019	0.00005 J
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.00011	0.0015	<0.000027	0.000053 J	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.000088 J	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005	<0.000051	<0.00005
Benzo(a)pyrene	0.0002	0.0002	0.000066 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.0001 J	<0.000037	0.000052 J	<0.000037	0.00017 J	<0.000037	<0.000037	0.000064 J	0.000065 J	<0.000037	0.00015 J
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.00011	<0.00002	<0.00002	0.000023 J	0.000056 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	0.000025 J	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.000038 J
Fluoranthene	0.98	2.9	<0.000086	0.00025	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	0.000093 J	0.00026	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	0.0011	<0.00002	<0.00002	0.00011	0.00055	0.00011	<0.000059	<0.00002	<0.00002	0.0009	<0.00002	0.00025
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.000079
Phenanthrene	0.73	2.2	<0.00028	<0.000021	<0.000021	0.00003 J	<0.000021	0.000035 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	0.00048	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	0.00055	<0.000035	<0.000035	<0.000035	0.00042
Pyrene	0.73	2.2	<0.000057	0.00053	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01	0.00135 J	0.00355	0.00166 J	0.000557 J	0.000653 J	0.00346	0.000864 J	<0.0004	0.000862 J	0.00138 J	0.000683 J	<0.0004

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-89B 01/16/2020	MW-90B 07/19/2018	MW-90B 01/22/2019	MW-90B 07/18/2019	MW-90B 01/20/2020	MW-91A 03/12/2020	MW-92B 03/12/2020	MW-93B 03/12/2020	MW-94A 03/12/2020	MW-95A 03/17/2020	MW-96B 03/17/2020	MW-96B 03/17/2020 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.00049 J	0.00092 J	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.002	0.002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.000021	<0.000021	<0.000022	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00004	<0.00004	<0.00004	<0.00004	0.000075 J	<0.000042	<0.00004	<0.000041	<0.000041	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.000061	<0.000059	<0.000059	<0.00006	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000044	<0.000042	<0.000043	<0.000043	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.000021	<0.000021	<0.000022	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	<0.000032	<0.000019	<0.000019	<0.000019	0.00034	<0.00002	<0.000019	<0.000019	<0.00002	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000049	<0.000047	<0.000048	<0.000048	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	<0.000083	<0.000027	<0.000027	<0.000027	0.00015	<0.000028	<0.000027	0.000039 J	<0.000028	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	0.000045 J	<0.000015	<0.000015	<0.000015	<0.000015	<0.000016	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.0002	<0.000014	<0.000014	<0.000014	0.00014	<0.000015	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.0091	0.02	0.00017	<0.00005	<0.00005	<0.00005	<0.00005	<0.000053	<0.000051	<0.000051	<0.000052	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	0.000079 J	<0.00002	<0.00002	<0.00002	0.000052 J	<0.000021	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.000032	<0.00003	<0.000031	<0.000031	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000099 J	0.000097 J	<0.000037	0.00011 J	<0.000037	<0.000039	<0.000037	<0.000038	0.000056 J	<0.000037	0.00035	0.00013 J
Chrysene	0.91	2	0.00024	<0.000021	<0.000021	<0.000021	<0.000021	<0.000022	<0.000021	<0.000021	<0.000022	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	<0.00011	<0.00002	<0.00002	<0.00002	0.00011	<0.000021	<0.00002	0.000071 J	<0.000021	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	0.000034 J	<0.00002	<0.00002	<0.00002	<0.00002	0.000031 J	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.012	<0.00001	<0.00001	<0.00001	<0.00012	<0.000011	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00018	<0.00003	<0.00003	<0.00003	0.00013	<0.000032	<0.00003	0.000058 J	<0.000031	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.00029	<0.00002	0.000045 J	0.000091 J	0.003	0.000083 J	<0.00002	<0.00002	<0.000021	<0.00002	<0.00002	<0.00002
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000025	<0.000024	<0.000024	<0.000025	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000026	<0.000025	<0.000026	<0.000026	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.000083	<0.00008	<0.000081	<0.000081	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.00083	<0.000021	<0.000021	<0.000021	0.0004	<0.000022	<0.000021	0.00016	<0.000022	<0.000021	<0.000021	<0.000021
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	0.00019 J	<0.000037	<0.000035	<0.000036	<0.000036	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.00081	<0.000019	<0.000019	<0.000019	<0.000077	<0.00002	<0.000019	<0.000019	<0.00002	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01	0.000463 J	0.00169 J	0.00346	<b>0.0135</b>	0.0029	0.00989	0.00201	0.00455	0.0054	0.000977 J	0.00312	0.00311

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	MW-97A 03/12/2020	MW-98A 03/12/2020	MW-98B 03/12/2020	MW-99C 03/11/2020	P-10 01/29/2008	P-10 01/29/2008 Duplicate	P-10 07/16/2008	P-10 01/22/2009	P-10 07/22/2009	P-10 07/22/2009 Duplicate	P-10 01/22/2010	P-10 01/22/2010 Duplicate
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002							<0.00052	
Benzene	0.005	0.005	<0.0002	<0.0002	<0.0002	<0.0002							<0.00025	
Chlorobenzene	0.1	0.1	<0.0003	<0.0003	<0.0003	<0.0003							<0.00047	
Ethylbenzene	0.7	0.7	<0.0003	<0.0003	<0.0003	<0.0003							<0.00025	
Methylene chloride	0.005	0.005	<0.001	<0.001	<0.001	<0.001							<0.00054	
Toluene	1	1	<0.0002	0.00065 J	0.002	0.00078 J							<0.00041	
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.0003	<0.0003	<0.0003	<0.0003							<0.00127	
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.000021	<0.000021	<0.000021								
2,4-Dimethylphenol	0.49	1.5	<0.00004	0.000078 J	<0.000041	<0.000041								
2,4-Dinitrotoluene	0.0013	0.003	<0.000059	<0.000059	<0.000059	<0.000059								
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.000043	<0.000043	<0.000043								
2-Chloronaphthalene	2	5.8	<0.000021	<0.000021	<0.000021	<0.000021								
2-Methylnaphthalene	0.098	0.29	<0.000019	<0.000019	<0.000019	0.000086 J								
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00002	<0.00002	<0.00002								
4-Nitrophenol	0.049	0.15	<0.000047	<0.000048	<0.000048	<0.000048								
Acenaphthene	1.5	4.4	<0.000027	0.00004 J	<0.000028	<0.000028	0.00373	0.00854	0.0106	<0.0008	<0.0009	<0.0009	<0.0009	<0.0009
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.00028	<0.00028	0.00053	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005
Anthracene	7.3	22	<0.000014	<0.000014	<0.000014	<0.000014	0.000703	0.00036 J	0.000747	<0.0007	<0.0006	<0.0006	<0.0006	<0.0006
Benzo(a)anthracene	0.0091	0.02	<0.000051	<0.000051	<0.000051	<0.000051								
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<0.00002	<0.00002	<0.00002								
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.000031	<0.000031	<0.000031								
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.000038	<0.000038	0.00007 J	0.00023 J	<0.00019	0.00022 J	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033
Chrysene	0.91	2	<0.000021	<0.000021	<0.000021	<0.000021								
Dibenzofuran	0.098	0.29	<0.00002	<0.00002	<0.00002	<0.00002	0.000713	0.00175	0.00176	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00002	<0.00002	<0.00002	<0.00019	<0.00019	0.00092 J	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005
Fluoranthene	0.98	2.9	<0.00001	0.000018 J	<0.00001	<0.00001	0.000506	0.00025 J	0.00022 J	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005
Fluorene	0.98	2.9	<0.00003	<0.000031	<0.000031	<0.000031	0.000668	0.00251	0.00245	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006
Naphthalene	0.49	1.5	<0.00002	0.000042 J	0.000036 J	0.00018	<0.00038	<0.00037	0.00079	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006
Nitrobenzene	0.049	0.15	<0.000024	<0.000024	<0.000024	<0.000024								
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.000026	<0.000026	<0.000026								
Pentachlorophenol	0.001	0.001	<0.00008	<0.000081	<0.000081	<0.000081								
Phenanthrene	0.73	2.2	<0.000021	<0.000021	<0.000021	0.000053 J								
Phenol	7.3	22	<0.000035	<0.000036	<0.000036	<0.000036	<0.00019	<0.00019	<0.00021	<0.0015	<0.0005	<0.0005	<0.0005	<0.0005
Pyrene	0.73	2.2	<0.000019	<0.000019	<0.000019	<0.000019	0.00039 J	<0.00019	<0.00021	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005
<b>Metals</b>														
Arsenic	0.01	0.01	0.00121 J	0.00963	0.00165 J	0.000866 J								

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
  8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	P-10 07/14/2010	P-10 07/14/2010 Duplicate	P-10 01/12/2011	P-10 07/12/2011	P-10 07/12/2011 Duplicate	P-10 01/31/2012	P-10 01/31/2012 Duplicate	P-10 07/11/2012	P-10 07/11/2012 Duplicate	P-10 01/10/2013	P-10 01/10/2013 Duplicate	P-10 07/11/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29												
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.0009	<0.0009	<0.0009	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000755	<0.0000808
Acenaphthylene	1.5	4.4	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000566	<0.0000566	<0.0000606
Anthracene	7.3	22	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	0.0000472 J	0.000148 J	0.000133 J
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0033	<0.0033	<0.0033	<0.0005	0.0015 J	<0.0005	<0.0005	<0.0005	<0.0005	0.000906	0.00112	0.000492 J
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0007	<0.0007	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000755	<0.0000808
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.000104	<0.000111
Fluoranthene	0.98	2.9	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000066	<0.000066	<0.0000707
Fluorene	0.98	2.9	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000066	<0.000066	<0.0000707
Naphthalene	0.49	1.5	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000755	<0.0000755	<0.0000808
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2												
Phenol	7.3	22	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000377	<0.0000377	<0.0000404
Pyrene	0.73	2.2	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.000104	<0.000111
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
- All values in milligrams per liter (mg/L).
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  - Concentrations > C/I AL and non-detects are highlighted dark gray.
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	P-10	P-10	P-10	P-10	P-10	P-10	P-10	P-10	P-10	P-10	P-10	P-10
			07/11/2013	01/09/2014	01/09/2014	07/02/2014	07/02/2014	01/07/2015	01/07/2015	07/08/2015	07/08/2015	01/13/2016	01/13/2016	07/07/2016
<b>Volatile Organic Compounds</b>			Duplicate		Duplicate		Duplicate		Duplicate		Duplicate		Duplicate	
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29				0.0000718 J	0.000891 J							
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.0000812	0.000102 J	0.000966	0.01 J	0.0169 J	<0.0000792	<0.0000792	0.023	0.022	<0.000027	<0.000027	<0.000027
Acenaphthylene	1.5	4.4	<0.0000609	<0.0000556	0.0000571 J	0.0000588 J	0.000265 J	<0.0000594	<0.0000594	0.00012	0.00012	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	0.000181 J	0.000323 J	0.000369 J	0.000375 J	0.000473 J	0.000122 J	0.000115 J	0.00039	0.00033	0.000014 J	0.00007 J	<0.000014
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.000575	<0.000343	<0.000343	0.00127	0.00133	0.000853 J	0.00155 J	0.0006 J	0.00032 J	0.0002 J	0.00035 J	0.00029
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.0000812	<0.0000741	0.000135 J	0.00205 J	0.00304 J	<0.0000792	<0.0000792	0.0024	0.0021	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000112	<0.000262	<0.000309	0.000108 J	0.000321 J	<0.000109	<0.000119	<0.000044	<0.000059	<0.000058	<0.00011	<0.000046
Fluoranthene	0.98	2.9	<0.0000711	<0.0000648	<0.0000648	0.00042 J	0.00039 J	0.000114 J	0.000113 J	0.0006	0.00047	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.0000711	<0.0000648	0.000262 J	0.00393	0.00514	<0.0000693	<0.0000693	0.0046	0.0038	<0.00003	<0.00003	<0.00003
Naphthalene	0.49	1.5	<0.0000812	<0.0000741	<0.0000741	0.0000784 J	0.000924 J	<0.0000792	<0.0000792	0.019	0.018	<0.00002	<0.00002	<0.00002
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2				0.000575 J	0.000808 J							
Phenol	7.3	22	<0.0000406	<0.000037	<0.000037	<0.0000392	<0.0000396	<0.0000396	<0.0000396	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000112	<0.000102	<0.000102	0.000318 J	0.000306 J	<0.000109	<0.000109	0.00038	0.00029	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
1. All values in milligrams per liter (mg/L).
  2. Concentrations > RAL and non-detects are highlighted light gray.
  3. Concentrations > C/I AL and non-detects are highlighted dark gray
  4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
  6. J = Estimated value, < = not detected at the specified detection limit.
  7. MW-32A was screened in the B-CZ & replaced with MW-32AR
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HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	P-10 07/07/2016	P-10 01/12/2017	P-10 01/12/2017	P-10 07/13/2017	P-10 07/13/2017	P-10 01/04/2018	P-10 01/04/2018	P-10 07/19/2018	P-10 07/19/2018	P-10 01/08/2019	P-10 01/08/2019	P-10 07/01/2019
			Duplicate		Duplicate		Duplicate		Duplicate		Duplicate		Duplicate	
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												
Benzene	0.005	0.005												
Chlorobenzene	0.1	0.1												
Ethylbenzene	0.7	0.7												
Methylene chloride	0.005	0.005												
Toluene	1	1												
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10												
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												
2,4-Dimethylphenol	0.49	1.5												
2,4-Dinitrotoluene	0.0013	0.003												
2,6-Dinitrotoluene	0.0013	0.003												
2-Chloronaphthalene	2	5.8												
2-Methylnaphthalene	0.098	0.29												
4,6-Dinitro-2-methylphenol	0.0024	0.0073												
4-Nitrophenol	0.049	0.15												
Acenaphthene	1.5	4.4	<0.000027	0.000027 J	0.000078 J	<0.000027	<0.000027	<0.000027	0.0033	0.041	0.043	<0.000027	<0.000027	0.0028
Acenaphthylene	1.5	4.4	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.00038	0.00032	<0.000015	<0.000015	<0.000015
Anthracene	7.3	22	<0.000014	0.000014 J	0.000081 J	<0.000014	<0.000014	<0.000014	0.00017	0.00058	0.0006	0.00025	0.00021	0.000065 J
Benzo(a)anthracene	0.0091	0.02												
Benzo(a)pyrene	0.0002	0.0002												
bis(2-Chloroethoxy)methane	0.00083	0.0019												
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.0003	0.00025 J	0.00018 J	0.000097 J	0.00016 J	<0.00024	<0.00017	<0.00019	<0.00022	<0.000061	<0.000057	0.00029
Chrysene	0.91	2												
Dibenzofuran	0.098	0.29	<0.00002	0.00002 J	0.000055 J	<0.00002	<0.00002	0.00029	0.00093	0.0047	0.0042	<0.00002	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000031	<0.00002	<0.00002	<0.00002	<0.00002	<0.000042	<0.000034	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	<0.00001	0.00001 J	0.000088 J	<0.00001	<0.00001	0.00017	0.00016	0.00099	0.001	<0.00001	<0.00001	<0.00001
Fluorene	0.98	2.9	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00042	0.0012	0.0073	0.007	<0.00003	0.00003 J	<0.00003
Naphthalene	0.49	1.5	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.0034	0.00014	0.000056 J	<0.00002	<0.00002	0.000063 J
Nitrobenzene	0.049	0.15												
N-Nitrosodiphenylamine	0.19	0.42												
Pentachlorophenol	0.001	0.001												
Phenanthrene	0.73	2.2												
Phenol	7.3	22	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	0.0014	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	<0.000019	<0.000019	0.000022 J	<0.000019	<0.000019	<0.000019	<0.000019	0.00049	0.00048	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	0.01	0.01												

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray.
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
  - J = Estimated value, < = not detected at the specified detection limit.
  - MW-32A was screened in the B-CZ & replaced with MW-32AR
  - Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	P-10 01/13/2020	P-11 01/30/2008	P-11 07/15/2008	P-11 02/04/2009	P-11 01/21/2010	P-11 06/22/2010	P-11 01/18/2011	P-11 07/27/2011	P-11 02/02/2012	P-11 07/26/2012	P-11 02/05/2013	P-11 08/01/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005		<0.00052	<0.00052	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014
Benzene	0.005	0.005		<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00008	<0.00008
Chlorobenzene	0.1	0.1		<0.00047	<0.00047	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012
Ethylbenzene	0.7	0.7		<0.00025	<0.00025	<0.0005	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011
Methylene chloride	0.005	0.005		<0.00054	<0.00054	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015
Toluene	1	1		<0.00041	<0.00041	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015
Vinyl chloride	0.002	0.002											<0.00011	<0.00011
Xylenes (total)	10	10		<0.00127	<0.00127	<0.001	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026		<0.00008	<0.00008	<0.0001	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105
2,4-Dimethylphenol	0.49	1.5		<0.00031	<0.00028	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	R	<0.000295
2,4-Dinitrotoluene	0.0013	0.003		<0.00021	<0.00019	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124
2,6-Dinitrotoluene	0.0013	0.003		<0.00021	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.0000762
2-Chloronaphthalene	2	5.8		<0.00042	<0.00038	<0.00012	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762
2-Methylnaphthalene	0.098	0.29		0.000783	<0.00038	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	0.00023	<0.000052	0.000127 J	<0.0000667
4,6-Dinitro-2-methylphenol	0.0024	0.0073		<0.00021	<0.00047	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	R	<0.00079
4-Nitrophenol	0.049	0.15		<0.00026	<0.00024	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	R	<0.000533
Acenaphthene	1.5	4.4	<0.000027	0.0776	<0.00028	0.0057	<0.00009	0.0037	<0.00009	0.00075	0.03	0.018	<0.0000755	<0.0000762
Acenaphthylene	1.5	4.4	<0.000015	<0.00031	<0.00028	<0.00006	<0.00007	<0.00007	<0.00007	<0.00005	0.0002	0.0001 J	<0.0000566	<0.0000571
Anthracene	7.3	22	<0.000014	0.00356	<0.00019	0.00015 J	<0.00007	0.00012 J	<0.00007	0.00012 J	0.0016	0.00039	0.00025 J	0.0000997 J
Benzo(a)anthracene	0.0091	0.02		<0.00021	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762
Benzo(a)pyrene	0.0002	0.0002		<0.00021	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762
bis(2-Chloroethoxy)methane	0.00083	0.0019		<0.00042	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.000124
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	0.00116 J	<0.00019	0.00022	<0.00051	<0.00021	0.0016	0.00018 J	<0.00013	0.00021	0.00036 J	0.000593 J
Chrysene	0.91	2		<0.00021	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762
Dibenzofuran	0.098	0.29	<0.00002	<0.00031	<0.00028	0.00024	<0.00008	0.000093 J	<0.00008	0.00013 J	0.0035	0.00059	0.000135 J	<0.0000762
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	<0.00021	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105
Fluoranthene	0.98	2.9	<0.00001	0.0061	<0.00019	<0.00007	<0.00007	0.00042	<0.00007	0.000081 J	0.0022	0.00048	<0.000066	<0.0000667
Fluorene	0.98	2.9	<0.00003	0.0219	<0.00019	0.0018	<0.00007	0.0016	<0.00007	0.000082 J	0.011	0.0044 J	0.0000769 J	<0.0000667
Naphthalene	0.49	1.5	0.00017	0.0324	<0.00038	0.0027	<0.0001	0.0027	<0.0001	0.00013 J	0.0017	<0.00026	0.000662	<0.0000762
Nitrobenzene	0.049	0.15		<0.00042	<0.00038	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105
N-Nitrosodiphenylamine	0.19	0.42		<0.00026	<0.00024	<0.00009	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000952
Pentachlorophenol	0.001	0.001		<0.00021	<0.00019	<0.00008	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	R	<0.000581
Phenanthrene	0.73	2.2		0.0196	<0.00019	0.00048	<0.00007	0.00053	<0.00007	0.000086 J	0.0045	0.00055	0.0000854 J	<0.0000571
Phenol	7.3	22	<0.000035	<0.00021	<0.00019	<0.00007	<0.00007	<0.00007	<0.00007	<0.00005	<0.00015	<0.00005	R	<0.0000381
Pyrene	0.73	2.2	<0.000019	0.00369	<0.00019	<0.00007	<0.00007	0.00015 J	<0.00007	<0.00005	0.0013	0.00023	<0.000104	<0.000105
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
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CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	P-11	P-11	P-11	P-11	P-11	P-11	P-11	P-11	P-12	P-12	P-12	P-12
			01/15/2014	07/29/2014	01/24/2018	03/23/2018	05/24/2018	01/09/2019	07/11/2019	01/14/2020	01/29/2008	07/16/2008	07/16/2008 Duplicate	01/22/2009
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002		<0.00109	<0.00052	
Benzene	0.005	0.005	0.000207 J	<0.00008	<0.0002	<0.0002	0.00021 J	<0.0002	<0.0002	<0.0002		<0.00112	<0.00025	
Chlorobenzene	0.1	0.1	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003		<0.0015	<0.00047	
Ethylbenzene	0.7	0.7	0.000253 J	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003		<0.00142	<0.00025	
Methylene chloride	0.005	0.005	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001		<0.00122	<0.00054	
Toluene	1	1	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002		<0.00138	<0.00041	
Vinyl chloride	0.002	0.002												
Xylenes (total)	10	10	<0.00058	<0.00026	<0.0003	<0.0003	0.00032 J	<0.0003	<0.0003	<0.0003		<0.00302	<0.00127	
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000104	<0.000108	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021				
2,4-Dimethylphenol	0.49	1.5	<0.000292	<0.000304	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004				
2,4-Dinitrotoluene	0.0013	0.003	<0.000123	<0.000127	<0.000058	<0.000058	<0.000059	<0.000058	<0.000058	<0.000058				
2,6-Dinitrotoluene	0.0013	0.003	<0.0000755	<0.0000784	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042				
2-Chloronaphthalene	2	5.8	<0.0000755	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021				
2-Methylnaphthalene	0.098	0.29	0.000257 J	<0.0000686	<0.000019	<0.000019	0.00015	<0.000019	0.00005 J	0.000076 J				
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.000783	<0.000814	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002				
4-Nitrophenol	0.049	0.15	<0.000528	<0.000549	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047				
Acenaphthene	1.5	4.4	0.00951	0.000653	<0.000027	0.00021	0.081	<0.000027	0.0019	0.0046	<0.00029	<0.0003	<0.0003	<0.0008
Acenaphthylene	1.5	4.4	<0.0000566	<0.0000588	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00029	<0.0003	<0.0003	<0.0007
Anthracene	7.3	22	0.000503	0.000119 J	<0.000014	<0.000014	0.0037	<0.000014	0.000063 J	0.000081 J	0.000645	0.000552	0.000566	<0.0007
Benzo(a)anthracene	0.0091	0.02	<0.0000755	<0.0000784	<0.00005	<0.00005	<0.000051	<0.00005	<0.00005	<0.00005				
Benzo(a)pyrene	0.0002	0.0002	<0.0000755	<0.0000784	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002				
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.000123	<0.000127	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003				
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	0.00403	<0.000711	0.00018 J	<0.000037	0.000063 J	<0.000037	0.000087 J	<0.000037	<0.00019	0.00034 J	0.00064 J	<0.0012
Chrysene	0.91	2	<0.0000755	<0.0000784	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021				
Dibenzofuran	0.098	0.29	0.00103	0.000176 J	<0.00002	<0.00002	0.0016	<0.00002	0.000079 J	0.00016	<0.00029	<0.0003	<0.0003	<0.0007
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000104	<0.000108	<0.00002	<0.00002	0.000025 J	<0.00002	<0.00002	<0.00002	<0.00019	0.00085 J	0.00087 J	<0.0007
Fluoranthene	0.98	2.9	0.000287 J	0.0000771 J	0.000038 J	<0.00001	0.0056	<0.00001	0.00006 J	<0.00001	<0.00019	<0.0002	<0.0002	<0.0006
Fluorene	0.98	2.9	0.00264	0.000344 J	0.000045 J	<0.00003	0.037	<0.00003	0.00074	0.0013	<0.00019	<0.0002	<0.0002	<0.0008
Naphthalene	0.49	1.5	0.0554	<0.0000784	<0.00002	<0.00002	0.0024	<0.00002	<0.0018	0.00048	<0.00038	0.000626	0.000639	<0.0008
Nitrobenzene	0.049	0.15	<0.000104	<0.000108	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024				
N-Nitrosodiphenylamine	0.19	0.42	<0.0000943	<0.000098	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025				
Pentachlorophenol	0.001	0.001	<0.000575	<0.000598	<0.000079	<0.000079	<0.00008	<0.000079	<0.000079	<0.000079				
Phenanthrene	0.73	2.2	0.00189	<0.000317	<0.000021	<0.000021	0.04	<0.000021	0.0002	0.00035				
Phenol	7.3	22	<0.0000377	<0.0000392	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00019	<0.0002	<0.0002	<0.0015
Pyrene	0.73	2.2	0.000274 J	<0.000108	0.000039 J	<0.000019	0.0032	<0.000019	0.000036 J	<0.000019	0.00932	0.00211	0.00166	0.0026 J
<b>Metals</b>														
Arsenic	0.01	0.01			<b>0.0374</b>	<b>0.016</b>	<b>0.0622</b>	<b>0.0183</b>	<b>0.0704</b>	<b>0.036</b>				

Notes:

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2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

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	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	P-12 01/22/2009	P-12 07/22/2009	P-12 01/22/2010	P-12 07/14/2010	P-12 01/12/2011	P-12 07/12/2011	P-12 01/31/2012	P-12 07/11/2012	P-12 01/09/2013	P-12 07/11/2013	P-12 01/09/2014	P-12 07/02/2014	
<b>Volatile Organic Compounds</b>															
1,2-Dichloroethane	0.005	0.005													
Benzene	0.005	0.005													
Chlorobenzene	0.1	0.1													
Ethylbenzene	0.7	0.7													
Methylene chloride	0.005	0.005													
Toluene	1	1													
Vinyl chloride	0.002	0.002													
Xylenes (total)	10	10													
<b>Semivolatile Organic Compounds</b>															
1,2-Diphenylhydrazine	0.0011	0.0026													
2,4-Dimethylphenol	0.49	1.5													
2,4-Dinitrotoluene	0.0013	0.003													
2,6-Dinitrotoluene	0.0013	0.003													
2-Chloronaphthalene	2	5.8													
2-Methylnaphthalene	0.098	0.29		<0.0009										<0.000066	
4,6-Dinitro-2-methylphenol	0.0024	0.0073													
4-Nitrophenol	0.049	0.15													
Acenaphthene	1.5	4.4	<0.0008	<0.0009	<0.0009	<0.0009	<0.0009	<0.0009	<0.0005	<0.0005	<0.0005	<0.0000755	<0.00008	<0.0000741	<0.0000755
Acenaphthylene	1.5	4.4	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000566	<0.00006	<0.0000556	<0.0000566
Anthracene	7.3	22	<0.0007	<0.0006	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.0000472	<0.00005	0.0002 J	0.000189 J
Benzo(a)anthracene	0.0091	0.02													
Benzo(a)pyrene	0.0002	0.0002													
bis(2-Chloroethoxy)methane	0.00083	0.0019													
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.0012	<0.0033	<0.0033	<0.0033	<0.0033	<0.0033	<0.0005	<0.0005	<0.0005	0.00142 J	0.00039 J	0.000515 J	0.000439 J
Chrysene	0.91	2													
Dibenzofuran	0.098	0.29	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007	<0.0007	<0.0005	<0.0005	<0.0005	<0.0000755	<0.00008	<0.0000741	<0.0000755
Di-n-butylphthalate (DBP)	2.4	7.3	<0.0007	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.00011	<0.000416	0.000144 J
Fluoranthene	0.98	2.9	<0.0006	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000066	<0.00007	<0.0000648	<0.000066
Fluorene	0.98	2.9	<0.0008	<0.0006	<0.0006	<0.0006	<0.0006	<0.0006	<0.0005	<0.0005	<0.0005	<0.000066	<0.00007	<0.0000648	<0.000066
Naphthalene	0.49	1.5	<0.0008	<0.0006	<0.0006	<0.0006	0.0006 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000755	<0.00008	<0.0000741	<0.0000755
Nitrobenzene	0.049	0.15													
N-Nitrosodiphenylamine	0.19	0.42													
Pentachlorophenol	0.001	0.001													
Phenanthrene	0.73	2.2		<0.0005											<0.0000566
Phenol	7.3	22	<0.0015	<0.0005	<0.0005	<0.0005	0.0005 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.0000377	<0.00004	<0.000037	<0.0000377
Pyrene	0.73	2.2	0.0012 J	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.000104	<0.00011	<0.000102	0.00189
<b>Metals</b>															
Arsenic	0.01	0.01													

- Notes:
- All values in milligrams per liter (mg/L).
  - Concentrations > RAL and non-detects are highlighted light gray.
  - Concentrations > C/I AL and non-detects are highlighted dark gray
  - TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
  - RAL = Residential Assessment Level, C/I = Commercial/Industrial
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ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	P-12 01/07/2015	P-12 07/08/2015	P-12 01/13/2016	P-12 07/07/2016	P-12 01/12/2017	P-12 07/13/2017	P-12 01/04/2018	P-12 07/19/2018	P-12 01/07/2019	P-12 07/01/2019	P-12 01/13/2020	TW-41B 01/19/2010
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.005	0.005												<0.0005
Benzene	0.005	0.005												<0.0005
Chlorobenzene	0.1	0.1												<0.0005
Ethylbenzene	0.7	0.7												<0.0005
Methylene chloride	0.005	0.005												<0.0005
Toluene	1	1												<0.0005
Vinyl chloride	0.002	0.002												<0.0005
Xylenes (total)	10	10												<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.0011	0.0026												<0.0001
2,4-Dimethylphenol	0.49	1.5												<0.00008
2,4-Dinitrotoluene	0.0013	0.003												<0.00009
2,6-Dinitrotoluene	0.0013	0.003												<0.00007
2-Chloronaphthalene	2	5.8												<0.0001
2-Methylnaphthalene	0.098	0.29												<0.00007
4,6-Dinitro-2-methylphenol	0.0024	0.0073												<0.00008
4-Nitrophenol	0.049	0.15												<0.00007
Acenaphthene	1.5	4.4	<0.0000792	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.00009
Acenaphthylene	1.5	4.4	<0.0000594	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.00007
Anthracene	7.3	22	<0.0000495	<0.000014	<0.000014	0.000026 J	<0.000014	<0.000014	0.000036 J	<0.000014	0.000063 J	0.000052 J	0.0001	<0.00007
Benzo(a)anthracene	0.0091	0.02												<0.00007
Benzo(a)pyrene	0.0002	0.0002												<0.00008
bis(2-Chloroethoxy)methane	0.00083	0.0019												<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000366	0.00055	0.00036	0.000086 J	0.00029	0.00026	<0.00012	<0.00009	<0.000037	<0.000037	<0.000037	0.0011
Chrysene	0.91	2												<0.00007
Dibenzofuran	0.098	0.29	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008
Di-n-butylphthalate (DBP)	2.4	7.3	<0.000109	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.000022	<0.00002	<0.00002	<0.00002	<0.00002	<0.00007
Fluoranthene	0.98	2.9	<0.0000693	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001	<0.00007
Fluorene	0.98	2.9	<0.0000693	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	0.00015 J
Naphthalene	0.49	1.5	<0.0000792	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.00016	0.00014 J
Nitrobenzene	0.049	0.15												<0.00009
N-Nitrosodiphenylamine	0.19	0.42												<0.00009
Pentachlorophenol	0.001	0.001												<0.00008
Phenanthrene	0.73	2.2												<0.00007
Phenol	7.3	22	<0.0000396	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.00007
Pyrene	0.73	2.2	0.00152 J	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.000097 J	<0.000019	<0.000019	<0.000019	0.00063	<0.00007
<b>Metals</b>														
Arsenic	0.01	0.01												

**Notes:**

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Residential Assessment Level</i>	<i>C/I Assessment Level</i>	TW-41B 07/27/2011	TW-41B 02/01/2012	TW-41B 07/26/2012	TW-41B 02/05/2013	TW-41B 07/31/2013	TW-41B 01/16/2014	TW-41B 07/25/2014	TW-41B 01/24/2018	TW-41B 03/20/2018	TW-41B 05/16/2018	TW-41B 01/09/2019	TW-41B 07/12/2019
<b><i>Volatile Organic Compounds</i></b>														
1,2-Dichloroethane	0.005	0.005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.002	<0.0002
Benzene	0.005	0.005	<0.001	<0.001	<0.0005	<0.00008	0.000347 J	<0.0002	0.000594 J	0.00065 J	0.001 J	0.0013	<0.002	<0.0013
Chlorobenzene	0.1	0.1	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.003	<0.0003
Ethylbenzene	0.7	0.7	0.0075	<0.0011	<0.0005	<0.00011	0.00115	<0.00019	0.00501	<0.0003	0.0036	0.0029	<0.003	<0.0011
Methylene chloride	0.005	0.005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1	1	0.0033 J	<0.001	<0.0005	<0.00015	<0.000279	<0.00017	0.00116	<0.0002	0.00068 J	0.0012	<0.002	<0.0011
Vinyl chloride	0.002	0.002				<0.00011			<0.00011					
Xylenes (total)	10	10	0.0052 J	<0.0031	<0.0015	<0.00026	0.000386 J	<0.00058	0.0101	0.0079	0.013	0.015	<0.003	0.019
<b><i>Semivolatile Organic Compounds</i></b>														
1,2-Diphenylhydrazine	0.0011	0.0026	<0.00005	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	0.49	1.5	<0.00005	0.00005 J	0.0014	<0.000292	<0.000292	<0.000301	<0.000041	<0.000041	<0.000041	<0.000041	<0.000041	<0.000041
2,4-Dinitrotoluene	0.0013	0.003	<0.00005	0.00005 J	<0.00005	<0.000123	<0.000123	<0.000126	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.0013	0.003	<0.00006	0.00006 J	<0.00006	<0.0000755	<0.0000762	<0.0000755	<0.0000777	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	2	5.8	<0.00005	0.00005 J	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	0.098	0.29	0.015	0.00005 J	<0.0001	<0.000066	0.000256 J	<0.0000846	0.0125	0.0003	0.011	0.026	0.0098	0.067
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00008	<0.00008	<0.00008	<0.000783	<0.00079	<0.000783	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	0.049	0.15	<0.00005	0.00005 J	<0.00005	<0.000528	<0.000533	<0.000528	<0.000544	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	1.5	4.4	0.041	0.00005 J	0.039	<0.0000755	0.0252	<0.0000755	0.142	0.087	0.072	0.08	0.058	0.15
Acenaphthylene	1.5	4.4	0.00053	0.00005 J	0.00041	0.0000751 J	0.000409 J	0.0000926 J	0.00185	0.0019	0.0017	0.0016	0.00091	0.0014
Anthracene	7.3	22	0.0022	0.00016 J	0.0011	0.000979	0.00161	0.00093	0.00697	0.0016	0.0034	0.0039	0.0023	0.0055
Benzo(a)anthracene	0.0091	0.02	<0.00005	<0.00005	<0.00005	<0.0000755	0.0000879 J	<0.0000755	<0.0000777	<0.000051	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.0002	0.0002	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000777	<0.00002	<0.00002	<0.00002	0.000097 J	<0.00002
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00005	0.00005 J	<0.00005	<0.000123	<0.000124	<0.000123	<0.000126	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.00022	<0.00023	<0.0001	<0.000349	<0.000352	<0.000349	<0.000359	0.000058 J	<0.000037	<0.000037	<0.000037	<0.000037
Chrysene	0.91	2	<0.00005	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	0.098	0.29	0.029	0.00005 J	0.016	<0.0000755	0.0104	<0.0000755	0.0845	0.022	0.026	0.034	0.026	0.074
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00005	<0.00005	<0.0001	<0.000104	<0.000138	0.000116 J	<0.000107	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	0.98	2.9	0.0022	<0.00005	0.0015	<0.000066	0.00153	0.000206 J	0.00475	0.0019	0.0026	0.0022	0.0014	0.0034
Fluorene	0.98	2.9	0.028	0.00005 J	0.0054 J	0.0000917 J	0.00386	<0.000066	<0.0811	0.035	0.037	0.045	0.035	0.085
Naphthalene	0.49	1.5	0.049	0.00051 J	<0.0013	0.000156 J	0.00309 J	<0.000259	0.149	0.027	0.12	0.28	0.061	<b>0.69</b>
Nitrobenzene	0.049	0.15	<0.00005	0.00005 J	<0.00005	<0.000104	<0.000105	<0.000104	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	0.19	0.42	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0000952	<0.0000943	<0.0000971	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.001	0.001	<0.00005	0.00005 J	<0.00005	<0.000575	<0.000581	<0.000575	<0.000592	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	0.73	2.2	0.019	<0.00005	<0.00005	<0.0000566	0.000659	<0.000176	0.0573	0.00048	0.0089	0.013	0.0035	0.027
Phenol	7.3	22	<0.00005	0.000057 J	0.0016	<0.0000377	<0.0000381	<0.0000377	<0.0000388	<0.000036	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	0.73	2.2	0.00095	<0.00005	0.00066	<0.000104	0.000702	0.000223 J	0.00209	0.00083	0.0011	0.001	0.00056	0.0014
<b><i>Metals</i></b>														
Arsenic	0.01	0.01								<b>0.0376</b>	<b>0.0953</b>	<b>0.0976</b>	<b>0.125</b>	<b>0.113</b>

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.
- MW-32A was screened in the B-CZ & replaced with MW-32AR
- Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 2  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 2 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Residential Assessment Level</b>	<b>C/I Assessment Level</b>	TW-41B 01/13/2020	TW-56A 01/20/2010	TW-56A 07/14/2011	TW-56A 02/02/2012	TW-56A 07/11/2012	TW-56A 01/31/2013
<b>Volatile Organic Compounds</b>								
1,2-Dichloroethane	0.005	0.005	<0.0002	<b>0.023 J</b>	<0.005	<b>&lt;0.01</b>	<0.0025	<0.0014
Benzene	0.005	0.005	0.00039 J	<b>0.26</b>	<b>0.27</b>	<b>0.15</b>	<b>0.26</b>	<b>0.238</b>
Chlorobenzene	0.1	0.1	<0.0003	<0.0025	<0.005	<0.01	<0.0025	0.00412 J
Ethylbenzene	0.7	0.7	<0.0003	0.36	0.16	0.068	0.14	0.202
Methylene chloride	0.005	0.005	<0.001	<0.0025	<b>&lt;0.0065</b>	<b>&lt;0.013</b>	<0.005	<0.0015
Toluene	1	1	<0.0002	0.32	0.14	0.028 J	0.069	0.0314
Vinyl chloride	0.002	0.002		<b>&lt;0.0025</b>	<b>0.0069 J</b>	<b>0.01 J</b>	<b>0.016</b>	<b>0.0126 J</b>
Xylenes (total)	10	10	0.0021	0.98	0.61	0.53	0.43	0.5
<b>Semivolatile Organic Compounds</b>								
1,2-Diphenylhydrazine	0.0011	0.0026	<0.000021	<0.0001	<0.00005	<0.0005	<0.00005	<b>&lt;0.055</b>
2,4-Dimethylphenol	0.49	1.5	0.00017 J	<b>2.9</b>	<b>6.8</b>	<b>4.2</b>	<b>3.8</b>	<b>4.81</b>
2,4-Dinitrotoluene	0.0013	0.003	<0.000058	<0.00009	<0.00005	<0.0005	<0.00005	<b>&lt;0.065</b>
2,6-Dinitrotoluene	0.0013	0.003	<0.000042	<0.00007	<0.00006	<0.0006	<0.00006	<b>&lt;0.04</b>
2-Chloronaphthalene	2	5.8	0.00012 J	<0.0001	<0.00005	<0.0005	<0.00005	<0.04
2-Methylnaphthalene	0.098	0.29	0.0071	<b>0.15</b>	<b>0.16</b>	<b>0.11</b>	0.052	<b>0.123 J</b>
4,6-Dinitro-2-methylphenol	0.0024	0.0073	<0.00002	<0.00008	<0.00008	<0.0008	<0.00008	<b>&lt;0.415</b>
4-Nitrophenol	0.049	0.15	<0.000047	<0.00007	<0.00005	<0.0005	<0.00005	<b>&lt;0.28</b>
Acenaphthene	1.5	4.4	0.12	0.077	0.18	0.19	0.095	0.25
Acenaphthylene	1.5	4.4	0.0011	0.0024	0.004	0.0038	0.0028	<0.03
Anthracene	7.3	22	0.0031	0.0035	0.021	0.02	0.0083	0.0338 J
Benzo(a)anthracene	0.0091	0.02	<0.00005	0.00099	0.0014	0.0016 J	0.0024	<b>&lt;0.04</b>
Benzo(a)pyrene	0.0002	0.0002	<0.00002	<b>0.00031</b>	<b>0.00047</b>	<b>0.00051 J</b>	<b>0.0008</b>	<b>&lt;0.04</b>
bis(2-Chloroethoxy)methane	0.00083	0.0019	<0.00003	<0.00009	<0.00005	<0.0005	<0.00005	<b>&lt;0.065</b>
bis(2-Ethylhexyl)phthalate (DEHP)	0.006	0.006	<0.000037	<0.00025	<0.0018	<0.001	<0.0001	<b>&lt;0.185</b>
Chrysene	0.91	2	<0.000021	0.00084	0.0014	0.0018 J	0.0022	<0.04
Dibenzofuran	0.098	0.29	0.055	0.043	0.09	0.049	0.038	<b>0.108 J</b>
Di-n-butylphthalate (DBP)	2.4	7.3	<0.00002	0.00045	<0.00005	<0.0005	<0.00005	<0.055
Fluoranthene	0.98	2.9	0.0021	0.0021	0.021	0.02	0.027	<0.035
Fluorene	0.98	2.9	0.078	0.033	0.09	0.058	0.047	0.12 J
Naphthalene	0.49	1.5	0.058	<b>2.5</b>	<b>2.3</b>	<b>2.2</b>	<b>0.81</b>	<b>1.75 J</b>
Nitrobenzene	0.049	0.15	<0.000024	<0.00009	<0.00005	<0.0005	<0.00005	<b>&lt;0.055</b>
N-Nitrosodiphenylamine	0.19	0.42	<0.000025	<0.00009	<0.00005	<0.0005	<0.00005	<0.05
Pentachlorophenol	0.001	0.001	<0.000079	0.00013 J	0.00076 J	<0.0005	0.00091	<b>&lt;0.305</b>
Phenanthrene	0.73	2.2	0.0084	0.06	0.17	0.2	0.073	0.217 J
Phenol	7.3	22	<0.000035	0.014	<0.00005	0.0063	<0.00005	<0.02
Pyrene	0.73	2.2	0.00082	0.0067	0.012	0.015	0.018	<0.055
<b>Metals</b>								
Arsenic	0.01	0.01	<b>0.0557</b>					

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray.
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.
7. MW-32A was screened in the B-CZ & replaced with MW-32AR
8. Based on historical data, MW-25A and MW-25C were likely mislabeled in March 2018 and results are provided correctly in this table.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-32B 02/09/2012 DNAPL	MW-32B 07/16/2012 DNAPL	MW-32B 02/06/2013 DNAPL	MW-32B 01/21/2014 DNAPL	MW-32B 01/20/2020 DNAPL	MW-33B 01/29/2008 DNAPL	MW-33B 07/14/2008 DNAPL	MW-33B 07/14/2008 Duplicate	MW-33B 02/03/2009 DNAPL	MW-33B 01/13/2010 DNAPL	MW-33B 06/29/2010 DNAPL	MW-33B 01/24/2011 DNAPL
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	0.025 J	<0.0005	<0.00014	<0.0002	<0.005	<0.00052	<0.00109	<0.00109	<0.005	<0.005	<0.005	<0.005
Benzene	0.5	0.5	<b>2.6 J</b>	<0.0005	0.00428	0.239	<b>1.8</b>	<b>1.92</b>	<b>2.73</b>	<b>2.69</b>	<b>2.4</b>	<b>1.2</b>	<b>2</b>	<b>1</b>
Chlorobenzene	10	10	0.025 J	<0.0005	<0.000343	<0.00018	<0.0075	<0.00047	<0.0015	<0.0015	<0.005	<0.005	<0.005	<0.005
Ethylbenzene	70	70	0.53 J	<0.0005	0.00561	0.254	0.73	0.491	0.626	0.598	0.47	0.41	0.62	0.36
Methylene chloride	0.5	0.5	0.032 J	<0.001	<0.00015	<0.00022	<0.025	<0.00054	<0.00122	<0.00122	0.0096 J	<0.005	<0.005	<0.005
Toluene	100	100	2.2 J	<0.0005	0.00261	0.541	2.3	0.1	0.136	0.13	0.084	0.019 J	0.016 J	0.0067 J
Vinyl chloride	0.2	0.2					<0.005							
Xylenes (total)	1000	1000	1.5 J	<0.0015	0.0203	0.749	2.1	1.24	1.63	1.58	1.4	1.2	1.5	0.85
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	0.0005 J	<0.00005	<0.0262	<0.00104	<0.00042	<0.01	<0.008	<0.008	<0.0001	<0.0001	<0.0001	<0.0001
2,4-Dimethylphenol	49	150	46	0.0014	<0.0738	0.178	26	<0.043	<0.028	<0.028	<0.00008	0.0035	<0.00008	0.0029
2,4-Dinitrotoluene	0.13	0.3	0.0005 J	<0.00005	<0.031	<0.00123	<0.0012	<0.029	<0.019	<0.019	<0.00009	<0.00009	<0.00009	<0.00009
2,6-Dinitrotoluene	0.13	0.3	0.0006 J	<0.00006	<0.019	<0.000755	<0.00084	<0.029	<0.019	<0.019	<0.00007	<0.00007	<0.00007	<0.00007
2-Chloronaphthalene	200	580	0.0005 J	<0.00005	<0.019	<0.000755	<0.00042	<0.057	<0.038	<0.038	<0.00012	<0.0001	<0.0001	<0.0001
2-Methylnaphthalene	9.8	29	0.53	0.00019 J	<0.0167	0.137	<b>50</b>	0.443	0.808	0.787	1.9	0.71	0.51	0.52
4,6-Dinitro-2-methylphenol	0.24	0.73	0.0008 J	<0.00008	<0.198	<0.00783	<0.0004	<0.029	<0.019	<0.019	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	4.9	15	0.0005 J	<0.00005	<0.133	<0.00528	<0.00094	<0.036	<0.024	<0.024	<0.00007	<0.00007	<0.00007	<0.00007
Acenaphthene	150	440	0.28 J	0.014	0.0416 J	0.0427	26	0.137	0.152	0.182	0.41	0.17	0.096	0.15
Acenaphthylene	150	440	0.0059 J	0.00085	<0.0143	<0.000566	<0.0003	<0.043	<0.028	<0.028	0.0037	0.0016	0.0011	0.0015
Anthracene	730	2200	0.059 J	0.0048	<0.0119	0.144	23	<0.029	<0.019	0.035	0.14	0.015	0.011	0.027
Benzo(a)anthracene	0.91	2	0.0033 J	0.0033	<0.019	0.0195	<b>2.2</b>	<0.029	<0.019	<0.019	0.022	0.00019 J	0.000073 J	0.0019
Benzo(a)pyrene	0.02	0.02	0.0005 J	0.00089	<0.019	0.00649	<b>0.74</b>	<0.029	<0.019	<0.019	0.0045	<0.00008	<0.00008	0.00073
bis(2-Chloroethoxy)methane	0.083	0.19	<0.0005	<0.00005	<0.031	<0.00123	<0.0006	<0.057	<0.038	<0.038	<0.00009	<0.00009	<0.00009	<0.00009
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	0.001 J	0.00079	<0.0881	<0.00349	<0.00074	<0.029	<0.019	<0.019	0.00031	0.008	<0.00054	0.00091
Chrysene	91	200	0.0042 J	0.0023	<0.019	0.018	2.4	<0.029	<0.019	<0.019	0.02	0.00018 J	0.000092 J	0.0018
Dibenzofuran	9.8	29	0.28 J	0.0012	<0.019	0.0428	<b>28</b>	0.118	0.17	0.205	0.46	0.18	0.13	0.17
Di-n-butylphthalate (DBP)	240	730	0.0005 J	<0.00005	<0.0262	<0.00104	<0.0004	<0.029	<0.019	<0.019	<0.00007	<0.00007	<0.00007	<0.00007
Fluoranthene	98	290	0.03 J	0.031	<0.0167	0.121	21	<0.029	<0.019	0.042	0.2	0.0033	0.0018	0.033
Fluorene	98	290	0.15 J	0.0021	<0.0167	0.0282	23	0.046	0.0683	0.098	0.26	0.068	0.048	0.069
Naphthalene	49	150	26	<0.00057	<0.019	2.17 J	<b>300</b>	12.5	16	13.1	20	10	2.2	7
Nitrobenzene	4.9	15	<0.0005	<0.00005	<0.0262	<0.00104	<0.00048	<0.057	<0.038	<0.038	<0.00009	<0.00009	<0.00009	<0.00009
N-Nitrosodiphenylamine	19	42	0.0005 J	<0.00005	<0.0238	<0.000943	<0.0005	<0.036	<0.024	<0.024	<0.00009	<0.00009	<0.00009	<0.00009
Pentachlorophenol	0.1	0.1	0.0005 J	<0.00005	<b>&lt;0.145</b>	<0.00575	<0.0016	<0.029	<0.019	<0.019	<0.00008	<0.00008	<0.00008	<0.00008
Phenanthrene	73	220	0.25 J	0.0012	<0.0143	0.0548	69	0.0903	0.0688	0.186	0.72	0.066	0.041	0.09
Phenol	730	2200	38 J	0.00066 J	<0.00952	0.0357	17	<0.029	<0.019	<0.019	0.003	<0.00007	0.0032	<0.00007
Pyrene	73	220	0.02 J	0.04	<0.0262	0.0841	14	0.045	<0.019	0.022	0.13	0.0016	0.00092	0.007
<b>Metals</b>														
Arsenic	1	1						0.00193 J						

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray.
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.



**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-33B 07/19/2011 DNAPL	MW-33B 02/15/2012	MW-33B 07/17/2012	MW-33BR 02/06/2013	MW-33BR 08/07/2013	MW-33BR 01/21/2014	MW-33BR 07/28/2014	MW-33BR 01/28/2018	MW-33BR 03/29/2018	MW-33BR 05/31/2018	MW-33BR 01/22/2019	MW-33BR 07/30/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.01	<0.005	<0.0005	<0.007	<0.0007	<0.0002	<0.0007	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.5	0.5	1.6	2	0.3	1.61	1.62	0.837	1.41	<0.0002	<0.0002	0.12	0.0025	0.25
Chlorobenzene	10	10	<0.01	<0.005	<0.0005	<0.006	<0.0006	0.000349 J	<0.0006	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	70	70	0.4	0.46	0.07	0.471	0.389	0.128	0.348	<0.0003	<0.0003	0.0058	0.013	0.065
Methylene chloride	0.5	0.5	<0.013	<0.0065	<0.001	0.011 J	<0.00075	<0.00022	<0.00075	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	100	100	<0.01	0.12	0.023	0.157	0.0645	0.00942	0.00638	<0.0002	<0.0002	<0.0002	<0.0002	0.0031
Vinyl chloride	0.2	0.2	<0.01	<0.005	<0.0005	<0.0055	<0.00055	<0.00018	<0.00055	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	1000	1000	1.2	0.82	0.15	0.924	0.182	0.128	0.0649	<0.0003	<0.0003	0.0058	<0.0003	0.016
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.00005	<0.00005	<0.0005	<0.529	<0.00519	<0.000529	<0.000107	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	49	150	0.0034	<0.00005	0.0074	<1.49	<0.0146	<0.00149	<0.000301	<0.00004	<0.00004	<0.00004	<0.00004	0.00028
2,4-Dinitrotoluene	0.13	0.3	<0.00005	<0.00005	<0.0005	<0.625	<0.00613	<0.000625	<0.000126	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.13	0.3	<0.00006	<0.00006	<0.0006	<0.385	<0.00377	<0.000385	<0.0000777	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	200	580	<0.00005	<0.00005	<0.0005	<0.385	<0.00377	<0.000385	<0.0000777	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	9.8	29	1.6	0.81	0.55	0.993 J	0.198	0.0558	0.277	<0.000019	<0.000019	0.0029	<0.000019	0.014
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00008	<0.00008	<0.0008	<3.99	<0.0392	<0.00399	<0.000806	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	4.9	15	<0.00005	<0.00005	<0.0005	<2.69	<0.0264	<0.00269	<0.000544	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	150	440	0.41	0.23	0.099	<0.385	0.0455	0.0625	0.0711	<0.000027	<0.000027	0.0019	0.0013	0.0078
Acenaphthylene	150	440	0.0033	<0.00005	0.0014 J	<0.288	<0.00283	<0.000679	0.00087	<0.000015	<0.000015	0.000068 J	<0.000015	<0.000015
Anthracene	730	2200	0.16	0.054	0.011	<0.24	<0.00236	0.0045	0.00564	<0.000014	<0.000014	0.00018	<0.000014	0.00055
Benzo(a)anthracene	0.91	2	0.032	0.000074 J	<0.0005	<0.385	<0.00377	<0.000385	0.000119 J	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.02	0.02	0.0077	<0.00005	<0.0005	<0.385	<0.00377	<0.000385	<0.0000777	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00005	<0.00005	<0.0005	<0.625	<0.00613	<0.000625	<0.000126	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.00046	<0.00018	<0.001	<1.78	<0.0175	<0.00178	0.000722	0.000062 J	<0.000037	0.0001 J	<0.000037	0.00007 J
Chrysene	91	200	0.026	0.000073 J	<0.0005	<0.385	<0.00377	<0.000385	0.000132 J	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	9.8	29	0.53	0.38	0.15	<0.385	0.0498	0.0769	0.0868	<0.000021	<0.00002	0.0019	0.000079 J	0.012
Di-n-butylphthalate (DBP)	240	730	<0.00005	<0.00005	<0.0005	<0.529	<0.00519	<0.000529	<0.000384	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	98	290	0.28	0.0049	0.01	<0.337	<0.0033	0.00107 J	0.00265	0.000049 J	<0.00001	0.0003	0.000053 J	0.00039
Fluorene	98	290	0.31	0.12	0.051	<0.337	0.0181 J	0.0287	0.035 J	<0.00003	<0.00003	0.00058	<0.00003	0.0042
Naphthalene	49	150	13	21	7.3	14.9 J	6.54	1.68 J	6.59	<0.00018	<0.00002	0.069	0.00004 J	0.53 J
Nitrobenzene	4.9	15	<0.00005	<0.00005	<0.0005	<0.529	<0.00519	<0.000529	<0.000107	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	19	42	<0.00005	0.00081	<0.0005	<0.481	<0.00472	<0.000481	<0.0000971	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.1	0.1	<0.00005	<0.00005	<0.0005	<2.93	<0.0288	<0.00293	<0.000592	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	73	220	0.79	0.17	0.091	<0.288	0.0141 J	0.0238	0.0313 J	<0.000047	<0.000021	0.0008	<0.000021	0.0048
Phenol	730	2200	0.001	0.0043	0.0014 J	<0.192	<0.00189	<0.000192	<0.0000388	<0.000035	<0.000035	<0.000035	<0.000035	0.00021
Pyrene	73	220	0.17	0.0025	0.0054	<0.529	<0.00519	0.000734 J	0.00126	<0.000019	<0.000019	0.00019	0.00003 J	0.00026
<b>Metals</b>														
Arsenic	1	1								0.00144 J	0.00187 J	0.00294	0.00143 J	0.00117 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-33BR 01/15/2020	MW-36B 07/15/2010	MW-36B 07/15/2010	MW-36B 01/20/2011	MW-36B 07/19/2011	MW-36B 02/08/2012	MW-36B 07/17/2012	MW-36B 01/31/2013	MW-36B 08/06/2013	MW-36B 01/16/2014	MW-36B 07/28/2014	MW-36B 01/25/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002
Benzene	0.5	0.5	0.12	<0.0005	<0.0005	0.0018 J	0.0014 J	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002
Chlorobenzene	10	10	<0.0003	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003
Ethylbenzene	70	70	0.043	<0.0005	<0.0005	<0.0005	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003
Methylene chloride	0.5	0.5	<0.001	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001
Toluene	100	100	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002
Vinyl chloride	0.2	0.2	<0.0002	<0.0002	<0.0002	<0.0002	<0.001	<0.001	<0.0005	<0.00011	<0.00011	<0.00018	<0.00011	<0.0002
Xylenes (total)	1000	1000	0.011	<0.001	<0.001	<0.001	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.000021	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.000104	<0.00011	<0.000104	<0.000107	<0.000021
2,4-Dimethylphenol	49	150	0.00018 J	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000292	<0.00031	<0.000292	<0.000301	<0.00004
2,4-Dinitrotoluene	0.13	0.3	<0.000058	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.00013	<0.000123	<0.000126	<0.000059
2,6-Dinitrotoluene	0.13	0.3	<0.000042	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.0000755	<0.00008	<0.0000755	<0.0000777	<0.000042
2-Chloronaphthalene	200	580	<0.000021	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.0000755	<0.00008	<0.0000755	<0.0000777	<0.000021
2-Methylnaphthalene	9.8	29	0.00025	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000762	<0.00007	<0.000271	<0.000068	<0.000019
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.000783	<0.00083	<0.000783	<0.000806	<0.00002
4-Nitrophenol	4.9	15	<0.000047	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000528	<0.00056	<0.000528	<0.000544	<0.000047
Acenaphthene	150	440	0.0051	<0.00009	<0.00009	0.00023	0.00014 J	0.00023	0.00016 J	<0.0000755	<0.00008	0.000463 J	<0.0000777	0.00014
Acenaphthylene	150	440	0.00015	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000566	<0.00006	<0.0000566	<0.0000583	<0.000015
Anthracene	730	2200	0.00046	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000472	<0.00005	0.00035 J	<0.0000485	<0.000014
Benzo(a)anthracene	0.91	2	0.0001	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000755	<0.00008	0.00012 J	<0.0000777	<0.000051
Benzo(a)pyrene	0.02	0.02	0.000039 J	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.0000755	<0.00008	<0.0000755	<0.0000777	<0.00002
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00003	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000123	<0.00013	<0.000123	<0.000126	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	0.00067	0.01	0.0024	<0.00048	<0.00068	<0.00033	0.00021	<0.000349	<0.00037	0.00044 J	<0.000359	0.00015 J
Chrysene	91	200	0.00011	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.0000755	<0.00008	0.000146 J	<0.0000777	<0.000021
Dibenzofuran	9.8	29	0.0039	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	0.00011 J	<0.000118	<0.00008	<0.000409	<0.0000777	<0.00002
Di-n-butylphthalate (DBP)	240	730	0.00099	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000104	<0.00011	<0.000104	<0.000107	<0.00002
Fluoranthene	98	290	0.00082	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000066	<0.00007	0.000756	<0.000068	<0.00001
Fluorene	98	290	0.0018	<0.00007	<0.00007	<0.00007	<0.00005	0.00011 J	<0.00005	<0.000066	<0.00007	0.000434 J	<0.000068	<0.00003
Naphthalene	49	150	0.0051	<0.0001	<0.0001	<0.0001	<0.00005	<0.00024	<0.00005	<0.000943	0.0000895 J	0.000825 J	<0.0000777	<0.00002
Nitrobenzene	4.9	15	<0.000024	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.000104	<0.00011	<0.000104	<0.000107	<0.000024
N-Nitrosodiphenylamine	19	42	<0.000025	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.0000943	<0.0001	<0.0000943	<0.0000971	<0.000025
Pentachlorophenol	0.1	0.1	<0.000079	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	<0.00005	<0.000575	<0.00061	<0.000575	<0.000592	<0.00008
Phenanthrene	73	220	0.0013	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00027	<0.0000566	<0.00006	0.00183	<0.0000583	<0.000021
Phenol	730	2200	0.00018 J	<0.00007	<0.00007	0.000089 J	<0.00005	0.00026	<0.00005	<0.0000377	<0.00004	<0.0000377	<0.0000388	<0.000035
Pyrene	73	220	0.00057	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.000104	<0.00011	0.00046 J	<0.000107	<0.000019
<b>Metals</b>														
Arsenic	1	1	0.000877 J											0.00116 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-36B 03/21/2018	MW-36B 05/31/2018	MW-36B 01/14/2019	MW-36B 07/16/2019	MW-36B 01/09/2020	MW-49B 02/04/2009	MW-49B 01/20/2010	MW-49B 06/24/2010	MW-49B 01/20/2011	MW-49B 01/20/2011 Duplicate	MW-49B 07/22/2011	MW-49B 02/07/2012
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Benzene	0.5	0.5	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0095	0.013	0.1	0.0057	0.0056	0.056	0.0056
Chlorobenzene	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.001	<0.001
Ethylbenzene	70	70	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.0081	0.024	0.019	0.004 J	0.0046 J	0.0091	0.0042 J
Methylene chloride	0.5	0.5	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013
Toluene	100	100	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.016	0.045	0.071	0.0072	0.0072	0.038	0.0057
Vinyl chloride	0.2	0.2	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002						<0.001	<0.001
Xylenes (total)	1000	1000	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	0.024	0.07	0.047	0.0066 J	0.0073 J	0.02	0.008 J
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	<0.0001	<0.0005	<0.0001	<0.0001	<0.00005	<0.00005
2,4-Dimethylphenol	49	150	<0.00004	<0.00004	<0.00004	0.000078 J	<0.000054	0.031	0.013	1.2	0.18	0.16	0.59	0.19
2,4-Dinitrotoluene	0.13	0.3	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058	<0.00009	<0.00009	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005
2,6-Dinitrotoluene	0.13	0.3	<0.000042	0.0073	<0.000042	<0.000042	<0.000042	<0.00007	<0.00007	<0.00035	<0.00007	<0.00007	<0.00006	<0.00006
2-Chloronaphthalene	200	580	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.00012	<0.0001	<0.0005	<0.0001	<0.0001	<0.00005	<0.00005
2-Methylnaphthalene	9.8	29	<0.000019	<0.000019	<0.000019	0.00058	<0.00013	0.14	<0.00007	0.0016	<0.00007	<0.00007	0.0029	0.0095
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.0004	<0.00008	<0.00008	<0.00008	<0.00008
4-Nitrophenol	4.9	15	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.00007	<0.00007	<0.00035	<0.00007	<0.00007	<0.00005	<0.00005
Acenaphthene	150	440	<0.000027	0.00014	<0.000027	0.00014	<0.00017	0.094	0.017	0.014	0.00067	0.00049	0.0051	0.034
Acenaphthylene	150	440	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	0.0016	0.0007	0.00063 J	<0.00007	0.000072 J	0.00019 J	0.0007
Anthracene	730	2200	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	0.019	0.00015 J	<0.00035	0.00031	0.000083 J	0.00093	0.0029
Benzo(a)anthracene	0.91	2	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	0.00035	<0.00007	<0.00035	<0.00007	<0.00007	0.00018 J	<0.00005
Benzo(a)pyrene	0.02	0.02	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00008	<0.00008	<0.0004	<0.00008	<0.00008	0.000057 J	<0.00005
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00009	<0.00009	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.000037	0.00017 J	0.00022	<0.000088	<0.000037	0.00029	<0.00053	<0.001	<0.00055	<0.0015	<0.00024	0.00069
Chrysene	91	200	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	0.00038	<0.00007	<0.00035	<0.00007	<0.00007	0.00016 J	<0.00005
Dibenzofuran	9.8	29	<0.00002	<0.00002	<0.00002	0.000087 J	<0.00012	0.071	0.0024	0.0026	0.00018 J	0.00012 J	0.0018	0.019
Di-n-butylphthalate (DBP)	240	730	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	0.0013	0.000083 J	<0.00035	<0.00007	<0.00007	<0.00005	<0.00005
Fluoranthene	98	290	<0.00001	<0.00001	<0.00001	0.000017 J	0.000051 J	0.014	0.00023	<0.00035	0.00019 J	<0.00007	0.0011	0.0015
Fluorene	98	290	<0.00003	<0.00003	<0.00003	0.000047 J	<0.00012	0.071	0.0036	0.0016	0.00018 J	<0.00007	0.0014	0.019
Naphthalene	49	150	<0.00002	<0.00002	<0.00002	0.006	<0.00093	1.4	0.00044	0.23	<0.00052	<0.0002	0.13	0.047
Nitrobenzene	4.9	15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.00009	<0.00009	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005
N-Nitrosodiphenylamine	19	42	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.00009	<0.00009	<0.00045	<0.00009	<0.00009	<0.00005	<0.00005
Pentachlorophenol	0.1	0.1	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079	<0.00008	<0.00008	<0.0004	<0.00008	<0.00008	<0.00005	<0.00005
Phenanthrene	73	220	<0.000021	<0.000021	<0.000021	<0.000021	<0.0001	0.11	0.00017 J	<0.00035	<0.00021	<0.0003	0.0025	0.0098
Phenol	730	2200	<0.000035	<0.000035	<0.000035	<0.00017	<0.000075	<0.00007	<0.00007	0.0053	0.00044	0.0004	0.00021	<0.00005
Pyrene	73	220	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	0.0074	0.0002	<0.00035	0.00024	0.00017 J	0.00066	0.00083
<b>Metals</b>														
Arsenic	1	1	0.000942 J	0.000817 J	0.00118 J	0.00127 J	0.001 J							

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-49B 07/23/2012	MW-49B 02/07/2013 DNAPL	MW-49B 08/01/2013 DNAPL	MW-49B 01/16/2014	MW-49B 07/16/2014	MW-49B 01/29/2018	MW-49B 03/21/2018	MW-49B 05/25/2018	MW-49B 01/27/2020 DNAPL	MW-54B 03/11/2020	MW-57B 02/15/2012 DNAPL	MW-57B 07/24/2012 DNAPL
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0005	<0.00014	<0.0028	<0.0002	<0.0028	<0.0002	<0.0002	<0.0002	<0.01	<0.0002	<0.01	<0.005
Benzene	0.5	0.5	0.11	0.0631	0.469	0.0691	0.346	0.0073	0.026	0.26	<b>0.6</b>	0.00043 J	<b>1.4</b>	<b>1.5</b>
Chlorobenzene	10	10	<0.0005	<0.00012	0.0103 J	<0.00018	<0.0024	<0.0003	<0.0003	<0.0003	<0.015	<0.0003	<0.01	<0.005
Ethylbenzene	70	70	0.023	0.0182	0.0825	0.0425	0.0847	<0.0003	<0.0003	0.048	0.26	0.0011	0.39	0.42
Methylene chloride	0.5	0.5	<0.001	<0.00015	<0.003	<0.00022	0.0212	<0.001	<0.001	<0.001	<0.05	<0.001	<0.013	0.017 J
Toluene	100	100	0.089	0.0633	0.345	0.091	0.31	0.0058	0.014	0.23	0.72	<0.0002	1.3	1.4
Vinyl chloride	0.2	0.2	<0.0005	<0.00011	<0.0022	<0.00018	<0.0022	<0.0002	<0.0002	<0.0002	<0.01	<0.0002		
Xylenes (total)	1000	1000	0.06	0.0527	0.222	0.112	0.249	0.0095	0.012	0.14	0.71	0.0018	1.2	1.1
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.00005	<0.00524	<0.0105	<0.0259	<0.00208	<0.000021	<0.00021	<0.00021	<0.0013	<0.000021	<0.0005	<0.0005
2,4-Dimethylphenol	49	150	6.3	1.09	21.4	4.96 J	13.6	0.2	0.34	3.4	25	<0.00004	6.3	16
2,4-Dinitrotoluene	0.13	0.3	<0.00005	<0.00619	<0.0124	<0.0307	<0.00245	<0.000058	<0.00058	<0.00058	<0.0035	<0.000059	<0.0005	<0.0005
2,6-Dinitrotoluene	0.13	0.3	<0.00006	<0.00381	<0.00762	<0.0189	<0.00151	0.002	<0.00042	<0.00042	<0.0025	<0.000042	<0.0006	<0.0006
2-Chloronaphthalene	200	580	<0.00005	<0.00381	<0.00762	<0.0189	<0.00151	<0.000021	<0.00021	<0.00021	<0.0013	<0.000021	<0.0005	<0.0005
2-Methylnaphthalene	9.8	29	0.18	0.297	0.223	0.691	0.276	0.00011	0.00084 J	0.072	<b>250</b>	0.00011	0.92	1.6
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00008	<0.0395	<0.079	<0.196	<0.0157	<0.00002	<0.0002	<0.0002	<0.0012	<0.00002	<0.0008	<0.0008
4-Nitrophenol	4.9	15	<0.00005	<0.0267	<0.0533	<0.132	<0.0106	<0.000047	<0.00047	<0.00047	<0.0028	<0.000047	<0.0005	<0.0005
Acenaphthene	150	440	0.14	0.248	0.0964	0.622	0.117	0.071	0.066	0.071	<b>190</b>	0.0093	0.35	0.44
Acenaphthylene	150	440	0.0013	<0.00286	<0.00571	<0.0142	0.00432 J	0.0017	0.003	0.026	2.1	0.000087 J	0.006	0.0087
Anthracene	730	2200	0.056	0.0876	<0.00476	0.221	0.013	0.003	0.0038	0.0073	87	0.00017	0.023	0.05
Benzo(a)anthracene	0.91	2	0.013	0.0228 J	<0.00762	0.0671 J	<0.00151	0.000088 J	<0.0005	<0.0005	<b>23</b>	<0.000051	0.0011 J	0.0012 J
Benzo(a)pyrene	0.02	0.02	0.0038	<0.00381	<0.00762	<0.0189	<0.00151	<0.00002	<0.0002	<0.0002	<b>7.5</b>	<0.00002	<0.0005	<0.0005
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00005	<0.00619	<0.0124	<0.0307	<0.00245	<0.00003	<0.0003	<0.0003	<0.0018	<0.00003	<0.0005	<0.0005
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.00055	<0.0176	<0.0352	<0.0873	<0.00698	0.00011 J	<0.00037	<0.00037	<0.0022	0.000089 J	0.0019 J	<0.001
Chrysene	91	200	0.015	0.0207 J	<0.00762	0.0737 J	<0.00151	0.00012	<0.00021	<0.00021	23	<0.000021	0.00099 J	0.0016 J
Dibenzofuran	9.8	29	0.12	0.2	<0.00762	0.484	0.08	0.0028	0.009	0.03	<b>160</b>	0.000076 J	0.28	0.38
Di-n-butylphthalate (DBP)	240	730	<0.00005	<0.00524	<0.0105	<0.0259	<0.00208	<0.00002	<0.0002	<0.0002	<0.0012	0.000033 J	<0.0005	<0.0005
Fluoranthene	98	290	0.093	0.167	<0.00667	0.415	0.00456 J	0.0027	0.0038	0.0036	<b>170</b>	0.000022 J	0.0081	0.016
Fluorene	98	290	0.13	0.217	0.049	0.464	0.0633	0.0087	0.018	0.031	<b>170</b>	0.00015	0.095	0.23
Naphthalene	49	150	2.3	1.58 J	9.38	6.75	5.57	0.00042	0.0075	2.5	<b>1200</b>	0.0021	24	27
Nitrobenzene	4.9	15	<0.00005	<0.00524	<0.0105	<0.0259	<0.00208	<0.000024	<0.00024	<0.00024	<0.0014	<0.000024	<0.0005	<0.0005
N-Nitrosodiphenylamine	19	42	<0.00005	<0.00476	<0.00952	<0.0236	<0.00189	<0.000025	<0.00025	<0.00025	<0.0015	<0.000025	<0.0005	<0.0005
Pentachlorophenol	0.1	0.1	<0.00005	<0.029	<0.0581	<b>&lt;0.144</b>	<0.0115	<0.000079	<0.00079	<0.00079	<0.0047	<0.00008	<0.0005	<0.0005
Phenanthrene	73	220	0.35	0.466	0.039 J	1.29	0.0458	0.0073	0.011	0.028	<b>500</b>	0.0028	0.16	0.24
Phenol	730	2200	0.0063	<0.0019	<0.00381	0.0445 J	0.0145	<0.000035	<0.00035	<0.00035	<0.0021	<0.000035	0.45	1
Pyrene	73	220	0.062	0.101	<0.0105	0.262	<0.00208	0.0014	0.002	0.0021	<b>96</b>	<0.000019	0.0075	0.011
<b>Metals</b>														
Arsenic	1	1						0.000564 J	0.000746 J	0.00146 J	0.0107	0.00117 J		

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-57B 01/31/2013 DNAPL	MW-57B 07/31/2013 DNAPL	MW-57B 01/15/2014 DNAPL	MW-57B 07/29/2014 DNAPL	MW-57B 01/31/2018	MW-57B 04/01/2018	MW-57B 05/25/2018	MW-57B 07/10/2019	MW-57B 01/08/2020	MW-59B 07/15/2010	MW-59B 01/20/2011	MW-59B 07/18/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0035	<0.014	<0.0002	<0.0007	<0.0002	<0.001	<0.0002	<0.0002	<0.002	<0.0005	<0.0005	<0.001
Benzene	0.5	0.5	<b>0.733</b>	<b>1.49</b>	<b>0.716</b>	<b>1.25</b>	0.012	0.01	<b>0.82</b>	<b>0.84</b>	<b>0.52</b>	<0.0005	<0.0005	<0.001
Chlorobenzene	10	10	<0.003	<0.012	<0.00018	<0.0006	<0.0003	<0.0015	<0.0003	<0.0003	<0.003	<0.0005	<0.0005	<0.001
Ethylbenzene	70	70	0.193	0.501	0.174	0.371	0.026	0.032	0.3	0.39	0.25	<0.0005	<0.0005	<0.0011
Methylene chloride	0.5	0.5	<0.00375	0.0405 J	<0.00022	<0.00075	<0.001	<0.005	<0.001	<0.001	<0.01	<0.0005	<0.0005	<0.0013
Toluene	100	100	0.692	1.62	0.63	1.33	0.0043	0.0019 J	0.84	0.93	0.64	<0.0005	<0.0005	<0.001
Vinyl chloride	0.2	0.2				0.00299 J				0.00048 J				<0.001
Xylenes (total)	1000	1000	0.589	1.4	0.574	1.16	0.057	0.055	0.84	1.2	0.7	<0.001	<0.001	<0.0031
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.055	<0.0267	<0.0519	<0.0109	<0.00021	<0.000021	<0.00021	<0.00021	<0.00021	<0.0001	<0.0001	<0.00005
2,4-Dimethylphenol	49	150	13.8	9.67	19.8 J	15	0.01	<0.00004	3.7	6.5	2.8	<0.00008	<0.00008	<0.00005
2,4-Dinitrotoluene	0.13	0.3	<0.065	<0.0316	<0.0613	<0.0129	<0.00058	<0.000058	<0.00058	<0.00058	<0.00058	<0.00009	<0.00009	<0.00005
2,6-Dinitrotoluene	0.13	0.3	<0.04	<0.0194	<0.0377	<0.00792	<0.00042	<0.000042	<0.00042	<0.00042	<0.00042	<0.00007	<0.00007	<0.00006
2-Chloronaphthalene	200	580	<0.04	<0.0194	<0.0377	<0.00792	<0.00021	<0.000021	<0.00021	<0.00021	<0.00021	<0.0001	<0.0001	<0.00005
2-Methylnaphthalene	9.8	29	1.75	1.07	0.892	0.945	0.17	0.029	0.61	0.93	1.1	<0.00007	<0.00007	<0.00005
4,6-Dinitro-2-methylphenol	0.24	0.73	<b>&lt;0.415</b>	<0.201	<b>&lt;0.392</b>	<0.0822	<0.0002	<0.00002	<0.0002	<0.0002	<0.0002	<0.00008	<0.00008	<0.00008
4-Nitrophenol	4.9	15	<0.28	<0.136	<0.264	<0.0554	<0.00047	<0.000047	<0.00047	<0.00047	<0.00047	<0.00007	<0.00007	<0.00005
Acenaphthene	150	440	0.93	0.423	0.524	0.267	0.13	0.13	0.3	0.38	0.55	<0.00009	<0.00009	<0.00005
Acenaphthylene	150	440	<0.03	<0.0146	<0.0283	<0.00594	0.0029	0.0011	0.0063	0.0051	0.0086	<0.00007	<0.00007	<0.00005
Anthracene	730	2200	0.292	0.0493 J	0.0844 J	0.0355 J	0.014	0.005	0.44	0.039	0.19	<0.00007	<0.00007	<0.00005
Benzo(a)anthracene	0.91	2	0.0543 J	<0.0194	<0.0377	<0.00792	0.001	<0.00005	0.03	0.0071	0.04	<0.00007	<0.00007	<0.00005
Benzo(a)pyrene	0.02	0.02	<b>&lt;0.04</b>	<0.0194	<b>&lt;0.0377</b>	<0.00792	0.00058 J	<0.00002	0.0094	0.0021	0.012	<0.00008	<0.00008	<0.00005
bis(2-Chloroethoxy)methane	0.083	0.19	<0.065	<0.0316	<0.0613	<0.0129	<0.0003	<0.00003	<0.0003	<0.0003	<0.0003	<0.00009	<0.00009	<0.00005
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.185	<0.0898	<0.175	<0.0366	0.00043 J	0.000095 J	<0.00037	<0.00037	<0.00037	0.002	<0.00021	<0.00031
Chrysene	91	200	0.0561 J	<0.0194	<0.0377	<0.00792	0.00091 J	<0.000021	0.024	0.0065	0.042	<0.00007	<0.00007	<0.00005
Dibenzofuran	9.8	29	0.814	0.322	0.392	0.226	0.11	0.081	0.29	0.32	0.49	<0.00008	<0.00008	<0.00005
Di-n-butylphthalate (DBP)	240	730	<0.055	<0.0267	<0.0519	<0.0109	<0.0002	<0.00002	<0.0002	<0.0002	<0.0002	<0.00007	<0.00007	<0.00005
Fluoranthene	98	290	0.387	0.0301 J	0.0752 J	0.0109 J	0.012	0.0044	0.19	0.047	0.34	<0.00007	<0.00007	<0.00005
Fluorene	98	290	0.65	0.208	0.298	0.138	0.081	0.096	0.27	0.21	0.46	<0.00007	<0.00007	<0.00005
Naphthalene	49	150	18.9 J	18.1	10.6	17	1.8	0.39	12	17	13	0.00014 J	<0.0001	<0.00005
Nitrobenzene	4.9	15	<0.055	<0.0267	<0.0519	<0.0109	<0.00024	<0.000024	<0.00024	<0.00024	<0.00024	<0.00009	<0.00009	<0.00005
N-Nitrosodiphenylamine	19	42	<0.05	<0.0243	<0.0472	<0.0099	<0.00025	<0.000025	<0.00025	<0.00025	<0.00025	<0.00009	<0.00009	<0.00005
Pentachlorophenol	0.1	0.1	<b>&lt;0.305</b>	<b>&lt;0.148</b>	<b>&lt;0.288</b>	<0.0604	<0.00079	<0.000079	<0.00079	<0.00079	<0.00079	<0.00008	<0.00008	<0.00005
Phenanthrene	73	220	1.39	0.242	0.456	0.127	0.094	0.033	0.63	0.29	1.1	<0.00007	<0.00007	<0.00005
Phenol	730	2200	1	0.645	1	0.495	<0.00035	<0.000035	0.22	1	0.32	0.0002	<0.00007	<0.00005
Pyrene	73	220	0.245 J	<0.0267	<0.0519	<0.0109	0.0068	0.003	0.13	0.03	0.21	<0.00007	<0.00007	<0.00005
<b>Metals</b>														
Arsenic	1	1					0.0419	0.00179 J	0.00285	0.00404	0.00404			

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-59B 02/06/2012	MW-59B 07/27/2012	MW-59B 01/31/2013	MW-59B 08/01/2013	MW-59B 01/16/2014	MW-59B 07/30/2014	MW-59B 01/29/2018	MW-59B 03/20/2018	MW-59B 05/25/2018	MW-59B 01/23/2019	MW-59B 07/17/2019	MW-59B 01/16/2020
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.5	0.5	<0.001	<0.0005	<0.00008	0.0000981 J	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	10	10	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	70	70	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.5	0.5	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	100	100	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.18
Vinyl chloride	0.2	0.2	<0.001	<0.0005	<0.00011	<0.00011	<0.00018	<0.00011	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Xylenes (total)	1000	1000	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	R	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	49	150	0.25	<0.00005	<0.000292	<0.000295	<0.000292	R	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.13	0.3	<0.00005	<0.00005	<0.000123	<0.000124	<0.000123	R	<0.000058	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	0.13	0.3	<0.00006	<0.00006	<0.0000755	<0.0000762	<0.0000755	R	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	200	580	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	R	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	9.8	29	<0.00005	<0.00005	<0.000066	<0.000067	<0.000066	R	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019	<0.000038
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00008	<0.00008	<0.0000783	<0.000079	<0.0000783	R	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	4.9	15	<0.00005	<0.00005	<0.000528	<0.000533	<0.000528	R	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	150	440	0.0017	<0.00005	<0.0000755	<0.0000762	<0.0000755	0.000621U	<0.000027	<0.000027	<0.000027	<0.000027	<0.000027	<0.000051
Acenaphthylene	150	440	0.00014 J	<0.00005	<0.0000566	<0.0000571	<0.0000566	R	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	730	2200	0.000054 J	<0.00005	<0.0000472	<0.0000476	<0.0000472	R	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014	<0.000031
Benzo(a)anthracene	0.91	2	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	R	<0.00005	<0.000051	<0.000051	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.02	0.02	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	R	<0.00002	<0.00002	<0.00002	0.000033 J	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00005	<0.00005	<0.000123	<0.000124	<0.000123	R	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	0.00068	0.00018 J	<0.000349	<0.000352	<0.000349	R	0.000058 J	<0.000037	<0.000037	<0.000056	<0.000039	<0.000037
Chrysene	91	200	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	R	<0.000021	<0.000021	<0.000021	0.000036 J	<0.000021	<0.000021
Dibenzofuran	9.8	29	<0.00005	<0.00005	<0.0000755	<0.0000762	<0.0000755	0.000201U	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00005
Di-n-butylphthalate (DBP)	240	730	<0.00005	<0.00005	<0.000104	0.000115 J	0.000105 J	R	<0.00002	<0.00002	<0.00002	<0.00002	0.000055 J	<0.00002
Fluoranthene	98	290	<0.00005	<0.00005	<0.000066	<0.000067	<0.000066	R	<0.00001	<0.00001	<0.00001	0.000051 J	<0.00001	<0.000025
Fluorene	98	290	<0.00005	<0.00005	<0.000066	<0.000067	<0.000066	0.000189U	<0.00003	<0.00003	<0.00003	<0.00003	<0.00003	<0.000042
Naphthalene	49	150	0.00012 J	0.00006 J	<0.000269	<0.000166	<0.000172	0.00627U	<0.00002	<0.00002	<0.00002	0.000072 J	<0.00015	<0.00024
Nitrobenzene	4.9	15	<0.00005	<0.00005	<0.000104	<0.000105	<0.000104	R	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
N-Nitrosodiphenylamine	19	42	<0.00005	<0.00005	<0.0000943	<0.0000952	<0.0000943	R	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.1	0.1	0.00011 J	<0.00005	<0.000575	<0.000581	<0.000575	R	<0.000079	<0.00008	<0.00008	<0.000079	<0.000079	<0.000079
Phenanthrene	73	220	0.00025	<0.00005	<0.0000566	<0.0000571	<0.000112	R	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000099
Phenol	730	2200	0.00033	<0.00005	<0.0000377	<0.0000381	<0.0000377	R	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	73	220	0.000062 J	<0.00005	<0.000104	<0.000105	<0.000104	R	<0.000019	<0.000019	<0.000019	0.000053 J	<0.000019	<0.000021
<b>Metals</b>														
Arsenic	1	1							<0.0004	<0.0004	<0.0004	0.000983 J	0.000542 J	0.000486 J

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray.
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/ Assessment Level</b>	MW60B 03/17/2020	MW61B 03/17/2020	MW-63B 01/13/2010	MW-63B 06/30/2010	MW-63B 01/27/2011	MW-63B 07/19/2011	MW-63B 07/19/2011 Duplicate	MW-63B 02/09/2012	MW-63B 02/09/2012 Duplicate	MW-63B 07/18/2012	MW-63B 07/18/2012 Duplicate	MW-63B 02/07/2013
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0002	<0.0002	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.00014
Benzene	0.5	0.5	<0.0002	<0.0002	0.21	0.015	0.019	0.019	0.024	<0.001	<0.001	0.0015 J	0.0016 J	0.00952
Chlorobenzene	10	10	<0.0003	<0.0003	<0.0005	<0.0005	<0.0005	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005	<0.00012
Ethylbenzene	70	70	<0.0003	<0.0003	0.2	0.072	0.071	0.04	0.045	0.0012 J	0.0012 J	0.0014 J	0.0014 J	0.0165
Methylene chloride	0.5	0.5	<0.001	<0.001	<0.0005	<0.0005	<0.0005	<0.0013	<0.0013	<0.0013	<0.0013	<0.001	<0.001	<0.00015
Toluene	100	100	<0.0002	<0.0002	0.015	0.0016 J	0.0018 J	0.0017 J	0.0019 J	<0.001	<0.001	0.0038 J	0.0038 J	0.00241
Vinyl chloride	0.2	0.2												
Xylenes (total)	1000	1000	<0.0003	<0.0003	0.082	0.02	0.016	0.013 J	0.014 J	<0.0031	<0.0031	<0.0015	<0.0015	0.00629
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00524
2,4-Dimethylphenol	49	150	<0.00004	<0.00004	<0.00008	<0.00008	<0.00008	0.000056 J	<0.00005	0.00005 J	0.00044 J	<0.00005	<0.00005	<0.0148
2,4-Dinitrotoluene	0.13	0.3	<0.000058	<0.000058	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00619
2,6-Dinitrotoluene	0.13	0.3	<0.000042	<0.000042	<0.00007	<0.00007	<0.00007	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.00006	<0.00381
2-Chloronaphthalene	200	580	<0.000021	<0.000021	<0.0001	<0.0001	<0.0001	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00381
2-Methylnaphthalene	9.8	29	<0.000019	<0.000019	0.11	0.031	0.025	0.014	0.013	0.0029	0.0029	0.0034	0.0032	0.0104 J
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00002	<0.00002	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.00008	<0.0395
4-Nitrophenol	4.9	15	<0.000047	<0.000047	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.0267
Acenaphthene	150	440	<0.000027	<0.000027	0.028	0.013	0.017	0.0053	0.0066	0.002	0.0019	0.0023	0.0022	0.00952 J
Acenaphthylene	150	440	<0.000015	<0.000015	0.00051	0.00018 J	<0.00007	0.000066 J	0.00011 J	0.00012 J	0.00016 J	<0.00005	<0.00005	<0.00286
Anthracene	730	2200	<0.000014	<0.000014	0.00068	0.00039	0.0011	0.00011 J	0.0002 J	0.00015 J	0.00015 J	<0.00013	<0.00011	0.00238 J
Benzo(a)anthracene	0.91	2	<0.00005	<0.00005	<0.00007	<0.00007	0.00087	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00381
Benzo(a)pyrene	0.02	0.02	<0.00002	<0.00002	<0.00008	<0.00008	0.00027	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00381
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00003	<0.00003	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00619
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	0.00092	0.00015 J	<0.00036	<0.00036	0.0006	<0.00051	<0.00029	0.00096	0.00073	0.00096 J	0.00051 J	<0.0176
Chrysene	91	200	<0.000021	<0.000021	<0.00007	<0.00007	0.00079	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00381
Dibenzofuran	9.8	29	<0.00002	<0.00002	0.022	0.008	0.013	0.0041	0.0049	0.0024	0.0019	0.0026	0.0023	0.00576 J
Di-n-butylphthalate (DBP)	240	730	0.000047 J	<0.00002	0.00019 J	<0.00007	<0.00007	<0.00005	<0.00005	0.00014 J	0.00005 J	<0.00005	<0.00005	0.00524 J
Fluoranthene	98	290	0.000015 J	<0.00001	<0.00007	<0.00007	0.0042	<0.00005	<0.00005	0.000091 J	0.00005 J	<0.00013	<0.00013	<0.00333
Fluorene	98	290	<0.00003	<0.00003	0.0078	0.0041	0.0054	0.0019	0.0024	0.00093	0.00078	0.0011	0.001	<0.00333
Naphthalene	49	150	<0.00002	<0.00002	3.1	0.67	0.76	0.36 J	0.49 J	0.027	0.026	0.044	0.04	0.251
Nitrobenzene	4.9	15	<0.000024	<0.000024	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00524
N-Nitrosodiphenylamine	19	42	<0.000025	<0.000025	<0.00009	<0.00009	<0.00009	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.00476
Pentachlorophenol	0.1	0.1	<0.000079	<0.000079	<0.00008	<0.00008	<0.00008	<0.00005	<0.00005	0.00018 J	0.00005 J	<0.00005	<0.00005	<0.029
Phenanthrene	73	220	<0.000021	<0.000021	0.0034	0.00076	0.0044	0.00075	0.00098	0.00072 J	0.0005 J	<0.001	<0.001	0.00286 J
Phenol	730	2200	<0.000035	<0.000035	<0.00007	<0.00007	<0.00007	<0.00005	<0.00005	0.00057 J	0.00095 J	<0.00005	<0.00005	<0.0019
Pyrene	73	220	<0.000019	<0.000019	<0.00007	<0.00007	0.0029	<0.00005	<0.00005	0.000063 J	<0.00005	<0.00005	<0.00005	<0.00524
<b>Metals</b>														
Arsenic	1	1	0.00234	0.00461										

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-63B 02/07/2013	MW-63B 08/07/2013	MW-63B 01/22/2014	MW-63B 07/24/2014	MW-63B 01/28/2018	MW-63B 03/26/2018	MW-63B 06/06/2018	MW-63B 01/14/2019	MW-63B 07/16/2019	MW-63B 01/16/2020	MW-67B 07/15/2010	MW-67B 01/27/2011
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0005	<0.0005
Benzene	0.5	0.5	0.00919	0.0869	0.0762	0.108	0.0033	0.0026	0.048	0.35	0.11	0.018	<0.0005	<0.0005
Chlorobenzene	10	10	<0.00012	<0.00012	<0.00018	0.000216 J	<0.0003	<0.0003	<0.0003	0.00073 J	<0.0003	<0.0003	<0.0005	<0.0005
Ethylbenzene	70	70	0.0163	0.0341	0.0418	0.151	0.012	0.0059	0.048	0.48	0.14	0.051	0.0015 J	<0.0005
Methylene chloride	0.5	0.5	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0005	<0.0005
Toluene	100	100	0.00231	<0.000434	0.000399 J	0.00257	<0.0002	<0.0002	0.00093 J	0.0071	0.00071 J	<0.0002	<0.0005	<0.0005
Vinyl chloride	0.2	0.2				<0.00011								
Xylenes (total)	1000	1000	0.00635	0.0113	0.0156	0.0535	0.0048	0.0016	0.011	0.11	0.041	0.013	0.0012 J	<0.001
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.000105	<0.000104	<0.00109	<0.0011	<0.000021	<0.000021	<0.00021	<0.00021	<0.000021	<0.000021	<0.0001	<0.0001
2,4-Dimethylphenol	49	150	<0.000295	<0.000292	<0.00307	<0.0031	<0.00004	<0.00004	<0.0004	<0.0004	<0.00004	<0.00004	<0.00008	<0.00008
2,4-Dinitrotoluene	0.13	0.3	<0.000124	<0.000123	<0.00129	<0.0013	<0.000058	<0.000058	<0.00058	<0.00059	<0.000058	<0.000058	<0.00009	<0.00009
2,6-Dinitrotoluene	0.13	0.3	<0.0000762	<0.0000755	<0.000792	<0.0008	<0.000042	<0.000042	<0.00042	<0.00042	<0.000042	<0.000042	<0.00007	<0.00007
2-Chloronaphthalene	200	580	<0.0000762	<0.0000755	<0.000792	<0.0008	<0.000021	<0.000021	<0.00021	<0.00021	<0.000021	<0.000021	<0.0001	<0.0001
2-Methylnaphthalene	9.8	29	0.00865	0.00242	0.00756	0.0302	0.000059 J	<0.000019	0.0016	0.042	0.037	<0.000056	<0.00007	0.00007 J
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00079	<0.000783	<0.00822	<0.0083	<0.00002	<0.00002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00008	<0.00008
4-Nitrophenol	4.9	15	<0.000533	<0.000528	<0.00554	<0.0056	<0.000047	<0.000047	<0.00047	<0.00047	<0.000047	<0.000047	<0.00007	<0.00007
Acenaphthene	150	440	0.0065	0.000833	0.00274 J	0.00754	0.00066	<0.000027	0.0027	<0.00027	0.012	<0.000059	0.00011 J	<0.00009
Acenaphthylene	150	440	<0.0000571	<0.0000566	<0.000594	<0.0006	<0.000015	<0.000015	<0.00015	0.0029	0.00016	0.000032 J	<0.00007	0.00007 J
Anthracene	730	2200	0.000124 J	<0.0000472	<0.000495	<0.0005	0.00011	0.00022	<0.00014	0.00017 J	0.00034	<0.000081	<0.00007	<0.00007
Benzo(a)anthracene	0.91	2	<0.0000762	<0.0000755	<0.000792	<0.0008	<0.00005	<0.00005	<0.0005	<0.00051	<0.00005	<0.00005	<0.00007	<0.00007
Benzo(a)pyrene	0.02	0.02	<0.0000762	<0.0000755	<0.000792	<0.0008	<0.00002	<0.00002	<0.0002	<0.0002	<0.00002	<0.00002	<0.00008	<0.00008
bis(2-Chloroethoxy)methane	0.083	0.19	<0.000124	<0.000123	<0.00129	<0.0013	<0.00003	<0.00003	<0.0003	<0.0003	<0.00003	<0.00003	<0.00009	0.00009 J
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.000376	0.000381 J	<0.00366	<0.0037	<0.000037	<0.000037	<0.00037	<0.00037	<0.000037	0.000065 J	0.0016	0.0022 J
Chrysene	91	200	<0.0000762	<0.0000755	<0.000792	<0.0008	<0.000021	<0.000021	<0.00021	<0.00021	<0.000021	<0.000021	<0.00007	<0.00007
Dibenzofuran	9.8	29	0.00514	0.00104	0.002 J	0.00663	0.00065	<0.00002	0.002	0.0087	0.011	<0.000063	<0.00008	0.00008 J
Di-n-butylphthalate (DBP)	240	730	0.000127 J	<0.000104	<0.00109	<0.0011	<0.00002	<0.00002	<0.0002	<0.0002	<0.00002	0.000034 J	<0.00007	0.000083 J
Fluoranthene	98	290	<0.0000667	<0.000066	<0.000693	<0.0007	0.0001	<0.00001	<0.0001	<0.0001	<0.00001	<0.000023	<0.00007	<0.00007
Fluorene	98	290	0.00208	0.000349 J	0.00102 J	0.00248 J	0.00033	<0.00003	0.00099 J	0.0029	0.0037	<0.00012	<0.00007	<0.00007
Naphthalene	49	150	0.162 J	0.146	0.374	1.69	<0.00066	<0.00002	0.29	2.1	1.5	<0.00035	<0.0001	0.00062
Nitrobenzene	4.9	15	<0.000105	<0.000104	<0.00109	<0.0011	<0.000024	<0.000024	<0.00024	<0.00024	<0.000024	<0.000024	<0.00009	0.00009 J
N-Nitrosodiphenylamine	19	42	<0.0000952	<0.0000943	<0.00099	<0.001	<0.000025	<0.000025	<0.00025	<0.00025	<0.000025	<0.000025	<0.00009	<0.00009
Pentachlorophenol	0.1	0.1	<0.000581	<0.000575	<0.00604	<0.0061	<0.000079	<0.000079	<0.00079	<0.0008	<0.000079	<0.000079	<0.00008	<0.00008
Phenanthrene	73	220	0.000776 J	<0.0000566	0.00133 J	<0.0006	<0.000075	<0.000021	<0.00021	0.00094 J	0.0013	<0.000072	<0.00007	<0.00007
Phenol	730	2200	<0.0000381	<0.0000377	0.000889 J	<0.0004	<0.000035	<0.000035	<0.00035	<0.00035	<0.000035	<0.000035	<0.00007	<0.00007
Pyrene	73	220	<0.000105	<0.000104	<0.00109	<0.0011	<0.000019	<0.000019	<0.00019	<0.00019	<0.000019	<0.000019	<0.00007	<0.00007
<b>Metals</b>														
Arsenic	1	1					0.00114 J	0.00211	0.000818 J	0.00338	0.00156 J	0.00204		

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.



**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-67B 07/20/2011	MW-67B 02/09/2012	MW-67B 07/17/2012	MW-67B 02/12/2013	MW-67B 08/08/2013	MW-67B 01/23/2014	MW-67B 07/24/2014	MW-67B 01/31/2018	MW-67B 03/27/2018	MW-67B 06/06/2018	MW-67B 01/24/2019	MW-67B 07/31/2019
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.001	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.5	0.5	<0.001	<0.001	<0.0005	<0.00008	<0.00008	<0.0002	<0.00008	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	10	10	<0.001	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	70	70	<0.0011	<0.0011	<0.0005	<0.00011	<0.00011	<0.00019	<0.00011	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.5	0.5	<0.0013	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	100	100	<0.001	<0.001	<0.0005	<0.00015	<0.00015	<0.00017	<0.00015	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.2	0.2					<0.00011	<0.00011	<0.00011					
Xylenes (total)	1000	1000	<0.0031	<0.0031	<0.0015	<0.00026	<0.00026	<0.00058	<0.00026	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000106	<0.000106	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	49	150	<0.00005	0.0005 J	<0.00005	<0.000295	<0.000292	<0.000298	<0.000298	<0.00004	<0.00004	<0.00004	<0.00004	<0.00004
2,4-Dinitrotoluene	0.13	0.3	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000125	<0.000125	<0.000058	<0.000058	0.000058 J	<0.000058	<0.000058
2,6-Dinitrotoluene	0.13	0.3	<0.00006	<0.00006	0.0022 J	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	200	580	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.000021	<0.000021	0.000021 J	<0.000021	<0.000021
2-Methylnaphthalene	9.8	29	<0.00005	0.00023 J	0.000062 J	<0.0000667	<0.000066	<0.0000673	<0.0000673	<0.000031	<0.000019	0.000061 J	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00008	<0.00008	<0.00008	<b>R</b>	<0.000783	<0.000798	<0.000798	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
4-Nitrophenol	4.9	15	<0.00005	<0.00005	<0.00005	<0.000533	<0.000528	<0.000538	<0.000538	<0.000047	<0.000047	0.000047 J	<0.000047	<0.000047
Acenaphthene	150	440	<0.00005	0.00012 J	<0.00005	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.000027	<0.000027	0.000027 J	<0.000027	<0.000027
Acenaphthylene	150	440	<0.00005	<0.00005	<0.00005	<0.0000571	<0.0000566	<0.0000577	<0.0000577	<0.000015	<0.000015	0.000015 J	<0.000015	<0.000015
Anthracene	730	2200	<0.00005	<0.00005	<0.00005	<0.0000476	<0.0000472	<0.0000424	<0.0000481	<0.000014	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	0.91	2	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Benzo(a)pyrene	0.02	0.02	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00005	<0.00005	<0.00005	<0.000124	<0.000123	<0.000125	<0.000125	<0.00003	<0.00003	0.00003 J	<0.00003	<0.00003
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.00094	0.00042	0.00012 J	<0.000352	<0.000349	<0.000356	0.00184	<0.000056	<0.000037	<0.000056	0.000051 J	<0.000037
Chrysene	91	200	<0.00005	<0.00005	<0.00005	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Dibenzofuran	9.8	29	<0.00005	0.00013 J	<0.00005	<0.0000762	<0.0000755	<0.0000769	<0.0000769	<0.00002	<0.00002	0.000069 J	<0.00002	<0.00002
Di-n-butylphthalate (DBP)	240	730	<0.00005	<0.00005	<0.00005	<0.000105	0.000119 J	<0.000834	<0.000106	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002
Fluoranthene	98	290	<0.00005	<0.00005	<0.00005	<0.0000667	<0.000066	<0.0000673	<0.0000673	<0.00001	<0.00001	<0.00001	<0.00001	<0.00001
Fluorene	98	290	<0.00005	0.0001 J	<0.00005	<0.0000667	<0.000066	<0.0000673	<0.0000673	<0.00003	<0.00003	0.00003 J	<0.00003	<0.00003
Naphthalene	49	150	<0.00005	0.0019	0.00049	<0.0000762	0.000433 J	<0.000711	0.000275 J	<0.00045	0.00013	0.00047 J	<0.00002	0.000079 J
Nitrobenzene	4.9	15	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000106	<0.000106	<0.000024	<0.000024	0.000024 J	<0.000024	<0.000024
N-Nitrosodiphenylamine	19	42	<0.00005	<0.00005	<0.00005	<0.0000952	<0.0000943	<0.0000962	<0.0000962	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Pentachlorophenol	0.1	0.1	<0.00005	<0.00005	<0.00005	0.000581 J	<0.000575	<0.000587	<0.000587	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	73	220	<0.00005	0.00011 J	<0.00005	<0.0000571	<0.0000566	<0.0000377	<0.0000577	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	730	2200	<0.00005	0.000089 J	<0.00005	<0.0000381	<0.0000377	<0.0000385	<0.0000385	0.00018 J	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	73	220	<0.00005	<0.00005	<0.00005	<0.000105	<0.000104	<0.000106	<0.000106	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>														
Arsenic	1	1								0.000751 J	0.000565 J	0.000416 J	<0.0004	0.000494 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3**  
**ANALYTICAL RESULTS SUMMARY (2008 - 2020)**  
**CLASS 3 GROUNDWATER**  
**HOUSTON, TX - WOOD PRESERVING WORKS**

	<i>Class 3 Residential Assessment</i>	<i>Class 3 C/I Assessment Level</i>	MW-67B 01/15/2020	MW-70B 07/17/2012 DNAPL	MW-70B 02/07/2013 DNAPL	MW-70B 01/22/2014 DNAPL	MW-70B 07/28/2014 DNAPL	MW-70B 01/20/2020 DNAPL	MW-71B 02/08/2012	MW-71B 07/18/2012	MW-71B 02/07/2013	MW-71B 08/07/2013	MW-71B 01/24/2014	MW-71B 07/28/2014
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0002	<0.0025	<0.007	<0.0002	<0.0014	<0.01	<0.001	<0.0005	<0.00014	<0.00014	<0.0002	<0.00014
Benzene	0.5	0.5	<0.0002	0.21	<b>2.01</b>	<b>2.39</b>	<b>2.55</b>	<b>1.9</b>	0.012	0.0014 J	0.0124	0.103	0.039	0.00155
Chlorobenzene	10	10	<0.0003	<0.0025	0.0317 J	0.000715	<0.0012	<0.015	<0.001	<0.0005	<0.00012	<0.00012	<0.00018	<0.00012
Ethylbenzene	70	70	<0.0003	0.058	0.524	0.621	0.742	0.57	0.0045 J	0.0075	0.00541	0.0354	0.00793	<0.00011
Methylene chloride	0.5	0.5	<0.001	<0.005	<0.0075	<0.00022	<0.0015	<0.05	<0.0013	<0.001	<0.00015	<0.00015	<0.00022	<0.00015
Toluene	100	100	<0.0002	0.22	1.65	2.31	2.76	2.1	0.0077	0.0078	0.0104	0.0355	0.00918	0.00423
Vinyl chloride	0.2	0.2						<0.01				<0.00011		
Xylenes (total)	1000	1000	<0.0003	0.19	1.51	1.68	2.11	1.5	0.016	0.033	0.0143	0.0615	0.0202	0.0126
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.000021	<0.0005	<0.075	<0.0157	<0.0107	<0.00042	<0.00005	<0.00005	<0.000105	<0.00519	<0.000106	<0.000107
2,4-Dimethylphenol	49	150	<0.00004	2.6	<2.11	<b>72</b>	<b>50.8</b>	49	0.0034	<0.00005	<0.000295	<0.0146	0.0225	<0.000301
2,4-Dinitrotoluene	0.13	0.3	<0.000058	<0.0005	<0.0886	<0.0186	<0.0126	<0.0012	<0.00005	<0.00005	<0.000124	<0.00613	<0.000125	<0.000126
2,6-Dinitrotoluene	0.13	0.3	<0.000042	<0.0006	<0.0545	<0.0114	<0.00777	<0.00084	<0.00006	<0.00006	<0.0000762	<0.00377	<0.0000769	<0.0000777
2-Chloronaphthalene	200	580	<0.000021	<0.0005	<0.0545	<0.0114	<0.00777	<0.00042	<0.00005	<0.00005	<0.0000762	<0.00377	<0.0000769	<0.0000777
2-Methylnaphthalene	9.8	29	<0.000019	0.94	1.21	1.4	1.31	<b>10</b>	0.0076	<0.0004	0.000377 J	0.114	0.0476	<0.000068
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00002	<0.0008	<b>&lt;0.566</b>	<0.119	<0.0806	<0.0004	<0.00008	<0.00008	<0.00079	<0.0392	<0.000798	<0.000806
4-Nitrophenol	4.9	15	<0.000047	<0.0005	<0.382	<0.08	<0.0544	<0.00094	<0.00005	<0.00005	<0.000533	<0.0264	<0.000538	<0.000544
Acenaphthene	150	440	<0.000027	0.91	0.515	0.454	0.374	7.1	0.0039	<0.00017	0.0044	0.0346	0.0212	0.000785
Acenaphthylene	150	440	<0.000015	0.011	0.0424 J	<0.00857	0.0114 J	0.079	0.00019 J	<0.00005	0.000135 J	<0.00283	0.00122	<0.0000583
Anthracene	730	2200	<0.000014	0.096	0.051 J	0.0423 J	0.0387 J	5.8	0.00056	<0.00013	0.000452 J	0.00383 J	0.00198	<0.0000485
Benzo(a)anthracene	0.91	2	<0.00005	0.016	<0.0545	<0.0114	<0.00777	0.52	0.000081 J	0.00011 J	<0.0000762	<0.00377	<0.0000769	<0.0000777
Benzo(a)pyrene	0.02	0.02	<0.00002	0.0041	<b>&lt;0.0545</b>	<0.0114	<0.00777	<b>0.13</b>	0.00012 J	0.00014 J	<0.0000762	<0.00377	<0.0000769	<0.0000777
bis(2-Chloroethoxy)methane	0.083	0.19	<0.00003	<0.0005	<b>&lt;0.0886</b>	<0.0186	<0.0126	<0.0006	<0.00005	<0.00005	<0.000124	<0.00613	<0.000125	<0.000126
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	0.00009 J	0.0068	<0.252	<0.0529	<0.0359	0.0039 J	<0.0013	0.00012 J	<0.000352	<0.0175	<0.000356	<0.000359
Chrysene	91	200	<0.000021	0.013	<0.0545	<0.0114	<0.00777	0.63	0.000089 J	0.00015 J	<0.0000762	<0.00377	<0.0000769	<0.0000777
Dibenzofuran	9.8	29	<0.00002	0.69	0.345	0.355	0.278	6.2	0.0031	<0.00016	0.00244	0.0292	0.0175	<0.0000777
Di-n-butylphthalate (DBP)	240	730	<0.00002	<0.0005	<0.075	<0.0157	<0.0107	<0.0004	<0.00005	<0.00005	<0.000105	<0.00519	<0.00085	<0.000107
Fluoranthene	98	290	<0.00001	0.28	<0.0477	0.0105 J	0.013 J	4.7	0.00053	<0.00026	0.000387 J	<0.0033	0.000712	0.000149 J
Fluorene	98	290	<0.00003	0.66	0.211 J	0.217	0.186	6.7	0.002	<0.00023	0.00168	0.0127 J	0.0104	<0.000068
Naphthalene	49	150	<0.00002	5.3	17.3 J	30.1	18.1	<b>71</b>	0.051	<0.0019	0.0000937 J	2.07	0.504	<0.000471
Nitrobenzene	4.9	15	<0.000024	<0.0005	<0.075	<0.0157	<0.0107	<0.00048	<0.00005	<0.00005	<0.000105	<0.00519	<0.000106	<0.000107
N-Nitrosodiphenylamine	19	42	<0.000025	<0.0005	<0.0682	<0.0143	<0.00971	<0.0005	<0.00005	<0.00005	<0.0000952	<0.00472	<0.0000962	<0.0000971
Pentachlorophenol	0.1	0.1	<0.000079	<0.0005	<b>&lt;0.416</b>	<0.0871	<0.0592	<0.0016	0.00022	<0.00005	<0.000581	<0.0288	<0.000587	<0.000592
Phenanthrene	73	220	<0.000021	0.93	0.227 J	0.175	0.162	14	0.0025	<0.00034	0.00127	0.0124 J	0.00677	<0.0000583
Phenol	730	2200	<0.000035	0.077	2.87	3.86	1.69	3.4	0.00037	<0.00005	<0.0000381	<0.00189	<0.0000385	<0.0000388
Pyrene	73	220	<0.000019	0.094	<0.075	<0.0157	<0.0107	3.2	0.00057	<0.00026	0.000253 J	<0.00519	0.000353 J	<0.000107
<b>Metals</b>														
Arsenic	1	1	0.000467 J					0.0012 J						

Notes:

- All values in milligrams per liter (mg/L).
- Concentrations > RAL and non-detects are highlighted light gray.
- Concentrations > C/I AL and non-detects are highlighted dark gray
- TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
- RAL = Residential Assessment Level, C/I = Commercial/Industrial
- J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/ Assessment Level</b>	MW-71B 01/25/2018	MW-71B 03/26/2018	MW-71B 06/06/2018	MW-71B 01/15/2019	MW-71B 07/17/2019	MW-71B 01/15/2020	MW-72B 07/12/2012	MW-72B 02/01/2013	MW-72B 07/29/2013	MW-72B 01/15/2014	MW-72B 02/08/2018	MW-72B 03/19/2018
<b>Volatile Organic Compounds</b>														
1,2-Dichloroethane	0.5	0.5	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.005	<0.007	<0.014	<0.0002	0.18	<0.0002
Benzene	0.5	0.5	<0.0002	0.0042	0.027	0.0024	0.13	0.0021	1.4	1.45	1.23	0.932	0.8	1.1
Chlorobenzene	10	10	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.005	<0.006	<0.012	0.00029 J	0.00033 J	<0.0003
Ethylbenzene	70	70	<0.0003	0.00065 J	0.0055	0.00093 J	0.031	0.0013	0.31	0.321	0.332	0.224	0.26	0.31
Methylene chloride	0.5	0.5	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.01	<0.0075	0.291	<0.00022	<0.001	<0.001
Toluene	100	100	<0.0002	0.00094 J	0.0033	<0.0002	0.023	<0.0002	1.1	1.18	1.12	0.724	0.72	0.99
Vinyl chloride	0.2	0.2												
Xylenes (total)	1000	1000	<0.0003	0.0044	0.013	0.00084 J	0.054	0.003	0.88	0.96	0.928	0.661	0.87	0.94
<b>Semivolatile Organic Compounds</b>														
1,2-Diphenylhydrazine	0.11	0.26	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0005	<0.0524	<0.0267	<0.156	<0.00021	<0.00021
2,4-Dimethylphenol	49	150	<0.000041	<0.000041	0.00044	<0.00004	<0.00004	<0.00004	20	98.1	29.9	182 J	10	16
2,4-Dinitrotoluene	0.13	0.3	<0.000059	<0.000059	<0.000058	<0.000058	<0.000058	<0.000058	<0.0005	<0.0619	<0.0316	<0.184	<0.00058	<0.00058
2,6-Dinitrotoluene	0.13	0.3	<0.000043	<0.000043	<0.000042	<0.000042	<0.000042	<0.000042	<0.0006	<0.0381	<0.0194	<0.113	<0.00042	<0.00042
2-Chloronaphthalene	200	580	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021	<0.0005	<0.0381	<0.0194	<0.113	<0.00021	<0.00021
2-Methylnaphthalene	9.8	29	<0.000019	0.0017	0.00031	<0.000019	0.024	0.006	0.74	1.39	1.19	3.37	0.33	0.42
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.00002	<0.0008	<0.395	<0.201	<1.17	<0.0002	<0.0002
4-Nitrophenol	4.9	15	<0.000048	<0.000048	<0.000047	<0.000047	<0.000047	<0.000047	<0.0005	<0.267	<0.136	<0.792	<0.00047	<0.00047
Acenaphthene	150	440	<0.000028	0.0043	0.0023	<0.000027	0.016	0.0053	0.23	0.584	0.476	1.6	0.07	0.15
Acenaphthylene	150	440	<0.000015	0.000066 J	0.00011	<0.000015	0.00033	0.00015	0.0073	<0.0286	<0.0146	<0.0849	0.0021	<0.00015
Anthracene	730	2200	0.000064 J	0.0022	0.00041	<0.000014	0.0066	0.0039	0.017	0.0646 J	0.033 J	0.179 J	0.0085	0.02
Benzo(a)anthracene	0.91	2	0.00015	0.00067	0.00013	<0.00005	0.0026	0.0011	<0.0005	<0.0381	<0.0194	<0.113	<0.0005	<0.0005
Benzo(a)pyrene	0.02	0.02	0.0002	0.00029	0.00018	0.000021 J	0.00083	0.00052	<0.0005	<0.0381	<0.0194	<0.113	<0.0002	<0.0002
bis(2-Chloroethoxy)methane	0.083	0.19	<0.000031	<0.000031	<0.00003	<0.00003	<0.00003	<0.00003	<0.0005	<0.0619	<0.0316	<0.184	<0.0003	<0.0003
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	0.00018 J	0.000081 J	<0.00056	<0.000037	<0.00014	0.00063	<0.001	<0.176	<0.0898	<0.524	<0.00037	<0.00037
Chrysene	91	200	0.00023	0.00077	0.00025	<0.000021	0.0024	0.0011	<0.0005	<0.0381	<0.0194	<0.113	<0.00021	<0.00021
Dibenzofuran	9.8	29	<0.00002	0.004	0.0019	<0.00002	0.016	0.0049	0.18	0.355	0.348	1.21	0.06	0.13
Di-n-butylphthalate (DBP)	240	730	<0.00002	<0.00002	<0.00002	0.000022 J	<0.00002	0.00098	<0.0005	<0.0524	<0.0267	<0.156	<0.0002	<0.0002
Fluoranthene	98	290	0.00033	0.0045	0.0005	0.000039 J	0.017	0.0057	0.0034	<0.0333	<0.017	<0.0991	<0.0001	0.0015
Fluorene	98	290	<0.000031	0.0032	0.0013	<0.00003	0.0093	0.0038	0.11	0.253	0.224	0.7 J	0.032	0.069
Naphthalene	49	150	<0.00002	0.00048	0.00011	<0.00002	0.44	0.0064	16	88.5 J	25	82.8	7.5	13
Nitrobenzene	4.9	15	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.0005	<0.0524	<0.0267	<0.156	<0.00024	<0.00024
N-Nitrosodiphenylamine	19	42	<0.000026	<0.000026	<0.000025	<0.000025	<0.000025	<0.000025	<0.0005	<0.0476	<0.0243	<0.142	<0.00025	<0.00025
Pentachlorophenol	0.1	0.1	<0.000081	<0.000081	<0.000079	<0.000079	<0.000079	<0.000079	<0.0005	<0.29	<0.148	<0.863	<0.00079	<0.00079
Phenanthrene	73	220	0.00012	0.012	0.0016	<0.000021	0.03	0.017	0.079	0.264	0.182	0.76	0.02	0.084
Phenol	730	2200	<0.000036	<0.000036	<0.000035	<0.000035	0.000051 J	<0.000035	3.4	7.51	6.31	31.4	4.2	4.2
Pyrene	73	220	0.00031	0.0028	0.00064	0.000037 J	0.0088	0.0042	0.0019 J	<0.0524	<0.0267	<0.156	<0.00019	0.0012
<b>Metals</b>														
Arsenic	1	1	0.00174 J	0.00214	0.000851 J	0.00158 J	0.000626 J	0.00279					0.00127 J	0.000624 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray.
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

**TABLE 3  
ANALYTICAL RESULTS SUMMARY (2008 - 2020)  
CLASS 3 GROUNDWATER  
HOUSTON, TX - WOOD PRESERVING WORKS**

	<b>Class 3 Residential Assessment</b>	<b>Class 3 C/I Assessment Level</b>	MW-72B 05/16/2018	MW-72B 01/24/2019	MW-72B 07/10/2019	MW-72B 01/09/2020
<b>Volatil Organic Compounds</b>						
1,2-Dichloroethane	0.5	0.5	<0.001	0.011	<0.0002	<0.002
Benzene	0.5	0.5	<b>1.2</b>	<b>0.63</b>	<b>1.1</b>	<b>0.7</b>
Chlorobenzene	10	10	<0.0015	<0.003	<0.0003	<0.003
Ethylbenzene	70	70	0.34	0.2	0.3	0.13
Methylene chloride	0.5	0.5	<0.005	<0.01	<0.001	<0.01
Toluene	100	100	0.95	0.58	0.96	0.58
Vinyl chloride	0.2	0.2				
Xylenes (total)	1000	1000	1.1	0.63	0.97	0.36
<b>Semivolatile Organic Compounds</b>						
1,2-Diphenylhydrazine	0.11	0.26	<0.00021	<0.00021	<0.00021	<0.00021
2,4-Dimethylphenol	49	150	14	2	19	13
2,4-Dinitrotoluene	0.13	0.3	<0.00058	<0.00058	<0.00058	<0.00058
2,6-Dinitrotoluene	0.13	0.3	<0.00042	<0.00042	<0.00042	<0.00042
2-Chloronaphthalene	200	580	<0.00021	<0.00021	<0.00021	<0.00021
2-Methylnaphthalene	9.8	29	0.23	0.071	0.57	0.22
4,6-Dinitro-2-methylphenol	0.24	0.73	<0.0002	<0.0002	<0.0002	<0.0002
4-Nitrophenol	4.9	15	<0.00047	0.0073 J	<0.00047	<0.00047
Acenaphthene	150	440	0.12	0.019	0.2	0.079
Acenaphthylene	150	440	0.003	0.00069 J	0.0044	0.0023
Anthracene	730	2200	<0.00014	0.0015	0.034	0.02
Benzo(a)anthracene	0.91	2	<0.0005	<0.0005	0.0069	0.0031
Benzo(a)pyrene	0.02	0.02	<0.0002	<0.0002	0.0019	0.0011
bis(2-Chloroethoxy)methane	0.083	0.19	<0.0003	<0.0003	<0.0003	<0.0003
bis(2-Ethylhexyl)phthalate (DEHP)	0.6	0.6	<0.00037	<0.00037	<0.00037	<0.00037
Chrysene	91	200	<0.00021	<0.00021	0.0069	0.0033
Dibenzofuran	9.8	29	0.082	0.017	0.16	0.061
Di-n-butylphthalate (DBP)	240	730	<0.0002	<0.0002	<0.0002	<0.0002
Fluoranthene	98	290	0.00095 J	<0.0001	0.046	0.024
Fluorene	98	290	0.051	0.0091	0.12	0.049
Naphthalene	49	150	12	1.2	12	2
Nitrobenzene	4.9	15	<0.00024	<0.00024	<0.00024	<0.00024
N-Nitrosodiphenylamine	19	42	<0.00025	<0.00025	<0.00025	<0.00025
Pentachlorophenol	0.1	0.1	<0.00079	<0.00079	<0.00079	<0.00079
Phenanthrene	73	220	0.045	0.0042	0.16	0.075
Phenol	730	2200	2.3	0.58	9.3	5.2
Pyrene	73	220	0.00074 J	<0.00019	0.029	0.013
<b>Metals</b>						
Arsenic	1	1	0.000951 J	0.00106 J	0.000861 J	0.00105 J

Notes:

1. All values in milligrams per liter (mg/L).
2. Concentrations > RAL and non-detects are highlighted light gray.
3. Concentrations > C/I AL and non-detects are highlighted dark gray
4. TRRP PCLs (30 TAC §350, Table 3), updated November 2019.
5. RAL = Residential Assessment Level, C/I = Commercial/Industrial
6. J = Estimated value, < = not detected at the specified detection limit.

TABLE 4

Analytical Results Summary - A-TZ Wells - VISL Comparison  
Houston, tx - Wood Preserving Works

Location ID:	Target Groundwater Concentration (TCR=1E-05 or THQ=0.1) Cgw,Target (mg/L)	MW-25A	MW-25A	MW-26A	MW-26A	MW-27A	MW-27A	MW-28A	MW-28A
Sample Date:		07/16/2019	01/15/2020	07/17/2019	01/16/2020	07/18/2019	01/15/2020	07/16/2019	01/16/2020
<b>Semivolatile Organic Compounds (mg/L)</b>									
Benzo(a)anthracene	0.344	<0.00005	0.00012	<0.00005	<0.00005	<0.00005	<0.00005	0.000055 J	0.000073 J
Naphthalene	0.0174	0.00023	0.0001	<0.00032	<0.00017	0.0026	0.000053 J	0.0013	<0.0006
Nitrobenzene	0.715	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
<b>Volatile Organic Compounds (mg/L)</b>									
1,2-Dichloroethane	0.0151	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.0138	<0.0002	<0.0002	0.00036 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.041	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.0349	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.471	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1.92	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.00147	<0.0002	<0.0002	NA	NA	NA	NA	NA	NA
Xylenes (total)	0.0385	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003

Notes: VISL Calculator output for target groundwater screening level with inputs: residential scenario, hazard quotient of 0.1, carcinogenic risk of 10-5, and groundwater temperature of 25 degree Celsius (EPA, 2019).

J = Estimated value

< = Compound not detected at the specified detection limit.

TABLE 4

Analytical Results Summary - A-TZ Wells - VISL Comparison  
Houston, tx - Wood Preserving Works

Location ID:	Target Groundwater Concentration (TCR=1E-05 or THQ=0.1)	MW-32AR	MW-32AR	MW-33A	MW-33A (DUP)	MW-33A	MW-33A (DUP)	MW-35A	MW-35A	MW-36A	MW-36A
Sample Date:	Cgw,Target (mg/L)	07/30/2019	01/20/2020	07/17/2019	07/17/2019	01/20/2020	01/20/2020	07/18/2019	01/10/2020	07/16/2019	01/09/2020
<b>Semivolatile Organic Compounds (mg/L)</b>											
Benzo(a)anthracene	0.344	0.000054 J	<0.00005	<0.00005	<0.00005	0.00012	0.00012	<0.00005	<0.00005	<0.00005	<0.00005
Naphthalene	0.0174	<0.00002	<0.00014	<0.000076	<0.000096	<0.0005	<0.00066	0.00083	<0.0002	0.00015	<0.00034
Nitrobenzene	0.715	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
<b>Volatile Organic Compounds (mg/L)</b>											
1,2-Dichloroethane	0.0151	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.0138	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.041	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.0349	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.471	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1.92	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.00147	NA	<0.0002	NA	NA	<0.0002	<0.0002	NA	<0.0002	<0.0002	<0.0002
Xylenes (total)	0.0385	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003

Notes: VISL Calculator output for target groundwater screening level with inputs: residential scenario, hazard quotient of 0.1, carcinogenic risk of 10-5, and groundwater temperature of 25 degree Celsius (EPA, 2019).

J = Estimated value

< = Compound not detected at the specified detection limit.

TABLE 4

Analytical Results Summary - A-TZ Wells - VISL Comparison  
Houston, tx - Wood Preserving Works

Location ID:	Target Groundwater Concentration (TCR=1E-05 or THQ=0.1) Cgw, Target (mg/L)	MW-38A	MW-38A	MW-44A	MW-44A	MW-47A	MW-59A	MW-59A	MW-60A	MW-60AR
Sample Date:		07/31/2019	01/21/2020	07/17/2019	01/09/2020	03/20/2020	07/17/2019	01/16/2020	07/17/2019	03/20/2020
<b>Semivolatile Organic Compounds (mg/L)</b>										
Benzo(a)anthracene	0.344	0.000087 J	<0.00005	0.000085 J	0.000054 J	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005
Naphthalene	0.0174	0.000068 J	0.0011	<0.00019	<0.003	<0.00002	<0.00012	<0.0012	<0.00013	<0.00002
Nitrobenzene	0.715	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
<b>Volatile Organic Compounds (mg/L)</b>										
1,2-Dichloroethane	0.0151	0.00031 J	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.0138	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.041	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.0349	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.471	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1.92	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.00147	NA	NA	<0.0002	NA	NA	<0.0002	<0.0002	<0.0002	NA
Xylenes (total)	0.0385	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003

Notes: VISL Calculator output for target groundwater screening level with inputs: residential scenario, hazard quotient of 0.1, carcinogenic risk of 10-5, and groundwater temperature of 25 degree Celsius (EPA, 2019).

J = Estimated value

< = Compound not detected at the specified detection limit.

TABLE 4

Analytical Results Summary - A-TZ Wells - VISL Comparison  
Houston, tx - Wood Preserving Works

Location ID:	Target Groundwater Concentration (TCR=1E-05 or THQ=0.1) Cgw, Target (mg/L)	MW-61A	MW-61A	MW-68A	MW-68A	MW-69A	MW-84A	MW-91A	MW-94A	MW-95A
Sample Date:		07/17/2019	01/16/2020	07/18/2019	01/17/2020	07/17/2019	03/12/2020	03/12/2020	03/12/2020	03/17/2020
<b>Semivolatile Organic Compounds (mg/L)</b>										
Benzo(a)anthracene	0.344	<0.00005	<0.00005	<0.00005	<0.00005	<0.00005	<0.000051	<0.000053	<0.000052	<0.00005
Naphthalene	0.0174	<0.00002	<0.00021	0.00035	<0.00022	<0.000078	0.000023 J	0.000083 J	<0.000021	<0.00002
Nitrobenzene	0.715	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024	<0.000025	<0.000025	<0.000024
<b>Volatile Organic Compounds (mg/L)</b>										
1,2-Dichloroethane	0.0151	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Benzene	0.0138	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Chlorobenzene	0.041	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Ethylbenzene	0.0349	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003
Methylene chloride	0.471	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Toluene	1.92	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Vinyl chloride	0.00147	<0.0002	<0.0002	NA	NA	<0.0002	NA	NA	NA	NA
Xylenes (total)	0.0385	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003	<0.0003

Notes: VISL Calculator output for target groundwater screening level with inputs: residential scenario, hazard quotient of 0.1, carcinogenic risk of 10-5, and groundwater temperature of 25 degree Celsius (EPA, 2019).

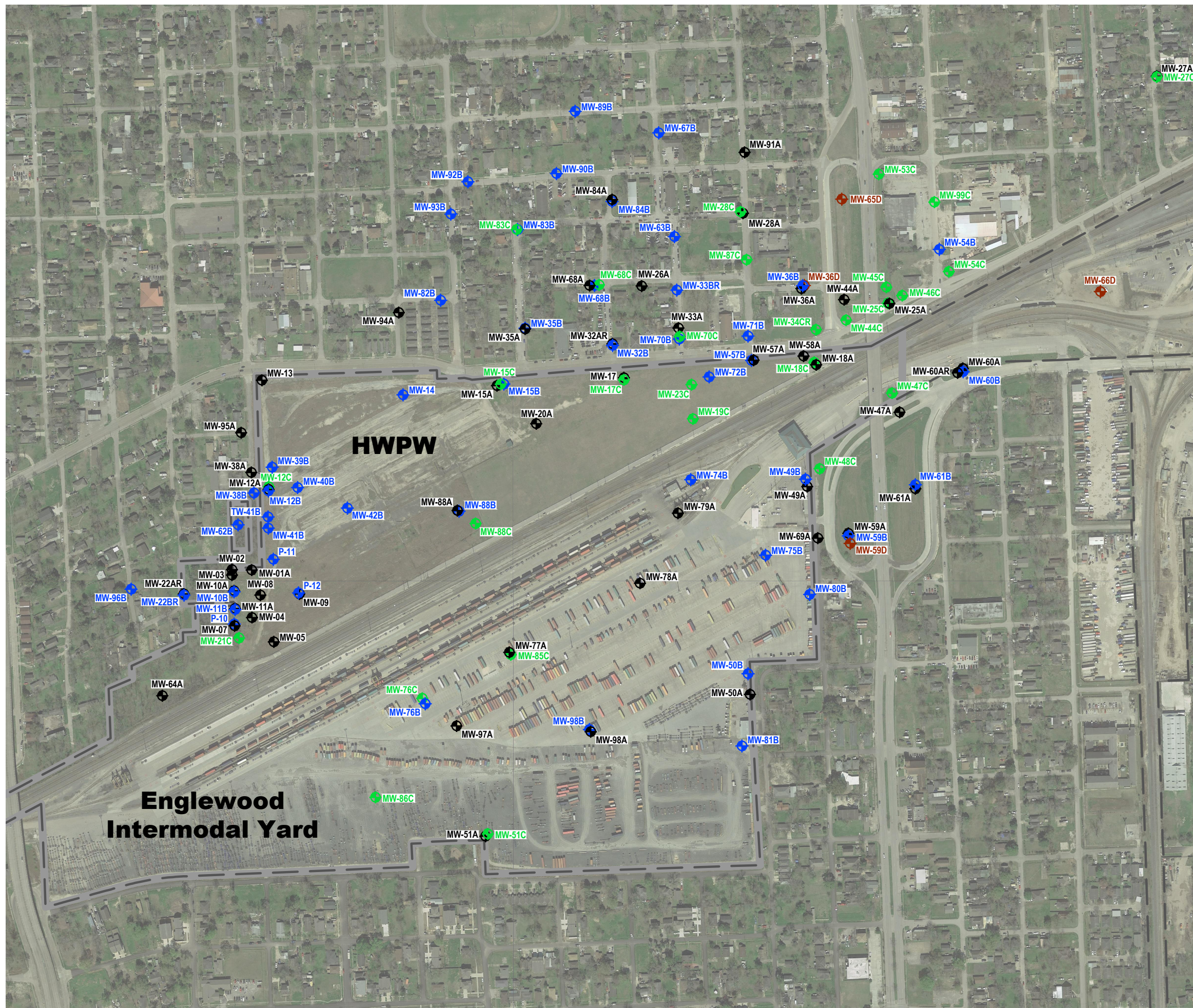
J = Estimated value

< = Compound not detected at the specified detection limit.



**FIGURES**





**LEGEND**

- UPRR PROPERTY BOUNDARY
- A-TZ MONITORING WELL LOCATION
- B-CZ/B-TZ MONITORING WELL LOCATION
- C-TZ MONITORING WELL LOCATION
- D-TZ MONITORING WELL LOCATION

**REFERENCE(S)**  
 PARCEL BOUNDARIES: CITY OF HOUSTON GEOGRAPHIC INFORMATION & MANAGEMENT SYSTEMS (GIMS).  
 AERIAL: GOOGLE EARTH, IMAGERY DATED 2/23/19.



UNION PACIFIC RAILROAD CO.

HOUSTON WOOD PRESERVING WORKS

**MONITORING WELL LOCATION MAP**

2020-03-18

AJD

AJD

MH

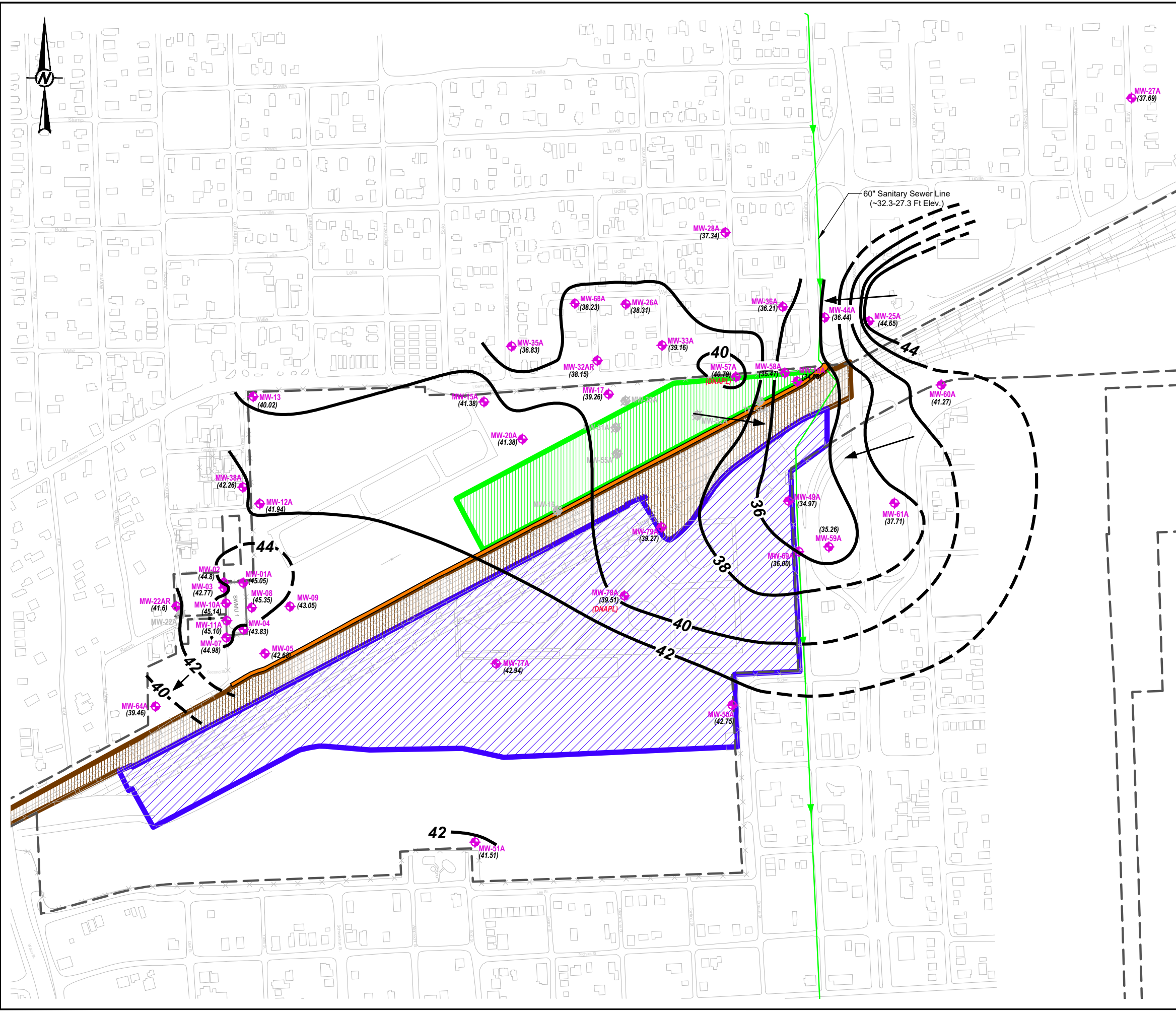
ECM

**GOLDER**

19119232 0 **FIGURE 1**



Path: \\uswest\erdm\dat\proj\19119232 - houston\2019-07-19\19119232 - houston\2019-07-19\Cad\1 - Groundwater Gradient Map A-TZ - July 2019.dwg | Last Edited By: rlsalazar | Date: 2020-04-30 | Printed By: rlsalazar | Date: 2020-04-30 | Time: 12:01 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- A-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- (37.71) GROUNDWATER ELEVATION (FT, HVD) (NM = NOT MEASURED)
- 39 GROUNDWATER ELEVATION CONTOUR (FT, HVD) C.I. = 2 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA

- NOTE(S)**
1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
  2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
  3. \*\* - NOT USED TO DEVELOP POTENTIOMETRIC SURFACE.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

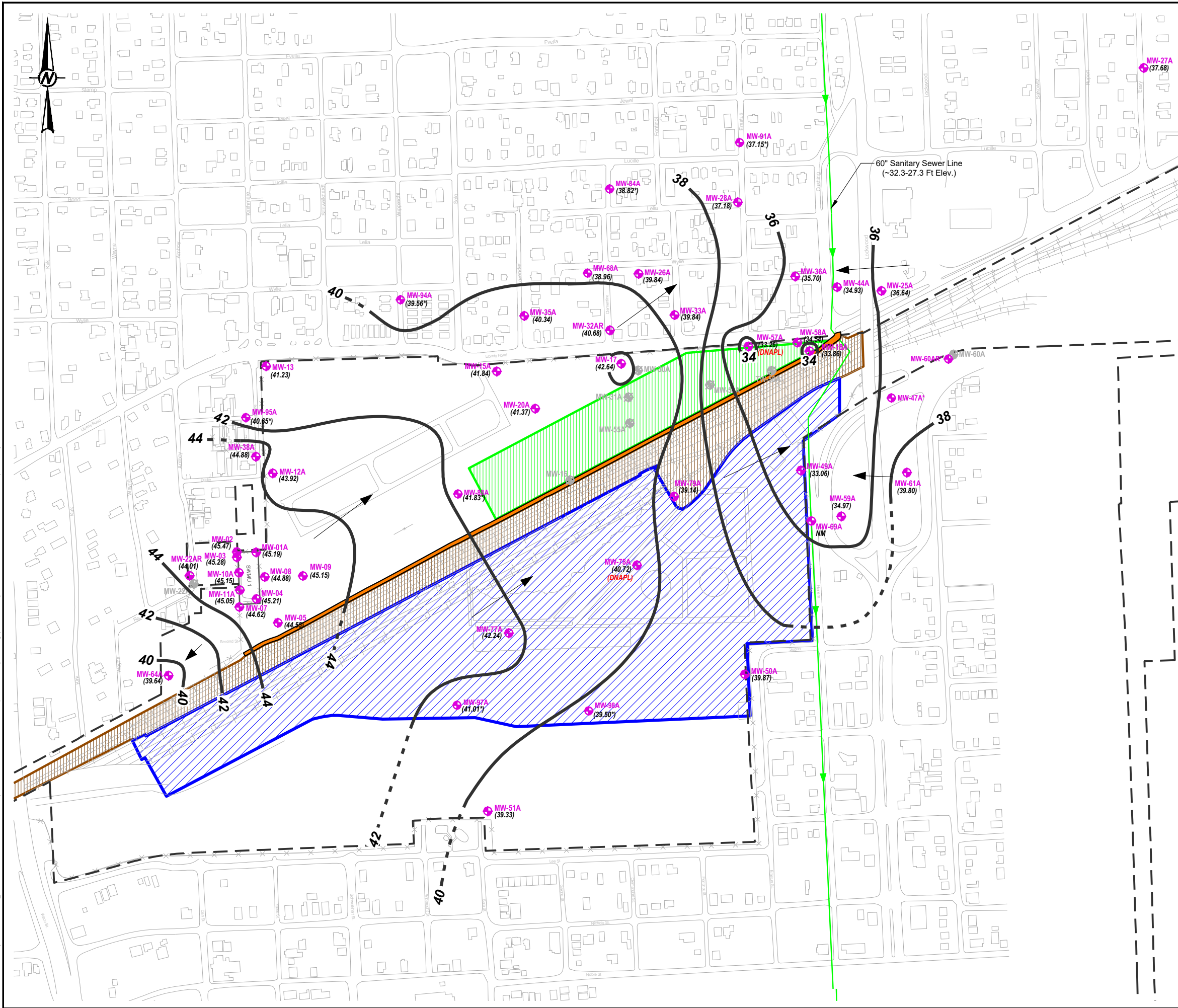
**TITLE**  
 GROUNDWATER GRADIENT MAP A-TZ  
 JULY 2019

CONSULTANT	DATE	BY
YYYY-MM-DD	2020-04-30	
DESIGNED		AJD
PREPARED		AJD
REVIEWED		MH
APPROVED		ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5A-1a

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\proj\19119232 - HWPW\2020-31\Month 1 - File Name: Groundwater Gradient Map A-TZ - January 2020.dwg | User: erand | Date: 2020-04-30 11:04:49 AM | Printed By: Rbalazar | Date: 2020-04-30 11:11:17 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- A-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- (37.71) GROUNDWATER ELEVATION (FT, HVD) (NM = NOT MEASURED)
- 39 GROUNDWATER ELEVATION CONTOUR (FT, HVD) C.I. = 2 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA

- NOTE(S)**
1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
  2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  3. \*\* - NOT USED TO DEVELOP POTENTIOMETRIC SURFACE.
  4. \* - WATER ELEVATION MEASURED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

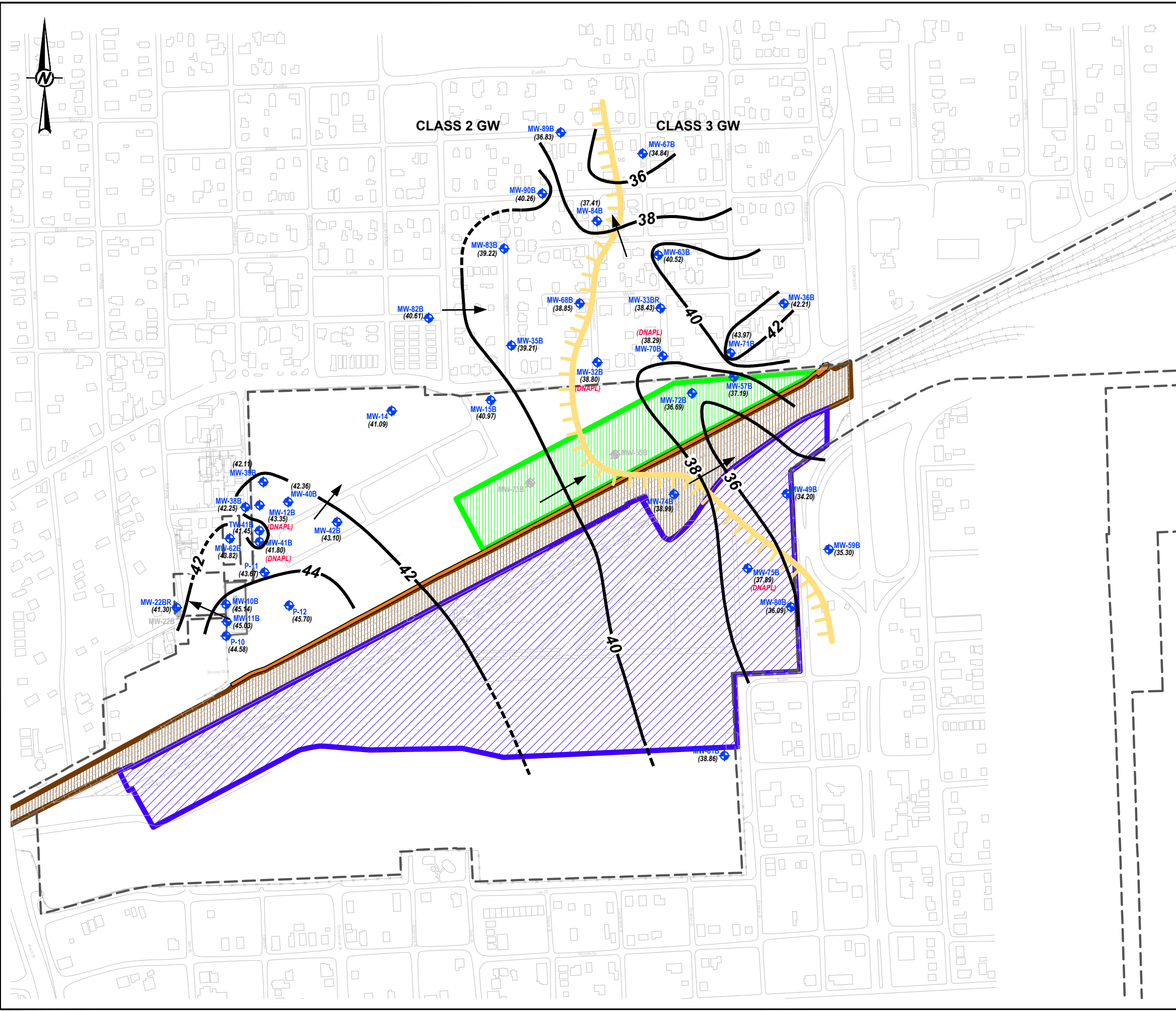


<b>CLIENT</b>		UNION PACIFIC RAILROAD CO.	
<b>PROJECT</b>		HOUSTON WOOD PRESERVING WORKS	
<b>TITLE</b>		GROUNDWATER GRADIENT MAP - A-TZ JANUARY-MARCH 2020	
<b>CONSULTANT</b>	YYYY-MM-DD	2020-04-30	
<b>GOLDER</b>	DESIGNED		
<small>TEXAS GEOSCIENCE FIRM NO. 50969 TEXAS ENGINEERING FIRM NO. 2578</small>	PREPARED	RS	
	REVIEWED	MH	
	APPROVED	ECM	
<b>PROJECT NO.</b>	<b>REV.</b>	<b>FIGURE</b>	
19119232	0	5A-1b	

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



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 Date: 2020-04-30 | Time: 1:22:50 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW) B-TZ/B-CZ BOUNDARY
- B-TZ (CLASS 2 GW) (39.98)
- GROUNDWATER ELEVATION (FT, HVD) (NM = NOT MEASURED)
- GROUNDWATER ELEVATION CONTOUR (FT, HVD) C.I. = 2 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA

**NOTE(S)**

1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
3. \*\* - NOT USED TO GENERATE CONTOURS.

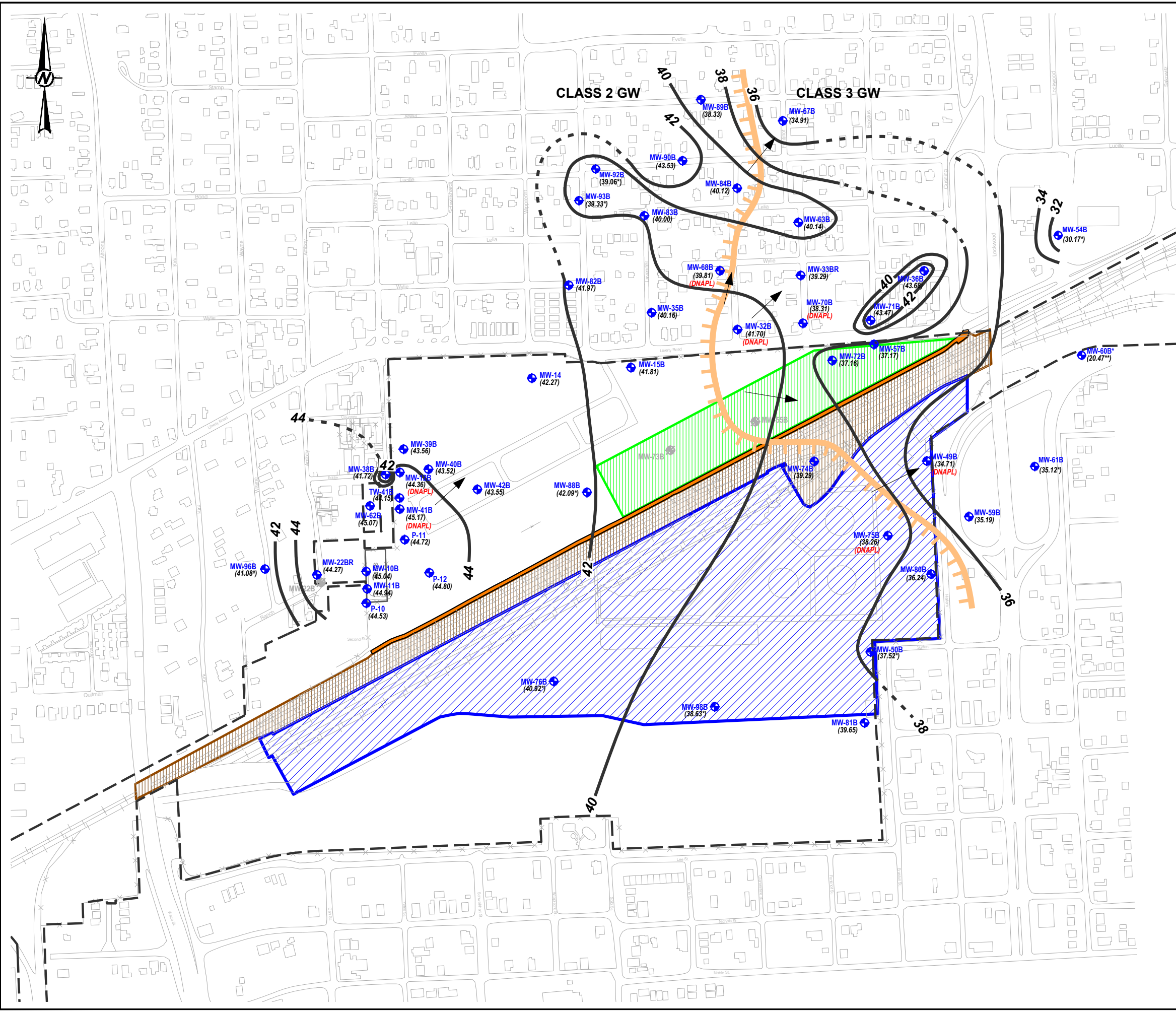
**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



<b>CLIENT</b> UNION PACIFIC RAILROAD CO.		
<b>PROJECT</b> HOUSTON WOOD PRESERVING WORKS		
<b>TITLE</b> GROUNDWATER GRADIENT MAP B-TZ AND B-CZ JULY 2019		
<b>CONSULTANT</b>	YYYY-MM-DD	2020-04-30
	DESIGNED	AJD
	PREPARED	AJD
	REVIEWED	MH
	APPROVED	ECM
<b>PROJECT NO.</b> 19119232	<b>REV.</b> 0	<b>FIGURE</b> 5A-2a

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\2019\01\19232 - HWPP\2020-31\Month 1 - File Name: Groundwater Gradient Map B-TZ/B-CZ - January 2020.dwg | Last Edited By: rsalazar | Date: 2020-04-30 | Printed By: RSalazar | Date: 2020-04-30 | Time: 11:04:15 AM | File Name: Groundwater Gradient Map B-TZ/B-CZ - January 2020.dwg | Last Edited By: rsalazar | Date: 2020-04-30 | Printed By: RSalazar | Date: 2020-04-30 | Time: 12:11:46 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW) B-TZ/B-CZ BOUNDARY
- B-TZ (CLASS 2 GW) (39.98)
- GROUNDWATER ELEVATION (FT, HVD) (NM = NOT MEASURED)
- GROUNDWATER ELEVATION CONTOUR (FT, HVD) C.I. = 2 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA

**NOTE(S)**

1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
3. \*\* - NOT USED TO GENERATE CONTOURS.
4. \* - WATER ELEVATION MEASURED IN MARCH 2020.

**REFERENCE(S)**

BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

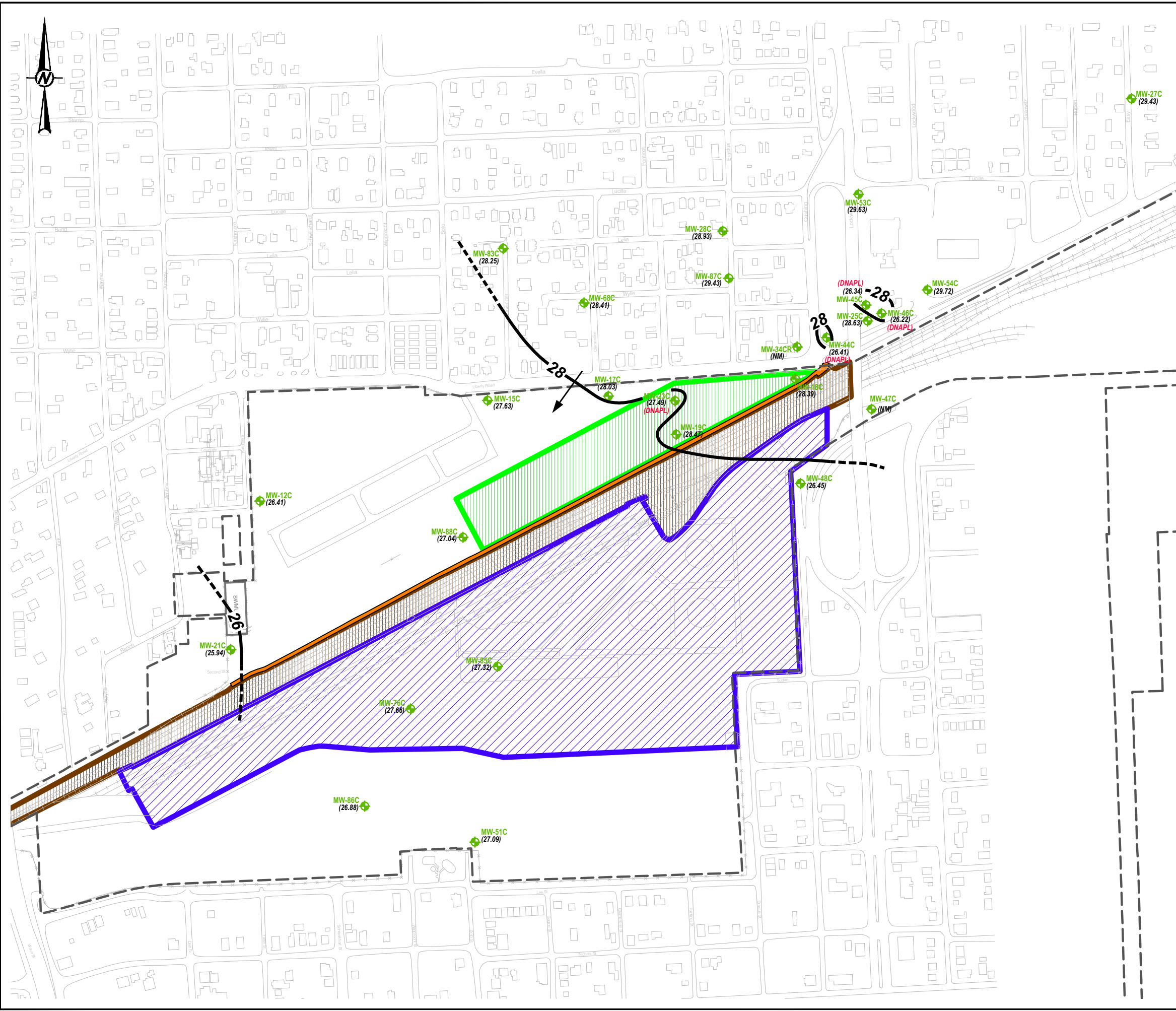


<p>CLIENT UNION PACIFIC RAILROAD CO.</p> <hr/> <p>PROJECT HOUSTON WOOD PRESERVING WORKS</p> <hr/> <p>TITLE <b>GROUNDWATER GRADIENT MAP - B-TZ/B-CZ</b> JANUARY-MARCH 2020</p> <hr/> <p>CONSULTANT <b>GOLDER</b> TEXAS GEOSCIENCE FIRM NO. 50369 TEXAS ENGINEERING FIRM NO. 2578</p>	<table border="0" style="width: 100%; border-collapse: collapse;"> <tr> <td>YYYY-MM-DD</td> <td>2020-04-30</td> </tr> <tr> <td>DESIGNED</td> <td></td> </tr> <tr> <td>PREPARED</td> <td>RS</td> </tr> <tr> <td>REVIEWED</td> <td>MH</td> </tr> <tr> <td>APPROVED</td> <td>ECM</td> </tr> </table> <hr/> <p>PROJECT NO. 19119232</p> <p style="text-align: right;">REV. 0</p> <p style="text-align: right;">FIGURE 5A-2b</p>	YYYY-MM-DD	2020-04-30	DESIGNED		PREPARED	RS	REVIEWED	MH	APPROVED	ECM
YYYY-MM-DD	2020-04-30										
DESIGNED											
PREPARED	RS										
REVIEWED	MH										
APPROVED	ECM										

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\msanet\share\data\projects\19119232\_19119232 - Inpwr\2019-10-10\Cad\1 - File Name: FIG 5A-3 - Groundwater Gradient Map C-TZ - July 2019.dwg | Last Edited By: rmlsaur | Date: 2020-04-30 | Time: 1:20:55 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- GROUNDWATER ELEVATION (FT, HVD)  
(NM = NOT MEASURED)
- GROUNDWATER ELEVATION CONTOUR (FT, HVD)  
C.I. = 2 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA

- NOTE(S)**
1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
  2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
  3. \*\* NOT USED TO CREATE POTENTIOMETRIC SURFACE.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

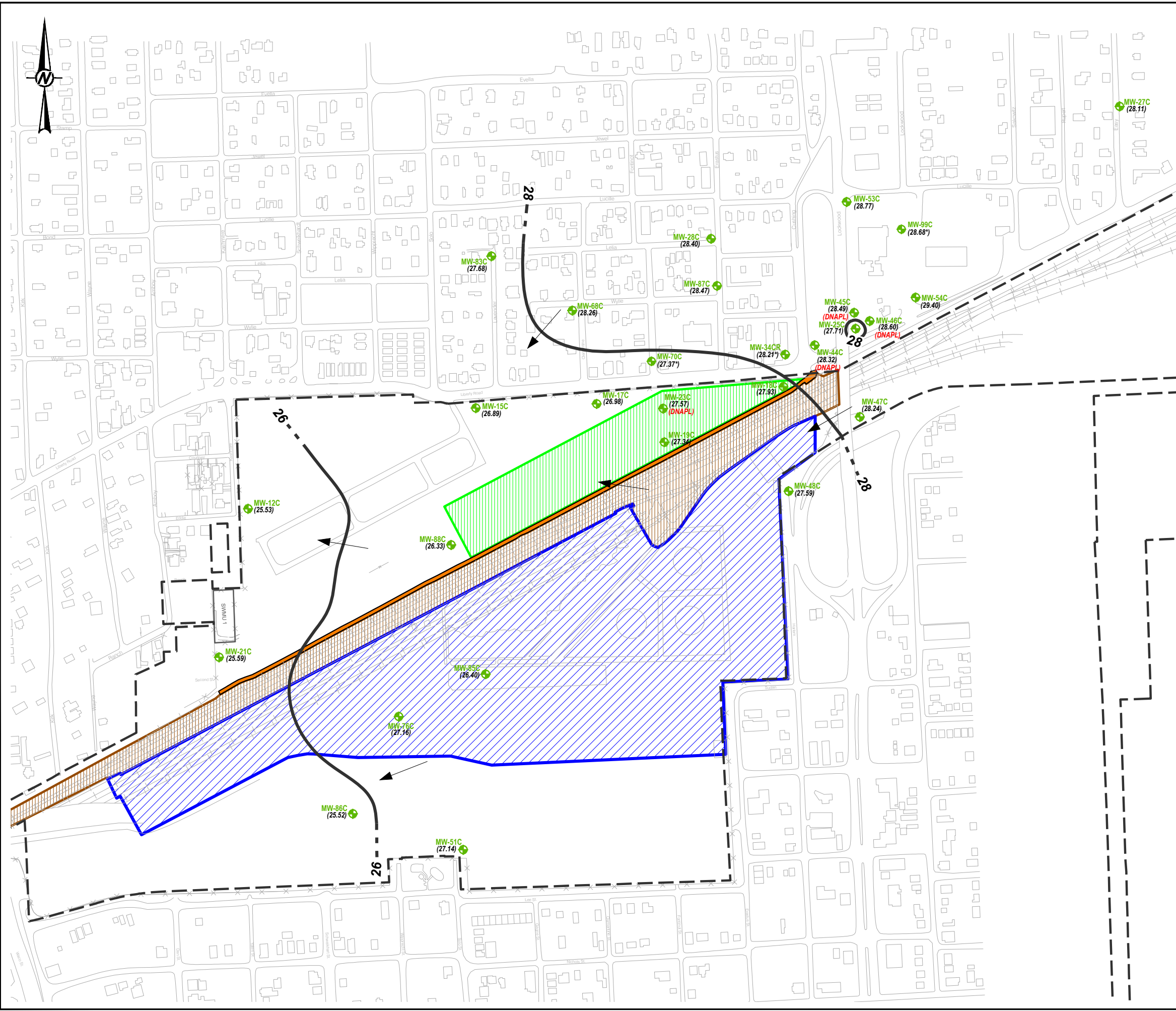
**TITLE**  
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 JULY 2019

<b>CONSULTANT</b>	YYYY-MM-DD	2020-04-30
	DESIGNED	AJD
	PREPARED	AJD
	REVIEWED	MH
	APPROVED	ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5A-3a

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\19119232 - HWPW\2020-31\Month 1 - File Name: Groundwater Gradient Map C-TZ - January 2020.dwg | Last Edited By: rmatzner | Date: 2020-04-30 | Time: 11:03:38 AM | Plotted By: RMatzner | Date: 2020-04-30 | Time: 12:14 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ▭ HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- ××× FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- (29.40) GROUNDWATER ELEVATION (FT, HVD)  
(NM = NOT MEASURED)
- 28 — GROUNDWATER ELEVATION CONTOUR (FT, HVD)  
C.I. = 2 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- ▨ RAILROAD BALLAST CAP AREA
- ▨ ASPHALT CAP AREA
- ▨ SOIL CAP
- ▨ CONCRETE CAP AREA

- NOTE(S)**
1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
  2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  3. \*\* - NOT USED TO DEVELOP POTENTIOMETRIC SURFACE.
  4. \* - WATER ELEVATION MEASURED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER GRADIENT MAP - C-TZ**  
**JANUARY-MARCH 2020**

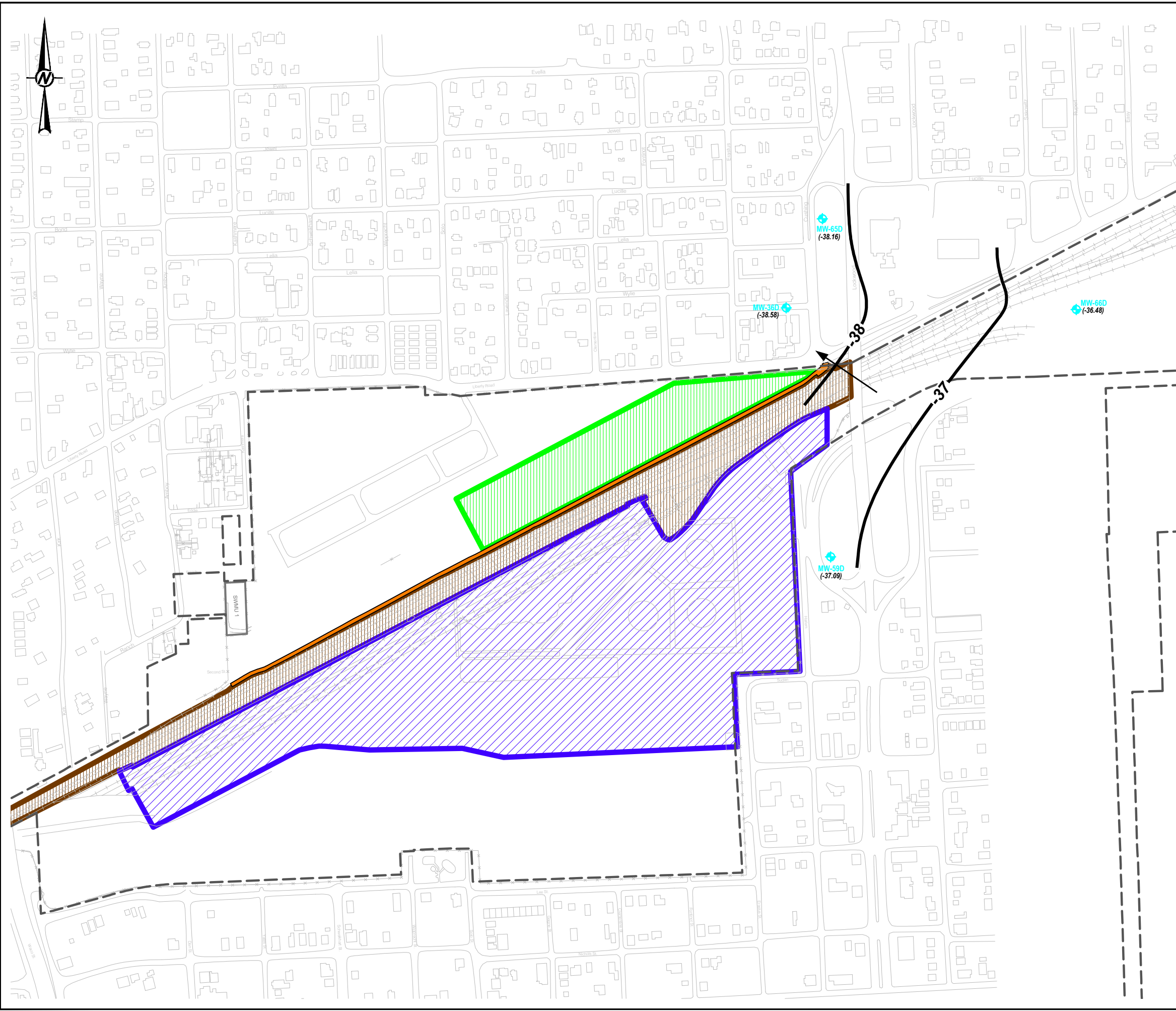
CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5A-3b

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erand\del\proj\del\2019\01\19232 - hwy\2019-10-10\Del\_1\_Files\Name: FIG 5A-4 - Groundwater Gradient Map D-TZ - July 2019.dwg | Last Edited By: rsmiller Date: 2020-04-30 Time: 1:20:00 PM  
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**LEGEND**

- UPRR PROPERTY BOUNDARY
- HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- D-TZ MONITORING WELL LOCATION
- GROUNDWATER ELEVATION (FT, HVD)  
(NM = NOT MEASURED)
- GROUNDWATER ELEVATION CONTOUR (FT, HVD)  
C.I. = 1 FT
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA

**NOTE(S)**

1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
2. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).

**REFERENCE(S)**

BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
UNION PACIFIC RAILROAD CO.

**PROJECT**  
HOUSTON WOOD PRESERVING WORKS

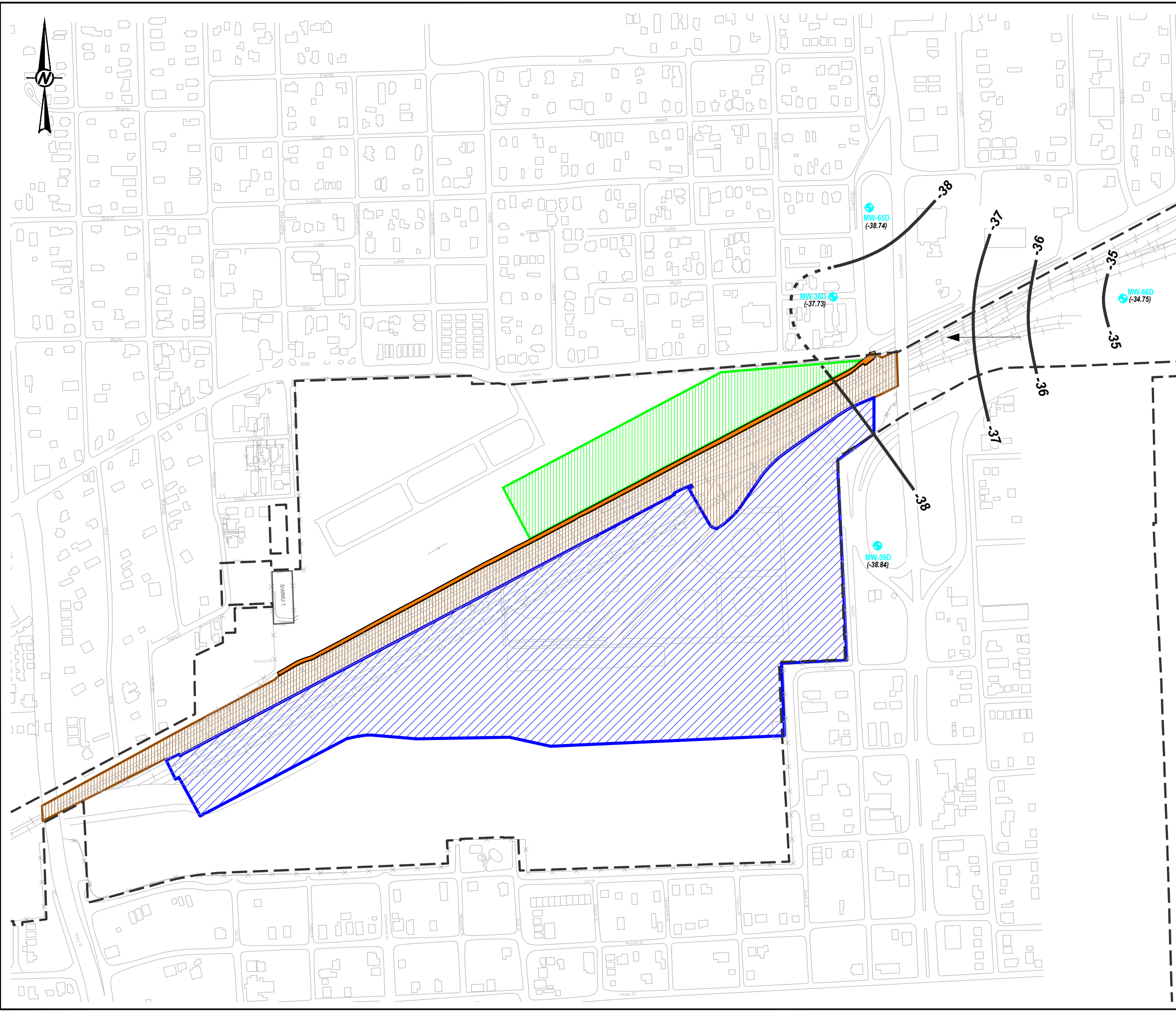
**TITLE**  
GROUNDWATER GRADIENT MAP D-TZ  
JULY 2019

CONSULTANT	YYYY-MM-DD	2020-04-30
<b>GOLDER</b> TEXAS GEOSCIENCE FIRM NO. 50969 TEXAS ENGINEERING FIRM NO. 2578	DESIGNED	AJD
	PREPARED	AJD
	REVIEWED	MH
	APPROVED	ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5A-4a

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\staff\Projects - Round Rock\2019\01\19232 - HWPP\2020-31\Month 1 - File Name: Groundwater Gradient Map D-TZ - January 2020.dwg | Last Edited By: rmatzner | Date: 2020-04-30 | Time: 11:03:14 AM | Printed By: RMatzner | Date: 2020-04-30 | Time: 12:45 PM



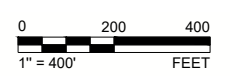
**LEGEND**

- UPRR PROPERTY BOUNDARY
- ▭ HISTORIC STRUCTURE AND FEATURE
- ROAD, PARKING LOT, SIDEWALK
- ××× FENCE
- RAILROAD
- ⊕ D-TZ MONITORING WELL LOCATION
- ⊙ PLUGGED AND ABANDONED MONITORING WELL
- (37.71) GROUNDWATER ELEVATION (FT, HVD)  
(NM = NOT MEASURED)
- 39 — GROUNDWATER ELEVATION CONTOUR (FT, HVD)  
C.I. = 2 FT
- ➔ INFERRED GROUNDWATER FLOW DIRECTION
- ▨ RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- ▨ SOIL CAP
- ▨ CONCRETE CAP AREA

**NOTE(S)**

1. VERTICAL DATUM BASED ON CITY OF HOUSTON VERTICAL DATUM (HVD).
2. DNAPL = DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).

**REFERENCE(S)**  
BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



<b>CLIENT</b> UNION PACIFIC RAILROAD CO.	
<b>PROJECT</b> HOUSTON WOOD PRESERVING WORKS	
<b>TITLE</b> GROUNDWATER GRADIENT MAP - D-TZ JANUARY-MARCH 2020	
<b>CONSULTANT</b>	YYYY-MM-DD 2020-04-30
<b>DESIGNED</b>	RS
<b>PREPARED</b>	RS
<b>REVIEWED</b>	MH
<b>APPROVED</b>	ECM
<b>PROJECT NO.</b> 19119232	<b>REV.</b> 0
<b>FIGURE</b> 5A-4b	

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

**ATTACHMENT A**

# Monitoring Well Construction Logs

**TABLE A-1  
SUMMARY OF GROUNDWATER MONITORING WELLS  
HOUSTON, TX - WOOD PRESERVING WORKS**

WELL NO.	DATE INSTALLED	NORTHING	EASTING	TOP OF CASING ELEVATION (FT HVD)	TOTAL DEPTH (FT BGS)	Top Screen Interval (FT BGS)	Bottom Screen Interval (FT BGS)	Zone	Plugged
<b>SWMU No. 1 Monitoring Wells</b>									
MW-01A <sup>1</sup>	1/0/1900	728,006	3,165,936	47.90	19	8.5	18.5	A-TZ	
MW-02 <sup>1</sup>	4/17/1984	728,007	3,165,857	47.89	18.5	8.5	18.5	A-TZ	
MW-07 <sup>1</sup>	3/27/1991	727,779	3,165,867	48.91	23	14.1	19.1	A-TZ	
MW-08 <sup>1</sup>	3/27/1991	727,903	3,165,972	49.33	24	14.2	19.2	A-TZ	
MW-10A <sup>1</sup>	9/13/1994	727,921	3,165,866	49.83	23	11	20.5	A-TZ	
MW-10B <sup>1</sup>	9/14/1994	727,916	3,165,866	49.96	46	27.1	41.6	B-TZ	
MW-11A <sup>1</sup>	9/15/1994	727,849	3,165,869	50.16	22	10	19.3	A-TZ	
MW-11B <sup>1</sup>	9/19/1994	727,845	3,165,869	50.24	44	27.5	41.2	B-TZ	
P-10 <sup>1</sup>	3/26/1991	727,786	3,165,866	47.71	50	36.2	38.2	B-TZ	
P-12 <sup>1</sup>	3/27/1991	727,912	3,166,127	48.76	50	36.3	38.3	B-TZ	
<b>Site-Wide Monitoring Wells</b>									
MW-03	4/17/1984	727,985	3,165,857	48.34	18.5	8.5	18.5	A-TZ	
MW-04	4/18/1984	727,813	3,165,938	49.85	21	11	21	A-TZ	
MW-05	----	727,715	3,166,026	49.24	26	10	25	A-TZ	
MW-09	3/26/1991	727,908	3,166,130	49.26	24	14.8	19.8	A-TZ	
MW-12A	2/27/1997	728,333	3,166,004	49.96	30	17.5	27.5	A-TZ	
MW-12B	2/27/1997	728,328	3,166,004	50.02	45	32.5	42.5	B-TZ	
MW-12C	4/21/1997	728,345	3,166,005	50.14	75.3	69	73.5	C-TZ	
MW-13	2/25/1997	728,777	3,165,977	50.65	25	9	22.5	A-TZ	
MW-14	2/27/1997	728,718	3,166,550	50.66	45	28	42.5	B-TZ	
MW-15A	2/25/1997	728,755	3,166,931	50.41	30	12	26.1	A-TZ	
MW-15B	12/19/2011	728,761	3,166,960	47.05	40	28	38	B-TZ	
MW-15C	4/25/1997	728,761	3,166,947	50.01	75	64	73.5	C-TZ	
MW-16	2/26/1997	728,305	3,167,235	51.51	30	12.5	27	A-TZ	P
MW-17	3/25/1997	728,787	3,167,447	50.92	35	18	32.5	A-TZ	
MW-17C	12/10/2003	728,779	3,167,446	50.17	70	59.5	69.5	C-TZ	
MW-18A	2/26/1997	728,839	3,168,227	51.57	35	18	32.5	A-TZ	P
MW-18C	4/25/1997	728,849	3,168,219	51.47	80.2	62	76.5	C-TZ	P
MW-19C	10/15/1998	728,620	3,167,727	53.05	73	63	73	C-TZ	
MW-20A	9/28/1998	728,600	3,167,091	50.43	30	15	25	A-TZ	
MW-21C	10/26/1998	727,730	3,165,884	49.05	72.5	62.5	72.5	C-TZ	
MW-22A	10/1/1998	727,876	3,165,677	46.07	25	10	20	A-TZ	P
MW-22AR	1/24/2018	727,909	3,165,660	45.56	25	9.6	19.6	A-TZ	
MW-22B	10/27/1998	727,871	3,165,678	45.86	38	27.5	37.5	B-TZ	P
MW-22BR	1/24/2018	727,904	3,165,660	45.71	39	27.6	37.6	B-TZ	
MW-23C	10/14/1998	728,759	3,167,722	54.16*	72.5	62.5	72.5	C-TZ	
MW-24AR	1/27/2009	727,531	3,165,207	45.65	21	11	21	A-TZ	
MW-24B	3/7/2000	727,534	3,165,208	46.06	50	38.5	48.5	B-TZ	
MW-24C	3/8/2000	727,542	3,165,206	46.05	74	63	73	C-TZ	
MW-25A	3/7/2000	729,089	3,168,524	44.65	29	18.5	28.5	A-TZ	
MW-25C	3/7/2000	729,089	3,168,518	44.49	74	58	68	C-TZ	
MW-26A	3/7/2000	729,159	3,167,519	44.62	26	14.5	24.5	A-TZ	
MW-27A	3/26/2001	730,002	3,169,610	44.90	30	17	27	A-TZ	
MW-27C	4/16/2001	730,009	3,169,610	45.04	73.5	60.5	70.5	C-TZ	
MW-28A	3/26/2001	729,462	3,167,926	43.86	28	16	26	A-TZ	
MW-28C	4/12/2001	729,461	3,167,920	43.96	88	75	85	C-TZ	
MW-29A	4/19/2001	727,310	3,164,239	46.59	23	9	19	A-TZ	P
MW-29B	4/12/2001	727,303	3,164,239	46.26	57	44	54	B-TZ	P
MW-29C	4/27/2001	727,293	3,164,240	46.46	75	62.5	72.5	C-TZ	P
MW-30A	12/8/2003	728,759	3,167,517	50.45	31	19.5	29.5	A-TZ	P
MW-31A	12/8/2003	728,648	3,167,477	52.08	33	21.5	31.5	A-TZ	P
MW-32A	12/29/203	728,914	3,167,401	43.77	33	20.5	30.5	A-TZ	P
MW-32AR	12/15/2011	728,925	3,167,400	44.74	22	10	20	A-TZ	
MW-32B	12/15/2011	728,918	3,167,400	44.73	40	26	36	B-TZ	
MW-33A	12/30/2003	728,989	3,167,668	44.25	25	13	23	A-TZ	
MW-33B	2/23/2007	729,150	3,167,661	44.35	75	32	42	B-CZ	P
MW-33BR	12/19/2011	729,142	3,167,662	44.86	40	28	38	B-CZ	
MW-33C	8/14/2006	729,092	3,167,663	NA	40	NA	NA	C-TZ	P
MW-34C	1/13/2004	728,934	3,168,160	45.31	72.5	60	70	C-TZ	P
MW-34CR	5/9/2014	728,982	3,168,227	46.47	70	60	70	C-TZ	
MW-35A	2/21/2007	728,985	3,167,045	44.75	28	13	28	A-TZ	
MW-35B	2/26/2007	728,988	3,167,045	44.83	42	32	42	B-CZ	
MW-35C	8/14/2006	729,021	3,167,040	NA	40	NA	NA	C-TZ	P
MW-36A	2/22/2007	729,148	3,168,167	44.53	28	18	28	A-TZ	
MW-36B	6/24/2010	729,161	3,168,172	44.07	43	38	43	B-CZ	
MW-36D	6/23/2010	729,162	3,168,180	44.33	110	100	110	D-TZ	
MW-37A	8/16/2006	728,896	3,166,210	NA	25	NA	NA	A-TZ	P
MW-38A	2/21/2007	728,402	3,165,934	46.39	22	12	22	A-TZ	
MW-38B	12/31/2003	728,319	3,165,945	45.51	37	25.5	35.5	B-TZ	
MW-39B	12/16/2003	728,424	3,166,019	49.58	40	29.5	39.5	B-TZ	

**TABLE A-1  
SUMMARY OF GROUNDWATER MONITORING WELLS  
HOUSTON, TX - WOOD PRESERVING WORKS**

WELL NO.	DATE INSTALLED	NORTHING	EASTING	TOP OF CASING ELEVATION (FT HVD)	TOTAL DEPTH (FT BGS)	Top Screen Interval (FT BGS)	Bottom Screen Interval (FT BGS)	Zone	Plugged
MW-40B	12/15/2004	728,341	3,166,122	49.59	40	29.5	39.5	B-TZ	
MW-41B	1/7/2003	728,176	3,166,003	49.37	40	29.5	39.5	B-TZ	
MW-42B	8/24/2006	728,257	3,166,324	50.52	42	30	40	B-TZ	
MW-44A	2/22/2007	729,021	3,168,349	45.11	28	18	28	A-TZ	
MW-44C	1/16/2004	729,021	3,168,349	45.13	70	57.5	67.5	C-TZ	
MW-45C	1/20/2004	729,155	3,168,512	44.73	70	58	68	C-TZ	
MW-46C	1/9/2004	729,121	3,168,576	44.94	72	60	70	C-TZ	
MW-47A	3/17/2020	728,644	3,168,558	43.92	26	16	26	A-TZ	
MW-47C	3/16/2007	728,725	3,168,535	45.52	71	61	71	C-TZ	
MW-48C	2/2/2004	728,417	3,168,241	44.68	72	60	70	C-TZ	
MW-49A	2/28/2007	728,345	3,168,191	46.18	30	20	30	A-TZ	
MW-49B	1/24/2009	728,375	3,168,184	46.43	35	30	35	B-CZ	
MW-50A	3/1/2007	727,501	3,167,958	46.96	25	15	25	A-TZ	
MW 50B	2/7/2020	727,585	3,167,950	45.89	40	34.5	39.5	B-TZ	
MW-51A	2/28/2007	726,925	3,166,885	47.80	25	15	25	A-TZ	
MW-51C	5/10/2014	726,935	3,166,894	47.48	80	62	72	C-TZ	
MW-52A	2/27/2007	728,699	3,167,814	51.91	30	20	30	A-TZ	P
MW-53C	8/15/2006	729,613	3,168,481	45.49	72	60	70	C-TZ	
MW 54B	2/22/2020	729,308	3,168,727	43.59	40	34.6	39.6	B-CZ	
MW-54C	8/15/2006	729,218	3,168,766	44.99	72	60	70	C-TZ	
MW-55A	1/12/2009	728,540	3,167,482	52.01	25	10	25	A-TZ	P
MW-55B	12/14/2011	728,538	3,167,474	49.15	40	32	37	B-TZ	P
MW-55C	8/22/2006	728,540	3,167,440	NA	30	NA	NA	C-TZ	P
MW-57A	1/22/2009	728,859	3,167,973	50.89*	27	12	27	A-TZ	
MW-57B	12/21/2011	728,857	3,167,965	47.93	40	34	39	B-TZ	
MW-58A	1/23/2009	728,875	3,168,176	47.76	29	14	29	A-TZ	
MW-59A	1/28/2009	728,155	3,168,358	44.18	21	11	21	A-TZ	
MW-59B	6/26/2010	728,145	3,168,358	44.36	33	28	33	B-CZ	
MW-59D	1/27/2009	728,114	3,168,365	44.22	118	108	118	D-TZ	
MW-60A	1/26/2009	728,825	3,168,823	46.79	28.5	18.5	28.5	A-TZ	
MW-60AR	3/17/2020	728,810	3,168,818	45.43	30	20	30	A-TZ	
MW-60B	2/27/2020	728,812	3,168,824	45.38	40	35	40	B-CZ	
MW-61A	1/26/2009	728,336	3,168,630	44.67	22	12	22	A-TZ	
MW-61B	2/23/2020	728,352	3,168,630	43.43	40	28	33	B-CZ	
MW-62B	1/21/2009	728,190	3,165,880	48.16	35	25	35	B-TZ	
MW-63B	1/28/2009	729,361	3,167,652	44.48	36	31	36	B-CZ	
MW-64A	1/26/2009	727,496	3,165,573	44.55	19.5	14.5	19.5	A-TZ	
MW-65D	1/17/2009	729,512	3,168,331	44.55	110	100	110	D-TZ	
MW-66D	1/20/2009	729,137	3,169,381	46.51	103	93	103	D-TZ	
MW-67B	6/26/2010	729,782	3,167,588	43.93	40	35	40	B-CZ	
MW-68A	5/21/2019	729,160	3,167,324	43.24	23.4	13	23	A-TZ	
MW-68B	12/15/2011	729,162	3,167,328	44.93	40	28	38	B-TZ	
MW-68C	6/25/2010	729,164	3,167,346	44.80	70	60	70	C-TZ	
MW-69A	6/23/2010	728,136	3,168,234	45.71	18.5	8.5	18.5	A-TZ	
MW-70B	12/14/2011	728,944	3,167,671	45.02	40	25	35	B-CZ	
MW 70C	2/21/2020	728,954	3,167,674	43.41	70	58	68	C-TZ	
MW-71B	12/13/2011	728,956	3,167,951	45.06	40	32	37	B-TZ	
MW-72B	12/21/2011	728,790	3,167,792	48.69	41	32	37	B-TZ	
MW-73B	12/13/2011	728,419	3,167,123	48.66	55	47	52	B-TZ	P
MW-74B	12/20/2011	728,373	3,167,718	47.83	40	26.5	36.5	B-TZ	
MW-75B	12/20/2011	728,066	3,168,022	47.18	40	32.2	37.2	B-TZ	P
MW 76B	2/12/2020	727,463	3,166,640	46.32	40	31	36	B-TZ	
MW-76C	5/7/2014	727,485	3,166,628	47.84	70	60	70	C-TZ	
MW-77A	5/7/2014	727,672	3,166,981	49.05	25	13	23	A-TZ	
MW-78A	5/6/2014	727,953	3,167,512	48.68	30	15	25	A-TZ	
MW-79A	5/7/2014	728,237	3,167,666	48.95	30	17	27	A-TZ	
MW-80B	5/8/2014	727,907	3,168,201	47.11	35	29	34	B-TZ	
MW-81B	5/11/2014	727,292	3,167,926	46.77	40	29	34	B-TZ	
MW-82B	1/22/2018	729,102	3,166,703	44.64	38	29.6	34.6	B-TZ	
MW-83B	1/30/2018	729,389	3,167,015	45.33	37	29.6	34.6	B-TZ	
MW-83C	2/2/2018	729,389	3,167,011	45.42	70	58.6	68.6	C-TZ	
MW 84A	2/19/2020	729,510	3,167,399	43.01	24	13.5	23.5	A-TZ	
MW-84B	1/23/2018	729,503	3,167,399	44.50	43	34.6	39.6	B-TZ	
MW-85C	1/20/2018	727,661	3,166,988	49.10	75	59.9	69.9	C-TZ	
MW-86C	1/18/2018	727,083	3,166,438	46.61	70	59.6	69.6	C-TZ	
MW-87C	1/31/2018	729,266	3,167,944	44.26	70	57.6	67.6	C-TZ	
MW-88A	2/22/2020	728,247	3,166,771	49.83	24	13	23	A-TZ	
MW-88B	2/19/2020	728,245	3,166,778	49.91	24	35	40	B-TZ	
MW-88C	1/21/2018	728,195	3,166,845	51.17	75	64.6	74.6	C-TZ	
MW-89B	7/12/2018	729,869	3,167,249	44.57	40.4	30	40	B-TZ	
MW-90B	7/13/2018	729,616	3,167,173	44.39	35.4	30	35	B-TZ	
MW-91A	2/19/2020	729,702	3,167,937	42.36	25	14.5	24.5	A-TZ	
MW-92B	2/20/2020	729,583	3,166,814	43.25	35.5	30	35	B-TZ	

**TABLE A-1  
SUMMARY OF GROUNDWATER MONITORING WELLS  
HOUSTON, TX - WOOD PRESERVING WORKS**

WELL NO.	DATE INSTALLED	NORTHING	EASTING	TOP OF CASING ELEVATION (FT HVD)	TOTAL DEPTH (FT BGS)	Top Screen Interval (FT BGS)	Bottom Screen Interval (FT BGS)	Zone	Plugged
MW-93B	2/26/2020	729,451	3,166,744	43.39	35.5	30	35	B-TZ	
MW-94A	2/18/2020	729,052	3,166,533	43.55	22	11.5	21.5	A-TZ	
MW-95A	2/19/2020	728,564	3,165,893	44.53	21	15	25	A-TZ	
MW-96B	2/24/2020	727,929	3,165,446	45.36	22	27	37	B-TZ	
MW-97A	2/13/2020	727,373	3,166,767	46.11	21	10.5	20.5	A-TZ	
MW-98A	2/10/2020	727,350	3,167,312	46.69	21	10.5	20.5	A-TZ	
MW-98B	2/11/2020	727,359	3,167,306	46.93	40.5	35	40	B-TZ	
MW-99C	2/23/2020	729,500	3,168,707	43.67	70	59.5	69.5	C-TZ	
TW-03	3/12/2007	727,734	3,167,007	NA	25	14	24	NA	
TW-41B	1/22/2009	728,222	3,166,002	49.67	40	30	40	B-TZ	
TW-56A	1/23/2009	728,758	3,168,070	51.89	31	21	31	A-TZ	P
GB-1	1/14/2009	729,146	3,169,447	NA	135	NA	NA	NA	
P-11	3/25/1991	728,049	3,166,025	48.98	50	36.2	38.2	B-TZ	

**Notes:**

1 - Point of Compliance Wells for SWMU No. 1 (resurveyed every 5 years)

BGS=Below Ground Surface

HVD = Elevations relative to Houston Vertical Datum, Houston Monument System

Northing/Easting = Coordinates based on NAD 1927 Texas State Plane, South Central Zone, US Survey Feet

\* Wells were resurveyed after casings were modified from above ground to flush grade during soil cap construction (2016).



**LOG OF BORING:  
MW-47A**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX		Drilling Company: Cascade Drilling	Borehole Diameter (in.): 8	8	
Project No. 19119232		Completion Date: 3/17/2020	Total Depth (ft): 26	26	
		Driller: Davis Ocana	Water Level (ft bgl): --	--	
		Driller's License: TX 59380	Northing: 728644.1	728644.1	
		Logged By: K. Worley	Easting: 3168558.3	3168558.3	
		Drilling Method: HSA	TOC Elev. (ft AMSL): 43.92	43.92	
		Sampling Method: 2" x 5' Core	Ground Elev. (ft AMSL): 43.80	43.80	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 15) No Recovery - Hydrovac to 15.0'
2					
4					
6					
8		0.0/15.0		NR	
10					
12					
14					
16		5.0/5.0	0.9	SP	(15 - 24.8) SAND, SP, tan to gray, fine, poorly sorted, soft, unconsolidated, interbedded tan silty clay lenses, no odor, color change to gray at 17.3', fine to medium grained at 17.9', interbedded gray fine sand and brown medium sand at 24.2'
18			0.7		
20			0.0		
22		5.0/5.0	0.1		
24			0.0		
26		1.0/1.0	0.0	CL	(24.8 - 26) Sandy, Silty, CLAY, CL, brown to reddish brown, medium density, wet, no odor, clay content increases with depth.

**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 15.0' drilled out with a hydrovac to clear for utilities.

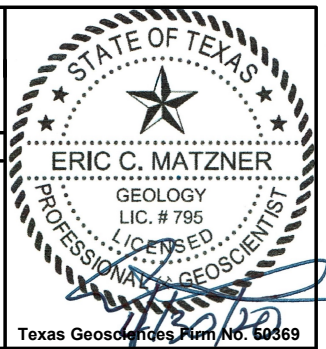
**Casing Materials**  
(0.0 - 16.0) 2" SCH-40 PVC Casing  
(16.0 - 26.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(26.0 - 26.5) End Cap

**Annular Materials**  
(0.0 - 1.0) Concrete  
(1.0 - 14.0) 3/8" Bentonite Chips  
(14.0 - 26.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-50B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/7/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Joey Lester	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	727585.1	
	Logged By:	K. Worley	Easting:	3167950.5	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	45.89	
	Sampling Method:	4" x 10' Core	Ground Elev. (ft AMSL):	46.16	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No Recovery - Hydrovac to 10.0'.
2					
4					
6					
8					
10		0.0/10.0		NR	
12			0.0		(10 - 14.8) Sandy CLAY, CL, tan with gray mottling, 20% fine sand, high plasticity clay, medium stiff to firm, slight odor.
14			0.0	CL	
16		10.0/10.0	0.0	SC	(14.8 - 16.5) Clayey SAND, SC, light brown to gray, soft, fine to very fine sand, sand content increases with depth, no odor, wet.
18			0.0		(16.5 - 22) Sand, SP, gray, poorly sorted, fine grained, unconsolidated, no odor, wet.
20			0.0	SP	
22		5.0/5.0	0.0	SC	(22 - 23.2) Sandy CLAY, SC, gray to brown, medium plasticity.
24			0.0	CL	(23.2 - 25) CLAY, CL, gray to brown with orange mottling, stiff, low plasticity, moist.
26			0.0		(25 - 40) Silty CLAY, CH, gray, becomes red brown at 27.2', stiff, medium plasticity at 31.2', no plasticity at 33', carbonate gravel abundant at 33' to 37.2', silty sand seam, moist, strong odor at 37.8' to 38.7'
28			0.0		
30		10.0/10.0	0.0		
32			0.0	CH	
34			0.0		
36		5.0/5.0	0.0		
38			0.0		
40			0.0		



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 27.0'. Removed during grouting.

**Casing Materials**

(0.0 - 34.5) 2" SCH-40 PVC Casing  
(34.5 - 39.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(39.5 - 40.0) End Cap

**Annular Materials**

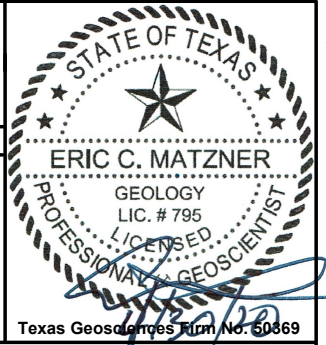
(0.0 - 2.0) Concrete  
(2.0 - 26.0) Grout  
(26.0 - 32.0) 3/8" Bentonite Chips  
(32.0 - 40.0) 16/30 Silica Sand



LOG OF BORING:  
**MW-54B**



# Union Pacific Railroad



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/22/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Joey Lester	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	729308	
	Logged By:	M. Hermiston	Easting:	3168726.9	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.59	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	--	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No Recovery - Hydrovac to 10.0'.
2					
4		0.0/10.0		NR	
6					
8					
10					
12		2.5/2.5	0.0	CL	(10 - 17.8) Silty CLAY, CL, light gray GLEY1 7/ with yellow 2.5Y 7/8 mottling, black staining, moist, no odor, white beds/mottling throughout 11.5' to 13.5', yellowish reddish 5YR 5/8 with reddish brown 5YR 5/4 mottling to 15.2', calcareous nodules abundant from 16' to 17.8', moderately firm, medium plasticity
14		2.5/2.5	0.0		
16		2.5/2.5	0.0		
18		2.5/2.5	0.0		
20		2.5/2.5	0.0	ML/CL	(17.8 - 20.2) Clayey SILT/ silty CLAY, ML/CL, light gray GLEY1 7/ with reddish brown 5YR 5/4 mottling, trace yellow 2.5Y 7/8 mottling, moist, no odor, cobble (2.4"x1"x1") at 19.8', sand content increases with depth.
22		2.5/2.5	0.0	ML/SC	(20.2 - 23.8) Sandy, clayey SILT / Clayey, silty SAND, ML/SC, reddish brown with some gray pockets, sand is very fine, wet, interbedded silty clay beds.
24		2.5/2.5	0.0		
26		2.5/2.5	0.0	CL	(23.8 - 25) Silty CLAY, CL, reddish brown 5YR 5/4, very firm, some calcareous nodules present.
28		2.5/2.5	0.0	SM	(25 - 27.8) Silty SAND, SM, reddish brown 5YR 5/4, very fine grained, wet.
30		2.5/2.5	0.0		
32		5.0/5.0		CL	(27.8 - 30.2) Silty CLAY, CL, reddish brown 5YR 5/4 and red 2.5 YR 4/8, thin sand lenses present.
34					
36		5.0/5.0			
38					(30.2 - 40) Silty CLAY, CL, light gray GLEY1 7/ with yellow 2.5Y 7/8 mottling, silt content decreases with depth, firm to hard, trace calcareous nodules throughout 32.4' to 40', calcareous gravel near 36', red mottling from 38' to 40'.
40					



## GOLDER

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 7" Sonic isolation casing to 25'. Removed during grouting.

**Casing Materials**

(0.0 - 34.5) 2" SCH-40 PVC Casing  
(34.5 - 39.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(39.5 - 40.0) End Cap

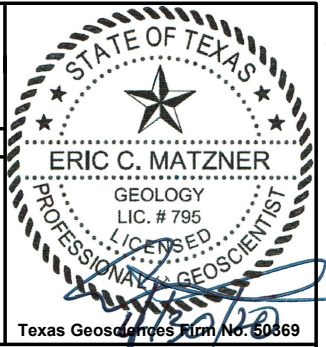
**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 30.0) Grout  
(30.0 - 33.0) 3/8" Bentonite Chips  
(33.0 - 40.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-60AR**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX		Drilling Company: Cascade Drilling	Borehole Diameter (in.): 8	Total Depth (ft): 30	
Project No. 19119232		Completion Date: 3/17/2020	Water Level (ft bgl): --	Northing: 728810.3	
		Driller: Davis Ocana	Eastings: 3168817.9	TOC Elev. (ft AMSL): 45.43	
		Driller's License: TX 59380	Ground Elev. (ft AMSL): 45.42		
		Logged By: K. Worley			
		Drilling Method: HSA			
		Sampling Method: 2" x 5' Core			

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 15) No Recovery - Hydrovac to 15.0'.
2					
4					
6					
8		0.0/15.0		NR	
10					
12					
14					
16			0.5	SC	(15 - 16.3) Clayey SAND, SC, gray, very fine to fine grained, saturated, no odor, soft.
18		5.0/5.0	0.9		(16.3 - 29.4) SAND, SP, greenish gray, poorly sorted, fine grained, wet, no odor, fine to medium grained, gray at 19.2' to 19.5', begins grading to reddish brown at 20.0', saturated at 22', brown at 27.3'.
20			1.0		
22		5.0/5.0	0.8	SP	
24			0.8		
26			0.6		
28		5.0/5.0	0.3		
30				CL	(29.4 - 30) Silty CLAY, CL, reddish brown, firm, moist, no odor.



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 15.0' drilled out with a hydrovac to clear for utilities.

**Casing Materials**

(0.0 - 20.0) 2" SCH-40 PVC Casing  
(20.0 - 30.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(30.0 - 30.5) End Cap

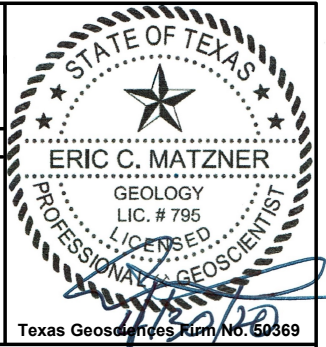
**Annular Materials**

(0.0 - 1.0) Concrete  
(1.0 - 16.0) 3/8" Bentonite Chips  
(16.0 - 30.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-60B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/27/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	728811.6	
	Logged By:	S. Ruiz	Easting:	3168823.6	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	45.38	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	45.64	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 11.2) No Recovery - Hydrovac to 10.0'.
2		0.0/10.0		NR	
4					
6		0.0/5.0		CL	(11.2 - 20.9) Sandy CLAY, CL, bluish gray with yellow-orange mottle, low plasticity, firm, wet. Abundant black streaks, very stiff from 18.1' - 18.8'.
8					
10		8.8/10.0	0.1	CL	
12			0.2		
14			0.3		
16		0.3		SP	(20.9 - 24.2) SAND, SP, fine, poorly sorted, light gray with brown streaks, wet, very soft.
18					
20		0.2		SM	(24.2 - 27.6) Silty SAND, SM, brown with light gray, wet, soft, slow dilatancy. Interbedded silt and clay lenses throughout, localized clay lenses with medium plasticity.
22					
24		10.0/10.0	0.1	CL	(27.6 - 40) Silty CLAY, CL, reddish brown with bluish gray mottle, some yellow-orange and black streaks, moist, firm to very firm, medium plasticity, laminated. Less cohesive from 30.0' - 32.5'. Color change to light brown with light gray mottle with brown-red streaks below 34.5'. Localized black gravel pockets, small to medium size, at 34.5', 35.1', 37.8', and 39.4'. Trace calcareous nodules below 38.0'.
26			0.2		
28			0.2		
30		5.0/5.0	0.3	CL	
32			0.2		
34			0.2		
36					
38					
40					



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 15.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 30.0'. Removed during grouting.

**Casing Materials**

(0.0 - 35.0) 2" SCH-40 PVC Casing  
(35.0 - 40.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(40.0 - 40.5) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 30.0) Grout  
(30.0 - 33.0) 3/8" Bentonite Chips  
(33.0 - 40.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-61B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/23/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Joey Lester	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	728352.3	
	Logged By:	M. Hermiston	Easting:	3168629.6	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.43	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	43.58	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No Recovery - Hydrovac to 10.0'.
2					
4					
6					
8					
10		0.0/10.0		NR	
12					
14		5.0/5.0	1.7	CL	(10 - 14) Sandy Silty CLAY, CL, light gray 7.5YR 7/1 with yellow 10YR 7/8 mottling and pinkish gray 7.5YR 7/2, no odor, sand content increases with depth.
16			0.4		
18					
20					
22					
24					
26					
28					
30					
32					
34					
36					
38					
40					



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 7.0' sonic isolation casing to 25.0'. Removed during grouting.  
3.) Photoionization detector (PID) readings were inconsistent.

**Casing Materials**  
(0.0 - 28.0) 2" SCH-40 PVC Casing  
(28.0 - 33.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(33.0 - 33.5) End Cap

**Annular Materials**  
(0.0 - 2.0) Concrete  
(2.0 - 22.5) Grout  
(22.5 - 26.0) 3/8" Bentonite Chips  
(26.0 - 33.0) 16/30 Silica Sand

LOG OF BORING:  
**MW-70C**



# Union Pacific Railroad



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	8	
	Completion Date:	2/21/2020	Total Depth (ft):	70	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	728954.390	
	Logged By:	S. Ruiz	Easting:	3167673.513	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.41	
	Sampling Method:	6" & 4" x 10' Core	Ground Elev. (ft AMSL):	--	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No Recovery - Hydrovac to 10.0'.
5		0.0/10.0		NR	
10			1.7		(10 - 15.7) Fine SAND, SP, bluish gray with light gray with localized brown discoloration, wet, very soft, rapid dilatancy.
15		10.0/10.0	2.2	SP	
			2.3	CL	(15.7 - 16.5) Sandy CLAY, CL, bluish gray with red-brown mottle, moist, firm, medium plasticity.
20			2.5		(16.5 - 22.9) Silty SAND, SM, light brown with bluish gray, wet, very soft, slow dilatancy, trace iron (Fe) staining at 17.5', odor.
			3.3	SM	
			7.4		
25		7.0/7.0	16.3	CH	(22.9 - 24.3) Silty CLAY, CH, red-brown with light gray mottle, dry, very firm, high plasticity.
			13.2		(24.3 - 37) CLAY, CL, light gray with light brown mottle, dry, very firm, medium plasticity, strong odor. Color transitions to tan with light gray mottle, streaks of black, red-brown, and yellow-orange, then to red-brown with light gray mottle below 32.0'. Strong odor throughout. Large calcareous gravel layers at 32.8' - 34.5', with thinner gravel lenses starting at 29.5' to 33.0'. Gravel lenses are moist and stained, with visible sheen on clay between 27.0' - 29.0'. Clay is very moist below 30.0', with thin, interbedded, stained silt lenses.
30		10.0/10.0	201.3	CL	
			234.2		
			42.3		
			21.3		
35			362.5		
40		7.0/7.0	59.5	CH	(37 - 45) CLAY, CH, red-brown with bluish gray mottle, moist, very firm and stiff, medium to high plasticity, trace black organic staining, slight odor.
			6.2		
			5.1		
			6.9		
45		5.0/5.0	12.0	ML	(45 - 46.5) SILT, ML, red-brown silt with interbedded clay nodules, moist, very soft, low plasticity.
			10.1		(46.5 - 58.6) Silty CLAY, CH, red-brown with trace black streaks, dry, very firm to hard, high plasticity, strong cementation, concretions from 47.6' - 47.8'. Calcareous lenses below 55.8'.
50			4.2		



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 12.0" Sonic isolation casing to 50.0'. Removed during grouting of 7.0" permanent casing. 7.0" permanent isolation casing to 52.0'.

**Casing Materials**

(2.0 - 52.0) 7" Permanent Steel Isolation Casing  
(0.0 - 58.0) 2" SCH-40 PVC Casing  
(58.0 - 68.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(68.0 - 68.5) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 50.0) Grout  
(50.0 - 58.0) 3/8" Bentonite Chips  
(58.0 - 68.0) 16/30 Silica Sand



**LOG OF BORING:  
MW-70C**



**Union Pacific Railroad**

**SITE INFORMATION**

**PROJECT AND DRILLING INFORMATION**

Houston Wood Preserving Works  
Houston, TX  
  
Project No. 19119232

Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	8
Completion Date:	2/21/2020	Total Depth (ft):	70
Driller:	Jesse Van Kirk	Water Level (ft bgl):	--
Driller's License:	TX 59618M	Northing:	728954.390
Logged By:	S. Ruiz	Easting:	3167673.513
Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.41
Sampling Method:	6" & 4" x 10' Core	Ground Elev. (ft AMSL):	--

Texas Geosciences Firm No. 50369

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
55		10.0/10.0	3.2	CH	
			1.8		
			2.1		
			3.4		
60		10.0/10.0	1.4	ML	(58.6 - 59.3) SILT, ML, red-brown silt, wet, very soft, low plasticity, increasing clay content with depth, odor.
			10.0	CH	(59.3 - 62.3) Silty CLAY, CH, red-brown, wet, firm, high plasticity, silt seams throughout.
65		10.0/10.0	23.5	SM	(62.3 - 67.5) Silty SAND, SM, red-brown, very fine, wet, very soft, slow dilatancy, trace black staining and odor throughout. Clay lenses from 64.8' - 65.5', small concretions with small to medium gravel from 65.5' - 66.1'.
			4.8	SM	
			2.0	SM	
70			1.6	CH	(67.5 - 70) Silty CLAY, CH, red-brown with black streaks, moist, very firm to hard, high plasticity, laminated, trace silt seams.



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 12.0" Sonic isolation casing to 50.0'. Removed during grouting of 7.0" permanent casing. 7.0" permanent isolation casing to 52.0'.

**Casing Materials**

(2.0 - 52.0) 7" Permanent Steel Isolation Casing  
(0.0 - 58.0) 2" SCH-40 PVC Casing  
(58.0 - 68.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(68.0 - 68.5) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 50.0) Grout  
(50.0 - 58.0) 3/8" Bentonite Chips  
(58.0 - 68.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-76B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/12/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	727463.2	
	Logged By:	S. Ruiz	Easting:	3166640.1	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	46.32	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	46.69	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'.
2		0.0/0.0		NR	
4					
6					
8					
10		12.5/15.0		CL	(10 - 19.9) Silty CLAY, CL, gray with yellow-orange mottle, moist, firm to very firm, medium plasticity. Calcareous lens from 12.9' - 13.3', black organic staining above 13.7', trace red mottle. Thin fine sand seams from 15.7' - 16.9'. Increasing orange mottle with depth.
12			0.2		
14			0.4		
16			0.4		
18					
20		10.0/10.0		SC/SM	(19.9 - 21.5) Silty SAND, SM/SC, tan with red-brown, fine to very fine with clay nodules throughout, some light gray, wet, soft, slow dilatancy, low plasticity. (21.5 - 40) Silty CLAY, CL, light gray with orange mottle, trace fine sand, moist, firm to very firm, medium to high plasticity. Increasing red-brown mottle with depth. Wet calcareous gravel, small to medium, mixed with silt layers from 29.2' - 29.8', 32.3' - 32.5', 32.9' - 33.2', 34.8' - 35.1'. Clay is dryer below 35', with dry calcareous nodule layers at 36.9' - 37.4' and 38.9' - 39.3'. Color change to red-brown with light gray mottle below 35.5, with increasing firmness, becoming very hard below 36.0'.
22			0.5		
24			0.4		
26			0.5		
28					
30		5.0/5.0		CL	
32			1.0		
34			0.9		
36			1.1		
38					
40					



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 27.0'. Removed during grouting.

**Casing Materials**

(0.0 - 31.0) 2" SCH-40 PVC Casing  
(31.0 - 36.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(36.0 - 36.5) End Cap

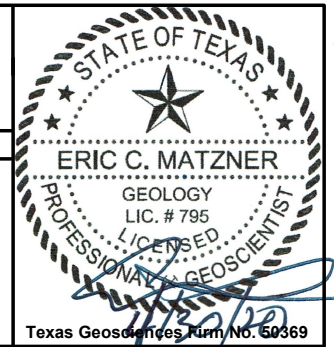
**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 26.0) Grout  
(26.0 - 29.0) 3/8" Bentonite Chips  
(29.0 - 36.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-84A**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX		Drilling Company: Cascade Drilling	Borehole Diameter (in.): 6	6	
Project No. 19119232		Completion Date: 2/19/2020	Total Depth (ft):	25	
		Driller: Jesse Van Kirk	Water Level (ft bgl):	--	
		Driller's License: TX 59618M	Northing:	729510.3	
		Logged By: S. Ruiz	Easting:	3167398.8	
		Drilling Method: Sonic	TOC Elev. (ft AMSL):	43.01	
		Sampling Method: 6" x 10' Core	Ground Elev. (ft AMSL):	--	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2					
4		0.0/10.0		NR	
6					
8					
10					(10 - 11.7) No recovery - lost
12					
14		8.3/10.0	0.8	SM	(11.7 - 18.3) Silty SAND, SM, light gray, very fine with clay nodules throughout, trace iron (Fe) staining at 13', wet, soft, slow dilatancy, low plasticity.
16			0.8		
18			0.6		
20			0.8	CL	(18.3 - 18.8) Sandy CLAY, CL, bluish gray with reddish brown mottle, wet, firm, medium plasticity.
22					(18.8 - 23.9) Silty SAND, SM, light gray, very fine, wet.
24		5.0/5.0	0.6	SM	
26			0.4	CL	(23.9 - 25) CLAY, CL, reddish brown with light gray mottle, moist, firm to very firm, medium to high plasticity.



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.

**Casing Materials**

(0.0 - 13.5) 2" SCH-40 PVC Casing  
(13.5 - 23.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(23.5 - 24) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 11.5) 3/8" Bentonite Chips  
(11.5 - 24) 16/30 Silica Sand



**LOG OF BORING:  
MW-88A**



**Union Pacific Railroad**



**SITE INFORMATION**

Houston Wood Preserving Works  
Houston, TX

Project No. 19119232

**PROJECT AND DRILLING INFORMATION**

Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6
Completion Date:	2/22/2020	Total Depth (ft):	25
Driller:	Joey Lester	Water Level (ft bgl):	--
Driller's License:	TX 59618M	Northing:	728247.4
Logged By:	M. Hermiston	Easting:	3166771
Drilling Method:	Sonic	TOC Elev. (ft AMSL):	49.83
Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	46.95

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2					
4					
6					
8					
10		0.0/10.0		NR	
12			0.0	CL	(10 - 15) Silty CLAY, CL, gray 7.5 YR 7/1 / greenish gray GLEY1 6/1 with yellow 2.5Y 7/8 mottling from 10' to 12.5', no mottling from 12.5' to 15', firm, moist to wet, no odor, med plasticity.
14		5.0/5.0	0.2		
16			0.0	ML	(15 - 17.8) Clayey SILT, ML, light bluish gray GLEY2 7/1, interbedded clay, mostly silt, wet, no odor, grades into sand unit below.
18		5.0/5.0	0.0		
20			0.0	SM	(17.8 - 21.8) Silty SAND, SM, light gray 7.5YR 7/1, wet
22		2.0/2.0	0.2		
24			0.2	CL	(21.8 - 25) Silty CLAY, CL, light gray with yellow mottling, moderate plasticity.
26		3.0/3.0	0.2		



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.

**Casing Materials**

(+2.9 - 13.0) 2" SCH-40 PVC Casing  
(13.0 - 23.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(23.0 - 23.5) End Cap

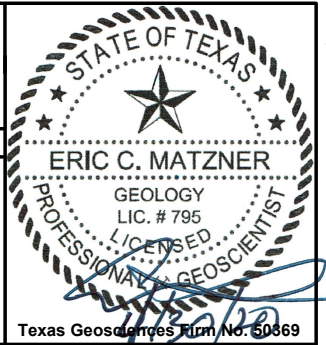
**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 11.0) 3/8" Bentonite Chips  
(11.0 - 23.0) 16/30 Silica Sand

LOG OF BORING:  
**MW-88B**



# Union Pacific Railroad



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/19/2020	Total Depth (ft):	45	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	728245	
	Logged By:	S. Ruiz	Easting:	3166777.6	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	49.91	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	46.86	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
4		0.0/10.0		NR	
8					
12			1.7	CL	(10 - 16.1) Silty CLAY, CL, bluish gray with reddish brown, orange yellow streaks with green and black streaks, moist, weak cementation, firm, medium plasticity, soft below 14' silt content increases with depth.
			1.4		
16		15.0/15.0	1.4	ML	(16.1 - 16.8) SILT, ML, light blueish gray with orange and some black organic staining, some clay, moist, soft, low plasticity.
			1.4	SM	
20			1.2	CH	(16.8 - 20.3) Silty SAND, SM, light gray with yellow and orange streaks, with blueish gray, at 18.5 transitions to light gray, wet, moderate cementation, very firm, high plasticity.
			1.4	ML/CL	
24			1.5		(20.3 - 20.7) Silty CLAY, CH, bluish gray clay with reddish brown streaks, clear boundary, moist, very firm, high plasticity.
			0.6		
28		5.0/5.0	0.8		(20.7 - 21.5) SILT with clay nodules, ML/CL, light gray, moist, very soft, medium to high plasticity.
			0.5		
32			0.6	CL	(21.5 - 40) Silty CLAY, CL, light grey with yellow streaks and reddish brown streaks below 24', moist, very soft, medium to high plasticity. Color transitions from light gray to red brown, with interbedded moist, soft, silt seams below 30'.
36		10.0/10.0	0.6		
40					(40 - 45) Silty CLAY, CH, reddish brown, with silt lenses, light gray mottle, calcareous nodules and gravel mixed throughout, wet, very firm to hard, moderate plasticity.
44		5.0/5.0	--	CH	
48					



## GOLDER

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0' Sonic isolation casing to 30.0'. Removed during sanding.

**Casing Materials**

(0.0 - 35.0) 2" SCH-40 PVC Casing  
(35.0 - 40.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(40.0 - 40.5) End Cap

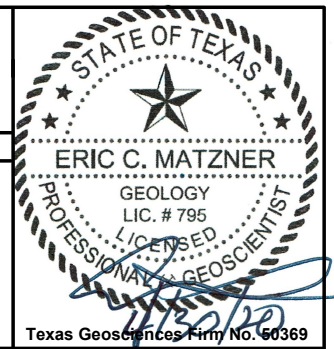
**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 30.0) Grout  
(30.0 - 33.0) 3/8" Bentonite Chips  
(33.0 - 40.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-91A**



**Union Pacific Railroad**



**SITE INFORMATION**

**PROJECT AND DRILLING INFORMATION**

Houston Wood Preserving Works  
Houston, TX

Drilling Company: Cascade Drilling  
Completion Date: 2/19/2020  
Driller: Jesse Van Kirk  
Driller's License: TX 59618M  
Logged By: S. Ruiz  
Drilling Method: Sonic  
Sampling Method: 6" x 10' Core

Borehole Diameter (in.): 6  
Total Depth (ft): 30  
Water Level (ft bgl): --  
Northing: 729702.2  
Easting: 3167936.8  
TOC Elev. (ft AMSL): 42.36  
Ground Elev. (ft AMSL): --

Project No. 19119232

Texas Geosciences Firm No. 50369

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2		0.0/10.0		NR	
4					
6					
8					
10					(10 - 10.7) No recovery - lost
12		6.3/7.0	1.3	CL	(10.7 - 17) Sandy CLAY, CL, light gray with reddish brown and yellow orange mottle with blueish streaks, moist, firm, medium to high plasticity, color change to light gray with just reddish brown mottle below 14', trace black streaks.
14			1.2	CL	
16			1.1	CL	
18		7.0/8.0	0.8	SM	(17 - 24.5) Silty SAND, SM, light orange brown with streaks of gray, trace clay nodules, color change to light brown below 21', with gray streaks, very fine, wet, very soft, rapid dilatancy.
20			0.8	SM	
22			1.0	SM	
24			0.9	SM	
26		5.0/5.0	1.5	CL	(24.5 - 30) Silty CLAY, CL, light brown with small light gray mottle, trace lenses of black staining at 26.5', moist, firm to very firm, medium to high plasticity.
28			--	CL	
30			--	CL	



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.

**Casing Materials**

(0.0 - 14.5) 2" SCH-40 PVC Casing  
(14.5 - 24.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(24.5 - 25) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 12.5) 3/8" Bentonite Chips  
(12.5 - 25) 16/30 Silica Sand

**LOG OF BORING:  
MW-92B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX		Drilling Company: Cascade Drilling	Borehole Diameter (in.): 6	6	
Project No. 19119232		Completion Date: 2/20/2020	Total Depth (ft): 40	40	
		Driller: Jesse Van Kirk	Water Level (ft bgl): --	--	
		Driller's License: TX 59618M	Northing: 729582.6	729582.6	
		Logged By: S. Ruiz	Easting: 3166813.6	3166813.6	
		Drilling Method: Sonic	TOC Elev. (ft AMSL): 43.25	43.25	
		Sampling Method: 6" x 10' Core	Ground Elev. (ft AMSL): --	--	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2					
4		0.0/10.0		NR	
6					
8					
10					(10 - 11.8) No recovery
12			0.7		
14			0.7	SP	(11.8 - 17.5) Fine SAND, SP, light gray with some Iron (Fe) staining, wet below 13', soft, low plasticity.
16			0.7		
18		13.2/15.0	0.7	CL	(17.5 - 19) Silty CLAY, CL, light gray with reddish brown mottle, trace calcareous nodules, clear boundary, moist, moderate cementation, very firm, medium plasticity.
20			6.7	SM	(19 - 21.9) Silty SAND, SM, light gray, wet, no cementation, very soft, rounded, no plasticity.
22			6.5		
24			6.6	CL	(21.9 - 26.8) Silty CLAY, CL, light gray with red mottle to 23', thin light gray with yellow mottling, clear boundary, sands increasing with depth, moist, moderate cementation, very firm, medium plasticity.
26			0.9		
28			0.8		
30			0.7	SM/SC	(26.8 - 33.2) Silty SAND Clayey SAND, SM/SC, tan with some light gray, moist to wet, weak cementation, soft, slow dilatancy, low plasticity.
32		15.0/15.0	1.0		
34			1.1		
36			1.0	CL	(33.2 - 40) CLAY, CL, brownish red with light brown, some blue streaks, moist, moderate cementation, very firm to hard, low to medium plasticity, calc gravel from 33.34' to 41.8' with trace calc nodules below.
38			0.6		
40			0.7		



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 25.0'. Removed during grouting.

**Casing Materials**

(0.0 - 30.0) 2" SCH-40 PVC Casing  
(30.0 - 35.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(35.0 - 35.5) End Cap

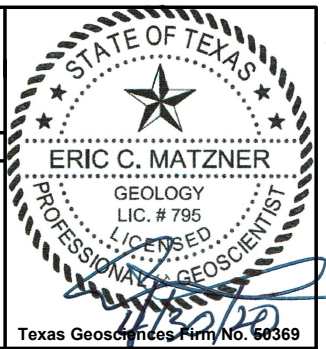
**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 24.0) Grout  
(24.0 - 27.0) 3/8" Bentonite Chips  
(27.0 - 35.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-93B**



**Union Pacific Railroad**



**SITE INFORMATION**

Houston Wood Preserving Works  
Houston, TX

Project No. 19119232

**PROJECT AND DRILLING INFORMATION**

Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6
Completion Date:	2/26/2020	Total Depth (ft):	40
Driller:	Jesse Van Kirk	Water Level (ft bgl):	--
Driller's License:	TX 59618M	Northing:	729451.5
Logged By:	S. Ruiz	Easting:	3166744.3
Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.39
Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	--

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2		0.0/10.0		NR	
4					
6					
8					
10		4.0/5.0		SP	(10 - 16.8) Fine SAND, SP, tan fine to medium grained, small unit of light gray sand at 14.3' to 14.6', wet, soft, subang, slow dilatancy.
12					
14					
16		8.0/8.0	0.3	SM	(16.8 - 21.6) Silty SAND, SM, light gray that transitions to bluish gray below 20.8', increasing clay content with depth, wet, weak cementation, very soft, slow dilatancy, low plasticity.
18			0.0		
20			0.1		
22			0.1	SC	(21.6 - 22.8) Clayey SAND, SC, light gray with tan, wet, weak cementation, soft, low plasticity.
24		2.0/2.0	0.1	CL	(22.8 - 26.8) Sandy CLAY, CL, bluish gray with tan mottle, sand grains above 25.5 are fine grained, light gray with tan and orange yellow mottle, moist, weak cementation, firm, low to medium plasticity.
26			0.2		
28		10.0/10.0	0.2	ML/SM	(26.8 - 31.5) SILT and Silty SAND, ML/SM, light gray with reddish brown, silty sand, clay nodules throughout, wet, weak cementation, soft, slow dilatancy, low plasticity.
30			0.0		
32			0.3		
34			0.2	CL	(31.5 - 34.8) Silty CLAY, CL, reddish brown with light gray, calc nodules from 32.5' to 34', some yellowish streaks, wet, moderate cementation, firm, subangular, high dry strength, medium plasticity.
36		0.1			
38		0.1			
40		5.0/5.0	0.0		(34.8 - 40) Silty Clay, CL, reddish brown with light bluish gray streaks, trace black streaks throughout, dry, strong cementation, very firm to hard, low plasticity.



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0' Sonic isolation casing to 25.0'. Removed during grouting.

**Casing Materials**

(0.0 - 30.0) 2" SCH-40 PVC Casing  
(30.0 - 35.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(35.0 - 35.5) End Cap

**Annular Materials**

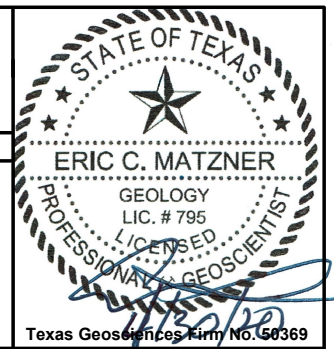
(0.0 - 2.0) Concrete  
(2.0 - 24.0) Grout  
(24.0 - 27.0) 3/8" Bentonite Chips  
(27.0 - 35.0) 16/30 Silica Sand



**LOG OF BORING:  
MW-94A**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/18/2020	Total Depth (ft):	22	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	729051.6	
	Logged By:	S. Ruiz	Easting:	3166533.1	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.55	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	--	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2					
4		0.0/10.0		NR	
6					
8					
10					(10 - 12) No recovery - lost
12		3.0/5.0	--	SC	(12 - 17) Clayey SAND, light gray with tan, fine grained, wet.
14					
16			--		
18		6.0/8.0	0.9	CL	(17 - 18.2) CLAY, CL, tan with light gray, mottled, moist firm, medium plasticity.
20			0.8	SM	(18.2 - 20.1) Silty SAND, SM, tan with light gray, mottled clay, moist, firm, medium plasticity.
22			0.8	CL	(20.1 - 20.8) CLAY, CL, tan with light gray, mottled, moist firm, medium plasticity.
				SM	(20.8 - 21.4) Silty SAND, SM, light gray, wet, soft, very fine grain.
				CL	(21.4 - 22) CLAY, CL, light gray, gravel, firm orange mottling with brownish red streaks below 22.5', dry, firm, medium plasticity.



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.

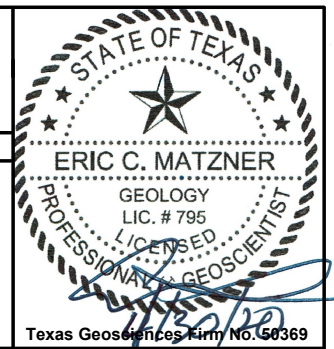
**Casing Materials**  
(0.0 - 11.5) 2" SCH-40 PVC Casing  
(11.5 - 21.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(21.5 - 22) End Cap

**Annular Materials**  
(0.0 - 2.0) Concrete  
(2.0 - 9.5) 3/8" Bentonite Chips  
(9.5 - 22) 16/30 Silica Sand

**LOG OF BORING:  
MW-95A**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX		Drilling Company: Cascade Drilling	Borehole Diameter (in.): 6	6	
Project No. 19119232		Completion Date: 2/19/2020	Total Depth (ft): 25	25	
		Driller: Joey Lester	Water Level (ft bgl): --	--	
		Driller's License: TX 59618M	Northing: 728563.9	728563.9	
		Logged By: K. Worley	Easting: 3165892.7	3165892.7	
		Drilling Method: Sonic	TOC Elev. (ft AMSL): 44.53	44.53	
		Sampling Method: 6" x 10' Core	Ground Elev. (ft AMSL): 45.57	45.57	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2					
4		0.0/10.0		NR	
6					
8					
10					
12		5.0/5.0	0.0	SC	(10 - 15.8) Clayey SAND, SC, gray to beige, orange mottling, iron staining, no odor, sand content increases with depth, moist.
14			0.0		
16			0.1		
18		5.0/5.0	0.0	SP	(15.8 - 21.5) SAND, SP, gray, poorly sorted, fine grained, unconsolidated, no odor, wet
20			0.0		
22		5.0/5.0	0.0	CL	(21.5 - 24.2) Sandy CLAY, CL, gray, orange mottling, medium plasticity, moist.
24			0.0		
26					(24.2 - 25) CLAY, CL, gray to brown, medium stiff, < 10% silt, no odor.



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 15.0'. Removed during sanding.

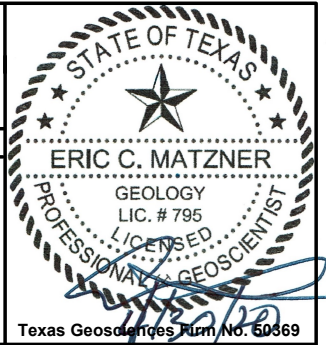
**Casing Materials**  
(0.0 - 15.0) 2" SCH-40 PVC Casing  
(15.0 - 25.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(25.0 - 25.5) End Cap

**Annular Materials**  
(0.0 - 2.0) Concrete  
(2.0 - 12.0) 3/8" Bentonite Chips  
(12.0 - 25.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-96B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/24/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Joey Lester	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	727928.5	
	Logged By:	M. Hermiston	Easting:	3165445.9	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	45.36	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	45.57	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION		
0					(0 - 10) No recovery - Hydrovac to 10.0'		
2							
4							
6							
8							
10		0.0/10.0		NR			
12			2.0	CL	(10 - 13.4) Silty CLAY, CL, light gray with reddish yellow mottling and black staining, trace white pockets, silt content increases with depth, no odor.		
14		5.0/5.0	6.3				
16			5.7	SM	(13.4 - 19.4) Silty SAND, SM, light gray with strong brown mottling from 16' to 19.4', no odor, wet.		
18		5.0/5.0	4.8				
20			4.3				
22			3.7	CL	(19.4 - 23) CLAY, CL, light gray with yellow mottling, becomes red mottling near 23' with black mottling near 23.5', more abundant red mottling with depth, no odor, firm, moderate plasticity.		
24		5.0/5.0	3.4				
26			4.4	CL	(23 - 37) Silty CLAY, CL, light gray with red mottling, calcareous nodules at 25.2', 28' to 29', 30.2' to 31.4', interbedded and friable clay from 33' to 34', moist from 35' to 37', some silt at 37', no odor.		
28		5.0/5.0	4.3				
30			3.6				
32		5.0/5.0	3.7				
34			5.7				
36		2.0/2.0	5.1				
38			4.7				
40		3.0/3.0	3.2				
							(37 - 40) CLAY, CL, red with some gray mottling, moist, very firm, no odor, moderate plasticity.



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 7.0' Sonic isolation casing to 25.0'. Removed during grouting.

**Casing Materials**

(0.0 - 27.0) 2" SCH-40 PVC Casing  
(27.0 - 37.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(37.0 - 37.5) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 22.0) Grout  
(22.0 - 25.0) 3/8" Bentonite Chips  
(25.0 - 37.0) 16/30 Silica Sand



**LOG OF BORING:  
MW-97A**



**Union Pacific Railroad**



**SITE INFORMATION**

**PROJECT AND DRILLING INFORMATION**

Houston Wood Preserving Works  
Houston, TX

Drilling Company: Cascade Drilling  
Completion Date: 2/13/2020  
Driller: Jesse Van Kirk  
Driller's License: TX 59618M  
Logged By: S. Ruiz  
Drilling Method: Sonic  
Sampling Method: 6" x 10' Core

Borehole Diameter (in.): 6  
Total Depth (ft): 23  
Water Level (ft bgl): --  
Northing: 727373  
Easting: 3166767.2  
TOC Elev. (ft AMSL): 46.21  
Ground Elev. (ft AMSL): 46.52

Project No. 19119232

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2					
4		0.0/10.0		NR	
6					
8					
10					(10 - 11.1) No recovery
12					(11.1 - 15.2) Sandy CLAY, CL, light gray with orange and yellow mottle and streaks of black, moist, firm, medium plasticity.
14		8.9/10.0	0.4	CL	
16			0.2	SC	(15.2 - 15.8) Clayey SAND, SC, light gray with yellow and orange mottling, trace black organic staining, wet, firm to soft, slow dilatancy, low plasticity.
18			0.3	CL	(15.8 - 17.8) CLAY, CL, light gray with reddish brown mottle and trace black organic staining, wet, very stiff.
20			0.4	SM/SC	(17.8 - 20.2) Silty SAND and Clayey SAND, SM/SC, light gray with tan, interbedded clay nodules, very fine, wet, no cementation, soft to very soft, slow dilatancy.
22		3.0/3.0	0.4	CL	(20.2 - 23) Sandy CLAY, CL, light gray, yellow and reddish brown mottle below 22', moist, very firm to stiff, medium to high plasticity.
24			0.4	CL	



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.

**Casing Materials**

(0.0 - 10.5) 2" SCH-40 PVC Casing  
(10.5 - 20.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(20.5 - 21) End Cap

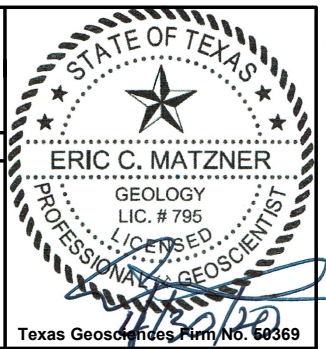
**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 8.5) 3/8" Bentonite Chips  
(8.5 - 21) 16/30 Silica Sand

**LOG OF BORING:  
MW-98A**



**Union Pacific Railroad**



**SITE INFORMATION**

**PROJECT AND DRILLING INFORMATION**

Houston Wood Preserving Works  
Houston, TX

Drilling Company: Cascade Drilling  
Completion Date: 2/10/2020  
Driller: Jesse Van Kirk  
Driller's License: TX 59618M  
Logged By: S. Ruiz  
Drilling Method: Sonic  
Sampling Method: 6" x 10' Core

Borehole Diameter (in.): 6  
Total Depth (ft): 25  
Water Level (ft bgl): --  
Northing: 727349.7  
Easting: 3167312  
TOC Elev. (ft AMSL): 46.69  
Ground Elev. (ft AMSL): 47.13

Project No. 19119232

Texas Geosciences Firm No. 59369

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 10) No recovery - Hydrovac to 10.0'
2		0.0/10.0		NR	
4					
6					
8					
10		10.0/10.0	0.1	CL	(10 - 19) CLAY, CL, light gray with tan and yellow orange mottle, abundant calcareous, color change to bluish gray below 15', gravel layer from 17.5' to 18.6', small to medium, moist, firm to very firm, medium to high plasticity.
12			0.0		
14			0.0		
16			0.2		
18			0.1		
20				SM	(19 - 20.6) Silty SAND, SM, light gray, gradual boundary, fine to very fine grain, wet, very soft, slow dilatancy.
22		5.0/5.0	0.1	CL	(20.6 - 25) Silty CLAY, CL, light gray, light brown mottling, increasing light brown color with depth, moist, very firm, medium to high plasticity.
24			0.0		
26			0.0		



**GOLDER**

2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

Notes: 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.

**Casing Materials**

(0.0 - 10.5) 2" SCH-40 PVC Casing  
(10.5 - 20.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(20.5 - 21) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 8.5) 3/8" Bentonite Chips  
(8.5 - 21) 16/30 Silica Sand

**LOG OF BORING:  
MW-98B**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/11/2020	Total Depth (ft):	40	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	727358.6	
	Logged By:	S. Ruiz	Easting:	3167306.4	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	46.93	
	Sampling Method:	6" x 10' Core	Ground Elev. (ft AMSL):	47.17	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0 - 10		0.0/10.0		NR	(0 - 10) No recovery - Hydrovac to 10.0'
10 - 14.4		8.0/10.0	0.3	CH	(10 - 14.4) CLAY, CH, gray with bluish gray and orange mottle, black organic staining from 15.0' to 15.5', small sand seams, moist, firm, medium to high plasticity.
14.4 - 16.8			0.3	SC	(14.4 - 16.8) Clayey SAND, SC, light gray with bluish gray and orange mottle, fine grain sand, moist, firm to soft, gradual boundary, low to medium plasticity.
16.8 - 21.1			0.2	SM	(16.8 - 21.1) Silty SAND, SM, light gray with light brown mottle above 17', clay nodules above 17.6', fine grain sand, wet at 17.6', very moist to wet, soft to very soft, slow dilatancy, diffused boundary, low plasticity.
21.1 - 33.8		5.0/5.0	0.6	CH	(21.1 - 33.8) CLAY, CH, gray with bluish gray and orange mottle, black organic staining from 15.0' to 15.5', small sand seams, calcareous nodules, moist, firm, medium to high plasticity.
33.8 - 34.7			0.2		
34.7 - 40		10.0/10.0	0.0	CH	(33.8 - 34.7) Clayey SAND, SC, gray with red below 31', calcareous nodules at 33', moist, soft, low to medium plasticity. (34.7 - 40) CLAY, CL, red to gray, color change to reddish brown below 34', 32.1' to 32.9' clayey sand, calcareous nodules at 33.3' with sandy clay below, silty sand seam at 39.3' to 39.61', dry, very hard, medium to high plasticity.
24 - 28			0.1		
28 - 32			0.7		
32 - 34			0.8		
34 - 36			1.1	CL	
36 - 38			1.3		
38 - 40			1.2		



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0' Sonic isolation casing to 28.0'. Removed during grouting.

**Casing Materials**

(0.0 - 35.0) 2" SCH-40 PVC Casing  
(35.0 - 40.0) 2" SCH-40 PVC 0.010" Slotted Screen  
(40.0 - 40.5) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 28.0) Grout  
(28.0 - 31.0) 3/8" Bentonite Chips  
(21.0 - 40.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-99C**



**Union Pacific Railroad**



SITE INFORMATION		PROJECT AND DRILLING INFORMATION			
Houston Wood Preserving Works Houston, TX	Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6	
	Completion Date:	2/23/2020	Total Depth (ft):	70	
Project No. 19119232	Driller:	Jesse Van Kirk	Water Level (ft bgl):	--	
	Driller's License:	TX 59618M	Northing:	729500.354	
	Logged By:	S. Ruiz	Easting:	3168707.132	
	Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.67	
	Sampling Method:	6" & 4" x 10' Core	Ground Elev. (ft AMSL):	--	

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
0					(0 - 11.5) No recovery - Hydrovac to 10.0'
5		0.0/10.0		NR	
10					
11.5			0.7	CL	(11.5 - 15.9) CLAY, CL, dark gray with greenish gray, calcareous nodules throughout, color change to light gray with reddish brown mottle below 18', some black streaks, increasing sand with depth, moist, weak cementation, soft, subangular grains, high plasticity.
13			0.1		
15		8.5/10.0	0.1		
15.9			0.0	SM	(15.9 - 22.5) Silty SAND, SM, clayey sand from 15.9 to 16.2, brown sand with lite gray interbedded clay lenses, wet, soft, slow dilatancy, low plasticity.
20			0.0		
22.5			0.0	CL	(22.5 - 23.6) CLAY, CL, silty sand, brown with light gray, calc nodules, thin silt seams, moist to wet, firm, low plasticity.
23.6			0.0		
23.6			0.3	SM	(23.6 - 26.1) Silty SAND, SM, brown with light gray above 24.5', very fine grain, wet, weak cementation, slow dilatancy.
26.1		10.0/10.0	0.3		
26.1			0.3	CH	(26.1 - 56.1) Silty CLAY, CH, reddish brown and gray with orange and yellow mottling below 30.51, abundant calcareous nodules below 35.51, some black streaks at 38.5' with gravel, orange and yellow mottle becomes reddish brown below 38', trace calcareous with localized black staining around black organic concentrations, dry silt seam with calcareous nodules from 52.6' to 54.8', dry, moderate cementation, firm, subangular grains, med dry strength.
30			0.2		
35			0.5		
35		10.0/10.0	0.7		
40			0.3		
40			0.4		
45			0.6		
45			0.0		
50			0.0		
50		15.0/15.0	0.1		
50			0.0		



**GOLDER**  
2201 Double Creek Dr., Suite 4004  
Round Rock, Texas 78664  
O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 30.0', 7" Sonic isolation casing to 45.0'. Both removed during grouting.

Casing Materials	Annular Materials
(0.0 - 59.5) 2" SCH-40 PVC Casing	(0.0 - 2.0) Concrete
(59.5 - 69.5) 2" SCH-40 PVC 0.010" Slotted Screen	(2.0 - 51.0) Grout
(69.5 - 70.0) End Cap	(51.0 - 56.0) 3/8" Bentonite Chips
	(56.0 - 70.0) 16/30 Silica Sand

**LOG OF BORING:  
MW-99C**



**Union Pacific Railroad**

**SITE INFORMATION**

**PROJECT AND DRILLING INFORMATION**

Houston Wood Preserving Works  
Houston, TX  
  
Project No. 19119232

Drilling Company:	Cascade Drilling	Borehole Diameter (in.):	6
Completion Date:	2/23/2020	Total Depth (ft):	70
Driller:	Jesse Van Kirk	Water Level (ft bgl):	--
Driller's License:	TX 59618M	Northing:	729500.354
Logged By:	S. Ruiz	Easting:	3168707.132
Drilling Method:	Sonic	TOC Elev. (ft AMSL):	43.67
Sampling Method:	6" & 4" x 10' Core	Ground Elev. (ft AMSL):	--

Texas Geosciences Firm No. 50369

Depth (ft)	Well Materials	Recovery (ft/ft)	PID (ppm)	USCS	LITHOLOGIC DESCRIPTION
55			0.0		
			0.3		
			0.1		
			0.3		(56.1 - 68.8) Silty SAND, SM, red, very fine grain, poorly sorted, clay nodules with calcareous nodules from 57' to 58', clay lenses from 60' to 61' with silt seams, clay nodules, calcareous nodules, wet, very soft
60		15.0/15.0	0.2	SM	
			0.5		
			0.5		
65			0.3		
			0.0		
70			0.4	CH	(68.8 - 70) Silty CLAY, CH, red with trace black streaks, moist, moderate cementation, very firm to hard, med toughness, high plasticity.



**GOLDER**

2201 Double Creek Dr., Suite 4004  
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O-512.671.3434 F-512.671.3446

**Notes:** 1.) Top 10.0' drilled out with a hydrovac to clear for utilities.  
2.) 8.0" Sonic isolation casing to 30.0', 7" Sonic isolation casing to 45.0'. Both removed during grouting.

**Casing Materials**

(0.0 - 59.5) 2" SCH-40 PVC Casing  
(59.5 - 69.5) 2" SCH-40 PVC 0.010" Slotted Screen  
(69.5 - 70.0) End Cap

**Annular Materials**

(0.0 - 2.0) Concrete  
(2.0 - 51.0) Grout  
(51.0 - 56.0) 3/8" Bentonite Chips  
(56.0 - 70.0) 16/30 Silica Sand

**ATTACHMENT B**

**Data Usability Summary and  
Analytical Reports from July 2019  
through March 2020 Site-Wide  
Sampling Events**

**JULY 2019**



# Memorandum

August 22, 2019

To: Eric Matzner Ref. No.: 11183954-1620

From: <sup>CK</sup> Chris G. Knight/eew/353-NF Tel: 512-506-8803

cc: Jesse Orth, Jon Lang; Julie Lidstone

**Subject: Data Usability Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works  
Houston, Texas  
July 2019**

## 1. Scope of Data Usability Study

This document details a Data Usability Summary (DUS) of analytical results for groundwater samples collected in support of the HWPW - Site-Wide Monitoring at the Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works site during July 2019. Samples were submitted to ALS Environmental (ALS), located in Houston, Texas and are reported in data packages HS19070575, HS19070609, HS19070735, HS19070864, HS19070994, HS19080069, and HS19080113. The intended use of the data is to support the HWPW - Site-Wide Monitoring at the site by providing current concentration of chemicals of concern.

Data were reviewed and validated by Chris G. Knight of GHD, in accordance with Title 30 of the Texas Administrative Code Section 350.54 (30 TAC 350.54) as described in the Texas Commission on Environmental Quality (TCEQ) Regulatory Guidance document entitled "Review and Reporting of COC Concentration Data under TRRP", (RG-366/TRRP-13), revised May 2010, herein referred to as "TRRP-13 Guidance". Evaluation of the data was based on information obtained from the chain of custody forms, the finished report forms, method blank data, recovery data from surrogate spikes/laboratory control samples (LCS)/matrix spikes (MS), duplicate data, field quality assurance/quality control (QA/QC) samples, the laboratory review checklists (LRC), and the laboratory exception reports (ER).

A sample collection and analysis summary is presented in Table 1. This summary provides a cross-reference of field sample identification numbers and location identification. Each sample is assigned a unique field identification number.

The validated sample results are presented in Table 2. A summary of the analytical methodology is presented in Table 3.





## 2. Laboratory Qualifications

The Laboratory's quality assurance program is consistent with the quality standards outlined in the National Environmental Laboratory Accreditation Program (NELAP). This laboratory was accredited under Texas Certification number # TX104704231 at the time the analysis was performed and the certificate is included in Attachment A.

## 3. Project Objectives

### 3.1 Sampling/Analytical QA/QC Objectives

The QA/QC program was designed to identify contamination resulting from the sampling, sample transport and analytical process through the analysis of trip blank samples, field blank samples, field duplicate sample sets, and method blanks. The QA/QC program was designed to evaluate the quality of the resulting data with respect to bias and precision through analysis of LCS, MS, and duplicate analyses.

## 4. Data Review/Validation Results

### 4.1 Sample Holding Time and Preservation

Samples were shipped with a chain of custody and the paper work was filled out properly with the following exceptions:

- i) HS19070864 – One of the sample vials for WG-1620-MW36A-20190716 mislabeled. It was logged in using the chain of custody. No further action was required.
- ii) HS19070994 – An additional trip blank sample was submitted but not listed on the chain of custody. The samples was logged in as WG-1620-TB06-20190719 for volatile organic compounds (VOCs) analysis. No further action was required.

All samples were properly preserved, delivered on ice, and stored by the laboratory at the required temperature (0-6°C).

The sample chain of custody documents and the analytical report were used to determine sample holding times. All samples were prepared and analyzed within the required holding times.

### 4.2 Sample Containers

Sample containers used were certified pre-cleaned glass and plastic containers provided by the laboratory. These containers meet or exceed analyte specifications established in the United States Environmental Protection Agency (USEPA) *Specifications and Guidance for Contaminant-free Sample Containers*.

### 4.3 Calibrations

According to the LRC, initial calibration and continuing calibration data met the criteria for the selected method.



#### **4.4 Laboratory Method Blank Analyses**

Method blanks are prepared from a purified matrix and analyzed with investigative samples to determine the existence and magnitude of sample contamination introduced during the analytical procedures. As these were not discrete samples handled in the field, these blanks are not listed on the sample identification cross-reference list found in the data packages.

For this study, laboratory method blanks were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch and results are reported in the laboratory data packages.

The method blank result was non-detect or below the method quantitation limit (MQL), indicating that laboratory contamination was not a factor for this investigation.

#### **4.5 Internal Standard and Surrogate Spike Recoveries**

Recoveries of internal standards are addressed in the LRC of the data packages. All internal standard recoveries associated with the compounds of interest were acceptable per the LRC.

In accordance with the methods employed, all samples, blanks, and QC samples analyzed for VOCs and semi-volatile organic compounds (SVOCs) are spiked with surrogate compounds prior to sample analysis. Surrogate recoveries provide a means to evaluate the effects of laboratory performance on individual sample matrices. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project. Each individual surrogate compound is expected to meet the laboratory control limits. According to the TRRP-13 Guidelines, one outlying surrogate is acceptable for methods with multiple surrogate spike compounds as long as the recovery is at least ten percent. Sample analyzed at elevated sample dilutions (five times or greater) were not assessed.

Surrogate recoveries were assessed against laboratory control limits and/or the guidance in TRRP-13. All surrogate recoveries met the above criteria.

#### **4.6 Laboratory Control Sample Analysis**

LCS or LCS/laboratory control sample duplicate (LCSD) are prepared and analyzed as samples to assess the analytical efficiencies of the methods employed, independent of sample matrix effects. The relative percent difference (RPD) of the LCS/LCSD recoveries is used to evaluate analytical precision. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project.

For this study, LCS or LCS/LCSD were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch.

The LCS or LCS/LCSD contained all analytes specified in the methods. All LCS recoveries and/or RPDs were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision (where applicable).



#### **4.7 Matrix Spike Analysis**

To evaluate the effects of sample matrices on the preparation process, measurement procedures, and accuracy of a particular analysis, samples are spiked with known concentrations of the analytes of interest and analyzed as MS/matrix spike duplicate (MSD) samples. The RPD between the MS and MSD is used to assess analytical precision.

MS/MSD analyses were performed as specified in Table 1. The recovery ranges established by the laboratory is adopted as the acceptance criteria for the project.

The MS/MSD samples were spiked with all analytes specified in the methods. All percent recoveries and the RPD value were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision with the following exceptions (see Table 4):

- i) Several MS/MSDs were reported with outlying recoveries and/or elevated RPDs for VOCs and/or SVOCs analyses due to possible matrix interferences and were not assessed. No further action was required.
- ii) One MS/MSD was reported with an elevated RPD for naphthalene. All associated detected sample results were qualified as estimated.

The laboratory also performed additional MS/MSD on non-site samples. These cannot be used to assess accuracy and precision for the site samples.

#### **4.8 Duplicate Sample Analyses**

Analytical precision is evaluated based on the analysis of laboratory duplicate samples. For this study, duplicate samples were prepared and analyzed by the laboratory as specified in Table 1 for arsenic. The RPDs established by the laboratory are adopted as the acceptance criteria for the project.

All duplicate analyses performed were acceptable, demonstrating acceptable analytical precision.

The laboratory also performed additional duplicate analyses on non-site samples. These cannot be used to assess precision for the site samples.

#### **4.9 Field QA/QC Samples**

The field QA/QC consisted of nine trip blank samples, five field blank samples, and five field duplicate sample sets.

##### ***Trip Blank Sample Analysis***

To evaluate contamination from sample collection, transportation, storage, and analytical activities, nine trip blank samples were submitted to the laboratory for VOCs analysis. All results were non-detect for the compounds of interest.



### *Field Blank Sample Analysis*

To assess ambient conditions at the site, five field blank samples were submitted for analysis, as identified in Table 1. All results were non-detect for the analytes of interest with the following exceptions (see Table 5):

- i) All field blanks submitted yielded low level detections for various VOCs and/or SVOCs. Associated sample results that were significantly greater than the concentrations found in the field blanks or were non-detect were not impacted. No further action was required. Associated sample results with comparable concentrations to the field blank detections were qualified as non-detect.

### *Field Duplicate Sample Analysis*

To assess the analytical and sampling protocol precision, two field duplicate sample sets were collected and submitted "blind" to the laboratory, as specified in Table 1. The RPDs associated with these duplicate samples must be less than thirty percent for water samples. The RPDs are only used when sample concentrations are above the estimated regions of detection.

Field duplicate summary data are presented in Table 2. All field duplicate results were within acceptable agreement, demonstrating acceptable sampling and analytical precision with the following exceptions (see Table 6):

- i) WG-1620-MW21C-20190716 and WG-1620-FD01-20190716 did show some variability in SVOCs results and were qualified as estimated.
- ii) WG-1620-MW68B-20190718 and WG-1620-FD03-20190718 did show some variability in SVOCs results and were qualified as estimated.

## **4.10 Field Procedures**

Golder Associates, Inc. collected groundwater samples in accordance with their Standard Operating Procedures (SOP) for sample collection.

## **4.11 Analyte Reporting**

The laboratory reported detected results for each analyte down to the sample detection limit (SDL), which is defined as the method detection limit (MDL) with sample-specific adjustments for dilutions, aliquot size, volumes, etc. Positive analyte detections less than the MQL but greater than the SDL were qualified as estimated (J) in Table 2 unless qualified otherwise in this memorandum.

The detectability check standard (DCS) results supported the laboratory MDLs.

## **5. Conclusion**

Based on the assessment detailed in the foregoing, the data summarized in Table 2 are usable for the purpose of supporting the HWPW - Site-Wide Monitoring at the site by providing current concentration of chemicals of concern with the specific qualifications noted herein.

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters			Comments
					VOCs	SVOCs	Arsenic	
WG-1620-MW18A-20190710	MW-18A	Water	07/10/2019	07:20	X	X	X	
WG-1620-MW18C-20190710	MW-18C	Water	07/10/2019	08:05	X	X	X	
WG-1620-MW58A-20190710	MW-58A	Water	07/10/2019	08:55	X	X	X	
WG-1620-MW57A-20190710	MW-57A	Water	07/10/2019	09:55	X	X	X	
WG-1620-MW57B-20190710	MW-57B	Water	07/10/2019	10:40	X	X	X	
WG-1620-MW72B-20190710	MW-72B	Water	07/10/2019	11:35	X	X	X	
WG-1620-MW19C-20190710	MW-19C	Water	07/10/2019	12:25	X	X	X	MS/MSD-P
WG-1620-MW17C-20190710	MW-17C	Water	07/10/2019	13:25	X	X	X	
WG-1620-MW17-20190710	MW-17	Water	07/10/2019	14:15	X	X	X	
WG-1620-MW20A-20190710	MW-20A	Water	07/10/2019	15:10	X	X	X	
WG-1620-MW15A-20190710	MW-15A	Water	07/10/2019	16:10	X	X	X	MS/MSD-P; DUP
WG-1620-MW15C-20190710	MW-15C	Water	07/10/2019	17:05	X	X	X	
WG-1620-MW15B-20190710	MW-15B	Water	07/10/2019	18:00	X	X	X	
WG-1620-FB01-20190710	-	Water	07/10/2019	18:15	X	X	X	Field Blank
WQ-1620-TB01-20190710	-	Water	07/10/2019	-	X			Trip Blank
WQ-1620-TB02-20190710	-	Water	07/10/2019	-	X			Trip Blank
WG-1620-MW14-20190711	MW-14	Water	07/11/2019	07:20	X	X	X	
WG-1620-MW13-20190711	MW-13	Water	07/11/2019	08:15	X	X	X	
WG-1620-MW39B-20190711	MW-39B	Water	07/11/2019	09:15	X	X	X	
WG-1620-MW12C-20190711	MW-12C	Water	07/11/2019	10:00	X	X	X	
WG-1620-MW12A-20190711	MW-12A	Water	07/11/2019	10:55	X	X	X	MS/MSD-P; DUP
WG-1620-MW40B-20190711	MW-40B	Water	07/11/2019	12:00	X	X	X	
WG-1620-MW42B-20190711	MW-42B	Water	07/11/2019	13:00	X	X	X	
WG-1620-P11-20190711	P-11	Water	07/11/2019	13:55	X	X	X	

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters			Comments
					VOCs	SVOCs	Arsenic	
WG-1620-MW05-20190711	MW-05	Water	07/11/2019	14:50	X	X	X	
WG-1620-MW64A-20190711	MW-64A	Water	07/11/2019	15:50	X	X	X	MS/MSD-P; DUP
WG-1620-FB02-20190711	-	Water	07/11/2019	16:05	X	X	X	Field Blank
WG-1620-TB02-20190711	-	Water	07/11/2019	-	X			Trip Blank
WG-1620-MW09-20190712	MW-09	Water	07/12/2019	07:35	X	X	X	MS/MSD-P; DUP
WG-1620-MW03-20190712	MW-03	Water	07/12/2019	08:35	X	X	X	
WG-1620-MW04-20190712	MW-04	Water	07/12/2019	09:25	X	X	X	
WG-1620-TW41B-20190712	TW-41B	Water	07/12/2019	10:40	X	X	X	
WG-1620-FB03-20190712	-	Water	07/12/2019	11:00	X	X	X	Field Blank
WG-1620-TB03-20190712	-	Water	07/12/2019	-	X			Trip Blank
WG-1620-MW21C-20190716	MW-21C	Water	07/16/2019	07:35	X	X	X	
WG-1620-FD01-20190716	MW-21C	Water	07/16/2019	07:35	X	X	X	Field duplicate of MW-21C
WG-1620-MW62B-20190716	MW-62B	Water	07/16/2019	08:25	X	X	X	
WG-1620-MW53C-20190716	MW-53C	Water	07/16/2019	10:15	X	X	X	
WG-1620-MW54C-20190716	MW-54C	Water	07/16/2019	11:15	X	X	X	
WG-1620-MW25C-20190716	MW-25C	Water	07/16/2019	12:15	X	X	X	
WG-1620-MW25A-20190716	MW-25A	Water	07/16/2019	13:10	X	X	X	
WG-1620-MW36B-20190716	MW-36B	Water	07/16/2019	14:20	X	X	X	MS/MSD-P
WG-1620-MW36A-20190716	MW-36A	Water	07/16/2019	15:20	X	X	X	
WG-1620-MW28C-20190716	MW-28C	Water	07/16/2019	16:15	X	X	X	MS/MSD; DUP
WG-1620-MW28A-20190716	MW-28A	Water	07/16/2019	17:00	X	X	X	
WG-1620-MW63B-20190716	MW-63B	Water	07/16/2019	17:55	X	X	X	
WG-1620-FB04-20190716	-	Water	07/16/2019	18:20	X	X	X	Field Blank
WG-1620-TB04-20190716	-	Water	07/16/2019	-	X			Trip Blank

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters			Comments
					VOCs	SVOCs	Arsenic	
WG-1620-MW61A-20190717	MW-61A	Water	07/17/2019	07:50	X	X	X	MS/MSD; DUP
WG-1620-MW60A-20190717	MW-60A	Water	07/17/2019	08:35	X	X	X	
WG-1620-MW69A-20190717	MW-69A	Water	07/17/2019	09:35	X	X	X	
WG-1620-MW47C-20190717	MW-47C	Water	07/17/2019	10:25	X	X	X	
WG-1620-MW48C-20190717	MW-48C	Water	07/17/2019	11:20	X	X	X	
WG-1620-MW59A-20190717	MW-59A	Water	07/17/2019	12:20	X	X	X	
WG-1620-MW59B-20190717	MW-59B	Water	07/17/2019	13:10	X	X	X	
WG-1620-MW44A-20190717	MW-44A	Water	07/17/2019	14:00	X	X	X	
WG-1620-MW87C-20190717	MW-87C	Water	07/17/2019	15:40	X	X	X	
WG-1620-MW71B-20190717	MW-71B	Water	07/17/2019	16:40	X	X	X	
WG-1620-MW33A-20190717	MW-33A	Water	07/17/2019	17:50	X	X	X	
WG-1620-FD02-20190717	MW-33A	Water	07/17/2019	17:50	X	X	X	Field duplicate of MW-33A
WG-1620-MW26A-20190717	MW-26A	Water	07/17/2019	18:45	X	X	X	
WG-1620-FB05-20190717	-	Water	07/17/2019	19:05	X	X	X	Field Blank
WG-1620-MW68A-20190718	MW-68A	Water	07/18/2019	07:55	X	X	X	
WG-1620-MW68B-20190718	MW-68B	Water	07/18/2019	08:55	X	X	X	
WG-1620-FD03-20190718	MW-68B	Water	07/18/2019	08:55	X	X	X	Field duplicate of MW-68B
WG-1620-MW68C-20190718	MW-68C	Water	07/18/2019	09:45	X	X	X	
WG-1620-MW83B-20190718	MW-83B	Water	07/18/2019	10:40	X	X	X	
WG-1620-MW83C-20190718	MW-83C	Water	07/18/2019	11:30	X	X	X	MS/MSD-P; DUP
WG-1620-MW35A-20190718	MW-35A	Water	07/18/2019	12:40	X	X	X	
WG-1620-MW35B-20190718	MW-35B	Water	07/18/2019	13:30	X	X	X	
WG-1620-MW90B-20190718	MW-90B	Water	07/18/2019	14:25	X	X	X	
WG-1620-MW89B-20190718	MW-89B	Water	07/18/2019	15:15	X	X	X	

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters			Comments
					VOCs	SVOCs	Arsenic	
WG-1620-MW38B-20190718	MW-38B	Water	07/18/2019	16:10	X	X	X	
WG-1620-MW27A-20190718	MW-27A	Water	07/18/2019	17:15	X	X	X	
WG-1620-MW27C-20190718	MW-27C	Water	07/18/2019	18:05	X	X	X	
WG-1620-MW51A-20190719	MW-51A	Water	07/19/2019	07:25	X	X	X	
WG-1620-MW51C-20190719	MW-51C	Water	07/19/2019	08:55	X	X	X	
WG-1620-MW81B-20190719	MW-81B	Water	07/19/2019	09:45	X	X	X	
WG-1620-MW50A-20190719	MW-50A	Water	07/19/2019	10:40	X	X	X	
WG-1620-TB05-20190719	-	Water	07/19/2019	-	X			Trip Blank
WG-1620-TB06-20190719	-	Water	07/19/2019	-	X			Trip Blank
WG-1620-MW32AR-20190730	MW-32AR	Water	07/30/2019	07:25	X	X	X	
WG-1620-MW33BR-20190730	MW-33BR	Water	07/30/2019	08:20	X	X	X	
WG-1620-MW84B-20190730	MW-84B	Water	07/30/2019	09:15	X	X	X	
WG-1620-MW82B-20190730	MW-82B	Water	07/30/2019	10:10	X	X	X	MS/MSD; DUP
WG-1620-MW80B-20190730	MW-80B	Water	07/30/2019	11:10	X	X	X	
WG-1620-MW77A-20190730	MW-77A	Water	07/30/2019	11:55	X	X	X	
WG-1620-MW85C-20190730	MW-85C	Water	07/30/2019	15:35	X	X	X	
WG-1620-MW86C-20190730	MW-86C	Water	07/30/2019	16:40	X	X	X	
WG-1620-FD04-20190730	MW-86C	Water	07/30/2019	16:40	X	X	X	Field duplicate of MW-86C
WG-1620-MW76C-20190730	MW-76C	Water	07/30/2019	17:30	X	X	X	
WG-1620-MW74B-20190730	MW-74B	Water	07/30/2019	18:20	X	X	X	
WG-1620-MW79A-20190730	MW-79A	Water	07/30/2019	19:10	X	X	X	
WG-1620-TB05-20190730	-	Water	07/30/2019	-	X			Trip Blank
WG-1620-MW36D-20190731	MW-36D	Water	07/31/2019	07:40	X	X	X	
WG-1620-MW38A-20190731	MW-38A	Water	07/31/2019	08:35	X	X	X	



Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters			Comments
					VOCs	SVOCs	Arsenic	
WG-1620-MW65D-20190731	MW-65D	Water	07/31/2019	09:45	X	X	X	MS/MSD-P; DUP
WG-1620-MW66D-20190731	MW-66D	Water	07/31/2019	10:45	X	X	X	
WG-1620-MW59D-20190731	MW-59D	Water	07/31/2019	11:50	X	X	X	
WG-1620-FD05-20190731	MW-59D	Water	07/31/2019	11:50	X	X	X	Field duplicate of MW-59D
WG-1620-MW22AR-20190731	MW-22AR	Water	07/31/2019	13:05	X	X	X	MS/MSD-P
WG-1620-MW22BR-20190731	MW-22BR	Water	07/31/2019	13:50	X	X	X	
WG-1620-MW88C-20190731	MW-88C	Water	07/31/2019	15:45	X	X	X	
WG-1620-MW67B-20190731	MW-67B	Water	07/31/2019	16:45	X	X	X	MS/MSD; DUP
WG-1620-MW49A-20190731	MW-49A	Water	07/31/2019	17:45	X	X	X	
WG-1620-TB06-20190731	-	Water	07/31/2019	-	X			Trip Blank

## Notes:

- VOCs - Volatile Organic Compounds  
SVOCs - Semi-volatile Organic Compounds  
MS/MSD - Matrix Spike/ Matrix Spike Duplicate  
MS/MSD-P - Matrix Spike/ Matrix Spike Duplicate (partial parameters)  
DUP - Laboratory Duplicate  
"- " - Not Applicable

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-03	MW-04	MW-05	MW-09	MW-12A
Sample Name:	WG-1620-MW03-20190712	WG-1620-MW04-20190712	WG-1620-MW05-20190711	WG-1620-MW09-20190712	WG-1620-MW12A-20190711
Sample Date:	07/12/2019	07/12/2019	07/11/2019	07/12/2019	07/11/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	0.00060 J
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000059	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000035	<0.000059	0.000087 J	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	0.013
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000020
Acenaphthene	mg/L	<0.000027	0.00052	<0.000047	<0.000047
Acenaphthylene	mg/L	<0.000015	<0.000015	0.00043	0.19
Anthracene	mg/L	<0.000015	<0.000015	<0.000015	0.0013
Benzo(a)anthracene	mg/L	0.000055 J	<0.000014	0.000097 J	0.00061 J
Benzo(a)pyrene	mg/L	<0.000050	<0.000050	<0.000050	0.00061 J
Benzo(a)pyrene	mg/L	<0.000050	<0.000050	<0.000050	0.00041
bis(2-Chloroethoxy)methane	mg/L	<0.000020	<0.000020	<0.000020	0.00012
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	<0.000037	0.00031
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	0.00034
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000058	0.000051 J	<0.000020
Fluoranthene	mg/L	<0.000010	<0.000010	<0.000010	0.14
Fluorene	mg/L	<0.000010	<0.000010	<0.000010	0.0097
Fluorene	mg/L	<0.000030	0.000053 J	0.000064 J	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	0.16
Naphthalene	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	<0.00029	<0.00053	<0.00071	<0.0015
Naphthalene	mg/L	<0.00029	<0.00053	<0.00071	<0.0012

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-03	MW-04	MW-05	MW-09	MW-12A
Sample Name:	WG-1620-MW03-20190712	WG-1620-MW04-20190712	WG-1620-MW05-20190711	WG-1620-MW09-20190712	WG-1620-MW12A-20190711
Sample Date:	07/12/2019	07/12/2019	07/11/2019	07/12/2019	07/11/2019

Parameters	Unit	MW-03	MW-04	MW-05	MW-09	MW-12A
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	0.063
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.000020 J	0.00016	0.00019	<0.000019	0.0047
<b>Metals</b>						
Arsenic	mg/L	0.000582 J	0.0127	0.0171	0.000901 J	0.00192 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-12C	MW-13	MW-14	MW-15A	MW-15B
Sample Name:	WG-1620-MW12C-20190711	WG-1620-MW13-20190711	WG-1620-MW14-20190711	WG-1620-MW15A-20190710	WG-1620-MW15B-20190710
Sample Date:	07/11/2019	07/11/2019	07/11/2019	07/10/2019	07/10/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	0.00074 J
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	0.00035 J
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	0.00035 J
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	0.0026
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.00043	<0.000063	<0.00013	<0.00024
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.00021	0.000077 J	0.00019	0.0077
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.000071 J	0.000066 J	0.00010	0.11
Acenaphthylene	mg/L	<0.000015	0.000033 J	<0.000015	0.00045
Anthracene	mg/L	0.000030 J	0.00047	0.000089 J	0.0026
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.000054 J	<0.000037	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.000070 J	0.000039 J	0.000087 J	0.035
Fluoranthene	mg/L	<0.000010	0.000032 J	0.000067 J	0.0018
Fluorene	mg/L	0.000052 J	0.000042 J	0.000069 J	0.054
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	<0.0031	<0.0011	<0.0020	<0.0066

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-12C	MW-13	MW-14	MW-15A	MW-15B
Sample Name:	WG-1620-MW12C-20190711	WG-1620-MW13-20190711	WG-1620-MW14-20190711	WG-1620-MW15A-20190710	WG-1620-MW15B-20190710
Sample Date:	07/11/2019	07/11/2019	07/11/2019	07/10/2019	07/10/2019

Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.000039 J	<0.000021	0.000059 J	0.0095	0.0065
Phenol	mg/L	0.00040	<0.000035	0.00019 J	0.00039	<0.00019
Pyrene	mg/L	<0.000019	0.000084 J	0.000032 J	0.00075	0.0018
<b>Metals</b>						
Arsenic	mg/L	0.00200	0.0715	0.00133 J	0.0251	0.0133

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-15C	MW-17	MW-17C	MW-18A	MW-18C
Sample Name:	WG-1620-MW15C-20190710	WG-1620-MW17-20190710	WG-1620-MW17C-20190710	WG-1620-MW18A-20190710	WG-1620-MW18C-20190710
Sample Date:	07/10/2019	07/10/2019	07/10/2019	07/10/2019	07/10/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.00050 J	0.54	0.0097	1.2
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	0.00052 J
Ethylbenzene	mg/L	<0.00030	0.23	0.027	0.32
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	0.83	0.0073	1.0
Vinyl chloride	mg/L	--	--	--	<0.00020
Xylenes (total)	mg/L	<0.00030	0.78	0.039	1.0
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.00021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.0052	5.9	0.21	0.0051
2,4-Dinitrotoluene	mg/L	<0.000058	<0.00058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.00042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.00021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.00066	0.39	0.026	0.28
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.00020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.00047	<0.000047	<0.000047
Acenaphthene	mg/L	0.017	0.16	0.035	0.058
Acenaphthylene	mg/L	0.0015	0.0042	0.00037	0.0050
Anthracene	mg/L	0.00064	0.0092	0.0013	0.0056
Benzo(a)anthracene	mg/L	<0.000050	<0.00050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.00020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.00030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.00037	0.00056	<0.000037
Chrysene	mg/L	<0.000021	<0.00021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.00020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.0039	0.12	0.027	0.054
Fluoranthene	mg/L	0.00086	0.0039	0.00090	0.0019
Fluorene	mg/L	0.00097	0.060	0.014	0.024
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.00025	<0.000025	<0.000025
Naphthalene	mg/L	0.019	8.9	0.97	9.9

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-15C	MW-17	MW-17C	MW-18A	MW-18C
Sample Name:	WG-1620-MW15C-20190710	WG-1620-MW17-20190710	WG-1620-MW17C-20190710	WG-1620-MW18A-20190710	WG-1620-MW18C-20190710
Sample Date:	07/10/2019	07/10/2019	07/10/2019	07/10/2019	07/10/2019

Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.00024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.00079	<0.000079	<0.000079	0.034
Phenanthrene	mg/L	<0.00039	0.045	0.0100	0.075	0.026
Phenol	mg/L	<0.00026	9.9	0.00073	0.00048	0.0050
Pyrene	mg/L	0.00043	0.0018	0.00046	0.0011	0.0014
<b>Metals</b>						
Arsenic	mg/L	0.000889 J	0.0460	0.000854 J	0.0248	0.00358

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-19C	MW-20A	MW-21C	MW-21C	MW-22AR
Sample Name:	WG-1620-MW19C-20190710	WG-1620-MW20A-20190710	WG-1620-MW21C-20190716	WG-1620-FD01-20190716	WG-1620-MW22AR-20190731
Sample Date:	07/10/2019	07/10/2019	07/16/2019	07/16/2019 Duplicate	07/31/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.00040 J	0.013	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.00096 J	0.0079	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.00086 J	0.0011	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	--	--
Xylenes (total)	mg/L	0.0022	0.022	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	0.000082 J	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.00056	0.0035	0.00010 J	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000071	0.072	0.00040 J	0.000050 J
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.0010	0.14	0.0057 J	0.00081 J
Acenaphthylene	mg/L	0.000030 J	0.00078	0.000074 J	<0.000015
Anthracene	mg/L	<0.000066	0.0062	0.00044 J	0.000050 J
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	<0.000040	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.00057	0.077	0.0032 J	0.00040 J
Fluoranthene	mg/L	0.00056	0.00058	0.00081 J	0.000087 J
Fluorene	mg/L	<0.00031	0.068	0.0019 J	0.00025 J
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	<0.0012	1.9	0.0044 J	0.00080 J



Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-19C	MW-20A	MW-21C	MW-21C	MW-22AR
Sample Name:	WG-1620-MW19C-20190710	WG-1620-MW20A-20190710	WG-1620-MW21C-20190716	WG-1620-FD01-20190716	WG-1620-MW22AR-20190731
Sample Date:	07/10/2019	07/10/2019	07/16/2019	07/16/2019 Duplicate	07/31/2019

Parameters	Unit	MW-19C	MW-20A	MW-21C	MW-21C	MW-22AR
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.00014	0.026	0.00087 J	0.00011 J	<0.000021
Phenol	mg/L	<0.00015	0.00096	<0.00018 J	0.00081 J	<0.000035
Pyrene	mg/L	0.00064	0.00026	0.00039 J	0.000033 J	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.00166 J	0.00574	0.00130 J	0.00326	0.000481 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-22BR	MW-25A	MW-25C	MW-26A	MW-27A
Sample Name:	WG-1620-MW22BR-20190731	WG-1620-MW25A-20190716	WG-1620-MW25C-20190716	WG-1620-MW26A-20190717	WG-1620-MW27A-20190718
Sample Date:	07/31/2019	07/16/2019	07/16/2019	07/17/2019	07/18/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	0.0013	0.00036 J
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	0.035	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	0.011	<0.00020
Vinyl chloride	mg/L	--	<0.00020	<0.00020	--
Xylenes (total)	mg/L	<0.00030	<0.00030	0.25	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.00021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	0.0075	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.00058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.00042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.00021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.000057 J	0.54	0.000059 J
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.00020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.00047	<0.000047
Acenaphthene	mg/L	<0.000027	<0.000027	0.20	0.028
Acenaphthylene	mg/L	<0.000015	<0.000015	0.0020	0.00013
Anthracene	mg/L	<0.000014	<0.000014	0.016	0.00047
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	0.0013	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	0.00037 J	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.00030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	<0.00037	<0.00019
Chrysene	mg/L	<0.000021	<0.000021	0.0014	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.00020	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	0.19	0.00021
Fluoranthene	mg/L	<0.000010	<0.000010	0.014	0.0029
Fluorene	mg/L	<0.000030	<0.000030	0.081	0.0017
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.00025	<0.000025
Naphthalene	mg/L	<0.000020	0.00023	4.2	<0.00032

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-22BR	MW-25A	MW-25C	MW-26A	MW-27A
Sample Name:	WG-1620-MW22BR-20190731	WG-1620-MW25A-20190716	WG-1620-MW25C-20190716	WG-1620-MW26A-20190717	WG-1620-MW27A-20190718
Sample Date:	07/31/2019	07/16/2019	07/16/2019	07/17/2019	07/18/2019

Parameters	Unit	MW-22BR	MW-25A	MW-25C	MW-26A	MW-27A
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.00024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.00079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	0.12	0.000083 J	0.00013
Phenol	mg/L	0.00069	<0.000035	<0.00035	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000019	0.0085	0.0015	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.000559 J	0.00285	0.00487	0.0933	0.000498 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-27C	MW-28A	MW-28C	MW-32AR	MW-33A
Sample Name:	WG-1620-MW27C-20190718	WG-1620-MW28A-20190716	WG-1620-MW28C-20190716	WG-1620-MW32AR-20190730	WG-1620-MW33A-20190717
Sample Date:	07/18/2019	07/16/2019	07/16/2019	07/30/2019	07/17/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.00020	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.000082 J	<0.000027	0.000098 J
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	0.000036 J	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	0.000055 J	<0.000050	0.000054 J
Benzo(a)pyrene	mg/L	<0.000020	0.000095 J	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000092 J	<0.00023	<0.00013	0.000074 J
Chrysene	mg/L	<0.000021	0.00010	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.000084 J	<0.000020	<0.000020
Fluoranthene	mg/L	<0.000010	0.00012	0.000015 J	<0.000010
Fluorene	mg/L	<0.000030	0.000053 J	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	0.000066 J	0.0013	0.000095 J	<0.000020

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-27C	MW-28A	MW-28C	MW-32AR	MW-33A
Sample Name:	WG-1620-MW27C-20190718	WG-1620-MW28A-20190716	WG-1620-MW28C-20190716	WG-1620-MW32AR-20190730	WG-1620-MW33A-20190717
Sample Date:	07/18/2019	07/16/2019	07/16/2019	07/30/2019	07/17/2019

Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	0.000049 J	<0.000021	<0.000021	<0.000021
Phenol	mg/L	0.00023	<0.000073	0.0025	<0.000035	0.00011 J
Pyrene	mg/L	<0.000019	0.00016	<0.000019	0.00045	0.00028
<b>Metals</b>						
Arsenic	mg/L	0.000428 J	0.00920	0.000456 J	0.0628	0.0155

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-33A	MW-33BR	MW-35A	MW-35B	MW-36A
Sample Name:	WG-1620-FD02-20190717	WG-1620-MW33BR-20190730	WG-1620-MW35A-20190718	WG-1620-MW35B-20190718	WG-1620-MW36A-20190716
Sample Date:	07/17/2019	07/30/2019	07/18/2019	07/18/2019	07/16/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	0.25	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	0.065	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	0.0031	<0.00020	0.00054 J
Vinyl chloride	mg/L	--	<0.00020	--	<0.00020
Xylenes (total)	mg/L	<0.00030	0.016	<0.00030	0.0062
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	0.00028	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.014	0.00012	0.041
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.0078	0.0028	0.029
Acenaphthylene	mg/L	<0.000015	<0.000015	0.000071 J	0.00022
Anthracene	mg/L	<0.000014	0.00055	0.00011	0.0022
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	0.000083 J
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	0.000053 J
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000084	0.000070 J	0.000057 J	0.00012 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	0.000091 J
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.012	0.00039	0.029
Fluoranthene	mg/L	0.00017	0.00039	0.00015	0.0017
Fluorene	mg/L	<0.000030	0.0042	0.00036	0.014
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	<0.000096	0.53 J	0.00083	1.1

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

<b>Location ID:</b>	<b>MW-33A</b>	<b>MW-33BR</b>	<b>MW-35A</b>	<b>MW-35B</b>	<b>MW-36A</b>
<b>Sample Name:</b>	<b>WG-1620-FD02-20190717</b>	<b>WG-1620-MW33BR-20190730</b>	<b>WG-1620-MW35A-20190718</b>	<b>WG-1620-MW35B-20190718</b>	<b>WG-1620-MW36A-20190716</b>
<b>Sample Date:</b>	<b>07/17/2019</b>	<b>07/30/2019</b>	<b>07/18/2019</b>	<b>07/18/2019</b>	<b>07/16/2019</b>
	<b>Duplicate</b>				

<b>Parameters</b>	<b>Unit</b>					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	0.0048	0.000042 J	0.019	<0.000021
Phenol	mg/L	0.000053 J	0.00021	<0.000035	0.00017 J	<0.00014
Pyrene	mg/L	0.00069	0.00026	0.000096 J	0.00079	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.0147	0.00117 J	0.0548	0.00120 J	0.00244

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-36B	MW-36D	MW-38A	MW-38B	MW-39B	
Sample Name:	WG-1620-MW36B-20190716	WG-1620-MW36D-20190731	WG-1620-MW38A-20190731	WG-1620-MW38B-20190718	WG-1620-MW39B-20190711	
Sample Date:	07/16/2019	07/31/2019	07/31/2019	07/18/2019	07/11/2019	
Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	0.00031 J	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.000078 J	<0.000040	<0.000040	0.000056 J	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.000058	<0.000019	<0.000019	0.000030	<0.000027
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.00014	<0.000027	<0.000027	0.00012	0.00065
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015	0.000021 J
Anthracene	mg/L	<0.000014	0.000017 J	0.000022 J	0.000037 J	0.00021
Benzo(a)anthracene	mg/L	<0.000050	0.000053 J	0.000087 J	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	0.000087 J	0.00012	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000088	0.000080 J	0.000084 J	0.00031	0.000042 J
Chrysene	mg/L	<0.000021	0.000058 J	0.000092 J	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.000087 J	<0.000020	<0.000020	0.00014	0.000037 J
Fluoranthene	mg/L	0.000017 J	0.00011	0.00018	<0.000010	0.000048 J
Fluorene	mg/L	0.000047 J	<0.000030	<0.000030	0.000060 J	0.000044 J
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	0.0060	<0.000020	0.000068 J	0.0026	<0.00041



**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-36B	MW-36D	MW-38A	MW-38B	MW-39B
Sample Name:	WG-1620-MW36B-20190716	WG-1620-MW36D-20190731	WG-1620-MW38A-20190731	WG-1620-MW38B-20190718	WG-1620-MW39B-20190711
Sample Date:	07/16/2019	07/31/2019	07/31/2019	07/18/2019	07/11/2019

Parameters	Unit	MW-36B	MW-36D	MW-38A	MW-38B	MW-39B
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	0.000066 J	0.00010	0.000079 J	<0.000021
Phenol	mg/L	<0.00017	<0.000035	<0.000035	0.000056 J	<0.000035
Pyrene	mg/L	<0.000019	0.000099 J	0.00014	<0.000019	0.000072 J
<b>Metals</b>						
Arsenic	mg/L	0.00127 J	<0.000400	<0.000400	0.000553 J	0.00144 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-40B	MW-42B	MW-44A	MW-47C	MW-48C
Sample Name:	WG-1620-MW40B-20190711	WG-1620-MW42B-20190711	WG-1620-MW44A-20190717	WG-1620-MW47C-20190717	WG-1620-MW48C-20190717
Sample Date:	07/11/2019	07/11/2019	07/17/2019	07/17/2019	07/17/2019

Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.0088	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.082	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.014	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	<0.00020	--	--
Xylenes (total)	mg/L	0.13	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.0013	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.18	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.21	0.000068 J	0.056	<0.000027	<0.000027
Acenaphthylene	mg/L	0.0021	<0.000015	0.00028	<0.000015	<0.000015
Anthracene	mg/L	0.011	0.000064 J	0.00089	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	0.000085 J	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.00015 J	<0.00025	<0.00027	<0.00030
Chrysene	mg/L	<0.000021	<0.000021	0.000052 J	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.13	0.000038 J	<0.000020	<0.000020	<0.000020
Fluoranthene	mg/L	0.0059	0.000040 J	0.0091	<0.000010	0.000035 J
Fluorene	mg/L	0.13	0.000057 J	0.014	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	3.6	<0.00043	<0.00019	<0.00020	<0.00020

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-40B	MW-42B	MW-44A	MW-47C	MW-48C
Sample Name:	WG-1620-MW40B-20190711	WG-1620-MW42B-20190711	WG-1620-MW44A-20190717	WG-1620-MW47C-20190717	WG-1620-MW48C-20190717
Sample Date:	07/11/2019	07/11/2019	07/17/2019	07/17/2019	07/17/2019

Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.093	0.000034 J	0.00012	<0.000021	<0.000021
Phenol	mg/L	0.000053 J	<0.000035	<0.000035	<0.000035	0.000074 J
Pyrene	mg/L	0.0024	0.000033 J	0.0057	<0.000019	0.000036 J
<b>Metals</b>						
Arsenic	mg/L	0.0520	0.00220	0.0303	0.000440 J	0.00167 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-49A	MW-50A	MW-51A	MW-51C	MW-53C
Sample Name:	WG-1620-MW49A-20190731	WG-1620-MW50A-20190719	WG-1620-MW51A-20190719	WG-1620-MW51C-20190719	WG-1620-MW53C-20190716
Sample Date:	07/31/2019	07/19/2019	07/19/2019	07/19/2019	07/16/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	0.0028
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	0.000067 J	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	0.00013 J	0.000079 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	0.000030 J	<0.000020
Fluoranthene	mg/L	<0.000010	<0.000010	<0.000010	<0.000010
Fluorene	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	<0.000020	0.00027	0.00055	0.00023

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-49A	MW-50A	MW-51A	MW-51C	MW-53C
Sample Name:	WG-1620-MW49A-20190731	WG-1620-MW50A-20190719	WG-1620-MW51A-20190719	WG-1620-MW51C-20190719	WG-1620-MW53C-20190716
Sample Date:	07/31/2019	07/19/2019	07/19/2019	07/19/2019	07/16/2019

Parameters	Unit	MW-49A	MW-50A	MW-51A	MW-51C	MW-53C
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	0.000026 J	<0.000021	<0.000021
Phenol	mg/L	0.00064	0.00016 J	0.00019 J	0.00027	<0.00016
Pyrene	mg/L	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.000658 J	0.000642 J	<0.000400	<0.000400	0.000569 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-54C	MW-57A	MW-57B	MW-58A	MW-59A
Sample Name:	WG-1620-MW54C-20190716	WG-1620-MW57A-20190710	WG-1620-MW57B-20190710	WG-1620-MW58A-20190710	WG-1620-MW59A-20190717
Sample Date:	07/16/2019	07/10/2019	07/10/2019	07/10/2019	07/17/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	0.00057 J	0.84	0.0049
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	0.00080 J
Ethylbenzene	mg/L	<0.00030	0.00085 J	0.39	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	0.93	<0.00020
Vinyl chloride	mg/L	--	<0.00020	0.00048 J	<0.00020
Xylenes (total)	mg/L	<0.00030	<0.00030	1.2	0.028
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.00021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.00059	6.5	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.00058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.00042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.00021	<0.000021
2-Methylnaphthalene	mg/L	0.0021	0.19	0.93	0.015
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.00020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.00047	<0.000047
Acenaphthene	mg/L	0.041	0.099	0.38	0.11
Acenaphthylene	mg/L	0.00037	0.0019	0.0051	0.0087
Anthracene	mg/L	0.0023	0.0084	0.039	0.062
Benzo(a)anthracene	mg/L	<0.000050	0.00100	0.0071	0.00082
Benzo(a)pyrene	mg/L	<0.000020	0.00054	0.0021	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.00030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.00021	<0.00037	<0.000037
Chrysene	mg/L	<0.000021	0.00100	0.0065	0.00084
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.00020	<0.000020
Dibenzofuran	mg/L	0.032	0.076	0.32	0.059
Fluoranthene	mg/L	0.0033	0.0091	0.047	0.020
Fluorene	mg/L	0.018	0.059	0.21	0.15
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.00025	<0.000025
Naphthalene	mg/L	0.032	0.61	17	0.46

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

<b>Location ID:</b>	<b>MW-54C</b>	<b>MW-57A</b>	<b>MW-57B</b>	<b>MW-58A</b>	<b>MW-59A</b>
<b>Sample Name:</b>	<b>WG-1620-MW54C-20190716</b>	<b>WG-1620-MW57A-20190710</b>	<b>WG-1620-MW57B-20190710</b>	<b>WG-1620-MW58A-20190710</b>	<b>WG-1620-MW59A-20190717</b>
<b>Sample Date:</b>	<b>07/16/2019</b>	<b>07/10/2019</b>	<b>07/10/2019</b>	<b>07/10/2019</b>	<b>07/17/2019</b>

<b>Parameters</b>	<b>Unit</b>					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.00024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	0.00059	<0.00079	<0.000079	<0.000079
Phenanthrene	mg/L	0.0065	0.059	0.29	0.036	<0.000021
Phenol	mg/L	0.00044	<0.00016	1.0	<0.000035	0.000058 J
Pyrene	mg/L	0.0016	0.0055	0.030	0.0099	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.00103 J	0.00447	0.00404	0.00748	0.00455

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-59B	MW-59D	MW-59D	MW-60A	MW-61A
Sample Name:	WG-1620-MW59B-20190717	WG-1620-MW59D-20190731	WG-1620-FD05-20190731	WG-1620-MW60A-20190717	WG-1620-MW61A-20190717
Sample Date:	07/17/2019	07/31/2019	07/31/2019 Duplicate	07/17/2019	07/17/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	<0.000019	0.000072 J
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.00039	<0.00037	<0.00037	<0.00055
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.000055 J	<0.000020	<0.000020	0.000075 J
Dibenzofuran	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Fluoranthene	mg/L	<0.000010	<0.000010	<0.000010	<0.000010
Fluorene	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	<0.00015	<0.000020	<0.000020	<0.00013



Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

	Location ID:	MW-59B	MW-59D	MW-59D	MW-60A	MW-61A
	Sample Name:	WG-1620-MW59B-20190717	WG-1620-MW59D-20190731	WG-1620-FD05-20190731	WG-1620-MW60A-20190717	WG-1620-MW61A-20190717
	Sample Date:	07/17/2019	07/31/2019	07/31/2019 Duplicate	07/17/2019	07/17/2019
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	mg/L	<0.000035	0.00011 J	<0.000035	<0.000035	0.00011 J
Pyrene	mg/L	<0.000019	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.000542 J	<0.000400	0.000502 J	0.000440 J	0.00117 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-62B	MW-63B	MW-64A	MW-65D	MW-66D
Sample Name:	WG-1620-MW62B-20190716	WG-1620-MW63B-20190716	WG-1620-MW64A-20190711	WG-1620-MW65D-20190731	WG-1620-MW66D-20190731
Sample Date:	07/16/2019	07/16/2019	07/11/2019	07/31/2019	07/31/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	0.11	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	0.14	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	0.00071 J	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	0.041	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.000096 J	0.037	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.000099 J	0.012	0.00030	<0.000027
Acenaphthylene	mg/L	<0.000015	0.00016	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	0.00034	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	<0.000037	0.00016 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.000048 J	0.011	<0.000020	<0.000020
Fluoranthene	mg/L	0.000017 J	<0.000010	<0.000010	<0.000010
Fluorene	mg/L	0.000036 J	0.0037	0.000031 J	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	0.00028	1.5	<0.00033	<0.000020

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-62B	MW-63B	MW-64A	MW-65D	MW-66D
Sample Name:	WG-1620-MW62B-20190716	WG-1620-MW63B-20190716	WG-1620-MW64A-20190711	WG-1620-MW65D-20190731	WG-1620-MW66D-20190731
Sample Date:	07/16/2019	07/16/2019	07/11/2019	07/31/2019	07/31/2019

Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.000021 J	0.0013	<0.000021	<0.000021	<0.000021
Phenol	mg/L	0.00056	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000019	0.00015	<0.000019	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.00194 J	0.00156 J	0.00939	0.00135 J	0.00124 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-67B	MW-68A	MW-68B	MW-68B	MW-68C
Sample Name:	WG-1620-MW67B-20190731	WG-1620-MW68A-20190718	WG-1620-MW68B-20190718	WG-1620-FD03-20190718	WG-1620-MW68C-20190718
Sample Date:	07/31/2019	07/18/2019	07/18/2019	07/18/2019 Duplicate	07/18/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.0020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	1.4	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.0030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	0.52	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	0.37	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	1.5	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.00021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	0.076	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.00058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.00042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.00021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.00012	0.72	0.000072 J
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.00020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.00047	<0.000047
Acenaphthene	mg/L	<0.000027	0.0019	0.20	0.00012
Acenaphthylene	mg/L	<0.000015	<0.000015	0.0027	0.000053 J
Anthracene	mg/L	<0.000014	0.000028 J	0.030 J	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	0.0051 J	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	0.0018	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.00030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.00024	<0.00037	0.00032
Chrysene	mg/L	<0.000021	<0.000021	0.0050 J	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	0.00015 J	<0.00020	0.000067 J
Dibenzofuran	mg/L	<0.000020	0.000082 J	0.21	0.000064 J
Fluoranthene	mg/L	<0.000010	<0.000010	0.043 J	<0.000010
Fluorene	mg/L	<0.000030	0.00012	0.11	0.000051 J
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.00025	<0.000025
Naphthalene	mg/L	0.000079 J	0.00035	9.6	0.0011

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-67B	MW-68A	MW-68B	MW-68B	MW-68C
Sample Name:	WG-1620-MW67B-20190731	WG-1620-MW68A-20190718	WG-1620-MW68B-20190718	WG-1620-FD03-20190718	WG-1620-MW68C-20190718
Sample Date:	07/31/2019	07/18/2019	07/18/2019	07/18/2019 Duplicate	07/18/2019

Parameters	Unit	MW-67B	MW-68A	MW-68B	MW-68B	MW-68C
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.00024	<0.00024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.00079	<0.00079	<0.000079
Phenanthrene	mg/L	<0.000021	0.000062 J	0.21	0.16	<0.000021
Phenol	mg/L	<0.000035	<0.000035	<0.00035	<0.00035	0.00026
Pyrene	mg/L	<0.000019	<0.000019	0.025 J	0.015 J	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.000494 J	0.0353	0.0134	0.0129	<0.000400

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-69A	MW-71B	MW-72B	MW-74B	MW-76C
Sample Name:	WG-1620-MW69A-20190717	WG-1620-MW71B-20190717	WG-1620-MW72B-20190710	WG-1620-MW74B-20190730	WG-1620-MW76C-20190730
Sample Date:	07/17/2019	07/17/2019	07/10/2019	07/30/2019	07/30/2019

Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.0010	<0.00020
Benzene	mg/L	<0.00020	0.13	1.1	0.59	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.0015	<0.00030
Ethylbenzene	mg/L	<0.00030	0.031	0.30	0.15	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0050	<0.0010
Toluene	mg/L	<0.00020	0.023	0.96	0.52	<0.00020
Vinyl chloride	mg/L	<0.00020	--	--	--	--
Xylenes (total)	mg/L	<0.00030	0.054	0.97	0.42	<0.00030
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.00021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	19	37	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.00058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.00042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.00021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.024	0.57	0.48	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.00020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.00047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.016	0.20	0.24	<0.000027
Acenaphthylene	mg/L	<0.000015	0.00033	0.0044	0.0037	<0.000015
Anthracene	mg/L	<0.000014	0.0066	0.034	0.0071	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	0.0026	0.0069	0.00022	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	0.00083	0.0019	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.00030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000034	<0.00014	<0.00037	<0.000037	<0.000037
Chrysene	mg/L	<0.000021	0.0024	0.0069	0.0016	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.00020	<0.000020	0.000054 J
Dibenzofuran	mg/L	<0.000020	0.016	0.16	0.19	<0.000020
Fluoranthene	mg/L	<0.000010	0.017	0.046	0.0035	<0.000010
Fluorene	mg/L	<0.000030	0.0093	0.12	0.14	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.00025	<0.000025	<0.000025
Naphthalene	mg/L	<0.000078	0.44	12	13 J	0.00018 J

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

<b>Location ID:</b>	<b>MW-69A</b>	<b>MW-71B</b>	<b>MW-72B</b>	<b>MW-74B</b>	<b>MW-76C</b>
<b>Sample Name:</b>	<b>WG-1620-MW69A-20190717</b>	<b>WG-1620-MW71B-20190717</b>	<b>WG-1620-MW72B-20190710</b>	<b>WG-1620-MW74B-20190730</b>	<b>WG-1620-MW76C-20190730</b>
<b>Sample Date:</b>	<b>07/17/2019</b>	<b>07/17/2019</b>	<b>07/10/2019</b>	<b>07/30/2019</b>	<b>07/30/2019</b>

<b>Parameters</b>	<b>Unit</b>					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.00024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.00079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	0.030	0.16	0.087	<0.000021
Phenol	mg/L	<0.000035	0.000051 J	9.3	33	<0.000035
Pyrene	mg/L	<0.000019	0.0088	0.029	0.0023	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.000642 J	0.000626 J	0.000861 J	0.00128 J	0.00216

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Location ID:	MW-77A	MW-79A	MW-80B	MW-81B	MW-82B
Sample Name:	WG-1620-MW77A-20190730	WG-1620-MW79A-20190730	WG-1620-MW80B-20190730	WG-1620-MW81B-20190719	WG-1620-MW82B-20190730
Sample Date:	07/30/2019	07/30/2019	07/30/2019	07/19/2019	07/30/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.063	0.013	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.072	0.0067	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.012	0.018	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	0.097	0.023	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.062	0.33	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.20	0.051	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.16	0.034	<0.000027	<0.000027
Acenaphthylene	mg/L	0.0013	0.00076	<0.000015	<0.000015
Anthracene	mg/L	0.0028	0.0020	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	0.00023	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.00055	0.000070 J	<0.000037
Chrysene	mg/L	<0.000021	0.00027	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.096	0.026	<0.000020	<0.000020
Fluoranthene	mg/L	0.00052	0.0012	<0.000010	<0.000010
Fluorene	mg/L	0.078	0.017	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	4.9 J	0.74 J	<0.000020	0.000099 J



**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

<b>Location ID:</b>	<b>MW-77A</b>	<b>MW-79A</b>	<b>MW-80B</b>	<b>MW-81B</b>	<b>MW-82B</b>
<b>Sample Name:</b>	<b>WG-1620-MW77A-20190730</b>	<b>WG-1620-MW79A-20190730</b>	<b>WG-1620-MW80B-20190730</b>	<b>WG-1620-MW81B-20190719</b>	<b>WG-1620-MW82B-20190730</b>
<b>Sample Date:</b>	<b>07/30/2019</b>	<b>07/30/2019</b>	<b>07/30/2019</b>	<b>07/19/2019</b>	<b>07/30/2019</b>

<b>Parameters</b>	<b>Unit</b>					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.032	0.0082	<0.000021	<0.000021	<0.000021
Phenol	mg/L	<0.000035	0.063	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.00041	0.0011	<0.000019	<0.000019	<0.000019
<b>Metals</b>						
Arsenic	mg/L	0.0231	0.00991	0.00162 J	0.000563 J	0.00873

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

	Location ID:	MW-83B	MW-83C	MW-84B	MW-85C
	Sample Name:	WG-1620-MW83B-20190718	WG-1620-MW83C-20190718	WG-1620-MW84B-20190730	WG-1620-MW85C-20190730
	Sample Date:	07/18/2019	07/18/2019	07/30/2019	07/30/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.021	<0.00020	0.010	0.00093 J
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.055	<0.00030	0.013	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.0042	<0.00020	0.0015	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	0.061	<0.00030	0.0041	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	0.00081	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.050	0.00012	0.0027	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.021	0.000094 J	0.0080	<0.000027
Acenaphthylene	mg/L	0.00025	<0.000015	0.00030	<0.000015
Anthracene	mg/L	0.00099	<0.000014	0.00042	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00036	0.00023	0.00026	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.00012 J	0.000040 J	0.000055 J	<0.000020
Dibenzofuran	mg/L	0.014	0.000070 J	0.0054	<0.000020
Fluoranthene	mg/L	0.00038	<0.000010	0.00015	<0.000010
Fluorene	mg/L	0.0076	<0.000030	0.0026	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	0.77	0.0014	0.060 J	0.0012 J

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

Location ID:	MW-83B	MW-83C	MW-84B	MW-85C
Sample Name:	WG-1620-MW83B-20190718	WG-1620-MW83C-20190718	WG-1620-MW84B-20190730	WG-1620-MW85C-20190730
Sample Date:	07/18/2019	07/18/2019	07/30/2019	07/30/2019

Parameters	Unit				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.0059	<0.000021	0.0021	<0.000021
Phenol	mg/L	0.00051	0.000068 J	<0.000035	<0.000035
Pyrene	mg/L	0.00020	<0.000019	0.00011	<0.000019
<b>Metals</b>					
Arsenic	mg/L	0.0648	0.00617	0.00838	0.000633 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

	Location ID:	MW-86C	MW-86C	MW-87C	MW-88C
	Sample Name:	WG-1620-MW86C-20190730	WG-1620-FD04-20190730	WG-1620-MW87C-20190717	WG-1620-MW88C-20190731
	Sample Date:	07/30/2019	07/30/2019 Duplicate	07/17/2019	07/31/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	<0.00020	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	<0.000027	<0.000027	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000060 J	0.000050 J	<0.00018	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.00011 J	0.000073 J	0.00017 J	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Fluoranthene	mg/L	<0.000010	<0.000010	<0.000010	<0.000010
Fluorene	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	0.00012 J	0.00031 J	<0.00017	<0.000020

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

	Location ID:	MW-86C	MW-86C	MW-87C	MW-88C
	Sample Name:	WG-1620-MW86C-20190730	WG-1620-FD04-20190730	WG-1620-MW87C-20190717	WG-1620-MW88C-20190731
	Sample Date:	07/30/2019	07/30/2019 Duplicate	07/17/2019	07/31/2019
Parameters	Unit				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	mg/L	<0.000035	<0.000035	0.00013 J	0.00055
Pyrene	mg/L	<0.000019	<0.000019	<0.000019	<0.000019
<b>Metals</b>					
Arsenic	mg/L	0.00236	0.00190 J	<0.000400	<0.000400

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

	Location ID:	MW-89B	MW-90B	P-11	TW-41B
	Sample Name:	WG-1620-MW89B-20190718	WG-1620-MW90B-20190718	WG-1620-P11-20190711	WG-1620-TW41B-20190712
	Sample Date:	07/18/2019	07/18/2019	07/11/2019	07/12/2019
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.0013
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.0011
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.0011
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	0.019
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.000050 J	<0.000019	0.000050 J	0.067
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	<0.000027	0.0019	0.15
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	0.0014
Anthracene	mg/L	<0.000014	<0.000014	0.000063 J	0.0055
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00015 J	0.00011 J	0.000087 J	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.000038 J	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	0.000079 J	0.074
Fluoranthene	mg/L	<0.000010	<0.000010	0.000060 J	0.0034
Fluorene	mg/L	<0.000030	<0.000030	0.00074	0.085
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025
Naphthalene	mg/L	0.00025	0.000091 J	<0.0018	0.69

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
July 2019**

	Location ID:	MW-89B	MW-90B	P-11	TW-41B
	Sample Name:	WG-1620-MW89B-20190718	WG-1620-MW90B-20190718	WG-1620-P11-20190711	WG-1620-TW41B-20190712
	Sample Date:	07/18/2019	07/18/2019	07/11/2019	07/12/2019
Parameters	Unit				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	0.00020	0.027
Phenol	mg/L	0.00042	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000019	0.000036 J	0.0014
<b>Metals</b>					
Arsenic	mg/L	<0.000400	0.0135	0.0704	0.113

Notes:

- < - Not detected at the associated reporting limit
- J - Estimated concentration
- "-" - Not applicable

Table 3

**Analytical Methods**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Method	Matrix	Holding Time	
			Collection to Extraction (Days)	Extraction to Analysis (Days)
VOCs	SW-846 8260C	Water	-	14
SVOCs	SW-846 8270D	Water	7	40
Arsenic	SW-846 6020A	Water	-	180

## Notes:

VOCs - Volatile Organic Compounds  
SVOCs - Semi-volatile Organic Compounds  
"- " - Not Applicable

## Method References:

SW-846 - "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846,  
Third Edition, 1986, with subsequent revisions



Table 4

**Qualified Sample Results Due to Outlying MS/MSD Results**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Sample ID	Analyte	MS	MSD	RPD (percent)	Control Limits		Qualified Result	Units
			% Recovery	% Recovery		% Recovery	RPD		
SVOCs	WG-1620-FD04-20190730	Naphthalene	65.1	104	46.1	70-127	20	0.00031 J	mg/L
	WG-1620-MW33BR-20190730							0.53 J	mg/L
	WG-1620-MW74B-20190730							13 J	mg/L
	WG-1620-MW76C-20190730							0.00018 J	mg/L
	WG-1620-MW77A-20190730							4.9 J	mg/L
	WG-1620-MW79A-20190730							0.74 J	mg/L
	WG-1620-MW84B-20190730							0.060 J	mg/L
	WG-1620-MW85C-20190730							0.0012 J	mg/L
WG-1620-MW86C-20190730	0.00012 J	mg/L							

## Notes:

- MS - Matrix Spike
- MSD - Matrix Spike Duplicate
- RPD - Relative Percent Difference
- SVOCs - Semi-volatile Organic Compounds
- J - Estimated concentration

Table 5

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB01-20190710	07/10/2019	2,4-Dimethylphenol	0.00037	WG-1620-MW15A-20190710	0.00024	<0.00024	mg/L
					WG-1620-MW15B-20190710	0.00016 J	<0.00016	mg/L
					WG-1620-MW19C-20190710	0.00056	<0.00056	mg/L
			2-Methylnaphthalene	0.00038	WG-1620-MW57A-20190710	0.00059	<0.00059	mg/L
					WG-1620-MW15C-20190710	0.00066	<0.00066	mg/L
					WG-1620-MW19C-20190710	0.00071 J	<0.00071	mg/L
			Anthracene	0.000032 J	WG-1620-MW19C-20190710	0.00066 J	<0.00066	mg/L
			Dibenzofuran	0.00018	WG-1620-MW19C-20190710	0.00057	<0.00057	mg/L
			Fluorene	0.00011	WG-1620-MW19C-20190710	0.00031	<0.00031	mg/L
			Naphthalene	0.0035	WG-1620-MW15A-20190710	0.0066	<0.0066	mg/L
					WG-1620-MW19C-20190710	0.0012	<0.0012	mg/L
			Phenanthrene	0.00013	WG-1620-MW15C-20190710	0.00039	<0.00039	mg/L
					WG-1620-MW19C-20190710	0.00014	<0.00014	mg/L
			Phenol	0.000071 J	WG-1620-MW15B-20190710	0.00019 J	<0.00019	mg/L
					WG-1620-MW15C-20190710	0.00026	<0.00026	mg/L
					WG-1620-MW19C-20190710	0.00015 J	<0.00015	mg/L
WG-1620-MW57A-20190710	0.00016 J	<0.00016			mg/L			
SVOCs	WG-1620-FB02-20190711	07/11/2019	2,4-Dimethylphenol	0.000064 J	WG-1620-MW05-20190711	0.000059 J	<0.000059	mg/L
					WG-1620-MW13-20190711	0.000063 J	<0.000063	mg/L
					WG-1620-MW14-20190711	0.00013 J	<0.00013	mg/L
			2-Methylnaphthalene	0.000077 J	WG-1620-MW39B-20190711	0.000027 J	<0.000027	mg/L

Table 5

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB02-20190711		Naphthalene	0.0062	WG-1620-MW05-20190711	0.00071	<0.00071	mg/L
					WG-1620-MW12A-20190711	0.0012	<0.0012	mg/L
					WG-1620-MW12C-20190711	0.0031	<0.0031	mg/L
					WG-1620-MW13-20190711	0.0011	<0.0011	mg/L
					WG-1620-MW14-20190711	0.0020	<0.0020	mg/L
					WG-1620-MW39B-20190711	0.00041	<0.00041	mg/L
					WG-1620-MW42B-20190711	0.00043	<0.00043	mg/L
					WG-1620-MW64A-20190711	0.00033	<0.00033	mg/L
	WG-1620-P11-20190711	0.0018	<0.0018	mg/L				
VOCs	WG-1620-FB03-20190712	07/12/2019	Benzene	0.00052 J	WG-1620-TW41B-20190712	0.0013	<0.0013	mg/L
			Ethylbenzene	0.00032 J	WG-1620-TW41B-20190712	0.0011	<0.0011	mg/L
VOCs	WG-1620-FB03-20190712	07/12/2019	Toluene	0.00037 J	WG-1620-TW41B-20190712	0.0011	<0.0011	mg/L
SVOCs			2-Methylnaphthalene	0.000052 J	WG-1620-MW03-20190712	0.000035 J	<0.000035	mg/L
					WG-1620-MW04-20190712	0.000059 J	<0.000059	mg/L
					WG-1620-MW04-20190712	0.000058 J	<0.000058	mg/L
					WG-1620-MW03-20190712	0.00029	<0.00029	mg/L
					WG-1620-MW04-20190712	0.00053	<0.00053	mg/L
	WG-1620-MW09-20190712	0.00015	<0.00015	mg/L				

Table 5

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units	
SVOCs	WG-1620-FB04-20190716	07/16/2019	bis(2-Ethylhexyl)phthalate (DEHP)	0.00045	WG-1620-MW21C-20190716	0.000040 J	<0.000040	mg/L	
					WG-1620-MW28A-20190716	0.00023	<0.00023	mg/L	
					WG-1620-MW28C-20190716	0.00013 J	<0.00013	mg/L	
					WG-1620-MW36B-20190716	0.000088 J	<0.000088	mg/L	
				Di-n-butylphthalate (DBP) Phenol	0.000043 J	WG-1620-FD01-20190716	0.000020 J	<0.000020	mg/L
			0.000084 J		WG-1620-MW21C-20190716	0.00018 J	<0.00018 J	mg/L	
					WG-1620-MW28A-20190716	0.000073 J	<0.000073	mg/L	
					WG-1620-MW36A-20190716	0.00014 J	<0.00014	mg/L	
					WG-1620-MW36B-20190716	0.00017 J	<0.00017	mg/L	
					WG-1620-MW53C-20190716	0.00016 J	<0.00016	mg/L	
SVOCs	WG-1620-FB05-20190717	07/17/2019	bis(2-Ethylhexyl)phthalate (DEHP)	0.00023	WG-1620-FD02-20190717	0.000084 J	<0.000084	mg/L	
					WG-1620-MW26A-20190717	0.00019 J	<0.00019	mg/L	
					WG-1620-MW33A-20190717	0.00022	<0.00022	mg/L	
					WG-1620-MW44A-20190717	0.00025	<0.00025	mg/L	
					WG-1620-MW47C-20190717	0.00027	<0.00027	mg/L	
					WG-1620-MW48C-20190717	0.00030	<0.00030	mg/L	
					WG-1620-MW59A-20190717	.000050	.000050	mg/L	
					WG-1620-MW59B-20190717	0.00039	<0.00039	mg/L	
					WG-1620-MW60A-20190717	0.00055	<0.00055	mg/L	

Table 5

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB05-20190717	07/17/2019	bis(2-Ethylhexyl)phthalate (DEHP)	0.00023	WG-1620-MW61A-20190717	0.00037	<0.00037	mg/L
					WG-1620-MW69A-20190717	0.00034	<0.00034	mg/L
					WG-1620-MW71B-20190717	0.00014 J	<0.00014	mg/L
					WG-1620-MW87C-20190717	0.00018 J	<0.00018	mg/L
			Naphthalene	0.00023	WG-1620-FD02-20190717	0.000096 J	<0.000096	mg/L
					WG-1620-MW26A-20190717	0.00032	<0.00032	mg/L
					WG-1620-MW33A-20190717	0.000076 J	<0.000076	mg/L
					WG-1620-MW44A-20190717	0.00019	<0.00019	mg/L
					WG-1620-MW59A-20190717	0.00012	<0.00012	mg/L
					WG-1620-MW59B-20190717	0.00015	<0.00015	mg/L
					WG-1620-MW60A-20190717	0.00013	<0.00013	mg/L
					WG-1620-MW69A-20190717	0.000078 J	<0.000078	mg/L
					WG-1620-MW87C-20190717	0.00017	<0.00017	mg/L

## Notes:

SVOCs - Semi-volatile Organic Compounds

VOCs - Volatile Organic Compounds

J - Estimated concentration

&lt; - Not detected at the associated reporting limit

Table 6

**Qualified Sample Data Due to Variability in Field Duplicate Results**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**July 2019**

Parameter	Analyte	RPD	Sample ID	Qualified Result	Field Duplicate Sample ID	Qualified Result	Units
SVOCs	2-Methylnaphthalene	155	WG-1620-MW21C-20190716	0.00040 J	WG-1620-FD01-20190716	0.000050 J	mg/L
	Acenaphthene	150		0.0057 J		0.00081 J	mg/L
	Anthracene	159		0.00044 J		0.000050 J	mg/L
	Dibenzofuran	155		0.0032 J		0.00040 J	mg/L
	Fluoranthene	161		0.00081 J		0.000087 J	mg/L
	Fluorene	153		0.0019 J		0.00025 J	mg/L
	Naphthalene	138		0.0044 J		0.00080 J	mg/L
	Phenanthrene	155		0.00087 J		0.00011 J	mg/L
	Phenol	120		<0.00020 J		0.00081 J	mg/L
Pyrene	168	0.00039 J	0.000033 J	mg/L			
SVOCs	Anthracene	30.8	WG-1620-MW68B-20190718	0.030 J	WG-1620-FD03-20190718	0.022 J	mg/L
	Benzo(a)anthracene	51.9		0.0051 J		0.0030 J	mg/L
	Chrysene	53.2		0.0050 J		0.0029 J	mg/L
	Fluoranthene	45.7		0.043 J		0.027 J	mg/L
	Pyrene	50.0		0.025 J		0.015 J	mg/L

## Notes:

RPD - Relative Percent Difference

SVOCs - Semi-volatile Organic Compounds

J - Estimated concentration

&lt; - Not detected at the associated reporting limit

# Attachment A

## Laboratory NELAP Certificate



# Texas Commission on Environmental Quality

## NELAP - Recognized Laboratory Fields of Accreditation



ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: *Drinking Water***

**Method** EPA 1613

Analyte	AB	Analyte ID	Method ID
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408

**Method** EPA 200.8

Analyte	AB	Analyte ID	Method ID
Copper	TX	1055	10014605
Lead	TX	1075	10014605





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**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 120.1			
Analyte Conductivity	TX	1610	10006403
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 160.4			
Analyte Residue-volatile	TX	1970	10010409
Method EPA 1613			
Analyte 1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10120408
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10120408
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10120408
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10120408
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10120408
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10120408
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10120408
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10120408
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,6,7,8-HxCDD)	TX	9456	10120408
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10120408
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10120408
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10120408
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10120408
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10120408



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**Matrix: Non-Potable Water**

2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10120408
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10120408
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10120408
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10120408
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10120408
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10120408
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10120408
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10120408
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10120408
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10120408
<b>Method EPA 1664</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10127807
<b>Method EPA 180.1</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Turbidity	TX	2055	10011606
<b>Method EPA 200.8</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Aluminum	TX	1000	10014605
Antimony	TX	1005	10014605
Arsenic	TX	1010	10014605
Barium	TX	1015	10014605
Beryllium	TX	1020	10014605
Boron	TX	1025	10014605
Cadmium	TX	1030	10014605
Calcium	TX	1035	10014605
Chromium	TX	1040	10014605
Cobalt	TX	1050	10014605
Copper	TX	1055	10014605
Iron	TX	1070	10014605



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**Matrix: Non-Potable Water**

Lead	TX	1075	10014605
Magnesium	TX	1085	10014605
Manganese	TX	1090	10014605
Molybdenum	TX	1100	10014605
Nickel	TX	1105	10014605
Potassium	TX	1125	10014605
Selenium	TX	1140	10014605
Silver	TX	1150	10014605
Sodium	TX	1155	10014605
Strontium	TX	1160	10014605
Thallium	TX	1165	10014605
Tin	TX	1175	10014605
Titanium	TX	1180	10014605
Uranium	TX	3035	10014605
Vanadium	TX	1185	10014605
Zinc	TX	1190	10014605

**Method EPA 245.1**

Analyte	AB	Analyte ID	Method ID
Mercury	TX	1095	10036609

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 325.1**

Analyte	AB	Analyte ID	Method ID
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**Matrix: Non-Potable Water**

Chloride	TX	1575	10056801
<b>Method</b> EPA 335.1			
<b>Analyte</b> Amenable cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1510	<b>Method ID</b> 10060001
<b>Method</b> EPA 335.2			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10278203
<b>Method</b> EPA 335.4			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10061402
<b>Method</b> EPA 350.3			
<b>Analyte</b> Ammonia as N	<b>AB</b> TX	<b>Analyte ID</b> 1515	<b>Method ID</b> 10064401
<b>Method</b> EPA 365.3			
<b>Analyte</b> Orthophosphate as P Phosphorus	<b>AB</b> TX TX	<b>Analyte ID</b> 1870 1910	<b>Method ID</b> 10070801 10070801
<b>Method</b> EPA 375.4			
<b>Analyte</b> Sulfate	<b>AB</b> TX	<b>Analyte ID</b> 2000	<b>Method ID</b> 10073800
<b>Method</b> EPA 376.1			
<b>Analyte</b> Sulfide	<b>AB</b> TX	<b>Analyte ID</b> 2005	<b>Method ID</b> 10074201
<b>Method</b> EPA 410.4			
<b>Analyte</b> Chemical oxygen demand (COD)	<b>AB</b> TX	<b>Analyte ID</b> 1565	<b>Method ID</b> 10077404
<b>Method</b> EPA 415.1			
<b>Analyte</b> Total Organic Carbon (TOC)	<b>AB</b> TX	<b>Analyte ID</b> 2040	<b>Method ID</b> 10078407
<b>Method</b> EPA 420.1			
<b>Analyte</b> Total phenolics	<b>AB</b> TX	<b>Analyte ID</b> 1905	<b>Method ID</b> 10079400



# Texas Commission on Environmental Quality



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**Matrix: Non-Potable Water**

**Method EPA 420.4**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10080203

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156419
Antimony	TX	1005	10156419
Arsenic	TX	1010	10156419
Barium	TX	1015	10156419
Beryllium	TX	1020	10156419
Boron	TX	1025	10156419
Cadmium	TX	1030	10156419
Calcium	TX	1035	10156419
Chromium	TX	1040	10156419
Cobalt	TX	1050	10156419
Copper	TX	1055	10156419
Iron	TX	1070	10156419
Lead	TX	1075	10156419
Lithium	TX	1080	10156419
Magnesium	TX	1085	10156419
Manganese	TX	1090	10156419
Molybdenum	TX	1100	10156419
Nickel	TX	1105	10156419
Potassium	TX	1125	10156419
Selenium	TX	1140	10156419
Silver	TX	1150	10156419
Sodium	TX	1155	10156419
Strontium	TX	1160	10156419
Thallium	TX	1165	10156419
Tin	TX	1175	10156419
Titanium	TX	1180	10156419



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**Matrix: Non-Potable Water**

Vanadium	TX	1185	10156419
Zinc	TX	1190	10156419
<b>Method EPA 608</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
4,4'-DDD	TX	7355	10103603
4,4'-DDE	TX	7360	10103603
4,4'-DDT	TX	7365	10103603
Aldrin	TX	7025	10103603
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10103603
alpha-Chlordane	TX	7240	10103603
Aroclor-1016 (PCB-1016)	TX	8880	10103603
Aroclor-1221 (PCB-1221)	TX	8885	10103603
Aroclor-1232 (PCB-1232)	TX	8890	10103603
Aroclor-1242 (PCB-1242)	TX	8895	10103603
Aroclor-1248 (PCB-1248)	TX	8900	10103603
Aroclor-1254 (PCB-1254)	TX	8905	10103603
Aroclor-1260 (PCB-1260)	TX	8910	10103603
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10103603
Chlordane (tech.)	TX	7250	10103603
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10103603
Dieldrin	TX	7470	10103603
Endosulfan I	TX	7510	10103603
Endosulfan II	TX	7515	10103603
Endosulfan sulfate	TX	7520	10103603
Endrin	TX	7540	10103603
Endrin aldehyde	TX	7530	10103603
Endrin ketone	TX	7535	10103603
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10103603
gamma-Chlordane	TX	7245	10103603
Heptachlor	TX	7685	10103603





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**Matrix: Non-Potable Water**

Heptachlor epoxide	TX	7690	10103603
Methoxychlor	TX	7810	10103603
Toxaphene (Chlorinated camphene)	TX	8250	10103603

**Method EPA 624**

Analyte	AB	Analyte ID	Method ID
1,1,1-Trichloroethane	TX	5160	10107207
1,1,2,2-Tetrachloroethane	TX	5110	10107207
1,1,2-Trichloroethane	TX	5165	10107207
1,1-Dichloroethane	TX	4630	10107207
1,1-Dichloroethylene	TX	4640	10107207
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10107207
1,2-Dichlorobenzene	TX	4610	10107207
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10107207
1,2-Dichloropropane	TX	4655	10107207
1,3-Dichlorobenzene	TX	4615	10107207
1,4-Dichlorobenzene	TX	4620	10107207
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10107207
2-Chloroethyl vinyl ether	TX	4500	10107207
Acetone (2-Propanone)	TX	4315	10107207
Acrolein (Propenal)	TX	4325	10107207
Acrylonitrile	TX	4340	10107207
Benzene	TX	4375	10107207
Bromodichloromethane	TX	4395	10107207
Bromoform	TX	4400	10107207
Carbon tetrachloride	TX	4455	10107207
Chlorobenzene	TX	4475	10107207
Chlorodibromomethane	TX	4575	10107207
Chloroethane (Ethyl chloride)	TX	4485	10107207
Chloroform	TX	4505	10107207
cis-1,2-Dichloroethylene	TX	4645	10107207



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**Matrix: Non-Potable Water**

cis-1,3-Dichloropropene	TX	4680	10107207
Ethylbenzene	TX	4765	10107207
m+p-xylene	TX	5240	10107207
Methyl bromide (Bromomethane)	TX	4950	10107207
Methyl chloride (Chloromethane)	TX	4960	10107207
Methyl tert-butyl ether (MTBE)	TX	5000	10107207
Methylene chloride (Dichloromethane)	TX	4975	10107207
Naphthalene	TX	5005	10107207
o-Xylene	TX	5250	10107207
Tetrachloroethylene (Perchloroethylene)	TX	5115	10107207
Toluene	TX	5140	10107207
trans-1,2-Dichloroethylene	TX	4700	10107207
trans-1,3-Dichloropropylene	TX	4685	10107207
Trichloroethene (Trichloroethylene)	TX	5170	10107207
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10107207
Vinyl chloride	TX	5235	10107207
Xylene (total)	TX	5260	10107207

**Method EPA 625**

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10107401
1,2,4-Trichlorobenzene	TX	5155	10107401
1,2-Dichlorobenzene	TX	4610	10107401
1,2-Diphenylhydrazine	TX	6220	10107401
1,3-Dichlorobenzene	TX	4615	10107401
1,4-Dichlorobenzene	TX	4620	10107401
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10107401
2,4,5-Trichlorophenol	TX	6835	10107401
2,4,6-Trichlorophenol	TX	6840	10107401
2,4-Dichlorophenol	TX	6000	10107401
2,4-Dimethylphenol	TX	6130	10107401





# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

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**Matrix: Non-Potable Water**

2,4-Dinitrophenol	TX	6175	10107401
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10107401
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10107401
2-Chloronaphthalene	TX	5795	10107401
2-Chlorophenol	TX	5800	10107401
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10107401
2-Methylphenol (o-Cresol)	TX	6400	10107401
2-Nitrophenol	TX	6490	10107401
3,3'-Dichlorobenzidine	TX	5945	10107401
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10107401
4-Chloro-3-methylphenol	TX	5700	10107401
4-Chlorophenyl phenylether	TX	5825	10107401
4-Methylphenol (p-Cresol)	TX	6410	10107401
4-Nitrophenol	TX	6500	10107401
Acenaphthene	TX	5500	10107401
Acenaphthylene	TX	5505	10107401
Anthracene	TX	5555	10107401
Benzidine	TX	5595	10107401
Benzo(a)anthracene	TX	5575	10107401
Benzo(a)pyrene	TX	5580	10107401
Benzo(b)fluoranthene	TX	5585	10107401
Benzo(g,h,i)perylene	TX	5590	10107401
Benzo(k)fluoranthene	TX	5600	10107401
bis(2-Chloroethoxy)methane	TX	5760	10107401
bis(2-Chloroethyl) ether	TX	5765	10107401
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10107401
Butyl benzyl phthalate	TX	5670	10107401
Chrysene	TX	5855	10107401
Dibenz(a,h) anthracene	TX	5895	10107401
Diethyl phthalate	TX	6070	10107401



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**Matrix: Non-Potable Water**

Dimethyl phthalate	TX	6135	10107401
Di-n-butyl phthalate	TX	5925	10107401
Di-n-octyl phthalate	TX	6200	10107401
Fluoranthene	TX	6265	10107401
Fluorene	TX	6270	10107401
Hexachlorobenzene	TX	6275	10107401
Hexachlorobutadiene	TX	4835	10107401
Hexachlorocyclopentadiene	TX	6285	10107401
Hexachloroethane	TX	4840	10107401
Indeno(1,2,3-cd) pyrene	TX	6315	10107401
Isophorone	TX	6320	10107401
Naphthalene	TX	5005	10107401
Nitrobenzene	TX	5015	10107401
n-Nitrosodiethylamine	TX	6525	10107401
n-Nitrosodimethylamine	TX	6530	10107401
n-Nitrosodi-n-butylamine	TX	5025	10107401
n-Nitrosodi-n-propylamine	TX	6545	10107401
n-Nitrosodiphenylamine	TX	6535	10107401
Pentachlorobenzene	TX	6590	10107401
Pentachlorophenol	TX	6605	10107401
Phenanthrene	TX	6615	10107401
Phenol	TX	6625	10107401
Pyrene	TX	6665	10107401
Pyridine	TX	5095	10107401
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603



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**Matrix: Non-Potable Water**

**Method EPA 8011**

Analyte	AB	Analyte ID	Method ID
1,2,3-Trichloropropane	TX	5180	10173009
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10173009
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10173009

**Method EPA 8015**

Analyte	AB	Analyte ID	Method ID
Diesel range organics (DRO)	TX	9369	10173203
Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402



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**Matrix: Non-Potable Water**

4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402
Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Hexachlorobenzene	TX	6275	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201



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**Matrix: Non-Potable Water**

Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8151**

Analyte	AB	Analyte ID	Method ID
2,4,5-T	TX	8655	10183003
2,4-D	TX	8545	10183003
2,4-DB	TX	8560	10183003
Dalapon	TX	8555	10183003
Dicamba	TX	8595	10183003
Dichloroprop (Dichloroprop, Weedone)	TX	8605	10183003
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10183003
MCPA	TX	7775	10183003
MCPP	TX	7780	10183003
Silvex (2,4,5-TP)	TX	8650	10183003

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404
1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404



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**Matrix: Non-Potable Water**

1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
2-Pentanone	TX	5045	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404
Allyl alcohol	TX	4350	10184404
Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404





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**Matrix: Non-Potable Water**

Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Di-isopropylether (DIPE)	TX	9375	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Ethyl-t-butylether (ETBE) (2-Ethoxy-2-methylpropane)	TX	4770	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404
Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404



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**Matrix: Non-Potable Water**

Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
T-amylmethylether (TAME)	TX	4370	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404
Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404





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**Matrix: Non-Potable Water**

Method EPA 8270

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203
2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203



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**Matrix: Non-Potable Water**

2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Dimethyl aminoazobenzene	TX	6105	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrobiphenyl	TX	6480	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Chloro-2-methylaniline	TX	5695	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203



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**Matrix: Non-Potable Water**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Captan	TX	7190	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203



# Texas Commission on Environmental Quality



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**Matrix: Non-Potable Water**

Chrysene	TX	5855	10185203
Coumaphos	TX	7315	10185203
Demeton	TX	7390	10185203
Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Dioxathion	TX	7495	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethion	TX	7565	10185203
Ethyl methanesulfonate	TX	6260	10185203
Famphur	TX	7580	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203



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**Matrix: Non-Potable Water**

Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203
Kepone	TX	7740	10185203
Maleic anhydride	TX	6335	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naled	TX	7905	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203



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**Matrix: Non-Potable Water**

Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Phosmet (Imidan)	TX	8000	10185203
Phthalic anhydride	TX	6640	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203
Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Resorcinol	TX	6680	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203
Trifluralin (Treflan)	TX	8295	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209





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**Matrix: Non-Potable Water**

1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209

**Method EPA 8316**

Analyte	AB	Analyte ID	Method ID
Acrylamide	TX	4330	10188202

**Method EPA 8330**

Analyte	AB	Analyte ID	Method ID
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807



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**Matrix: Non-Potable Water**

Methyl-2,4,6-trinitrophenylamine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807
<b>Method EPA 9014</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803
<b>Method EPA 9038</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Sulfate	TX	2000	10196608
<b>Method EPA 9040</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
pH	TX	1900	10196802
<b>Method EPA 9050</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	10198604
<b>Method EPA 9056</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209
<b>Method EPA 9060</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	10200201





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**Matrix: Non-Potable Water**

Method	Analyte	AB	Analyte ID	Method ID
EPA 9065	Total phenolics	TX	1905	10200405
EPA 9066	Total phenolics	TX	1905	10200609
EPA 9250	Chloride	TX	1575	10207202
EPA RSK 175	2-methylpropane (Isobutane)	TX	4942	10212905
	Ethane	TX	4747	10212905
	Ethene	TX	4752	10212905
	Methane	TX	4926	10212905
	n-Butane	TX	5007	10212905
	n-Propane	TX	5029	10212905
HACH 8000	Chemical oxygen demand (COD)	TX	1565	60003001
SM 2120 B	Color	TX	1605	20223807
SM 2310 B (4a)	Acidity, as CaCO <sub>3</sub>	TX	1500	20002806
SM 2320 B	Alkalinity as CaCO <sub>3</sub>	TX	1505	20045005
SM 2340 B		AB	Analyte ID	Method ID



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**Matrix: Non-Potable Water**

Total hardness as CaCO <sub>3</sub>	TX	1755	20046008
<b>Method</b> SM 2510 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	20048004
<b>Method</b> SM 2540 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-total (total solids)	TX	1950	20004608
<b>Method</b> SM 2540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-filterable (TDS)	TX	1955	20049803
<b>Method</b> SM 2540 D			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-nonfilterable (TSS)	TX	1960	20004802
<b>Method</b> SM 3500-Cr B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	20065809
<b>Method</b> SM 4500-Cl F			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total residual chlorine	TX	1940	20080482
<b>Method</b> SM 4500-Cl <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chloride	TX	1575	20019209
<b>Method</b> SM 4500-CN <sup>-</sup> C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20020808
<b>Method</b> SM 4500-CN <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20021209
<b>Method</b> SM 4500-CN <sup>-</sup> G			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	20021607



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**Matrix: Non-Potable Water**

Method	Analyte	AB	Analyte ID	Method ID
SM 4500-H+ B				
	pH	TX	1900	20104603
SM 4500-NH3 D				
	Ammonia as N	TX	1515	20108809
	Kjeldahl Nitrogen (Total Kjeldahl Nitrogen-TKN)	TX	1790	20108809
SM 4500-NH3 F				
	Ammonia as N	TX	1515	20023001
SM 4500-O G				
	Oxygen, dissolved	TX	1880	20025405
SM 4500-P E				
	Orthophosphate as P	TX	1870	20025803
	Phosphorus	TX	1910	20025803
SM 4500-S2 <sup>-</sup> F				
	Sulfide	TX	2005	20126209
SM 4500-SiO2 D				
	Silica as SiO2	TX	1990	20127202
SM 4500-SO3 <sup>-</sup> B				
	Sulfite	TX	2015	20026806
SM 5210 B				
	Biochemical oxygen demand (BOD)	TX	1530	20027401
	Carbonaceous BOD, CBOD	TX	1555	20027401
SM 5310 B				
	Analyte	AB	Analyte ID	Method ID



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**Matrix: Non-Potable Water**

Total Organic Carbon (TOC)	TX	2040	20137206
<b>Method</b> SM 5310 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	20138209
<b>Method</b> SM 5540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Surfactants - MBAS	TX	2025	20144405
<b>Method</b> TCEQ 1005			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
Method ASTM D2216			
Analyte Moisture	TX	10337	ASTM D2216-05
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 1030			
Analyte Ignitability	TX	1780	10117201
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 1668			
Analyte Decachlorobiphenyls	TX	10332	10262007
Dichlorobiphenyls	TX	464	10262007
Heptachlorobiphenyls	TX	486	10262007
Hexachlorobiphenyls	TX	487	10262007
Monochlorobiphenyls	TX	501	10262007
Nonachlorobiphenyls	TX	507	10262007
Octachlorobiphenyls	TX	508	10262007
Pentachlorobiphenyls	TX	515	10262007
Tetrachlorobiphenyls	TX	528	10262007
Trichlorobiphenyls	TX	541	10262007
Method EPA 200.8			
Analyte Uranium	TX	3035	10014605



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**Matrix: Solid & Chemical Materials**

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 310.1**

Analyte	AB	Analyte ID	Method ID
Alkalinity as CaCO3	TX	1505	10054805

**Method EPA 350.3**

Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	10064401

**Method EPA 365.3**

Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	10070801
Phosphorus	TX	1910	10070801

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156204
Antimony	TX	1005	10156204
Arsenic	TX	1010	10156204
Barium	TX	1015	10156204
Beryllium	TX	1020	10156204
Boron	TX	1025	10156204
Cadmium	TX	1030	10156204
Calcium	TX	1035	10156204
Chromium	TX	1040	10156204



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**Matrix: Solid & Chemical Materials**

Cobalt	TX	1050	10156204
Copper	TX	1055	10156204
Iron	TX	1070	10156204
Lead	TX	1075	10156204
Lithium	TX	1080	10156204
Magnesium	TX	1085	10156204
Manganese	TX	1090	10156204
Molybdenum	TX	1100	10156204
Nickel	TX	1105	10156204
Potassium	TX	1125	10156204
Selenium	TX	1140	10156204
Silver	TX	1150	10156204
Sodium	TX	1155	10156204
Strontium	TX	1160	10156204
Thallium	TX	1165	10156204
Tin	TX	1175	10156204
Titanium	TX	1180	10156204
Vanadium	TX	1185	10156204
Zinc	TX	1190	10156204
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603
<b>Method EPA 7471</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10166004
<b>Method EPA 8015</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Diesel range organics (DRO)	TX	9369	10173203





# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
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Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

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**Matrix: Solid & Chemical Materials**

Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402
4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402





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**Matrix: Solid & Chemical Materials**

Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201
Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404



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**Matrix: Solid & Chemical Materials**

1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404



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**Matrix: Solid & Chemical Materials**

Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404
Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404



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**Matrix: Solid & Chemical Materials**

Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404
Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404



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**Matrix: Solid & Chemical Materials**

Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404
<b>Method EPA 8270</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203



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**Matrix: Solid & Chemical Materials**

2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203
2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203





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**Matrix: Solid & Chemical Materials**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203
Chrysene	TX	5855	10185203



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**Matrix: Solid & Chemical Materials**

Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzo(a,e) pyrene	TX	5890	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethyl methanesulfonate	TX	6260	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203
Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203





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**Matrix: Solid & Chemical Materials**

Kepone	TX	7740	10185203
Malathion	TX	7770	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203
Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203



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**Matrix: Solid & Chemical Materials**

Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: Solid & Chemical Materials**

Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209
<b>Method EPA 8316</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Acrylamide	TX	4330	10188202
<b>Method EPA 8330</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807
Methyl-2,4,6-trinitrophenylnitramine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807
<b>Method EPA 9014</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803
<b>Method EPA 9038</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Sulfate	TX	2000	10196608



# Texas Commission on Environmental Quality

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**Matrix: Solid & Chemical Materials**

**Method EPA 9040**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197203
pH	TX	1900	10196802

**Method EPA 9045**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197805
pH	TX	1900	10197805

**Method EPA 9050**

Analyte	AB	Analyte ID	Method ID
Conductivity	TX	1610	10198604

**Method EPA 9056**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209

**Method EPA 9060**

Analyte	AB	Analyte ID	Method ID
Total Organic Carbon (TOC)	TX	2040	10200201

**Method EPA 9065**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10200405

**Method EPA 9071**

Analyte	AB	Analyte ID	Method ID
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10201204



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
Method EPA 9095			
Analyte Paint Filter Liquids Test	TX	10312	10204009
Method EPA 9250			
Analyte Chloride	TX	1575	10207202
Method SM 2320 B			
Analyte Alkalinity as CaCO3	TX	1505	20045005
Method SM 2510 B			
Analyte Conductivity	TX	1610	20048004
Method SM 2540 G			
Analyte Residue-total (total solids)	TX	1950	20005203
Method SSA/ASA Part 3:34			
Analyte Carbon, organic (Walkley-Black)	TX	10340	SSA/ASA Pt 3:34
Method TCEQ 1005			
Analyte Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

July 19, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19070575**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 21 sample(s) on Jul 11, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey



Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/19/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070575			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143014,143067,143129,R342400,R342410,R342474			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			2
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?	X				
		Were analytical duplicates analyzed at the appropriate frequency?	X				
		Were RPDs or relative standard deviations within the laboratory QC limits?	X				
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?			X		
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?			X		
		Are unadjusted MQLs and DCSs included in the laboratory data package?			X		
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				3
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/19/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070575			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143014,143067,143129,R342400,R342410,R342474			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 07/19/2019
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS19070575
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 143014,143067,143129,R342400,R342410,R342474
ER# <sup>5</sup>	Description	
1	Semivolatile Organics Method SW8270, samples WG-1620-MW18A-20190710, WG-1620-MW18C-20190710, WG-1620-MW58A-20190710, WG-1620-MW57A-20190710, WG-1620-MW57B-20190710, WG-1620-MW72B-20190710, WG-1620-MW17C-20190710, WG-1620-MW17-20190710, WG-1620-MW20A-20190710, WG-1620-MW15B-20190710, WG-1620-MW12A-20190711, the surrogate recoveries could not be determined due to dilution below the calibration range.	
2	Batch 143014, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.	
3	Batch 143014, Semivolatile Organics Method SW8270, samples WG-1620-MW57B-20190710, WG-1620-MW72B-20190710, WG-1620-MW17-20190710, the GCMS semi-volatile extract of the samples were run at a dilution due to a high level of matrix interference.	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.                      O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);                      NA = Not Applicable;                      NR = Not Reviewed;                      R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070575

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19070575-01	WG-1620-MW18A-20190710	Water		10-Jul-2019 07:20	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-02	WG-1620-MW18C-20190710	Water		10-Jul-2019 08:05	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-03	WG-1620-MW58A-20190710	Water		10-Jul-2019 08:55	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-04	WG-1620-MW57A-20190710	Water		10-Jul-2019 09:55	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-05	WG-1620-MW57B-20190710	Water		10-Jul-2019 10:40	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-06	WG-1620-MW72B-20190710	Water		10-Jul-2019 11:35	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-07	WG-1620-MW19C-20190710	Water		10-Jul-2019 12:25	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-08	WG-1620-MW17C-20190710	Water		10-Jul-2019 13:25	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-09	WG-1620-MW17-20190710	Water		10-Jul-2019 14:15	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-10	WG-1620-MW20A-20190710	Water		10-Jul-2019 15:10	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-11	WG-1620-MW15A-20190710	Water		10-Jul-2019 16:10	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-12	WG-1620-MW15C-20190710	Water		10-Jul-2019 17:05	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-13	WG-1620-MW15B-20190710	Water		10-Jul-2019 18:00	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-14	WG-1620-FB01-20190710	Water		10-Jul-2019 18:15	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-15	WG-1620-MW14-20190711	Water		11-Jul-2019 07:20	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-16	WG-1620-MW13-20190711	Water		11-Jul-2019 08:15	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-17	WG-1620-MW39B-20190711	Water		11-Jul-2019 09:15	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-18	WG-1620-MW12C-20190711	Water		11-Jul-2019 10:00	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-19	WG-1620-MW12A-20190711	Water		11-Jul-2019 10:55	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-20	WQ-1620-TB01-20190710	Water	C&G-062119-76	10-Jul-2019 00:00	11-Jul-2019 16:45	<input type="checkbox"/>
HS19070575-21	WQ-1620-TB02-20190710	Water	C&G-062119-84	10-Jul-2019 00:00	11-Jul-2019 16:45	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18A-20190710  
 Collection Date: 10-Jul-2019 07:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:56
<b>Benzene</b>	<b>0.16</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 04:56
<b>Chlorobenzene</b>	<b>0.0017</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 04:56
<b>Ethylbenzene</b>	<b>0.28</b>		<b>0.0015</b>	<b>0.0050</b>	<b>mg/L</b>	5	15-Jul-2019 19:23
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 04:56
<b>Toluene</b>	<b>0.057</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 04:56
<b>Vinyl chloride</b>	<b>0.0013</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 04:56
<b>Xylenes, Total</b>	<b>0.47</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 04:56
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>83.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:56</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.9</i>			<i>70-126</i>	<i>%REC</i>	<i>5</i>	<i>15-Jul-2019 19:23</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:56</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.8</i>			<i>81-113</i>	<i>%REC</i>	<i>5</i>	<i>15-Jul-2019 19:23</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:56</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.2</i>			<i>77-123</i>	<i>%REC</i>	<i>5</i>	<i>15-Jul-2019 19:23</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:56</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>5</i>	<i>15-Jul-2019 19:23</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18A-20190710  
 Collection Date: 10-Jul-2019 07:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>	<b>Method:SW8270</b>			Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 13:05
<b>2,4-Dimethylphenol</b>	<b>0.72</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	100	15-Jul-2019 19:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 13:05
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 13:05
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 13:05
<b>2-Methylnaphthalene</b>	<b>0.47</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 19:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 13:05
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 13:05
<b>Acenaphthene</b>	<b>0.30</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 19:37
<b>Acenaphthylene</b>	<b>0.0056</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:05
<b>Anthracene</b>	<b>0.0063</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:05
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 13:05
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 13:05
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 13:05
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 13:05
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 13:05
<b>Dibenzofuran</b>	<b>0.19</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 19:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 13:05
<b>Fluoranthene</b>	<b>0.0022</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:05
<b>Fluorene</b>	<b>0.091</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 19:17
<b>Naphthalene</b>	<b>5.6</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	15-Jul-2019 19:56
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 13:05
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 13:05
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 13:05
<b>Phenanthrene</b>	<b>0.075</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 19:17
<b>Phenol</b>	<b>0.00048</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 13:05
<b>Pyrene</b>	<b>0.0011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:05
<i>Surr: 2,4,6-Tribromophenol</i>	<i>77.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 13:05</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>89.8</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:17</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 19:37</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>15-Jul-2019 19:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>15-Jul-2019 19:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>49.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 13:05</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>62.1</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 19:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 13:05</i>
<i>Surr: 2-Fluorophenol</i>	<i>74.7</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:17</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 19:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>15-Jul-2019 19:56</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18A-20190710  
 Collection Date: 10-Jul-2019 07:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	15-Jul-2019 19:56
Surr: 4-Terphenyl-d14	69.4			40-135	%REC	1	15-Jul-2019 13:05
Surr: 4-Terphenyl-d14	92.2			40-135	%REC	10	15-Jul-2019 19:17
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	15-Jul-2019 19:37
Surr: Nitrobenzene-d5	56.3			41-120	%REC	1	15-Jul-2019 13:05
Surr: Nitrobenzene-d5	71.1			41-120	%REC	10	15-Jul-2019 19:17
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	15-Jul-2019 19:37
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	15-Jul-2019 19:56
Surr: Phenol-d6	71.2			20-120	%REC	1	15-Jul-2019 13:05
Surr: Phenol-d6	65.2			20-120	%REC	10	15-Jul-2019 19:17
Surr: Phenol-d6	0	JS		20-120	%REC	100	15-Jul-2019 19:37
Surr: Phenol-d6	0	JS		20-120	%REC	1000	15-Jul-2019 19:56
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.0248		0.000400	0.00200	mg/L	1	16-Jul-2019 22:38

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18C-20190710  
 Collection Date: 10-Jul-2019 08:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 06:09
<b>Benzene</b>	<b>1.2</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jul-2019 19:50
<b>Chlorobenzene</b>	<b>0.00052</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 06:09
<b>Ethylbenzene</b>	<b>0.32</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jul-2019 19:50
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 06:09
<b>Toluene</b>	<b>1.0</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jul-2019 19:50
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jul-2019 06:09
<b>Xylenes, Total</b>	<b>1.0</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jul-2019 19:50
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>81.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:09</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.1</i>			<i>70-126</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:50</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:09</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.3</i>			<i>81-113</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:50</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:09</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.9</i>			<i>77-123</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:50</i>
<i>Surr: Toluene-d8</i>	<i>100.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:09</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 19:50</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18C-20190710  
 Collection Date: 10-Jul-2019 08:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 13:25
<b>2,4-Dimethylphenol</b>	<b>0.0051</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 13:25
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 13:25
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 13:25
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 13:25
<b>2-Methylnaphthalene</b>	<b>0.28</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 20:36
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 13:25
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 13:25
<b>Acenaphthene</b>	<b>0.058</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 20:16
<b>Acenaphthylene</b>	<b>0.0050</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:25
<b>Anthracene</b>	<b>0.0056</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:25
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 13:25
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 13:25
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 13:25
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 13:25
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 13:25
<b>Dibenzofuran</b>	<b>0.054</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 20:16
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 13:25
<b>Fluoranthene</b>	<b>0.0019</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:25
<b>Fluorene</b>	<b>0.024</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 20:16
<b>Naphthalene</b>	<b>9.9</b>		<b>0.10</b>	<b>0.50</b>	<b>mg/L</b>	5000	16-Jul-2019 13:47
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 13:25
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 13:25
<b>Pentachlorophenol</b>	<b>0.034</b>		<b>0.00079</b>	<b>0.0020</b>	<b>mg/L</b>	10	15-Jul-2019 20:16
<b>Phenanthrene</b>	<b>0.026</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 20:16
<b>Phenol</b>	<b>0.0050</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 13:25
<b>Pyrene</b>	<b>0.0014</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:25
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	5000	16-Jul-2019 13:47
Surr: 2,4,6-Tribromophenol	38.0			34-129	%REC	1	15-Jul-2019 13:25
Surr: 2,4,6-Tribromophenol	91.7			34-129	%REC	10	15-Jul-2019 20:16
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	15-Jul-2019 20:36
Surr: 2-Fluorobiphenyl	68.1			40-125	%REC	10	15-Jul-2019 20:16
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	15-Jul-2019 20:36
Surr: 2-Fluorobiphenyl	42.6			40-125	%REC	1	15-Jul-2019 13:25
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	5000	16-Jul-2019 13:47
Surr: 2-Fluorophenol	0	JS		20-120	%REC	5000	16-Jul-2019 13:47
Surr: 2-Fluorophenol	43.8			20-120	%REC	1	15-Jul-2019 13:25
Surr: 2-Fluorophenol	114			20-120	%REC	10	15-Jul-2019 20:16
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	15-Jul-2019 20:36

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18C-20190710  
 Collection Date: 10-Jul-2019 08:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	74.9			40-135	%REC	10	15-Jul-2019 20:16
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	15-Jul-2019 20:36
Surr: 4-Terphenyl-d14	68.0			40-135	%REC	1	15-Jul-2019 13:25
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	5000	16-Jul-2019 13:47
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	5000	16-Jul-2019 13:47
Surr: Nitrobenzene-d5	57.2			41-120	%REC	1	15-Jul-2019 13:25
Surr: Nitrobenzene-d5	44.0			41-120	%REC	10	15-Jul-2019 20:16
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	15-Jul-2019 20:36
Surr: Phenol-d6	45.4			20-120	%REC	1	15-Jul-2019 13:25
Surr: Phenol-d6	71.7			20-120	%REC	10	15-Jul-2019 20:16
Surr: Phenol-d6	0	JS		20-120	%REC	100	15-Jul-2019 20:36
Surr: Phenol-d6	0	JS		20-120	%REC	5000	16-Jul-2019 13:47
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.00358		0.000400	0.00200	mg/L	1	16-Jul-2019 22:40

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW58A-20190710  
 Collection Date: 10-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	14-Jul-2019 05:20
<b>Benzene</b>	<b>0.0049</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 05:20
<b>Chlorobenzene</b>	<b>0.00080</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 05:20
<b>Ethylbenzene</b>	<b>0.017</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 05:20
Methylene chloride		U	0.0010	0.0020	mg/L	1	14-Jul-2019 05:20
Toluene		U	0.00020	0.0010	mg/L	1	14-Jul-2019 05:20
Vinyl chloride		U	0.00020	0.0010	mg/L	1	14-Jul-2019 05:20
<b>Xylenes, Total</b>	<b>0.028</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 05:20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>86.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:20</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:20</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:20</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW58A-20190710  
 Collection Date: 10-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 13:45
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jul-2019 13:45
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 13:45
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 13:45
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 13:45
<b>2-Methylnaphthalene</b>	<b>0.015</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:15
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 13:45
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 13:45
<b>Acenaphthene</b>	<b>0.11</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 21:34
<b>Acenaphthylene</b>	<b>0.00087</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:45
<b>Anthracene</b>	<b>0.062</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:15
<b>Benz(a)anthracene</b>	<b>0.00082</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:45
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 13:45
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 13:45
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 13:45
<b>Chrysene</b>	<b>0.00084</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:45
<b>Dibenzofuran</b>	<b>0.059</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:15
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 13:45
<b>Fluoranthene</b>	<b>0.020</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:15
<b>Fluorene</b>	<b>0.15</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 21:34
<b>Naphthalene</b>	<b>0.46</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 21:34
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 13:45
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 13:45
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 13:45
<b>Phenanthrene</b>	<b>0.036</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:15
Phenol	U		0.000035	0.00020	mg/L	1	15-Jul-2019 13:45
<b>Pyrene</b>	<b>0.0099</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 13:45
Surr: 2,4,6-Tribromophenol	93.8			34-129	%REC	10	15-Jul-2019 21:15
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	15-Jul-2019 21:34
Surr: 2,4,6-Tribromophenol	74.0			34-129	%REC	1	15-Jul-2019 13:45
Surr: 2-Fluorobiphenyl	47.7			40-125	%REC	1	15-Jul-2019 13:45
Surr: 2-Fluorobiphenyl	68.2			40-125	%REC	10	15-Jul-2019 21:15
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	15-Jul-2019 21:34
Surr: 2-Fluorophenol	76.7			20-120	%REC	10	15-Jul-2019 21:15
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	15-Jul-2019 21:34
Surr: 2-Fluorophenol	59.4			20-120	%REC	1	15-Jul-2019 13:45
Surr: 4-Terphenyl-d14	67.0			40-135	%REC	1	15-Jul-2019 13:45
Surr: 4-Terphenyl-d14	86.4			40-135	%REC	10	15-Jul-2019 21:15
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	15-Jul-2019 21:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW58A-20190710  
 Collection Date: 10-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	57.0			41-120	%REC	1	15-Jul-2019 13:45
Surr: Nitrobenzene-d5	78.8			41-120	%REC	10	15-Jul-2019 21:15
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	15-Jul-2019 21:34
Surr: Phenol-d6	0	JS		20-120	%REC	100	15-Jul-2019 21:34
Surr: Phenol-d6	62.6			20-120	%REC	1	15-Jul-2019 13:45
Surr: Phenol-d6	69.9			20-120	%REC	10	15-Jul-2019 21:15
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.00748		0.000400	0.00200	mg/L	1	16-Jul-2019 22:43

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57A-20190710  
 Collection Date: 10-Jul-2019 09:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:45
<b>Benzene</b>	<b>0.00057</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 05:45
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:45
<b>Ethylbenzene</b>	<b>0.00085</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 05:45
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 05:45
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:45
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:45
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:45
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:45</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:45</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:45</i>
<i>Surr: Toluene-d8</i>	<i>99.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:45</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57A-20190710  
 Collection Date: 10-Jul-2019 09:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jul-2019 14:04
<b>2,4-Dimethylphenol</b>	<b>0.00059</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jul-2019 14:04
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jul-2019 14:04
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jul-2019 14:04
<b>2-Methylnaphthalene</b>	<b>0.19</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:13
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jul-2019 14:04
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jul-2019 14:04
<b>Acenaphthene</b>	<b>0.099</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:54
<b>Acenaphthylene</b>	<b>0.0019</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Anthracene</b>	<b>0.0084</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Benz(a)anthracene</b>	<b>0.00100</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Benzo(a)pyrene</b>	<b>0.00054</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jul-2019 14:04
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00021</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Chrysene</b>	<b>0.00100</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Dibenzofuran</b>	<b>0.076</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:54
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jul-2019 14:04
<b>Fluoranthene</b>	<b>0.0091</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Fluorene</b>	<b>0.059</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:54
<b>Naphthalene</b>	<b>0.61</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:13
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jul-2019 14:04
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jul-2019 14:04
<b>Pentachlorophenol</b>	<b>0.00059</b>		<b>0.000079</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Phenanthrene</b>	<b>0.059</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 21:54
<b>Phenol</b>	<b>0.00016</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<b>Pyrene</b>	<b>0.0055</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 14:04
<i>Surr: 2,4,6-Tribromophenol</i>	<i>117</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 21:54</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:13</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>74.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:04</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>54.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:04</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>80.3</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 21:54</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:13</i>
<i>Surr: 2-Fluorophenol</i>	<i>61.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:04</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.0</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 21:54</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:13</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>65.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:04</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>97.9</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 21:54</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57A-20190710  
 Collection Date: 10-Jul-2019 09:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	58.0			41-120	%REC	1	15-Jul-2019 14:04
Surr: Nitrobenzene-d5	89.3			41-120	%REC	10	15-Jul-2019 21:54
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	15-Jul-2019 22:13
Surr: Phenol-d6	0	JS		20-120	%REC	100	15-Jul-2019 22:13
Surr: Phenol-d6	70.1			20-120	%REC	1	15-Jul-2019 14:04
Surr: Phenol-d6	87.2			20-120	%REC	10	15-Jul-2019 21:54
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.00447		0.000400	0.00200	mg/L	1	16-Jul-2019 22:45

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57B-20190710  
 Collection Date: 10-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 06:19
<b>Benzene</b>	<b>0.84</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 06:47
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 06:19
<b>Ethylbenzene</b>	<b>0.39</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 06:47
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 06:19
<b>Toluene</b>	<b>0.93</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 06:47
<b>Vinyl chloride</b>	<b>0.00048</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 06:19
<b>Xylenes, Total</b>	<b>1.2</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 06:47
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.4</i>			<i>70-126</i>	<i>%REC</i>	<i>10</i>	<i>14-Jul-2019 06:47</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>96.3</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:19</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:19</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>103</i>			<i>81-113</i>	<i>%REC</i>	<i>10</i>	<i>14-Jul-2019 06:47</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:19</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.2</i>			<i>77-123</i>	<i>%REC</i>	<i>10</i>	<i>14-Jul-2019 06:47</i>
<i>Surr: Toluene-d8</i>	<i>99.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 06:19</i>
<i>Surr: Toluene-d8</i>	<i>99.1</i>			<i>82-127</i>	<i>%REC</i>	<i>10</i>	<i>14-Jul-2019 06:47</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57B-20190710  
 Collection Date: 10-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	15-Jul-2019 14:24
<b>2,4-Dimethylphenol</b>	<b>6.5</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	1000	16-Jul-2019 14:06
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	15-Jul-2019 14:24
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	15-Jul-2019 14:24
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	15-Jul-2019 14:24
<b>2-Methylnaphthalene</b>	<b>0.93</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:33
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	15-Jul-2019 14:24
4-Nitrophenol	U		0.00047	0.010	mg/L	10	15-Jul-2019 14:24
<b>Acenaphthene</b>	<b>0.38</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:33
<b>Acenaphthylene</b>	<b>0.0051</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
<b>Anthracene</b>	<b>0.039</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
<b>Benz(a)anthracene</b>	<b>0.0071</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
<b>Benzo(a)pyrene</b>	<b>0.0021</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	15-Jul-2019 14:24
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	15-Jul-2019 14:24
<b>Chrysene</b>	<b>0.0065</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
<b>Dibenzofuran</b>	<b>0.32</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:33
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	15-Jul-2019 14:24
<b>Fluoranthene</b>	<b>0.047</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
<b>Fluorene</b>	<b>0.21</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:33
<b>Naphthalene</b>	<b>17</b>		<b>0.10</b>	<b>0.50</b>	<b>mg/L</b>	5000	16-Jul-2019 14:26
Nitrobenzene	U		0.00024	0.0020	mg/L	10	15-Jul-2019 14:24
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	15-Jul-2019 14:24
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	15-Jul-2019 14:24
<b>Phenanthrene</b>	<b>0.29</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jul-2019 22:33
<b>Phenol</b>	<b>1.0</b>		<b>0.035</b>	<b>0.20</b>	<b>mg/L</b>	1000	16-Jul-2019 14:06
<b>Pyrene</b>	<b>0.030</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:24
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 14:06</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>5000</i>	<i>16-Jul-2019 14:26</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>72.0</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 14:24</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:33</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>63.8</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 14:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:33</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 14:06</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>5000</i>	<i>16-Jul-2019 14:26</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 14:06</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>5000</i>	<i>16-Jul-2019 14:26</i>
<i>Surr: 2-Fluorophenol</i>	<i>70.5</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 14:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>15-Jul-2019 22:33</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57B-20190710  
 Collection Date: 10-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	77.3			40-135	%REC	10	15-Jul-2019 14:24
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	15-Jul-2019 22:33
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	16-Jul-2019 14:06
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	5000	16-Jul-2019 14:26
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	16-Jul-2019 14:06
Surr: Nitrobenzene-d5	78.7			41-120	%REC	10	15-Jul-2019 14:24
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	15-Jul-2019 22:33
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	5000	16-Jul-2019 14:26
Surr: Phenol-d6	0	JS		20-120	%REC	1000	16-Jul-2019 14:06
Surr: Phenol-d6	0	JS		20-120	%REC	5000	16-Jul-2019 14:26
Surr: Phenol-d6	75.3			20-120	%REC	10	15-Jul-2019 14:24
Surr: Phenol-d6	0	JS		20-120	%REC	100	15-Jul-2019 22:33
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.00404		0.000400	0.00200	mg/L	1	16-Jul-2019 22:47

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW72B-20190710  
 Collection Date: 10-Jul-2019 11:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 07:11
<b>Benzene</b>	<b>1.1</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 07:38
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 07:11
<b>Ethylbenzene</b>	<b>0.30</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 07:38
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 07:11
<b>Toluene</b>	<b>0.96</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 07:38
<b>Xylenes, Total</b>	<b>0.97</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 07:38
<i>Surr: 1,2-Dichloroethane-d4</i>	95.3			70-126	%REC	1	14-Jul-2019 07:11
<i>Surr: 1,2-Dichloroethane-d4</i>	96.8			70-126	%REC	10	14-Jul-2019 07:38
<i>Surr: 4-Bromofluorobenzene</i>	103			81-113	%REC	1	14-Jul-2019 07:11
<i>Surr: 4-Bromofluorobenzene</i>	101			81-113	%REC	10	14-Jul-2019 07:38
<i>Surr: Dibromofluoromethane</i>	97.2			77-123	%REC	1	14-Jul-2019 07:11
<i>Surr: Dibromofluoromethane</i>	98.2			77-123	%REC	10	14-Jul-2019 07:38
<i>Surr: Toluene-d8</i>	99.5			82-127	%REC	1	14-Jul-2019 07:11
<i>Surr: Toluene-d8</i>	98.4			82-127	%REC	10	14-Jul-2019 07:38

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW72B-20190710  
 Collection Date: 10-Jul-2019 11:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	15-Jul-2019 14:43
<b>2,4-Dimethylphenol</b>	<b>19</b>		<b>0.20</b>	<b>1.0</b>	<b>mg/L</b>	5000	17-Jul-2019 14:18
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	15-Jul-2019 14:43
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	15-Jul-2019 14:43
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	15-Jul-2019 14:43
<b>2-Methylnaphthalene</b>	<b>0.57</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 19:00
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	15-Jul-2019 14:43
4-Nitrophenol	U		0.00047	0.010	mg/L	10	15-Jul-2019 14:43
<b>Acenaphthene</b>	<b>0.20</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 19:00
<b>Acenaphthylene</b>	<b>0.0044</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
<b>Anthracene</b>	<b>0.034</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
<b>Benz(a)anthracene</b>	<b>0.0069</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
<b>Benzo(a)pyrene</b>	<b>0.0019</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	15-Jul-2019 14:43
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	15-Jul-2019 14:43
<b>Chrysene</b>	<b>0.0069</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
<b>Dibenzofuran</b>	<b>0.16</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 19:00
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	15-Jul-2019 14:43
<b>Fluoranthene</b>	<b>0.046</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
<b>Fluorene</b>	<b>0.12</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 19:00
<b>Naphthalene</b>	<b>12</b>		<b>0.10</b>	<b>0.50</b>	<b>mg/L</b>	5000	17-Jul-2019 14:18
Nitrobenzene	U		0.00024	0.0020	mg/L	10	15-Jul-2019 14:43
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	15-Jul-2019 14:43
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	15-Jul-2019 14:43
<b>Phenanthrene</b>	<b>0.16</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 19:00
<b>Phenol</b>	<b>9.3</b>		<b>0.035</b>	<b>0.20</b>	<b>mg/L</b>	1000	16-Jul-2019 19:19
<b>Pyrene</b>	<b>0.029</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 14:43
<i>Surr: 2,4,6-Tribromophenol</i>	<i>72.3</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 14:43</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>5000</i>	<i>17-Jul-2019 14:18</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 19:00</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 19:19</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 19:00</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>5000</i>	<i>17-Jul-2019 14:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 19:19</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>83.3</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 14:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>118</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 14:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 19:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>5000</i>	<i>17-Jul-2019 14:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 19:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW72B-20190710  
 Collection Date: 10-Jul-2019 11:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	16-Jul-2019 19:00
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	16-Jul-2019 19:19
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	5000	17-Jul-2019 14:18
Surr: 4-Terphenyl-d14	86.9			40-135	%REC	10	15-Jul-2019 14:43
Surr: Nitrobenzene-d5	42.9			41-120	%REC	10	15-Jul-2019 14:43
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	5000	17-Jul-2019 14:18
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	16-Jul-2019 19:00
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	16-Jul-2019 19:19
Surr: Phenol-d6	0	JS		20-120	%REC	100	16-Jul-2019 19:00
Surr: Phenol-d6	0	JS		20-120	%REC	1000	16-Jul-2019 19:19
Surr: Phenol-d6	0	JS		20-120	%REC	5000	17-Jul-2019 14:18
Surr: Phenol-d6	52.2			20-120	%REC	10	15-Jul-2019 14:43
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.000861	J	0.000400	0.00200	mg/L	1	17-Jul-2019 12:21

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW19C-20190710  
 Collection Date: 10-Jul-2019 12:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:59
<b>Benzene</b>	<b>0.00040</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 00:59
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:59
<b>Ethylbenzene</b>	<b>0.00096</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 00:59
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 00:59
<b>Toluene</b>	<b>0.00086</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 00:59
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:59
<b>Xylenes, Total</b>	<b>0.0022</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 00:59
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:59</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>103</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:59</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW19C-20190710  
 Collection Date: 10-Jul-2019 12:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	0.000082	J	0.000021	0.00020	mg/L	1	15-Jul-2019 15:03
2,4-Dimethylphenol	0.00056		0.000040	0.00020	mg/L	1	15-Jul-2019 15:03
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 15:03
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 15:03
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 15:03
2-Methylnaphthalene	0.000071	J	0.000019	0.00010	mg/L	1	15-Jul-2019 15:03
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 15:03
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 15:03
Acenaphthene	0.0010		0.000027	0.00010	mg/L	1	15-Jul-2019 15:03
Acenaphthylene	0.000030	J	0.000015	0.00010	mg/L	1	15-Jul-2019 15:03
Anthracene	0.000066	J	0.000014	0.00010	mg/L	1	15-Jul-2019 15:03
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 15:03
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 15:03
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 15:03
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 15:03
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 15:03
Dibenzofuran	0.00057		0.000020	0.00010	mg/L	1	15-Jul-2019 15:03
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 15:03
Fluoranthene	0.00056		0.000010	0.00010	mg/L	1	15-Jul-2019 15:03
Fluorene	0.00031		0.000030	0.00010	mg/L	1	15-Jul-2019 15:03
Naphthalene	0.0012		0.000020	0.00010	mg/L	1	15-Jul-2019 15:03
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 15:03
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 15:03
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 15:03
Phenanthrene	0.00014		0.000021	0.00010	mg/L	1	15-Jul-2019 15:03
Phenol	0.00015	J	0.000035	0.00020	mg/L	1	15-Jul-2019 15:03
Pyrene	0.00064		0.000019	0.00010	mg/L	1	15-Jul-2019 15:03
Surr: 2,4,6-Tribromophenol	90.9			34-129	%REC	1	15-Jul-2019 15:03
Surr: 2-Fluorobiphenyl	59.4			40-125	%REC	1	15-Jul-2019 15:03
Surr: 2-Fluorophenol	53.6			20-120	%REC	1	15-Jul-2019 15:03
Surr: 4-Terphenyl-d14	70.2			40-135	%REC	1	15-Jul-2019 15:03
Surr: Nitrobenzene-d5	60.0			41-120	%REC	1	15-Jul-2019 15:03
Surr: Phenol-d6	61.1			20-120	%REC	1	15-Jul-2019 15:03
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.00166	J	0.000400	0.00200	mg/L	1	16-Jul-2019 22:52

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17C-20190710  
 Collection Date: 10-Jul-2019 13:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 02:13
<b>Benzene</b>	<b>0.0097</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:13
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 02:13
<b>Ethylbenzene</b>	<b>0.027</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:13
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 02:13
<b>Toluene</b>	<b>0.0073</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:13
<b>Xylenes, Total</b>	<b>0.039</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:13
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:13</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:13</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:13</i>
<i>Surr: Toluene-d8</i>	<i>98.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17C-20190710  
 Collection Date: 10-Jul-2019 13:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 15:22
<b>2,4-Dimethylphenol</b>	<b>0.21</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	100	16-Jul-2019 19:58
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 15:22
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 15:22
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 15:22
<b>2-Methylnaphthalene</b>	<b>0.026</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 19:39
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 15:22
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 15:22
<b>Acenaphthene</b>	<b>0.035</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 19:39
<b>Acenaphthylene</b>	<b>0.00037</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
<b>Anthracene</b>	<b>0.0013</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 15:22
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 15:22
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 15:22
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00056</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 15:22
<b>Dibenzofuran</b>	<b>0.027</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 19:39
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 15:22
<b>Fluoranthene</b>	<b>0.00090</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
<b>Fluorene</b>	<b>0.014</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 19:39
<b>Naphthalene</b>	<b>0.97</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 19:58
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 15:22
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 15:22
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 15:22
<b>Phenanthrene</b>	<b>0.0100</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
<b>Phenol</b>	<b>0.00073</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
<b>Pyrene</b>	<b>0.00046</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 15:22
Surr: 2,4,6-Tribromophenol	61.3			34-129	%REC	10	16-Jul-2019 19:39
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	16-Jul-2019 19:58
Surr: 2,4,6-Tribromophenol	69.6			34-129	%REC	1	15-Jul-2019 15:22
Surr: 2-Fluorobiphenyl	62.4			40-125	%REC	10	16-Jul-2019 19:39
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	16-Jul-2019 19:58
Surr: 2-Fluorobiphenyl	45.2			40-125	%REC	1	15-Jul-2019 15:22
Surr: 2-Fluorophenol	52.5			20-120	%REC	1	15-Jul-2019 15:22
Surr: 2-Fluorophenol	76.5			20-120	%REC	10	16-Jul-2019 19:39
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	16-Jul-2019 19:58
Surr: 4-Terphenyl-d14	93.5			40-135	%REC	10	16-Jul-2019 19:39
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	16-Jul-2019 19:58
Surr: 4-Terphenyl-d14	68.8			40-135	%REC	1	15-Jul-2019 15:22

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17C-20190710  
 Collection Date: 10-Jul-2019 13:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	45.3			41-120	%REC	1	15-Jul-2019 15:22
Surr: Nitrobenzene-d5	62.8			41-120	%REC	10	16-Jul-2019 19:39
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	16-Jul-2019 19:58
Surr: Phenol-d6	66.3			20-120	%REC	10	16-Jul-2019 19:39
Surr: Phenol-d6	0	JS		20-120	%REC	100	16-Jul-2019 19:58
Surr: Phenol-d6	51.8			20-120	%REC	1	15-Jul-2019 15:22
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.000854	J	0.000400	0.00200	mg/L	1	16-Jul-2019 22:54

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17-20190710  
 Collection Date: 10-Jul-2019 14:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 08:03
<b>Benzene</b>	<b>0.54</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 08:30
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 08:03
<b>Ethylbenzene</b>	<b>0.23</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 08:30
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 08:03
<b>Toluene</b>	<b>0.83</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 08:30
<b>Xylenes, Total</b>	<b>0.78</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jul-2019 08:30
<i>Surr: 1,2-Dichloroethane-d4</i>	93.9			70-126	%REC	1	14-Jul-2019 08:03
<i>Surr: 1,2-Dichloroethane-d4</i>	96.0			70-126	%REC	10	14-Jul-2019 08:30
<i>Surr: 4-Bromofluorobenzene</i>	108			81-113	%REC	1	14-Jul-2019 08:03
<i>Surr: 4-Bromofluorobenzene</i>	107			81-113	%REC	10	14-Jul-2019 08:30
<i>Surr: Dibromofluoromethane</i>	95.4			77-123	%REC	1	14-Jul-2019 08:03
<i>Surr: Dibromofluoromethane</i>	96.9			77-123	%REC	10	14-Jul-2019 08:30
<i>Surr: Toluene-d8</i>	98.2			82-127	%REC	1	14-Jul-2019 08:03
<i>Surr: Toluene-d8</i>	97.1			82-127	%REC	10	14-Jul-2019 08:30

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17-20190710  
 Collection Date: 10-Jul-2019 14:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	15-Jul-2019 15:42
<b>2,4-Dimethylphenol</b>	<b>5.9</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	1000	16-Jul-2019 20:37
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	15-Jul-2019 15:42
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	15-Jul-2019 15:42
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	15-Jul-2019 15:42
<b>2-Methylnaphthalene</b>	<b>0.39</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 20:18
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	15-Jul-2019 15:42
4-Nitrophenol	U		0.00047	0.010	mg/L	10	15-Jul-2019 15:42
<b>Acenaphthene</b>	<b>0.16</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 20:18
<b>Acenaphthylene</b>	<b>0.0042</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 15:42
<b>Anthracene</b>	<b>0.0092</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 15:42
Benz(a)anthracene	U		0.00050	0.0010	mg/L	10	15-Jul-2019 15:42
Benzo(a)pyrene	U		0.00020	0.0010	mg/L	10	15-Jul-2019 15:42
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	15-Jul-2019 15:42
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	15-Jul-2019 15:42
Chrysene	U		0.00021	0.0010	mg/L	10	15-Jul-2019 15:42
<b>Dibenzofuran</b>	<b>0.12</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 20:18
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	15-Jul-2019 15:42
<b>Fluoranthene</b>	<b>0.0039</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 15:42
<b>Fluorene</b>	<b>0.060</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 15:42
<b>Naphthalene</b>	<b>8.9</b>		<b>0.10</b>	<b>0.50</b>	<b>mg/L</b>	5000	16-Jul-2019 20:57
Nitrobenzene	U		0.00024	0.0020	mg/L	10	15-Jul-2019 15:42
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	15-Jul-2019 15:42
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	15-Jul-2019 15:42
<b>Phenanthrene</b>	<b>0.045</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 15:42
<b>Phenol</b>	<b>9.9</b>		<b>0.18</b>	<b>1.0</b>	<b>mg/L</b>	5000	16-Jul-2019 20:57
<b>Pyrene</b>	<b>0.0018</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jul-2019 15:42
<i>Surr: 2,4,6-Tribromophenol</i>	<i>118</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 15:42</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>5000</i>	<i>16-Jul-2019 20:57</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 20:18</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 20:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 20:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 20:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>5000</i>	<i>16-Jul-2019 20:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>89.4</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 15:42</i>
<i>Surr: 2-Fluorophenol</i>	<i>72.9</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>15-Jul-2019 15:42</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>5000</i>	<i>16-Jul-2019 20:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 20:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>16-Jul-2019 20:37</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17-20190710  
 Collection Date: 10-Jul-2019 14:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	16-Jul-2019 20:18
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	16-Jul-2019 20:37
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	5000	16-Jul-2019 20:57
Surr: 4-Terphenyl-d14	86.5			40-135	%REC	10	15-Jul-2019 15:42
Surr: Nitrobenzene-d5	65.4			41-120	%REC	10	15-Jul-2019 15:42
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	16-Jul-2019 20:18
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	16-Jul-2019 20:37
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	5000	16-Jul-2019 20:57
Surr: Phenol-d6	0	JS		20-120	%REC	100	16-Jul-2019 20:18
Surr: Phenol-d6	0	JS		20-120	%REC	1000	16-Jul-2019 20:37
Surr: Phenol-d6	0	JS		20-120	%REC	5000	16-Jul-2019 20:57
Surr: Phenol-d6	117			20-120	%REC	10	15-Jul-2019 15:42
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.0460		0.000400	0.00200	mg/L	1	17-Jul-2019 12:23

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW20A-20190710  
 Collection Date: 10-Jul-2019 15:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 02:38
<b>Benzene</b>	<b>0.013</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:38
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 02:38
<b>Ethylbenzene</b>	<b>0.0079</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:38
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 02:38
<b>Toluene</b>	<b>0.0011</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:38
<b>Xylenes, Total</b>	<b>0.022</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 02:38
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:38</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>106</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:38</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:38</i>
<i>Surr: Toluene-d8</i>	<i>96.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 02:38</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW20A-20190710  
 Collection Date: 10-Jul-2019 15:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jul-2019 16:01
<b>2,4-Dimethylphenol</b>	<b>0.0035</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 16:01
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jul-2019 16:01
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jul-2019 16:01
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jul-2019 16:01
<b>2-Methylnaphthalene</b>	<b>0.072</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 21:16
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jul-2019 16:01
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jul-2019 16:01
<b>Acenaphthene</b>	<b>0.14</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	16-Jul-2019 21:36
<b>Acenaphthylene</b>	<b>0.00078</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:01
<b>Anthracene</b>	<b>0.0062</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:01
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	15-Jul-2019 16:01
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	15-Jul-2019 16:01
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jul-2019 16:01
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	15-Jul-2019 16:01
Chrysene		U	0.000021	0.00010	mg/L	1	15-Jul-2019 16:01
<b>Dibenzofuran</b>	<b>0.077</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 21:16
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jul-2019 16:01
<b>Fluoranthene</b>	<b>0.00058</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:01
<b>Fluorene</b>	<b>0.068</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 21:16
<b>Naphthalene</b>	<b>1.9</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	500	17-Jul-2019 15:16
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jul-2019 16:01
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jul-2019 16:01
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jul-2019 16:01
<b>Phenanthrene</b>	<b>0.026</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 21:16
<b>Phenol</b>	<b>0.00096</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 16:01
<b>Pyrene</b>	<b>0.00026</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:01
<i>Surr: 2,4,6-Tribromophenol</i>	<i>76.1</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 21:16</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 21:36</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>500</i>	<i>17-Jul-2019 15:16</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>76.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:01</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>60.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:01</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>500</i>	<i>17-Jul-2019 15:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>68.6</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 21:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 21:36</i>
<i>Surr: 2-Fluorophenol</i>	<i>96.5</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 21:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>16-Jul-2019 21:36</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>500</i>	<i>17-Jul-2019 15:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>60.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:01</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW20A-20190710  
 Collection Date: 10-Jul-2019 15:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	63.0			40-135	%REC	1	15-Jul-2019 16:01
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	500	17-Jul-2019 15:16
Surr: 4-Terphenyl-d14	98.3			40-135	%REC	10	16-Jul-2019 21:16
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	16-Jul-2019 21:36
Surr: Nitrobenzene-d5	67.4			41-120	%REC	10	16-Jul-2019 21:16
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	16-Jul-2019 21:36
Surr: Nitrobenzene-d5	66.5			41-120	%REC	1	15-Jul-2019 16:01
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	500	17-Jul-2019 15:16
Surr: Phenol-d6	0	JS		20-120	%REC	500	17-Jul-2019 15:16
Surr: Phenol-d6	68.2			20-120	%REC	1	15-Jul-2019 16:01
Surr: Phenol-d6	90.7			20-120	%REC	10	16-Jul-2019 21:16
Surr: Phenol-d6	0	JS		20-120	%REC	100	16-Jul-2019 21:36
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.00574		0.000400	0.00200	mg/L	1	16-Jul-2019 23:03

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15A-20190710  
 Collection Date: 10-Jul-2019 16:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 03:02
<b>Benzene</b>	<b>0.00074</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:02
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 03:02
<b>Ethylbenzene</b>	<b>0.00035</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:02
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 03:02
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 03:02
<b>Xylenes, Total</b>	<b>0.0026</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:02
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>96.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:02</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:02</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:02</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:02</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15A-20190710  
 Collection Date: 10-Jul-2019 16:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jul-2019 16:21
<b>2,4-Dimethylphenol</b>	<b>0.00024</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jul-2019 16:21
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jul-2019 16:21
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jul-2019 16:21
<b>2-Methylnaphthalene</b>	<b>0.0077</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jul-2019 16:21
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jul-2019 16:21
<b>Acenaphthene</b>	<b>0.11</b>		<b>0.0014</b>	<b>0.0050</b>	<b>mg/L</b>	50	17-Jul-2019 15:35
<b>Acenaphthylene</b>	<b>0.00045</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
<b>Anthracene</b>	<b>0.0026</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	15-Jul-2019 16:21
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	15-Jul-2019 16:21
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jul-2019 16:21
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	15-Jul-2019 16:21
Chrysene		U	0.000021	0.00010	mg/L	1	15-Jul-2019 16:21
<b>Dibenzofuran</b>	<b>0.035</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 21:55
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jul-2019 16:21
<b>Fluoranthene</b>	<b>0.0018</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
<b>Fluorene</b>	<b>0.054</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 21:55
<b>Naphthalene</b>	<b>0.0066</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jul-2019 16:21
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jul-2019 16:21
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jul-2019 16:21
<b>Phenanthrene</b>	<b>0.0095</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
<b>Phenol</b>	<b>0.00039</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
<b>Pyrene</b>	<b>0.00075</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:21
Surr: 2,4,6-Tribromophenol	71.3			34-129	%REC	10	16-Jul-2019 21:55
Surr: 2,4,6-Tribromophenol	59.6	J		34-129	%REC	50	17-Jul-2019 15:35
Surr: 2,4,6-Tribromophenol	65.0			34-129	%REC	1	15-Jul-2019 16:21
Surr: 2-Fluorobiphenyl	44.5			40-125	%REC	1	15-Jul-2019 16:21
Surr: 2-Fluorobiphenyl	60.0	J		40-125	%REC	50	17-Jul-2019 15:35
Surr: 2-Fluorobiphenyl	70.9			40-125	%REC	10	16-Jul-2019 21:55
Surr: 2-Fluorophenol	52.8	J		20-120	%REC	50	17-Jul-2019 15:35
Surr: 2-Fluorophenol	97.0			20-120	%REC	10	16-Jul-2019 21:55
Surr: 2-Fluorophenol	39.2			20-120	%REC	1	15-Jul-2019 16:21
Surr: 4-Terphenyl-d14	52.3			40-135	%REC	1	15-Jul-2019 16:21
Surr: 4-Terphenyl-d14	73.2			40-135	%REC	10	16-Jul-2019 21:55
Surr: 4-Terphenyl-d14	65.2	J		40-135	%REC	50	17-Jul-2019 15:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15A-20190710  
 Collection Date: 10-Jul-2019 16:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	65.2	J		41-120	%REC	50	17-Jul-2019 15:35
Surr: Nitrobenzene-d5	59.8			41-120	%REC	10	16-Jul-2019 21:55
Surr: Nitrobenzene-d5	44.0			41-120	%REC	1	15-Jul-2019 16:21
Surr: Phenol-d6	43.2			20-120	%REC	1	15-Jul-2019 16:21
Surr: Phenol-d6	83.2			20-120	%REC	10	16-Jul-2019 21:55
Surr: Phenol-d6	62.3	J		20-120	%REC	50	17-Jul-2019 15:35
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.0251		0.000400	0.00200	mg/L	1	16-Jul-2019 22:18

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15C-20190710  
 Collection Date: 10-Jul-2019 17:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 03:27
<b>Benzene</b>	<b>0.00050</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:27
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 03:27
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 03:27
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 03:27
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 03:27
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 03:27
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:27</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:27</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:27</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:27</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15C-20190710  
 Collection Date: 10-Jul-2019 17:05

**ANALYTICAL REPORT**

WorkOrder:HS19070575  
 Lab ID:HS19070575-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 16:41
<b>2,4-Dimethylphenol</b>	<b>0.0052</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 16:41
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 16:41
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 16:41
<b>2-Methylnaphthalene</b>	<b>0.00066</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 16:41
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 16:41
<b>Acenaphthene</b>	<b>0.017</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 22:15
<b>Acenaphthylene</b>	<b>0.0015</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
<b>Anthracene</b>	<b>0.00064</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 16:41
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 16:41
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 16:41
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 16:41
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 16:41
<b>Dibenzofuran</b>	<b>0.0039</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 16:41
<b>Fluoranthene</b>	<b>0.00086</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
<b>Fluorene</b>	<b>0.00097</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
<b>Naphthalene</b>	<b>0.019</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	16-Jul-2019 22:15
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 16:41
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 16:41
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 16:41
<b>Phenanthrene</b>	<b>0.00039</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
<b>Phenol</b>	<b>0.00026</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
<b>Pyrene</b>	<b>0.00043</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 16:41
<i>Surr: 2,4,6-Tribromophenol</i>	<i>103</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:41</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>73.5</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 22:15</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>65.7</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 22:15</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>70.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:41</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:41</i>
<i>Surr: 2-Fluorophenol</i>	<i>107</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 22:15</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.6</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 22:15</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:41</i>
<i>Surr: Nitrobenzene-d5</i>	<i>79.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:41</i>
<i>Surr: Nitrobenzene-d5</i>	<i>76.1</i>			<i>41-120</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 22:15</i>
<i>Surr: Phenol-d6</i>	<i>93.8</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>16-Jul-2019 22:15</i>
<i>Surr: Phenol-d6</i>	<i>65.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 16:41</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15C-20190710  
 Collection Date: 10-Jul-2019 17:05

**ANALYTICAL REPORT**

WorkOrder:HS19070575  
 Lab ID:HS19070575-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.000889	J	0.000400	0.00200	mg/L	1	16-Jul-2019 23:05

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15B-20190710  
 Collection Date: 10-Jul-2019 18:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 03:52
<b>Benzene</b>	<b>0.0023</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:52
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 03:52
<b>Ethylbenzene</b>	<b>0.0031</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:52
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 03:52
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 03:52
<b>Xylenes, Total</b>	<b>0.0025</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jul-2019 03:52
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:52</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>104</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:52</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:52</i>
<i>Surr: Toluene-d8</i>	<i>97.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 03:52</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15B-20190710  
 Collection Date: 10-Jul-2019 18:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jul-2019 17:00
<b>2,4-Dimethylphenol</b>	<b>0.00016</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jul-2019 17:00
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jul-2019 17:00
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jul-2019 17:00
<b>2-Methylnaphthalene</b>	<b>0.016</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jul-2019 12:42
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jul-2019 17:00
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jul-2019 17:00
<b>Acenaphthene</b>	<b>0.029</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jul-2019 12:42
<b>Acenaphthylene</b>	<b>0.00034</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<b>Anthracene</b>	<b>0.0016</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<b>Benz(a)anthracene</b>	<b>0.00016</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	15-Jul-2019 17:00
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jul-2019 17:00
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	15-Jul-2019 17:00
<b>Chrysene</b>	<b>0.00013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<b>Dibenzofuran</b>	<b>0.013</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jul-2019 12:42
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jul-2019 17:00
<b>Fluoranthene</b>	<b>0.0042</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<b>Fluorene</b>	<b>0.010</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jul-2019 12:42
<b>Naphthalene</b>	<b>0.33</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jul-2019 13:01
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jul-2019 17:00
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jul-2019 17:00
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jul-2019 17:00
<b>Phenanthrene</b>	<b>0.0065</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<b>Phenol</b>	<b>0.00019</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<b>Pyrene</b>	<b>0.0018</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:00
<i>Surr: 2,4,6-Tribromophenol</i>	<i>107</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:00</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>95.6</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 12:42</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:01</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>77.5</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 12:42</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:01</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:00</i>
<i>Surr: 2-Fluorophenol</i>	<i>58.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:00</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.4</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 12:42</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:01</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>88.1</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 12:42</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:01</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15B-20190710  
 Collection Date: 10-Jul-2019 18:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	70.9			41-120	%REC	1	15-Jul-2019 17:00
Surr: Nitrobenzene-d5	81.5			41-120	%REC	10	17-Jul-2019 12:42
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	17-Jul-2019 13:01
Surr: Phenol-d6	0	JS		20-120	%REC	100	17-Jul-2019 13:01
Surr: Phenol-d6	78.0			20-120	%REC	10	17-Jul-2019 12:42
Surr: Phenol-d6	66.4			20-120	%REC	1	15-Jul-2019 17:00
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	0.0133		0.000400	0.00200	mg/L	1	16-Jul-2019 23:08

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB01-20190710  
 Collection Date: 10-Jul-2019 18:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	13-Jul-2019 23:46
Benzene	U		0.00020	0.0010	mg/L	1	13-Jul-2019 23:46
Chlorobenzene	U		0.00030	0.0010	mg/L	1	13-Jul-2019 23:46
Ethylbenzene	U		0.00030	0.0010	mg/L	1	13-Jul-2019 23:46
Methylene chloride	U		0.0010	0.0020	mg/L	1	13-Jul-2019 23:46
Toluene	U		0.00020	0.0010	mg/L	1	13-Jul-2019 23:46
Vinyl chloride	U		0.00020	0.0010	mg/L	1	13-Jul-2019 23:46
Xylenes, Total	U		0.00030	0.0010	mg/L	1	13-Jul-2019 23:46
<i>Surr: 1,2-Dichloroethane-d4</i>		99.6		70-126	%REC	1	13-Jul-2019 23:46
<i>Surr: 4-Bromofluorobenzene</i>		96.9		81-113	%REC	1	13-Jul-2019 23:46
<i>Surr: Dibromofluoromethane</i>		100		77-123	%REC	1	13-Jul-2019 23:46
<i>Surr: Toluene-d8</i>		100		82-127	%REC	1	13-Jul-2019 23:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB01-20190710  
 Collection Date: 10-Jul-2019 18:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 17:20
<b>2,4-Dimethylphenol</b>	<b>0.00037</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 17:20
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 17:20
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 17:20
<b>2-Methylnaphthalene</b>	<b>0.00038</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 17:20
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 17:20
<b>Acenaphthene</b>	<b>0.00011</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jul-2019 17:20
<b>Anthracene</b>	<b>0.000032</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 17:20
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 17:20
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 17:20
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 17:20
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 17:20
<b>Dibenzofuran</b>	<b>0.00018</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
<b>Di-n-butyl phthalate</b>	<b>0.000024</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
<b>Fluoranthene</b>	<b>0.000025</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
<b>Fluorene</b>	<b>0.00011</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
<b>Naphthalene</b>	<b>0.0035</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 17:20
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 17:20
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 17:20
<b>Phenanthrene</b>	<b>0.00013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
<b>Phenol</b>	<b>0.000071</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:20
Pyrene	U		0.000019	0.00010	mg/L	1	15-Jul-2019 17:20
<i>Surr: 2,4,6-Tribromophenol</i>	<i>121</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:20</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>110</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>103</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:20</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>99.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:20</i>
<i>Surr: Nitrobenzene-d5</i>	<i>106</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:20</i>
<i>Surr: Phenol-d6</i>	<i>112</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:20</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	16-Jul-2019 23:10

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW14-20190711  
 Collection Date: 11-Jul-2019 07:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:16
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:16
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 04:16
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 04:16
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 04:16
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:16
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 04:16
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:16</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:16</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:16</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:16</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW14-20190711  
 Collection Date: 11-Jul-2019 07:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 17:39
<b>2,4-Dimethylphenol</b>	<b>0.00013</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 17:39
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 17:39
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 17:39
<b>2-Methylnaphthalene</b>	<b>0.00019</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 17:39
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 17:39
<b>Acenaphthene</b>	<b>0.00010</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jul-2019 17:39
<b>Anthracene</b>	<b>0.000089</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 17:39
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 17:39
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 17:39
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jul-2019 17:39
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 17:39
<b>Dibenzofuran</b>	<b>0.000087</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 17:39
<b>Fluoranthene</b>	<b>0.000067</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
<b>Fluorene</b>	<b>0.000069</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
<b>Naphthalene</b>	<b>0.0020</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 17:39
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 17:39
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 17:39
<b>Phenanthrene</b>	<b>0.000059</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
<b>Phenol</b>	<b>0.00019</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
<b>Pyrene</b>	<b>0.000032</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:39
Surr: 2,4,6-Tribromophenol	87.6			34-129	%REC	1	15-Jul-2019 17:39
Surr: 2-Fluorobiphenyl	77.2			40-125	%REC	1	15-Jul-2019 17:39
Surr: 2-Fluorophenol	79.3			20-120	%REC	1	15-Jul-2019 17:39
Surr: 4-Terphenyl-d14	76.6			40-135	%REC	1	15-Jul-2019 17:39
Surr: Nitrobenzene-d5	89.4			41-120	%REC	1	15-Jul-2019 17:39
Surr: Phenol-d6	86.2			20-120	%REC	1	15-Jul-2019 17:39
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00133</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	16-Jul-2019 23:12

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW13-20190711  
 Collection Date: 11-Jul-2019 08:15

**ANALYTICAL REPORT**

WorkOrder:HS19070575  
 Lab ID:HS19070575-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:41
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:41
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 04:41
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 04:41
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 04:41
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 04:41
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 04:41
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:41</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:41</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:41</i>
<i>Surr: Toluene-d8</i>	<i>99.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 04:41</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW13-20190711  
 Collection Date: 11-Jul-2019 08:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 17:59
<b>2,4-Dimethylphenol</b>	<b>0.000063</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 17:59
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 17:59
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 17:59
<b>2-Methylnaphthalene</b>	<b>0.000077</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 17:59
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 17:59
<b>Acenaphthene</b>	<b>0.000066</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
<b>Acenaphthylene</b>	<b>0.000033</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
<b>Anthracene</b>	<b>0.00047</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 17:59
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 17:59
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 17:59
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000054</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 17:59
<b>Dibenzofuran</b>	<b>0.000039</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 17:59
<b>Fluoranthene</b>	<b>0.000032</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
<b>Fluorene</b>	<b>0.000042</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
<b>Naphthalene</b>	<b>0.0011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 17:59
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 17:59
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 17:59
Phenanthrene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 17:59
Phenol	U		0.000035	0.00020	mg/L	1	15-Jul-2019 17:59
<b>Pyrene</b>	<b>0.000084</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 17:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>93.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>65.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>54.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>87.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.7</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:59</i>
<i>Surr: Phenol-d6</i>	<i>68.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 17:59</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0715</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	16-Jul-2019 23:14

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW39B-20190711  
 Collection Date: 11-Jul-2019 09:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:06
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:06
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:06
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:06
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 05:06
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:06
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:06
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:06</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:06</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:06</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:06</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW39B-20190711  
 Collection Date: 11-Jul-2019 09:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 18:19
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jul-2019 18:19
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 18:19
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 18:19
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 18:19
<b>2-Methylnaphthalene</b>	<b>0.000027</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 18:19
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 18:19
<b>Acenaphthene</b>	<b>0.00065</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
<b>Acenaphthylene</b>	<b>0.000021</b>	<b>J</b>	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
<b>Anthracene</b>	<b>0.00021</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jul-2019 18:19
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jul-2019 18:19
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 18:19
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000042</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 18:19
<b>Dibenzofuran</b>	<b>0.000037</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 18:19
<b>Fluoranthene</b>	<b>0.000048</b>	<b>J</b>	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
<b>Fluorene</b>	<b>0.000044</b>	<b>J</b>	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
<b>Naphthalene</b>	<b>0.00041</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 18:19
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 18:19
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 18:19
Phenanthrene	U		0.000021	0.00010	mg/L	1	15-Jul-2019 18:19
Phenol	U		0.000035	0.00020	mg/L	1	15-Jul-2019 18:19
<b>Pyrene</b>	<b>0.000072</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>15-Jul-2019 18:19</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>73.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:19</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>51.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>48.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:19</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>71.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:19</i>
<i>Surr: Nitrobenzene-d5</i>	<i>59.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:19</i>
<i>Surr: Phenol-d6</i>	<i>55.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:19</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00144</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>16-Jul-2019 23:16</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12C-20190711  
 Collection Date: 11-Jul-2019 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>					Analyst: AKP
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:30
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:30
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:30
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:30
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 05:30
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:30
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:30
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:30</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:30</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:30</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12C-20190711  
 Collection Date: 11-Jul-2019 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jul-2019 18:38
<b>2,4-Dimethylphenol</b>	<b>0.00043</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jul-2019 18:38
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jul-2019 18:38
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jul-2019 18:38
<b>2-Methylnaphthalene</b>	<b>0.00021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jul-2019 18:38
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jul-2019 18:38
<b>Acenaphthene</b>	<b>0.000071</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
Acenaphthylene		U	0.000015	0.00010	mg/L	1	15-Jul-2019 18:38
<b>Anthracene</b>	<b>0.000030</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	15-Jul-2019 18:38
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	15-Jul-2019 18:38
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jul-2019 18:38
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	15-Jul-2019 18:38
Chrysene		U	0.000021	0.00010	mg/L	1	15-Jul-2019 18:38
<b>Dibenzofuran</b>	<b>0.000070</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jul-2019 18:38
Fluoranthene		U	0.000010	0.00010	mg/L	1	15-Jul-2019 18:38
<b>Fluorene</b>	<b>0.000052</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
<b>Naphthalene</b>	<b>0.0031</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jul-2019 18:38
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jul-2019 18:38
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jul-2019 18:38
<b>Phenanthrene</b>	<b>0.000039</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
<b>Phenol</b>	<b>0.00040</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 18:38
Pyrene		U	0.000019	0.00010	mg/L	1	15-Jul-2019 18:38
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>50.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>54.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>71.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:38</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:38</i>
<i>Surr: Phenol-d6</i>	<i>62.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:38</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 16-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00200</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	16-Jul-2019 23:19

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12A-20190711  
 Collection Date: 11-Jul-2019 10:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:55
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:55
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:55
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 05:55
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 05:55
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 05:55
<b>Xylenes, Total</b>	<b>0.00060</b>	<b>J</b>	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	<b>1</b>	<b>14-Jul-2019 05:55</b>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:55</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:55</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:55</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 05:55</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12A-20190711  
 Collection Date: 11-Jul-2019 10:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jul-2019 18:58
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jul-2019 18:58
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jul-2019 18:58
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jul-2019 18:58
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jul-2019 18:58
<b>2-Methylnaphthalene</b>	<b>0.013</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jul-2019 13:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jul-2019 18:58
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jul-2019 18:58
<b>Acenaphthene</b>	<b>0.19</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jul-2019 13:59
<b>Acenaphthylene</b>	<b>0.0013</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<b>Anthracene</b>	<b>0.0087</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<b>Benz(a)anthracene</b>	<b>0.00041</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<b>Benzo(a)pyrene</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jul-2019 18:58
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00031</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<b>Chrysene</b>	<b>0.00034</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<b>Dibenzofuran</b>	<b>0.14</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jul-2019 13:59
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jul-2019 18:58
<b>Fluoranthene</b>	<b>0.0097</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<b>Fluorene</b>	<b>0.16</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jul-2019 13:59
<b>Naphthalene</b>	<b>0.0012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jul-2019 18:58
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jul-2019 18:58
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jul-2019 18:58
<b>Phenanthrene</b>	<b>0.063</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jul-2019 13:40
Phenol	U		0.000035	0.00020	mg/L	1	15-Jul-2019 18:58
<b>Pyrene</b>	<b>0.0047</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jul-2019 18:58
<i>Surr: 2,4,6-Tribromophenol</i>	<i>103</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 13:40</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:59</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>76.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>49.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.8</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 13:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>51.5</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 13:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>48.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:58</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>84.1</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>17-Jul-2019 13:40</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>17-Jul-2019 13:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>76.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 18:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12A-20190711  
 Collection Date: 11-Jul-2019 10:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	54.7			41-120	%REC	1	15-Jul-2019 18:58
Surr: Nitrobenzene-d5	70.5			41-120	%REC	10	17-Jul-2019 13:40
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	17-Jul-2019 13:59
Surr: Phenol-d6	68.7			20-120	%REC	10	17-Jul-2019 13:40
Surr: Phenol-d6	0	JS		20-120	%REC	100	17-Jul-2019 13:59
Surr: Phenol-d6	54.0			20-120	%REC	1	15-Jul-2019 18:58
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jul-2019		Analyst: JHD	
Arsenic	0.00192	J	0.000400	0.00200	mg/L	1	17-Jul-2019 22:23

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB01-20190710  
 Collection Date: 10-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:10
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:10
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:10
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:10
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 00:10
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:10
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:10
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:10
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:10</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:10</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:10</i>
<i>Surr: Toluene-d8</i>	<i>98.1</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:10</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB02-20190710  
 Collection Date: 10-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070575  
 Lab ID:HS19070575-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:35
Benzene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:35
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:35
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:35
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jul-2019 00:35
Toluene	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:35
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jul-2019 00:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jul-2019 00:35
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>101</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:35</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>96.6</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:35</i>
<i>Surr: Dibromofluoromethane</i>		<i>103</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:35</i>
<i>Surr: Toluene-d8</i>		<i>98.4</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jul-2019 00:35</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**Batch ID:** 143014      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070575-01	1	1000	1 (mL)	0.001
HS19070575-02	1	1000	1 (mL)	0.001
HS19070575-03	1	1000	1 (mL)	0.001
HS19070575-04	1	1000	1 (mL)	0.001
HS19070575-05	1	1000	1 (mL)	0.001
HS19070575-06	1	1000	1 (mL)	0.001
HS19070575-07	1	1000	1 (mL)	0.001
HS19070575-08	1	1000	1 (mL)	0.001
HS19070575-09	1	1000	1 (mL)	0.001
HS19070575-10	1	1000	1 (mL)	0.001
HS19070575-11	1	1000	1 (mL)	0.001
HS19070575-12	1	1000	1 (mL)	0.001
HS19070575-13	1	1000	1 (mL)	0.001
HS19070575-14	1	1000	1 (mL)	0.001
HS19070575-15	1	1000	1 (mL)	0.001
HS19070575-16	1	1000	1 (mL)	0.001
HS19070575-17	1	1000	1 (mL)	0.001
HS19070575-18	1	1000	1 (mL)	0.001
HS19070575-19	1	1000	1 (mL)	0.001

**Batch ID:** 143067      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070575-01	1	10	10 (mL)	1
HS19070575-02	1	10	10 (mL)	1
HS19070575-03	1	10	10 (mL)	1
HS19070575-04	1	10	10 (mL)	1
HS19070575-05	1	10	10 (mL)	1
HS19070575-06	1	10	10 (mL)	1
HS19070575-07	1	10	10 (mL)	1
HS19070575-08	1	10	10 (mL)	1
HS19070575-09	1	10	10 (mL)	1
HS19070575-10	1	10	10 (mL)	1
HS19070575-11	1	10	10 (mL)	1
HS19070575-12	1	10	10 (mL)	1
HS19070575-13	1	10	10 (mL)	1
HS19070575-14	1	10	10 (mL)	1
HS19070575-15	1	10	10 (mL)	1
HS19070575-16	1	10	10 (mL)	1
HS19070575-17	1	10	10 (mL)	1
HS19070575-18	1	10	10 (mL)	1

**Batch ID:** 143129      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070575-19	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143014 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20		15 Jul 2019 10:03	15 Jul 2019 19:37	100
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20		15 Jul 2019 10:03	15 Jul 2019 19:56	1000
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20		15 Jul 2019 10:03	15 Jul 2019 19:17	10
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20		15 Jul 2019 10:03	15 Jul 2019 13:05	1
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05		15 Jul 2019 10:03	15 Jul 2019 20:36	100
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05		15 Jul 2019 10:03	16 Jul 2019 13:47	5000
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05		15 Jul 2019 10:03	15 Jul 2019 20:16	10
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05		15 Jul 2019 10:03	15 Jul 2019 13:25	1
HS19070575-03	WG-1620-MW58A-20190710	10 Jul 2019 08:55		15 Jul 2019 10:03	15 Jul 2019 21:34	100
HS19070575-03	WG-1620-MW58A-20190710	10 Jul 2019 08:55		15 Jul 2019 10:03	15 Jul 2019 21:15	10
HS19070575-03	WG-1620-MW58A-20190710	10 Jul 2019 08:55		15 Jul 2019 10:03	15 Jul 2019 13:45	1
HS19070575-04	WG-1620-MW57A-20190710	10 Jul 2019 09:55		15 Jul 2019 10:03	15 Jul 2019 22:13	100
HS19070575-04	WG-1620-MW57A-20190710	10 Jul 2019 09:55		15 Jul 2019 10:03	15 Jul 2019 21:54	10
HS19070575-04	WG-1620-MW57A-20190710	10 Jul 2019 09:55		15 Jul 2019 10:03	15 Jul 2019 14:04	1
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40		15 Jul 2019 10:03	16 Jul 2019 14:26	5000
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40		15 Jul 2019 10:03	16 Jul 2019 14:06	1000
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40		15 Jul 2019 10:03	15 Jul 2019 22:33	100
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40		15 Jul 2019 10:03	15 Jul 2019 14:24	10
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35		15 Jul 2019 10:03	17 Jul 2019 14:18	5000
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35		15 Jul 2019 10:03	16 Jul 2019 19:19	1000
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35		15 Jul 2019 10:03	16 Jul 2019 19:00	100
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35		15 Jul 2019 10:03	15 Jul 2019 14:43	10
HS19070575-07	WG-1620-MW19C-20190710	10 Jul 2019 12:25		15 Jul 2019 10:03	15 Jul 2019 15:03	1
HS19070575-08	WG-1620-MW17C-20190710	10 Jul 2019 13:25		15 Jul 2019 10:03	16 Jul 2019 19:58	100
HS19070575-08	WG-1620-MW17C-20190710	10 Jul 2019 13:25		15 Jul 2019 10:03	16 Jul 2019 19:39	10
HS19070575-08	WG-1620-MW17C-20190710	10 Jul 2019 13:25		15 Jul 2019 10:03	15 Jul 2019 15:22	1
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15		15 Jul 2019 10:03	16 Jul 2019 20:37	1000
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15		15 Jul 2019 10:03	16 Jul 2019 20:57	5000
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15		15 Jul 2019 10:03	16 Jul 2019 20:18	100
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15		15 Jul 2019 10:03	15 Jul 2019 15:42	10
HS19070575-10	WG-1620-MW20A-20190710	10 Jul 2019 15:10		15 Jul 2019 10:03	17 Jul 2019 15:16	500
HS19070575-10	WG-1620-MW20A-20190710	10 Jul 2019 15:10		15 Jul 2019 10:03	16 Jul 2019 21:36	100
HS19070575-10	WG-1620-MW20A-20190710	10 Jul 2019 15:10		15 Jul 2019 10:03	16 Jul 2019 21:16	10
HS19070575-10	WG-1620-MW20A-20190710	10 Jul 2019 15:10		15 Jul 2019 10:03	15 Jul 2019 16:01	1
HS19070575-11	WG-1620-MW15A-20190710	10 Jul 2019 16:10		15 Jul 2019 10:03	17 Jul 2019 15:35	50
HS19070575-11	WG-1620-MW15A-20190710	10 Jul 2019 16:10		15 Jul 2019 10:03	16 Jul 2019 21:55	10
HS19070575-11	WG-1620-MW15A-20190710	10 Jul 2019 16:10		15 Jul 2019 10:03	15 Jul 2019 16:21	1
HS19070575-12	WG-1620-MW15C-20190710	10 Jul 2019 17:05		15 Jul 2019 10:03	16 Jul 2019 22:15	10
HS19070575-12	WG-1620-MW15C-20190710	10 Jul 2019 17:05		15 Jul 2019 10:03	15 Jul 2019 16:41	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
HS19070575-13	WG-1620-MW15B-20190710	10 Jul 2019 18:00		15 Jul 2019 10:03	17 Jul 2019 13:01	100
HS19070575-13	WG-1620-MW15B-20190710	10 Jul 2019 18:00		15 Jul 2019 10:03	17 Jul 2019 12:42	10
HS19070575-13	WG-1620-MW15B-20190710	10 Jul 2019 18:00		15 Jul 2019 10:03	15 Jul 2019 17:00	1
HS19070575-14	WG-1620-FB01-20190710	10 Jul 2019 18:15		15 Jul 2019 10:03	15 Jul 2019 17:20	1
HS19070575-15	WG-1620-MW14-20190711	11 Jul 2019 07:20		15 Jul 2019 10:03	15 Jul 2019 17:39	1
HS19070575-16	WG-1620-MW13-20190711	11 Jul 2019 08:15		15 Jul 2019 10:03	15 Jul 2019 17:59	1
HS19070575-17	WG-1620-MW39B-20190711	11 Jul 2019 09:15		15 Jul 2019 10:03	15 Jul 2019 18:19	1
HS19070575-18	WG-1620-MW12C-20190711	11 Jul 2019 10:00		15 Jul 2019 10:03	15 Jul 2019 18:38	1
HS19070575-19	WG-1620-MW12A-20190711	11 Jul 2019 10:55		15 Jul 2019 10:03	17 Jul 2019 13:59	100
HS19070575-19	WG-1620-MW12A-20190711	11 Jul 2019 10:55		15 Jul 2019 10:03	17 Jul 2019 13:40	10
HS19070575-19	WG-1620-MW12A-20190711	11 Jul 2019 10:55		15 Jul 2019 10:03	15 Jul 2019 18:58	1
<b>Batch ID:</b> 143067 ( 0 )		<b>Test Name :</b> ICP-MS METALS BY SW6020A			<b>Matrix:</b> Water	
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20		16 Jul 2019 11:30	16 Jul 2019 22:38	1
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05		16 Jul 2019 11:30	16 Jul 2019 22:40	1
HS19070575-03	WG-1620-MW58A-20190710	10 Jul 2019 08:55		16 Jul 2019 11:30	16 Jul 2019 22:43	1
HS19070575-04	WG-1620-MW57A-20190710	10 Jul 2019 09:55		16 Jul 2019 11:30	16 Jul 2019 22:45	1
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40		16 Jul 2019 11:30	16 Jul 2019 22:47	1
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35		16 Jul 2019 11:30	17 Jul 2019 12:21	1
HS19070575-07	WG-1620-MW19C-20190710	10 Jul 2019 12:25		16 Jul 2019 11:30	16 Jul 2019 22:52	1
HS19070575-08	WG-1620-MW17C-20190710	10 Jul 2019 13:25		16 Jul 2019 11:30	16 Jul 2019 22:54	1
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15		16 Jul 2019 11:30	17 Jul 2019 12:23	1
HS19070575-10	WG-1620-MW20A-20190710	10 Jul 2019 15:10		16 Jul 2019 11:30	16 Jul 2019 23:03	1
HS19070575-11	WG-1620-MW15A-20190710	10 Jul 2019 16:10		16 Jul 2019 11:30	16 Jul 2019 22:18	1
HS19070575-12	WG-1620-MW15C-20190710	10 Jul 2019 17:05		16 Jul 2019 11:30	16 Jul 2019 23:05	1
HS19070575-13	WG-1620-MW15B-20190710	10 Jul 2019 18:00		16 Jul 2019 11:30	16 Jul 2019 23:08	1
HS19070575-14	WG-1620-FB01-20190710	10 Jul 2019 18:15		16 Jul 2019 11:30	16 Jul 2019 23:10	1
HS19070575-15	WG-1620-MW14-20190711	11 Jul 2019 07:20		16 Jul 2019 11:30	16 Jul 2019 23:12	1
HS19070575-16	WG-1620-MW13-20190711	11 Jul 2019 08:15		16 Jul 2019 11:30	16 Jul 2019 23:14	1
HS19070575-17	WG-1620-MW39B-20190711	11 Jul 2019 09:15		16 Jul 2019 11:30	16 Jul 2019 23:16	1
HS19070575-18	WG-1620-MW12C-20190711	11 Jul 2019 10:00		16 Jul 2019 11:30	16 Jul 2019 23:19	1
<b>Batch ID:</b> 143129 ( 0 )		<b>Test Name :</b> ICP-MS METALS BY SW6020A			<b>Matrix:</b> Water	
HS19070575-19	WG-1620-MW12A-20190711	11 Jul 2019 10:55		17 Jul 2019 13:00	17 Jul 2019 22:23	1
<b>Batch ID:</b> R342400 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Water	
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20			14 Jul 2019 04:56	1
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05			14 Jul 2019 06:09	1
HS19070575-03	WG-1620-MW58A-20190710	10 Jul 2019 08:55			14 Jul 2019 05:20	1
HS19070575-04	WG-1620-MW57A-20190710	10 Jul 2019 09:55			14 Jul 2019 05:45	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: R342410 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40			14 Jul 2019 06:47	10
HS19070575-05	WG-1620-MW57B-20190710	10 Jul 2019 10:40			14 Jul 2019 06:19	1
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35			14 Jul 2019 07:38	10
HS19070575-06	WG-1620-MW72B-20190710	10 Jul 2019 11:35			14 Jul 2019 07:11	1
HS19070575-07	WG-1620-MW19C-20190710	10 Jul 2019 12:25			14 Jul 2019 00:59	1
HS19070575-08	WG-1620-MW17C-20190710	10 Jul 2019 13:25			14 Jul 2019 02:13	1
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15			14 Jul 2019 08:30	10
HS19070575-09	WG-1620-MW17-20190710	10 Jul 2019 14:15			14 Jul 2019 08:03	1
HS19070575-10	WG-1620-MW20A-20190710	10 Jul 2019 15:10			14 Jul 2019 02:38	1
HS19070575-11	WG-1620-MW15A-20190710	10 Jul 2019 16:10			14 Jul 2019 03:02	1
HS19070575-12	WG-1620-MW15C-20190710	10 Jul 2019 17:05			14 Jul 2019 03:27	1
HS19070575-13	WG-1620-MW15B-20190710	10 Jul 2019 18:00			14 Jul 2019 03:52	1
HS19070575-14	WG-1620-FB01-20190710	10 Jul 2019 18:15			13 Jul 2019 23:46	1
HS19070575-15	WG-1620-MW14-20190711	11 Jul 2019 07:20			14 Jul 2019 04:16	1
HS19070575-16	WG-1620-MW13-20190711	11 Jul 2019 08:15			14 Jul 2019 04:41	1
HS19070575-17	WG-1620-MW39B-20190711	11 Jul 2019 09:15			14 Jul 2019 05:06	1
HS19070575-18	WG-1620-MW12C-20190711	11 Jul 2019 10:00			14 Jul 2019 05:30	1
HS19070575-19	WG-1620-MW12A-20190711	11 Jul 2019 10:55			14 Jul 2019 05:55	1
HS19070575-20	WQ-1620-TB01-20190710	10 Jul 2019 00:00			14 Jul 2019 00:10	1
HS19070575-21	WQ-1620-TB02-20190710	10 Jul 2019 00:00			14 Jul 2019 00:35	1
<b>Batch ID: R342474 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS19070575-01	WG-1620-MW18A-20190710	10 Jul 2019 07:20			15 Jul 2019 19:23	5
HS19070575-02	WG-1620-MW18C-20190710	10 Jul 2019 08:05			15 Jul 2019 19:50	10

WorkOrder: HS19070575  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200

WorkOrder: HS19070575  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00010	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000096	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00012	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00011	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.00010	0.00012	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000013	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00013	0.000047	0.0010
A	Acenaphthene	83-32-9	0.00010	0.00011	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.00010	0.00010	0.000015	0.00010
A	Anthracene	120-12-7	0.00010	0.00011	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.00010	0.000099	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.00010	0.000076	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000086	0.000037	0.00020
A	Chrysene	218-01-9	0.00010	0.00011	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.00010	0.000094	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.000099	0.000020	0.00020
A	Fluoranthene	206-44-0	0.00010	0.00012	0.000010	0.00010
A	Fluorene	86-73-7	0.00010	0.00012	0.000030	0.00010
A	Naphthalene	91-20-3	0.00010	0.00010	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00013	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000085	0.000079	0.00020
A	Phenanthrene	85-01-8	0.00010	0.00011	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000094	0.000035	0.00020
A	Pyrene	129-00-0	0.00010	0.00011	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19070575  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00065	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00060	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00062	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00067	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00066	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00051	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00076	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010



WorkOrder: HS19070575  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00064	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00060	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00061	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00077	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00058	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00070	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00062	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

**Batch ID:** 143067 ( 0)      **Instrument:** ICPMS05      **Method:** ICP-MS METALS BY SW6020A

**MBLK**      Sample ID: **MBLK-143067**      Units: **mg/L**      Analysis Date: **16-Jul-2019 22:14**  
 Client ID:      Run ID: **ICPMS05\_342486**      SeqNo: **5169331**      PrepDate: **16-Jul-2019**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      RPD %RPD Limit Qual

Arsenic      U      0.00200

**LCS**      Sample ID: **LCS-143067**      Units: **mg/L**      Analysis Date: **16-Jul-2019 22:16**  
 Client ID:      Run ID: **ICPMS05\_342486**      SeqNo: **5169332**      PrepDate: **16-Jul-2019**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      RPD %RPD Limit Qual

Arsenic      0.05122      0.00200      0.05      0      102      80 - 120

**MS**      Sample ID: **HS19070575-11MS**      Units: **mg/L**      Analysis Date: **16-Jul-2019 22:23**  
 Client ID: **WG-1620-MW15A-20190710**      Run ID: **ICPMS05\_342486**      SeqNo: **5169335**      PrepDate: **16-Jul-2019**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      RPD %RPD Limit Qual

Arsenic      0.07665      0.00200      0.05      0.02509      103      80 - 120

**MSD**      Sample ID: **HS19070575-11MSD**      Units: **mg/L**      Analysis Date: **16-Jul-2019 22:25**  
 Client ID: **WG-1620-MW15A-20190710**      Run ID: **ICPMS05\_342486**      SeqNo: **5169336**      PrepDate: **16-Jul-2019**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      RPD %RPD Limit Qual

Arsenic      0.07628      0.00200      0.05      0.02509      102      80 - 120      0.07665      0.475      20

**PDS**      Sample ID: **HS19070575-11PDS**      Units: **mg/L**      Analysis Date: **16-Jul-2019 22:27**  
 Client ID: **WG-1620-MW15A-20190710**      Run ID: **ICPMS05\_342486**      SeqNo: **5169337**      PrepDate: **16-Jul-2019**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      RPD %RPD Limit Qual

Arsenic      0.1292      0.00200      0.1      0.02509      104      75 - 125

**SD**      Sample ID: **HS19070575-11SD**      Units: **mg/L**      Analysis Date: **16-Jul-2019 22:20**  
 Client ID: **WG-1620-MW15A-20190710**      Run ID: **ICPMS05\_342486**      SeqNo: **5169334**      PrepDate: **16-Jul-2019**      DF: **5**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %D %D Limit Qual

Arsenic      0.02454      0.0100      0.02509      2.21      10

**The following samples were analyzed in this batch:**

HS19070575-01	HS19070575-02	HS19070575-03	HS19070575-04
HS19070575-05	HS19070575-06	HS19070575-07	HS19070575-08
HS19070575-09	HS19070575-10	HS19070575-11	HS19070575-12
HS19070575-13	HS19070575-14	HS19070575-15	HS19070575-16
HS19070575-17	HS19070575-18		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

Batch ID: 143129 ( 0 )		Instrument: ICPMS05			Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-143129</b>	Units: <b>mg/L</b>			Analysis Date: <b>17-Jul-2019 22:19</b>					
Client ID:		Run ID: <b>ICPMS05_342573</b>			SeqNo: <b>5171278</b>		PrepDate: <b>17-Jul-2019</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-143129</b>	Units: <b>mg/L</b>			Analysis Date: <b>17-Jul-2019 22:21</b>					
Client ID:		Run ID: <b>ICPMS05_342573</b>			SeqNo: <b>5171279</b>		PrepDate: <b>17-Jul-2019</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.04949	0.00200	0.05	0	99.0	80 - 120				
<b>MS</b>	Sample ID: <b>HS19070575-19MS</b>	Units: <b>mg/L</b>			Analysis Date: <b>17-Jul-2019 22:28</b>					
Client ID: <b>WG-1620-MW12A-20190711</b>		Run ID: <b>ICPMS05_342573</b>			SeqNo: <b>5171282</b>		PrepDate: <b>17-Jul-2019</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.05141	0.00200	0.05	0.001917	99.0	80 - 120				
<b>MSD</b>	Sample ID: <b>HS19070575-19MSD</b>	Units: <b>mg/L</b>			Analysis Date: <b>17-Jul-2019 22:30</b>					
Client ID: <b>WG-1620-MW12A-20190711</b>		Run ID: <b>ICPMS05_342573</b>			SeqNo: <b>5171283</b>		PrepDate: <b>17-Jul-2019</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.05143	0.00200	0.05	0.001917	99.0	80 - 120	0.05141	0.0486	20	
<b>PDS</b>	Sample ID: <b>HS19070575-19PDS</b>	Units: <b>mg/L</b>			Analysis Date: <b>17-Jul-2019 22:32</b>					
Client ID: <b>WG-1620-MW12A-20190711</b>		Run ID: <b>ICPMS05_342573</b>			SeqNo: <b>5171284</b>		PrepDate: <b>17-Jul-2019</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.1245	0.00200	0.1	0.001917	123	75 - 125				
<b>SD</b>	Sample ID: <b>HS19070575-19SD</b>	Units: <b>mg/L</b>			Analysis Date: <b>17-Jul-2019 22:26</b>					
Client ID: <b>WG-1620-MW12A-20190711</b>		Run ID: <b>ICPMS05_342573</b>			SeqNo: <b>5171281</b>		PrepDate: <b>17-Jul-2019</b>		DF: <b>5</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual
Arsenic	0.002209	0.0100					0.001917	0	10	J

The following samples were analyzed in this batch: HS19070575-19

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

Batch ID: 143014 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143014	Units: ug/L			Analysis Date: 15-Jul-2019 12:07					
Client ID:	Run ID: SV-7_342456	SeqNo: 5167140		PrepDate: 15-Jul-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.701</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.828</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.6</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.803</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.1</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>3.47</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.4</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.43</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.183</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.7</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

Batch ID: 143014 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143014	Units: ug/L			Analysis Date: 15-Jul-2019 12:26					
Client ID:	Run ID: SV-7_342456	SeqNo: 5167141		PrepDate: 15-Jul-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.902	0.20	5	0	78.0	39 - 127				
2,4-Dimethylphenol	3.394	0.20	5	0	67.9	35 - 120				
2,4-Dinitrotoluene	4.229	0.20	5	0	84.6	50 - 122				
2,6-Dinitrotoluene	3.892	0.20	5	0	77.8	50 - 120				
2-Chloronaphthalene	3.393	0.20	5	0	67.9	50 - 120				
2-Methylnaphthalene	3.647	0.10	5	0	72.9	50 - 120				
4,6-Dinitro-2-methylphenol	3.043	0.20	5	0	60.9	25 - 121				
4-Nitrophenol	3.955	1.0	5	0	79.1	30 - 130				
Acenaphthene	3.452	0.10	5	0	69.0	45 - 120				
Acenaphthylene	3.423	0.10	5	0	68.5	47 - 120				
Anthracene	3.186	0.10	5	0	63.7	45 - 120				
Benz(a)anthracene	3.587	0.10	5	0	71.7	40 - 120				
Benzo(a)pyrene	2.95	0.10	5	0	59.0	45 - 120				
Bis(2-chloroethoxy)methane	3.517	0.20	5	0	70.3	45 - 120				
Bis(2-ethylhexyl)phthalate	3.572	0.20	5	0	71.4	40 - 139				
Chrysene	3.538	0.10	5	0	70.8	43 - 120				
Dibenzofuran	3.567	0.10	5	0	71.3	50 - 120				
Di-n-butyl phthalate	3.477	0.20	5	0	69.5	45 - 123				
Fluoranthene	3.475	0.10	5	0	69.5	45 - 125				
Fluorene	3.341	0.10	5	0	66.8	49 - 120				
Naphthalene	3.401	0.10	5	0	68.0	45 - 120				
Nitrobenzene	4.121	0.20	5	0	82.4	44 - 120				
N-Nitrosodiphenylamine	3.473	0.20	5	0	69.5	40 - 125				
Pentachlorophenol	2.714	0.20	5	0	54.3	19 - 121				
Phenanthrene	3.125	0.10	5	0	62.5	45 - 121				
Phenol	3.76	0.20	5	0	75.2	20 - 124				
Pyrene	3.463	0.10	5	0	69.3	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.938</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.8</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.839</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.8</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.753</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>75.1</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>3.464</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.3</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.278</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.197</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.9</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

Batch ID: 143014 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-143014		Units: ug/L		Analysis Date: 15-Jul-2019 12:46				
Client ID:		Run ID: SV-7_342456		SeqNo: 5167142		PrepDate: 15-Jul-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.957	0.20	5	0	79.1	39 - 127	3.902	1.41	20	
2,4-Dimethylphenol	3.566	0.20	5	0	71.3	35 - 120	3.394	4.94	20	
2,4-Dinitrotoluene	4.08	0.20	5	0	81.6	50 - 122	4.229	3.59	20	
2,6-Dinitrotoluene	3.82	0.20	5	0	76.4	50 - 120	3.892	1.87	20	
2-Chloronaphthalene	3.286	0.20	5	0	65.7	50 - 120	3.393	3.21	20	
2-Methylnaphthalene	3.596	0.10	5	0	71.9	50 - 120	3.647	1.41	20	
4,6-Dinitro-2-methylphenol	3.247	0.20	5	0	64.9	25 - 121	3.043	6.47	30	
4-Nitrophenol	3.728	1.0	5	0	74.6	30 - 130	3.955	5.92	20	
Acenaphthene	3.494	0.10	5	0	69.9	45 - 120	3.452	1.22	20	
Acenaphthylene	3.398	0.10	5	0	68.0	47 - 120	3.423	0.726	20	
Anthracene	3.226	0.10	5	0	64.5	45 - 120	3.186	1.26	20	
Benz(a)anthracene	3.638	0.10	5	0	72.8	40 - 120	3.587	1.4	20	
Benzo(a)pyrene	2.952	0.10	5	0	59.0	45 - 120	2.95	0.0403	20	
Bis(2-chloroethoxy)methane	3.497	0.20	5	0	69.9	45 - 120	3.517	0.59	20	
Bis(2-ethylhexyl)phthalate	3.575	0.20	5	0	71.5	40 - 139	3.572	0.0967	20	
Chrysene	3.526	0.10	5	0	70.5	43 - 120	3.538	0.344	20	
Dibenzofuran	3.479	0.10	5	0	69.6	50 - 120	3.567	2.5	20	
Di-n-butyl phthalate	3.423	0.20	5	0	68.5	45 - 123	3.477	1.58	20	
Fluoranthene	3.476	0.10	5	0	69.5	45 - 125	3.475	0.0122	20	
Fluorene	3.303	0.10	5	0	66.1	49 - 120	3.341	1.14	20	
Naphthalene	3.377	0.10	5	0	67.5	45 - 120	3.401	0.71	20	
Nitrobenzene	4.097	0.20	5	0	81.9	44 - 120	4.121	0.591	20	
N-Nitrosodiphenylamine	3.473	0.20	5	0	69.5	40 - 125	3.473	0.0165	20	
Pentachlorophenol	2.83	0.20	5	0	56.6	19 - 121	2.714	4.18	20	
Phenanthrene	3.152	0.10	5	0	63.0	45 - 121	3.125	0.835	20	
Phenol	3.643	0.20	5	0	72.9	20 - 124	3.76	3.19	20	
Pyrene	3.488	0.10	5	0	69.8	40 - 130	3.463	0.698	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.744</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.9</i>	<i>34 - 129</i>	<i>3.938</i>	<i>5.07</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.773</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>75.5</i>	<i>40 - 125</i>	<i>3.839</i>	<i>1.75</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.729</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.6</i>	<i>20 - 120</i>	<i>3.753</i>	<i>0.655</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>3.457</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.1</i>	<i>40 - 135</i>	<i>3.464</i>	<i>0.215</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>4.223</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>84.5</i>	<i>41 - 120</i>	<i>4.278</i>	<i>1.28</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.186</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.7</i>	<i>20 - 120</i>	<i>4.197</i>	<i>0.273</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

<b>Batch ID:</b> 143014 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D
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The following samples were analyzed in this batch:

HS19070575-01	HS19070575-02	HS19070575-03	HS19070575-04
HS19070575-05	HS19070575-06	HS19070575-07	HS19070575-08
HS19070575-09	HS19070575-10	HS19070575-11	HS19070575-12
HS19070575-13	HS19070575-14	HS19070575-15	HS19070575-16
HS19070575-17	HS19070575-18	HS19070575-19	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

<b>Batch ID:</b> R342400 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-190713</b>	Units: <b>ug/L</b>			Analysis Date: <b>13-Jul-2019 23:38</b>				
Client ID:	Run ID: <b>VOA2_342400</b>	SeqNo: <b>5166127</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>42.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>84.6</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.6</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-190713</b>	Units: <b>ug/L</b>			Analysis Date: <b>13-Jul-2019 23:13</b>				
Client ID:	Run ID: <b>VOA2_342400</b>	SeqNo: <b>5166126</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	19.36	1.0	20	0	96.8	70 - 124			
Benzene	20.6	1.0	20	0	103	74 - 120			
Chlorobenzene	19.47	1.0	20	0	97.3	76 - 113			
Ethylbenzene	19.57	1.0	20	0	97.8	77 - 117			
Methylene chloride	20.05	2.0	20	0	100	70 - 127			
Toluene	22.21	1.0	20	0	111	77 - 118			
Vinyl chloride	23.29	1.0	20	0	116	70 - 130			
Xylenes, Total	64.58	1.0	60	0	108	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.0</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>47.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.6</i>	<i>81 - 120</i>			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

**Batch ID:** R342400 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS19070509-10MS			Units: ug/L		Analysis Date: 14-Jul-2019 00:51			
Client ID:		Run ID: VOA2_342400			SeqNo: 5166130		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.54	1.0	20	0	82.7	70 - 127				
Benzene	17.76	1.0	20	0	88.8	70 - 127				
Chlorobenzene	17.58	1.0	20	0	87.9	70 - 114				
Ethylbenzene	17.7	1.0	20	0	88.5	70 - 124				
Methylene chloride	17.93	2.0	20	0	89.7	70 - 128				
Toluene	20.17	1.0	20	0	101	70 - 123				
Vinyl chloride	20.19	1.0	20	0	101	70 - 130				
Xylenes, Total	57.61	1.0	60	0	96.0	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>90.4</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.16</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.3</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.56</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

MSD		Sample ID: HS19070509-10MSD			Units: ug/L		Analysis Date: 14-Jul-2019 01:16			
Client ID:		Run ID: VOA2_342400			SeqNo: 5166131		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.17	1.0	20	0	80.8	70 - 127	16.54	2.3	20	
Benzene	17.18	1.0	20	0	85.9	70 - 127	17.76	3.31	20	
Chlorobenzene	16.7	1.0	20	0	83.5	70 - 114	17.58	5.17	20	
Ethylbenzene	16.96	1.0	20	0	84.8	70 - 124	17.7	4.27	20	
Methylene chloride	16.55	2.0	20	0	82.8	70 - 128	17.93	8.02	20	
Toluene	18.64	1.0	20	0	93.2	70 - 123	20.17	7.89	20	
Vinyl chloride	19.45	1.0	20	0	97.2	70 - 130	20.19	3.77	20	
Xylenes, Total	55.71	1.0	60	0	92.9	70 - 130	57.61	3.35	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>44.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>88.7</i>	<i>70 - 126</i>	<i>45.2</i>	<i>1.85</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.0</i>	<i>81 - 113</i>	<i>48.16</i>	<i>1.41</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>77 - 123</i>	<i>48.56</i>	<i>2.44</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.77</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>82 - 127</i>	<i>50.25</i>	<i>0.966</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070575-01    HS19070575-02    HS19070575-03    HS19070575-04

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

**Batch ID:** R342410 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190713</b>			Units: <b>ug/L</b>		Analysis Date: <b>13-Jul-2019 23:21</b>			
Client ID:		Run ID: <b>VOA4_342410</b>			SeqNo: <b>5166422</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Vinyl chloride	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.93</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>52.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-190713</b>			Units: <b>ug/L</b>		Analysis Date: <b>13-Jul-2019 22:31</b>			
Client ID:		Run ID: <b>VOA4_342410</b>			SeqNo: <b>5166421</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.33	1.0	20	0	91.6	70 - 124				
Benzene	18.78	1.0	20	0	93.9	74 - 120				
Chlorobenzene	19.89	1.0	20	0	99.4	76 - 113				
Ethylbenzene	19.73	1.0	20	0	98.7	77 - 117				
Methylene chloride	19.5	2.0	20	0	97.5	70 - 127				
Toluene	19.33	1.0	20	0	96.7	77 - 118				
Vinyl chloride	19.32	1.0	20	0	96.6	70 - 130				
Xylenes, Total	63.31	1.0	60	0	106	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.4</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.19</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.4</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

**Batch ID:** R342410 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS19070575-07MS			Units: ug/L		Analysis Date: 14-Jul-2019 01:24			
Client ID: WG-1620-MW19C-20190710		Run ID: VOA4_342410			SeqNo: 5166427		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.77	1.0	20	0	83.9	70 - 127				
Benzene	18.56	1.0	20	0.3991	90.8	70 - 127				
Chlorobenzene	18.71	1.0	20	0	93.5	70 - 114				
Ethylbenzene	19.94	1.0	20	0.9597	94.9	70 - 124				
Methylene chloride	18.1	2.0	20	0	90.5	70 - 128				
Toluene	19.42	1.0	20	0.8601	92.8	70 - 123				
Vinyl chloride	19.21	1.0	20	0	96.1	70 - 130				
Xylenes, Total	63.56	1.0	60	2.208	102	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.1</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.2</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.66</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>82 - 127</i>				

MSD		Sample ID: HS19070575-07MSD			Units: ug/L		Analysis Date: 14-Jul-2019 01:49			
Client ID: WG-1620-MW19C-20190710		Run ID: VOA4_342410			SeqNo: 5166428		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.01	1.0	20	0	85.0	70 - 127	16.77	1.39	20	
Benzene	18.15	1.0	20	0.3991	88.8	70 - 127	18.56	2.23	20	
Chlorobenzene	18.19	1.0	20	0	90.9	70 - 114	18.71	2.83	20	
Ethylbenzene	19.39	1.0	20	0.9597	92.2	70 - 124	19.94	2.78	20	
Methylene chloride	17.68	2.0	20	0	88.4	70 - 128	18.1	2.38	20	
Toluene	19.25	1.0	20	0.8601	92.0	70 - 123	19.42	0.87	20	
Vinyl chloride	18.26	1.0	20	0	91.3	70 - 130	19.21	5.06	20	
Xylenes, Total	60.19	1.0	60	2.208	96.6	70 - 130	63.56	5.45	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.27</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.5</i>	<i>70 - 126</i>	<i>48.1</i>	<i>1.73</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>	<i>51.66</i>	<i>2.67</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.0</i>	<i>77 - 123</i>	<i>48.64</i>	<i>0.291</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>49.6</i>	<i>1.44</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS19070575-05	HS19070575-06	HS19070575-07	HS19070575-08
HS19070575-09	HS19070575-10	HS19070575-11	HS19070575-12
HS19070575-13	HS19070575-14	HS19070575-15	HS19070575-16
HS19070575-17	HS19070575-18	HS19070575-19	HS19070575-20
HS19070575-21			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

Batch ID: R342474 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MBLK</b>	Sample ID: <b>VBLKW-190715</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jul-2019 12:40</b>					
Client ID:	Run ID: <b>VOA2_342474</b>	SeqNo: <b>5167694</b>		PrepDate:			DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
Benzene	U	1.0								
Ethylbenzene	U	1.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	44.39	1.0	50	0	88.8	70 - 123				
<i>Surr: 4-Bromofluorobenzene</i>	46.81	1.0	50	0	93.6	82 - 115				
<i>Surr: Dibromofluoromethane</i>	48.92	1.0	50	0	97.8	73 - 126				
<i>Surr: Toluene-d8</i>	50.4	1.0	50	0	101	81 - 120				
<b>LCS</b>	Sample ID: <b>VLCSW-190715</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jul-2019 11:51</b>					
Client ID:	Run ID: <b>VOA2_342474</b>	SeqNo: <b>5167693</b>		PrepDate:			DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
Benzene	19.79	1.0	20	0	98.9	74 - 120				
Ethylbenzene	20.53	1.0	20	0	103	77 - 117				
Toluene	22.56	1.0	20	0	113	77 - 118				
Xylenes, Total	66.83	1.0	60	0	111	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	46.06	1.0	50	0	92.1	70 - 130				
<i>Surr: 4-Bromofluorobenzene</i>	47.86	1.0	50	0	95.7	82 - 115				
<i>Surr: Dibromofluoromethane</i>	47.97	1.0	50	0	95.9	73 - 126				
<i>Surr: Toluene-d8</i>	50.17	1.0	50	0	100	81 - 120				
<b>MS</b>	Sample ID: <b>HS19070637-01MS</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jul-2019 14:47</b>					
Client ID:	Run ID: <b>VOA2_342474</b>	SeqNo: <b>5167699</b>		PrepDate:			DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
Benzene	16.39	1.0	20	0	82.0	70 - 127				
Ethylbenzene	16.9	1.0	20	0	84.5	70 - 124				
Toluene	18.59	1.0	20	0	93.0	70 - 123				
Xylenes, Total	55.82	1.0	60	0	93.0	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	45.39	1.0	50	0	90.8	70 - 126				
<i>Surr: 4-Bromofluorobenzene</i>	48.72	1.0	50	0	97.4	81 - 113				
<i>Surr: Dibromofluoromethane</i>	47.25	1.0	50	0	94.5	77 - 123				
<i>Surr: Toluene-d8</i>	50.44	1.0	50	0	101	82 - 127				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QC BATCH REPORT**

**Batch ID:** R342474 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MSD</b>		Sample ID: <b>HS19070637-01MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jul-2019 15:11</b>			
Client ID:		Run ID: <b>VOA2_342474</b>			SeqNo: <b>5167700</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	15.74	1.0	20	0	78.7	70 - 127	16.39	4.06	20	
Ethylbenzene	16.05	1.0	20	0	80.2	70 - 124	16.9	5.16	20	
Toluene	16.99	1.0	20	0	84.9	70 - 123	18.59	9.02	20	
Xylenes, Total	50.24	1.0	60	0	83.7	70 - 130	55.82	10.5	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.08</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.2</i>	<i>70 - 126</i>	<i>45.39</i>	<i>1.5</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.43</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.9</i>	<i>81 - 113</i>	<i>48.72</i>	<i>2.69</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.53</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.1</i>	<i>77 - 123</i>	<i>47.25</i>	<i>0.583</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>82 - 127</i>	<i>50.44</i>	<i>3.32</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070575-01      HS19070575-02

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070575

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kansas	E-10352 2018-2019	31-Jul-2019
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020







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# Chain of Custody Form

Page 1 of     

COC ID: 196131

## HS19070575

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UFRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold	
1	WG-1620-1801-20190710	7-10-19	-	Water	1	2		X										
2	WG-1620-MW18A-20190710	7-10-19	0720	W		6		X	X	X								
3	WG-1620-MW18C-20190710		0805	W		6		X	X	X								
4	WG-1620-MW58A-20190710		0855	W		6		X	X	X								
5	WG-1620-MW57A-20190710		0955	W		6		X	X	X								
6	WG-1620-MW57B-20190710		1040	W		6		X	X	X								
7	WG-1620-MW72B-20190710		1135	W		6		X	X	X								
8	WG-1620-MW19C-20190710		1225	W		6		X	X	X								
9	WG-1620-MW17C-20190710		1325	W		6		X	X	X								
10	WG-1620-MW-17-20190710		1415	W		6		X	X	X								

Sampler(s) Please Print & Sign  
**JOHN BRAYTON**      **John Br**

Relinquished by: **John Br**      Date: **7-11-19**      Time: **16:45**

Relinquished by: **John Br**      Date: **7/11/19**      Time: **16:45**

Logged by (Laboratory): \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

Shipment Method: **HAND DELIVERED**      Required Turnaround Time: (Check Box)  STD 10 Wk Drys       5 Wk Days       2 Wk Days       24 Hour

Notes: **UPRR Houston MWPW**

Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)
44145	0.7	<input type="checkbox"/> Level II Std OC <input checked="" type="checkbox"/> TRRP Checklist
43623	1.2	<input type="checkbox"/> Level III Std OC/Raw Data <input type="checkbox"/> TRRP Level IV
44608	1.0	<input type="checkbox"/> Level IV SW/MS/CLP
44381	0.9	<input type="checkbox"/> Other

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

ote: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
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 3. The Chain of Custody is a legal document. All information must be completed accurately.

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(1211      CF 0-1)



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# Chain of Custody Form

Page 2 of 2

COC ID: 196140

HS19070575

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UFRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92888	B	8200_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Malzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652848 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.malzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-T80-20190710</del>			Water	1	2		X									
2	WG-1620-MW20A-20190710	7-10-19	1510	W		6	X		X	X							
3	WG-1620-MW1SA-20190710		1610	W		6	X		X	X							
4	WG-1620-MW1SC-20190710		1705	W		6	X		X	X							
5	WG-1620-MW1SB-20190710		1800	W		6	X		X	X							
6	WG-1620-FB01-20190710		1815	W		6		X	X	X							
7	WG-1620-MW14-20190710	7-11-19	0720	W		6	X		X	X							
8	WG-1620-MW13-20190711		0815	W		6	X		X	X							
9	WG-1620-MW39B-20190711		0915	W		6	X		X	X							
10	WG-1620-MW12C-20190711		1000	W		6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BEASLEY</b>	Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour	Other: _____	Results Due Date: _____
Relinquished by: <b>John B</b>	Date: <b>7/11/19</b> Time: <b>16:45</b>	Received by: <b>J. M...</b>	Notes: <b>UPRR Houston MWPW</b>	
Relinquished by: <b>John B</b>	Date: <b>7/11/19</b> Time: <b>16:45</b>	Received by (Laboratory): <b>J. M...</b>	QC Package: (Check One Box Below)	
Logged by (Laboratory):	Date: _____ Time: _____	Checked by (Laboratory):	<input type="checkbox"/> Level II Std QC	<input checked="" type="checkbox"/> TRRP Checklist
			<input type="checkbox"/> Level III Std QC/RW Data	<input type="checkbox"/> TRRP Level IV
			<input type="checkbox"/> Level IV SW846/CLP	
			<input type="checkbox"/> Other: _____	

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

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# Chain of Custody Form

## HS19070575

W

Page 1 of 1

COC ID: 196123

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Malzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652846 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.malzner@pbwlc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TD0-201804</del>							*									
2	WG-1620-MW12A-20190711	7-11-19	1055	W		6	X		X	X							
3																	
4																	
5																	
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24-hour				Results Due Date:			
Relinquished by: <b>John Brayton</b>		Date: <b>7/12/19</b> Time: <b>16:45</b>		Received by:		Notes: <b>UPRR Houston MWPW</b>					
Relinquished by: <b>John Brayton</b>		Date: <b>7/11/19</b> Time: <b>16:45</b>		Received by (Laboratory): <b>J. MURPHY</b>		Cooler ID		Cooler Temp.		QC Package: (Check One Box Below)	
Logged by (Laboratory):		Date: Time:		Checked by (Laboratory):						<input type="checkbox"/> Level II Std GC <input checked="" type="checkbox"/> TTRP Checklist	
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035										<input type="checkbox"/> Level III Std GC/Row Data <input type="checkbox"/> TTRP Level IV	
										<input type="checkbox"/> Level IV SW/642CLP	
										<input type="checkbox"/> Other	

ote: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
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July 22, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19070609**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 8 sample(s) on Jul 12, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

---

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

---

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/22/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070609			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143051,143162,R342484			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			2
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/22/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070609			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143051,143162,R342484			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).



**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 07/22/2019
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS19070609
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 143051,143162,R342484

ER# <sup>5</sup>	Description
1	Semivolatile Organics Method SW8270, sample WG-1620-MW40B-20190711, the surrogate recoveries could not be determined due to dilution below the calibration range.
2	Batch 143051, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
 O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
 NA = Not Applicable;  
 NR = Not Reviewed;  
 R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070609

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19070609-01	WG-1620-MW40B-20190711	Water		11-Jul-2019 12:00	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-02	WG-1620-MW42B-20190711	Water		11-Jul-2019 13:00	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-03	WG-1620-P11-20190711	Water		11-Jul-2019 13:55	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-04	WG-1620-MW05-20190711	Water		11-Jul-2019 14:50	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-05	WG-1620-MW64A-20190711	Water		11-Jul-2019 15:50	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-06	WG-1620-FB02-20190711	Water		11-Jul-2019 16:05	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-07	WG-1620-TB02-20190711	Water	C&G-062119-78	11-Jul-2019 00:00	12-Jul-2019 12:15	<input type="checkbox"/>
HS19070609-08	WQ-1620-TB01-20190711	Water	C&G-062119-78	11-Jul-2019 00:00	12-Jul-2019 12:15	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW40B-20190711  
 Collection Date: 11-Jul-2019 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:00
<b>Benzene</b>	<b>0.0088</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jul-2019 14:00
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:00
<b>Ethylbenzene</b>	<b>0.082</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jul-2019 14:00
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 14:00
<b>Toluene</b>	<b>0.014</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jul-2019 14:00
<b>Xylenes, Total</b>	<b>0.13</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jul-2019 14:00
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:00</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>107</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:00</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:00</i>
<i>Surr: Toluene-d8</i>	<i>98.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW40B-20190711  
 Collection Date: 11-Jul-2019 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jul-2019 21:21
<b>2,4-Dimethylphenol</b>	<b>0.0013</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jul-2019 21:21
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jul-2019 21:21
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jul-2019 21:21
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jul-2019 21:21
<b>2-Methylnaphthalene</b>	<b>0.18</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	18-Jul-2019 15:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jul-2019 21:21
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jul-2019 21:21
<b>Acenaphthene</b>	<b>0.21</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	18-Jul-2019 15:40
<b>Acenaphthylene</b>	<b>0.0021</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:21
<b>Anthracene</b>	<b>0.011</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Jul-2019 15:21
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jul-2019 21:21
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jul-2019 21:21
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jul-2019 21:21
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jul-2019 21:21
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jul-2019 21:21
<b>Dibenzofuran</b>	<b>0.13</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	18-Jul-2019 15:40
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jul-2019 21:21
<b>Fluoranthene</b>	<b>0.0059</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:21
<b>Fluorene</b>	<b>0.13</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	18-Jul-2019 15:40
<b>Naphthalene</b>	<b>3.6</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	18-Jul-2019 15:59
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jul-2019 21:21
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jul-2019 21:21
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jul-2019 21:21
<b>Phenanthrene</b>	<b>0.093</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Jul-2019 15:21
<b>Phenol</b>	<b>0.000053</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jul-2019 21:21
<b>Pyrene</b>	<b>0.0024</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:21
Surr: 2,4,6-Tribromophenol	53.1			34-129	%REC	10	18-Jul-2019 15:21
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	18-Jul-2019 15:40
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	1000	18-Jul-2019 15:59
Surr: 2,4,6-Tribromophenol	88.2			34-129	%REC	1	17-Jul-2019 21:21
Surr: 2-Fluorobiphenyl	65.6			40-125	%REC	1	17-Jul-2019 21:21
Surr: 2-Fluorobiphenyl	59.9			40-125	%REC	10	18-Jul-2019 15:21
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	18-Jul-2019 15:40
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	1000	18-Jul-2019 15:59
Surr: 2-Fluorophenol	61.1			20-120	%REC	10	18-Jul-2019 15:21
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	18-Jul-2019 15:40
Surr: 2-Fluorophenol	0	JS		20-120	%REC	1000	18-Jul-2019 15:59
Surr: 2-Fluorophenol	65.2			20-120	%REC	1	17-Jul-2019 21:21

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW40B-20190711  
 Collection Date: 11-Jul-2019 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
Surr: 4-Terphenyl-d14	76.8			40-135	%REC	10	18-Jul-2019 15:21
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	18-Jul-2019 15:40
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	18-Jul-2019 15:59
Surr: 4-Terphenyl-d14	71.7			40-135	%REC	1	17-Jul-2019 21:21
Surr: Nitrobenzene-d5	66.2			41-120	%REC	10	18-Jul-2019 15:21
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	18-Jul-2019 15:40
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	18-Jul-2019 15:59
Surr: Nitrobenzene-d5	50.5			41-120	%REC	1	17-Jul-2019 21:21
Surr: Phenol-d6	54.1			20-120	%REC	1	17-Jul-2019 21:21
Surr: Phenol-d6	68.9			20-120	%REC	10	18-Jul-2019 15:21
Surr: Phenol-d6	0	JS		20-120	%REC	100	18-Jul-2019 15:40
Surr: Phenol-d6	0	JS		20-120	%REC	1000	18-Jul-2019 15:59
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Jul-2019		Analyst: JHD	
Arsenic	0.0520		0.000400	0.00200	mg/L	1	18-Jul-2019 18:11

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW42B-20190711  
 Collection Date: 11-Jul-2019 13:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:26
Benzene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:26
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:26
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:26
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 14:26
Toluene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:26
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:26
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:26</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:26</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:26</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:26</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW42B-20190711  
 Collection Date: 11-Jul-2019 13:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jul-2019 21:40
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jul-2019 21:40
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jul-2019 21:40
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jul-2019 21:40
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jul-2019 21:40
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	17-Jul-2019 21:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jul-2019 21:40
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jul-2019 21:40
<b>Acenaphthene</b>	<b>0.000068</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jul-2019 21:40
<b>Anthracene</b>	<b>0.000064</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jul-2019 21:40
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jul-2019 21:40
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jul-2019 21:40
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00015</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jul-2019 21:40
<b>Dibenzofuran</b>	<b>0.000038</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jul-2019 21:40
<b>Fluoranthene</b>	<b>0.000040</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
<b>Fluorene</b>	<b>0.000057</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
<b>Naphthalene</b>	<b>0.00043</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jul-2019 21:40
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jul-2019 21:40
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jul-2019 21:40
<b>Phenanthrene</b>	<b>0.000034</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
Phenol	U		0.000035	0.00020	mg/L	1	17-Jul-2019 21:40
<b>Pyrene</b>	<b>0.000033</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jul-2019 21:40
<i>Surr: 2,4,6-Tribromophenol</i>	78.4			34-129	%REC	1	17-Jul-2019 21:40
<i>Surr: 2-Fluorobiphenyl</i>	63.5			40-125	%REC	1	17-Jul-2019 21:40
<i>Surr: 2-Fluorophenol</i>	66.7			20-120	%REC	1	17-Jul-2019 21:40
<i>Surr: 4-Terphenyl-d14</i>	77.9			40-135	%REC	1	17-Jul-2019 21:40
<i>Surr: Nitrobenzene-d5</i>	63.9			41-120	%REC	1	17-Jul-2019 21:40
<i>Surr: Phenol-d6</i>	70.7			20-120	%REC	1	17-Jul-2019 21:40
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00220</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Jul-2019 18:14

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-P11-20190711  
 Collection Date: 11-Jul-2019 13:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:51
Benzene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:51
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:51
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:51
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 14:51
Toluene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 14:51
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jul-2019 14:51
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:51</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:51</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:51</i>
<i>Surr: Toluene-d8</i>	<i>98.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 14:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-P11-20190711  
 Collection Date: 11-Jul-2019 13:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 14:04
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 14:04
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 14:04
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 14:04
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 14:04
<b>2-Methylnaphthalene</b>	<b>0.000050</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 14:04
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 14:04
<b>Acenaphthene</b>	<b>0.0019</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 14:04
<b>Anthracene</b>	<b>0.000063</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 14:04
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 14:04
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 14:04
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000087</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 14:04
<b>Dibenzofuran</b>	<b>0.000079</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 14:04
<b>Fluoranthene</b>	<b>0.000060</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
<b>Fluorene</b>	<b>0.00074</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
<b>Naphthalene</b>	<b>0.0018</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 14:04
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 14:04
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 14:04
<b>Phenanthrene</b>	<b>0.00020</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 14:04
<b>Pyrene</b>	<b>0.000036</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:04
<i>Surr: 2,4,6-Tribromophenol</i>	83.5			34-129	%REC	1	18-Jul-2019 14:04
<i>Surr: 2-Fluorobiphenyl</i>	60.1			40-125	%REC	1	18-Jul-2019 14:04
<i>Surr: 2-Fluorophenol</i>	61.8			20-120	%REC	1	18-Jul-2019 14:04
<i>Surr: 4-Terphenyl-d14</i>	85.9			40-135	%REC	1	18-Jul-2019 14:04
<i>Surr: Nitrobenzene-d5</i>	61.2			41-120	%REC	1	18-Jul-2019 14:04
<i>Surr: Phenol-d6</i>	64.0			20-120	%REC	1	18-Jul-2019 14:04
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0704</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Jul-2019 18:16

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW05-20190711  
 Collection Date: 11-Jul-2019 14:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 15:17
Benzene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 15:17
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 15:17
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 15:17
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 15:17
Toluene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 15:17
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jul-2019 15:17
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:17</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>104</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:17</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:17</i>
<i>Surr: Toluene-d8</i>	<i>98.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:17</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW05-20190711  
 Collection Date: 11-Jul-2019 14:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 14:23
<b>2,4-Dimethylphenol</b>	<b>0.000059</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 14:23
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 14:23
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 14:23
<b>2-Methylnaphthalene</b>	<b>0.000087</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 14:23
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 14:23
<b>Acenaphthene</b>	<b>0.00043</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 14:23
<b>Anthracene</b>	<b>0.000097</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 14:23
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 14:23
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 14:23
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 14:23
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 14:23
<b>Dibenzofuran</b>	<b>0.000051</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 14:23
Fluoranthene	U		0.000010	0.00010	mg/L	1	18-Jul-2019 14:23
<b>Fluorene</b>	<b>0.000064</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
<b>Naphthalene</b>	<b>0.00071</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 14:23
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 14:23
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 14:23
Phenanthrene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 14:23
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 14:23
<b>Pyrene</b>	<b>0.00019</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:23
<i>Surr: 2,4,6-Tribromophenol</i>	<i>94.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:23</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>88.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:23</i>
<i>Surr: 2-Fluorophenol</i>	<i>82.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:23</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>96.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:23</i>
<i>Surr: Nitrobenzene-d5</i>	<i>90.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:23</i>
<i>Surr: Phenol-d6</i>	<i>87.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:23</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0171</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Jul-2019 18:18

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW64A-20190711  
 Collection Date: 11-Jul-2019 15:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 15:42
Benzene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 15:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 15:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 15:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 15:42
Toluene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 15:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jul-2019 15:42
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:42</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:42</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:42</i>
<i>Surr: Toluene-d8</i>	<i>98.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jul-2019 15:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW64A-20190711  
 Collection Date: 11-Jul-2019 15:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 14:42
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 14:42
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 14:42
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 14:42
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 14:42
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	18-Jul-2019 14:42
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 14:42
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 14:42
<b>Acenaphthene</b>	<b>0.00030</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:42
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 14:42
Anthracene	U		0.000014	0.00010	mg/L	1	18-Jul-2019 14:42
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 14:42
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 14:42
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 14:42
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 14:42
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 14:42
Dibenzofuran	U		0.000020	0.00010	mg/L	1	18-Jul-2019 14:42
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 14:42
Fluoranthene	U		0.000010	0.00010	mg/L	1	18-Jul-2019 14:42
<b>Fluorene</b>	<b>0.000031</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:42
<b>Naphthalene</b>	<b>0.00033</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:42
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 14:42
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 14:42
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 14:42
Phenanthrene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 14:42
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 14:42
<b>Pyrene</b>	<b>0.00015</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 14:42
<i>Surr: 2,4,6-Tribromophenol</i>	97.5			34-129	%REC	1	18-Jul-2019 14:42
<i>Surr: 2-Fluorobiphenyl</i>	90.7			40-125	%REC	1	18-Jul-2019 14:42
<i>Surr: 2-Fluorophenol</i>	84.9			20-120	%REC	1	18-Jul-2019 14:42
<i>Surr: 4-Terphenyl-d14</i>	80.5			40-135	%REC	1	18-Jul-2019 14:42
<i>Surr: Nitrobenzene-d5</i>	93.0			41-120	%REC	1	18-Jul-2019 14:42
<i>Surr: Phenol-d6</i>	93.3			20-120	%REC	1	18-Jul-2019 14:42
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00939</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Jul-2019 18:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB02-20190711  
 Collection Date: 11-Jul-2019 16:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 13:09
Benzene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 13:09
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 13:09
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 13:09
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 13:09
Toluene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 13:09
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jul-2019 13:09
<i>Surr: 1,2-Dichloroethane-d4</i>	99.3			70-126	%REC	1	15-Jul-2019 13:09
<i>Surr: 4-Bromofluorobenzene</i>	98.7			81-113	%REC	1	15-Jul-2019 13:09
<i>Surr: Dibromofluoromethane</i>	98.5			77-123	%REC	1	15-Jul-2019 13:09
<i>Surr: Toluene-d8</i>	100			82-127	%REC	1	15-Jul-2019 13:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB02-20190711  
 Collection Date: 11-Jul-2019 16:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	0.00026		0.00021	0.00020	mg/L	1	18-Jul-2019 15:02
2,4-Dimethylphenol	0.000064	J	0.00040	0.00020	mg/L	1	18-Jul-2019 15:02
2,4-Dinitrotoluene		U	0.00058	0.00020	mg/L	1	18-Jul-2019 15:02
2,6-Dinitrotoluene		U	0.00042	0.00020	mg/L	1	18-Jul-2019 15:02
2-Chloronaphthalene		U	0.00021	0.00020	mg/L	1	18-Jul-2019 15:02
2-Methylnaphthalene	0.000077	J	0.00019	0.00010	mg/L	1	18-Jul-2019 15:02
4,6-Dinitro-2-methylphenol		U	0.00020	0.00020	mg/L	1	18-Jul-2019 15:02
4-Nitrophenol		U	0.00047	0.0010	mg/L	1	18-Jul-2019 15:02
Acenaphthene		U	0.00027	0.00010	mg/L	1	18-Jul-2019 15:02
Acenaphthylene		U	0.00015	0.00010	mg/L	1	18-Jul-2019 15:02
Anthracene		U	0.00014	0.00010	mg/L	1	18-Jul-2019 15:02
Benz(a)anthracene		U	0.00050	0.00010	mg/L	1	18-Jul-2019 15:02
Benzo(a)pyrene		U	0.00020	0.00010	mg/L	1	18-Jul-2019 15:02
Bis(2-chloroethoxy)methane		U	0.00030	0.00020	mg/L	1	18-Jul-2019 15:02
Bis(2-ethylhexyl)phthalate		U	0.00037	0.00020	mg/L	1	18-Jul-2019 15:02
Chrysene		U	0.00021	0.00010	mg/L	1	18-Jul-2019 15:02
Dibenzofuran		U	0.00020	0.00010	mg/L	1	18-Jul-2019 15:02
Di-n-butyl phthalate		U	0.00020	0.00020	mg/L	1	18-Jul-2019 15:02
Fluoranthene		U	0.00010	0.00010	mg/L	1	18-Jul-2019 15:02
Fluorene		U	0.00030	0.00010	mg/L	1	18-Jul-2019 15:02
Naphthalene	0.00062		0.00020	0.00010	mg/L	1	18-Jul-2019 15:02
Nitrobenzene		U	0.00024	0.00020	mg/L	1	18-Jul-2019 15:02
N-Nitrosodiphenylamine		U	0.00025	0.00020	mg/L	1	18-Jul-2019 15:02
Pentachlorophenol		U	0.00079	0.00020	mg/L	1	18-Jul-2019 15:02
Phenanthrene		U	0.00021	0.00010	mg/L	1	18-Jul-2019 15:02
Phenol		U	0.00035	0.00020	mg/L	1	18-Jul-2019 15:02
Pyrene		U	0.00019	0.00010	mg/L	1	18-Jul-2019 15:02
Surr: 2,4,6-Tribromophenol	79.7			34-129	%REC	1	18-Jul-2019 15:02
Surr: 2-Fluorobiphenyl	103			40-125	%REC	1	18-Jul-2019 15:02
Surr: 2-Fluorophenol	98.2			20-120	%REC	1	18-Jul-2019 15:02
Surr: 4-Terphenyl-d14	88.2			40-135	%REC	1	18-Jul-2019 15:02
Surr: Nitrobenzene-d5	99.2			41-120	%REC	1	18-Jul-2019 15:02
Surr: Phenol-d6	104			20-120	%REC	1	18-Jul-2019 15:02
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Jul-2019		Analyst: JC	
Arsenic		U	0.000400	0.00200	mg/L	1	19-Jul-2019 15:24

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB02-20190711  
 Collection Date: 11-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070609  
 Lab ID:HS19070609-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jul-2019 13:35
Benzene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 13:35
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 13:35
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jul-2019 13:35
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jul-2019 13:35
Toluene	U		0.00020	0.0010	mg/L	1	15-Jul-2019 13:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jul-2019 13:35
<i>Surr: 1,2-Dichloroethane-d4</i>	99.6			70-126	%REC	1	15-Jul-2019 13:35
<i>Surr: 4-Bromofluorobenzene</i>	98.6			81-113	%REC	1	15-Jul-2019 13:35
<i>Surr: Dibromofluoromethane</i>	102			77-123	%REC	1	15-Jul-2019 13:35
<i>Surr: Toluene-d8</i>	99.0			82-127	%REC	1	15-Jul-2019 13:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.



## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**Batch ID:** 143051      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070609-01	1	1000	1 (mL)	0.001
HS19070609-02	1	1000	1 (mL)	0.001
HS19070609-03	1	1000	1 (mL)	0.001
HS19070609-04	1	1000	1 (mL)	0.001
HS19070609-05	1	1000	1 (mL)	0.001
HS19070609-06	1	1000	1 (mL)	0.001

**Batch ID:** 143162      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070609-01	1	10	10 (mL)	1
HS19070609-02	1	10	10 (mL)	1
HS19070609-03	1	10	10 (mL)	1
HS19070609-04	1	10	10 (mL)	1
HS19070609-05	1	10	10 (mL)	1
HS19070609-06	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143051 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS19070609-01	WG-1620-MW40B-20190711	11 Jul 2019 12:00		16 Jul 2019 09:08	18 Jul 2019 15:40	100
HS19070609-01	WG-1620-MW40B-20190711	11 Jul 2019 12:00		16 Jul 2019 09:08	18 Jul 2019 15:59	1000
HS19070609-01	WG-1620-MW40B-20190711	11 Jul 2019 12:00		16 Jul 2019 09:08	18 Jul 2019 15:21	10
HS19070609-01	WG-1620-MW40B-20190711	11 Jul 2019 12:00		16 Jul 2019 09:08	17 Jul 2019 21:21	1
HS19070609-02	WG-1620-MW42B-20190711	11 Jul 2019 13:00		16 Jul 2019 09:08	17 Jul 2019 21:40	1
HS19070609-03	WG-1620-P11-20190711	11 Jul 2019 13:55		16 Jul 2019 09:08	18 Jul 2019 14:04	1
HS19070609-04	WG-1620-MW05-20190711	11 Jul 2019 14:50		16 Jul 2019 09:08	18 Jul 2019 14:23	1
HS19070609-05	WG-1620-MW64A-20190711	11 Jul 2019 15:50		16 Jul 2019 09:08	18 Jul 2019 14:42	1
HS19070609-06	WG-1620-FB02-20190711	11 Jul 2019 16:05		16 Jul 2019 09:08	18 Jul 2019 15:02	1
<b>Batch ID: 143162 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19070609-01	WG-1620-MW40B-20190711	11 Jul 2019 12:00		18 Jul 2019 12:00	18 Jul 2019 18:11	1
HS19070609-02	WG-1620-MW42B-20190711	11 Jul 2019 13:00		18 Jul 2019 12:00	18 Jul 2019 18:14	1
HS19070609-03	WG-1620-P11-20190711	11 Jul 2019 13:55		18 Jul 2019 12:00	18 Jul 2019 18:16	1
HS19070609-04	WG-1620-MW05-20190711	11 Jul 2019 14:50		18 Jul 2019 12:00	18 Jul 2019 18:18	1
HS19070609-05	WG-1620-MW64A-20190711	11 Jul 2019 15:50		18 Jul 2019 12:00	18 Jul 2019 18:20	1
HS19070609-06	WG-1620-FB02-20190711	11 Jul 2019 16:05		18 Jul 2019 12:00	19 Jul 2019 15:24	1
<b>Batch ID: R342484 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19070609-01	WG-1620-MW40B-20190711	11 Jul 2019 12:00			15 Jul 2019 14:00	1
HS19070609-02	WG-1620-MW42B-20190711	11 Jul 2019 13:00			15 Jul 2019 14:26	1
HS19070609-03	WG-1620-P11-20190711	11 Jul 2019 13:55			15 Jul 2019 14:51	1
HS19070609-04	WG-1620-MW05-20190711	11 Jul 2019 14:50			15 Jul 2019 15:17	1
HS19070609-05	WG-1620-MW64A-20190711	11 Jul 2019 15:50			15 Jul 2019 15:42	1
HS19070609-06	WG-1620-FB02-20190711	11 Jul 2019 16:05			15 Jul 2019 13:09	1
HS19070609-07	WG-1620-TB02-20190711	11 Jul 2019 00:00			15 Jul 2019 13:35	1

WorkOrder: HS19070609  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200

WorkOrder: HS19070609  
 InstrumentID: ICPMS04  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.00100	0.000579	0.000400	0.00200

WorkOrder: HS19070609  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00010	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000096	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00012	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00011	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.00010	0.00012	0.000019	0.00010
A	2-Methylnaphthalene	91-57-6	0.000050	0.000049	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000013	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00013	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000051	0.000027	0.00010
A	Acenaphthene	83-32-9	0.00010	0.00011	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.00010	0.00010	0.000015	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.00010	0.00011	0.000014	0.00010
A	Anthracene	120-12-7	0.000050	0.000055	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000068	0.000050	0.00010
A	Benz(a)anthracene	56-55-3	0.00010	0.000099	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.00010	0.000076	0.000020	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000053	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000086	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000066	0.000021	0.00010
A	Chrysene	218-01-9	0.00010	0.00011	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.00010	0.000094	0.000020	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000052	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.000099	0.000020	0.00020
A	Fluoranthene	206-44-0	0.00010	0.00012	0.000010	0.00010
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.00010	0.00012	0.000030	0.00010
A	Fluorene	86-73-7	0.000050	0.000052	0.000030	0.00010
A	Naphthalene	91-20-3	0.00010	0.00010	0.000020	0.00010
A	Naphthalene	91-20-3	0.000050	0.000061	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00013	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000085	0.000079	0.00020
A	Phenanthrene	85-01-8	0.00010	0.00011	0.000021	0.00010
A	Phenanthrene	85-01-8	0.000050	0.000055	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000094	0.000035	0.00020
A	Pyrene	129-00-0	0.00010	0.00011	0.000019	0.00010
A	Pyrene	129-00-0	0.000050	0.000061	0.000019	0.00010

WorkOrder: HS19070609  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19070609  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00064	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00060	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00061	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00077	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00058	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00062	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QC BATCH REPORT**

Batch ID: 143162 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A						
<b>MBLK</b>	Sample ID: <b>MBLK-143162</b>	Units: <b>mg/L</b>		Analysis Date: <b>18-Jul-2019 17:57</b>						
Client ID:		Run ID: <b>ICPMS05_342669</b>	SeqNo: <b>5172670</b>	PrepDate: <b>18-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-143162</b>	Units: <b>mg/L</b>		Analysis Date: <b>18-Jul-2019 17:59</b>						
Client ID:		Run ID: <b>ICPMS05_342669</b>	SeqNo: <b>5172671</b>	PrepDate: <b>18-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.05156	0.00200	0.05	0	103	80 - 120				
<b>MS</b>	Sample ID: <b>HS19070609-05MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>18-Jul-2019 19:07</b>						
Client ID: <b>WG-1620-MW64A-20190711</b>		Run ID: <b>ICPMS05_342669</b>	SeqNo: <b>5172775</b>	PrepDate: <b>18-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.06093	0.00200	0.05	0.009394	103	80 - 120				
<b>MSD</b>	Sample ID: <b>HS19070609-05MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>18-Jul-2019 19:10</b>						
Client ID: <b>WG-1620-MW64A-20190711</b>		Run ID: <b>ICPMS05_342669</b>	SeqNo: <b>5172776</b>	PrepDate: <b>18-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.06097	0.00200	0.05	0.009394	103	80 - 120	0.06093	0.0689	20	
<b>PDS</b>	Sample ID: <b>HS19070609-05PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>18-Jul-2019 19:12</b>						
Client ID: <b>WG-1620-MW64A-20190711</b>		Run ID: <b>ICPMS05_342669</b>	SeqNo: <b>5172783</b>	PrepDate: <b>18-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.1102	0.00200	0.1	0.009394	101	75 - 125				
<b>SD</b>	Sample ID: <b>HS19070609-05SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>18-Jul-2019 19:05</b>						
Client ID: <b>WG-1620-MW64A-20190711</b>		Run ID: <b>ICPMS05_342669</b>	SeqNo: <b>5172774</b>	PrepDate: <b>18-Jul-2019</b>	DF: <b>5</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual
Arsenic	0.009487	0.0100					0.009394	0	10	J

The following samples were analyzed in this batch: HS19070609-01 HS19070609-02 HS19070609-03 HS19070609-04  
 HS19070609-05 HS19070609-06



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QC BATCH REPORT**

Batch ID: 143051 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143051	Units: ug/L			Analysis Date: 17-Jul-2019 11:25					
Client ID:	Run ID: SV-7_342594	SeqNo: 5170297	PrepDate: 16-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.164</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.71</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.02</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>100</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.296</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.9</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.224</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.169</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QC BATCH REPORT**

Batch ID: 143051 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143051	Units: ug/L			Analysis Date: 17-Jul-2019 11:44					
Client ID:	Run ID: SV-7_342594	SeqNo: 5170298	PrepDate: 16-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.891	0.20	5	0	97.8	39 - 127				
2,4-Dimethylphenol	4.456	0.20	5	0	89.1	35 - 120				
2,4-Dinitrotoluene	5.665	0.20	5	0	113	50 - 122				
2,6-Dinitrotoluene	5.12	0.20	5	0	102	50 - 120				
2-Chloronaphthalene	4.568	0.20	5	0	91.4	50 - 120				
2-Methylnaphthalene	4.785	0.10	5	0	95.7	50 - 120				
4,6-Dinitro-2-methylphenol	4.433	0.20	5	0	88.7	25 - 121				
4-Nitrophenol	5.291	1.0	5	0	106	30 - 130				
Acenaphthene	4.568	0.10	5	0	91.4	45 - 120				
Acenaphthylene	4.495	0.10	5	0	89.9	47 - 120				
Anthracene	4.502	0.10	5	0	90.0	45 - 120				
Benz(a)anthracene	5.224	0.10	5	0	104	40 - 120				
Benzo(a)pyrene	4.184	0.10	5	0	83.7	45 - 120				
Bis(2-chloroethoxy)methane	4.62	0.20	5	0	92.4	45 - 120				
Bis(2-ethylhexyl)phthalate	5.383	0.20	5	0	108	40 - 139				
Chrysene	5.78	0.10	5	0	116	43 - 120				
Dibenzofuran	4.573	0.10	5	0	91.5	50 - 120				
Di-n-butyl phthalate	4.964	0.20	5	0	99.3	45 - 123				
Fluoranthene	5.235	0.10	5	0	105	45 - 125				
Fluorene	4.312	0.10	5	0	86.2	49 - 120				
Naphthalene	4.422	0.10	5	0	88.4	45 - 120				
Nitrobenzene	5.307	0.20	5	0	106	44 - 120				
N-Nitrosodiphenylamine	4.846	0.20	5	0	96.9	40 - 125				
Pentachlorophenol	3.998	0.20	5	0	80.0	19 - 121				
Phenanthrene	4.511	0.10	5	0	90.2	45 - 121				
Phenol	4.734	0.20	5	0	94.7	20 - 124				
Pyrene	4.84	0.10	5	0	96.8	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.277</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>106</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.715</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.401</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.682</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.128</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.86</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.2</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QC BATCH REPORT**

**Batch ID:** 143051 ( 0 )      **Instrument:** SV-7      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

LCSD		Sample ID: LCSD-143051			Units: ug/L		Analysis Date: 17-Jul-2019 12:04			
Client ID:		Run ID: SV-7_342594			SeqNo: 5170299		PrepDate: 16-Jul-2019		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.027	0.20	5	0	101	39 - 127	4.891	2.75	20	
2,4-Dimethylphenol	4.508	0.20	5	0	90.2	35 - 120	4.456	1.17	20	
2,4-Dinitrotoluene	6.009	0.20	5	0	120	50 - 122	5.665	5.91	20	
2,6-Dinitrotoluene	5.413	0.20	5	0	108	50 - 120	5.12	5.56	20	
2-Chloronaphthalene	4.685	0.20	5	0	93.7	50 - 120	4.568	2.52	20	
2-Methylnaphthalene	4.646	0.10	5	0	92.9	50 - 120	4.785	2.96	20	
4,6-Dinitro-2-methylphenol	4.386	0.20	5	0	87.7	25 - 121	4.433	1.06	30	
4-Nitrophenol	5.432	1.0	5	0	109	30 - 130	5.291	2.62	20	
Acenaphthene	4.65	0.10	5	0	93.0	45 - 120	4.568	1.77	20	
Acenaphthylene	4.597	0.10	5	0	91.9	47 - 120	4.495	2.24	20	
Anthracene	4.626	0.10	5	0	92.5	45 - 120	4.502	2.72	20	
Benz(a)anthracene	5.314	0.10	5	0	106	40 - 120	5.224	1.7	20	
Benzo(a)pyrene	4.37	0.10	5	0	87.4	45 - 120	4.184	4.35	20	
Bis(2-chloroethoxy)methane	4.471	0.20	5	0	89.4	45 - 120	4.62	3.27	20	
Bis(2-ethylhexyl)phthalate	5.015	0.20	5	0	100	40 - 139	5.383	7.09	20	
Chrysene	5.4	0.10	5	0	108	43 - 120	5.78	6.79	20	
Dibenzofuran	4.789	0.10	5	0	95.8	50 - 120	4.573	4.61	20	
Di-n-butyl phthalate	4.9	0.20	5	0	98.0	45 - 123	4.964	1.29	20	
Fluoranthene	5.005	0.10	5	0	100	45 - 125	5.235	4.5	20	
Fluorene	4.585	0.10	5	0	91.7	49 - 120	4.312	6.13	20	
Naphthalene	4.38	0.10	5	0	87.6	45 - 120	4.422	0.956	20	
Nitrobenzene	5.058	0.20	5	0	101	44 - 120	5.307	4.82	20	
N-Nitrosodiphenylamine	4.98	0.20	5	0	99.6	40 - 125	4.846	2.72	20	
Pentachlorophenol	3.749	0.20	5	0	75.0	19 - 121	3.998	6.45	20	
Phenanthrene	4.452	0.10	5	0	89.0	45 - 121	4.511	1.31	20	
Phenol	4.892	0.20	5	0	97.8	20 - 124	4.734	3.28	20	
Pyrene	4.992	0.10	5	0	99.8	40 - 130	4.84	3.09	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.252</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>34 - 129</i>	<i>5.277</i>	<i>0.469</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.793</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.9</i>	<i>40 - 125</i>	<i>4.715</i>	<i>1.64</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.505</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.1</i>	<i>20 - 120</i>	<i>4.401</i>	<i>2.33</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.494</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.9</i>	<i>40 - 135</i>	<i>4.682</i>	<i>4.08</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>5.029</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>41 - 120</i>	<i>5.128</i>	<i>1.95</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.973</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.5</i>	<i>20 - 120</i>	<i>4.86</i>	<i>2.3</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070609-01    HS19070609-02    HS19070609-03    HS19070609-04  
 HS19070609-05    HS19070609-06

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QC BATCH REPORT**

**Batch ID:** R342484 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190715</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jul-2019 12:44</b>			
Client ID:		Run ID: <b>VOA4_342484</b>			SeqNo: <b>5167898</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>48.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.0</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-190715</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jul-2019 11:53</b>			
Client ID:		Run ID: <b>VOA4_342484</b>			SeqNo: <b>5167897</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.98	1.0	20	0	84.9	70 - 124				
Benzene	18.25	1.0	20	0	91.2	74 - 120				
Chlorobenzene	19.01	1.0	20	0	95.0	76 - 113				
Ethylbenzene	19.38	1.0	20	0	96.9	77 - 117				
Methylene chloride	18.63	2.0	20	0	93.1	70 - 127				
Toluene	18.74	1.0	20	0	93.7	77 - 118				
Xylenes, Total	63.1	1.0	60	0	105	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.6</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>52.69</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.8</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>48.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QC BATCH REPORT**

**Batch ID:** R342484 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS19070645-01MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jul-2019 16:33</b>			
Client ID:		Run ID: <b>VOA4_342484</b>			SeqNo: <b>5167907</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.64	1.0	20	0	73.2	70 - 127				
Benzene	54.83	1.0	20	39.27	77.8	70 - 127				
Chlorobenzene	15.71	1.0	20	0	78.5	70 - 114				
Ethylbenzene	16.09	1.0	20	0	80.5	70 - 124				
Methylene chloride	15.72	2.0	20	0	78.6	70 - 128				
Toluene	16.11	1.0	20	0	80.5	70 - 123				
Xylenes, Total	50.86	1.0	60	0	84.8	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.5</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.19</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.3</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.7</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS19070645-01MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jul-2019 16:58</b>			
Client ID:		Run ID: <b>VOA4_342484</b>			SeqNo: <b>5167908</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.23	1.0	20	0	71.2	70 - 127	14.64	2.81	20	
Benzene	54.47	1.0	20	39.27	76.0	70 - 127	54.83	0.657	20	
Chlorobenzene	15.64	1.0	20	0	78.2	70 - 114	15.71	0.441	20	
Ethylbenzene	16.03	1.0	20	0	80.2	70 - 124	16.09	0.374	20	
Methylene chloride	15.2	2.0	20	0	76.0	70 - 128	15.72	3.37	20	
Toluene	16.04	1.0	20	0	80.2	70 - 123	16.11	0.413	20	
Xylenes, Total	50.75	1.0	60	0	84.6	70 - 130	50.86	0.216	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>70 - 126</i>	<i>48.75</i>	<i>0.579</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.34</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>	<i>50.19</i>	<i>0.298</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>77 - 123</i>	<i>49.67</i>	<i>0.381</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 127</i>	<i>49.7</i>	<i>2.54</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070609-01    HS19070609-02    HS19070609-03    HS19070609-04  
 HS19070609-05    HS19070609-06    HS19070609-07

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070609

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kansas	E-10352 2018-2019	31-Jul-2019
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070609

**SAMPLE TRACKING**

---

Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS19070609-01	WG-1620-MW40B-20190711	Login	12/07/2019 15:13:02	NDR	MET044
HS19070609-01	WG-1620-MW40B-20190711	Login	12/07/2019 15:13:02	NDR	EXT133
HS19070609-01	WG-1620-MW40B-20190711	Login	12/07/2019 15:13:02	NDR	VOA244



Sample Receipt Checklist

Client Name: PBW
Work Order: HS19070609

Date/Time Received: 12-Jul-2019 12:15
Received by: JRM

Checklist completed by: Nilesh D. Ranchod
eSignature
Date: 12-Jul-2019

Reviewed by: Dane J. Wacasey
eSignature
Date: 15-Jul-2019

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:196133

Temperature(s)/Thermometer(s): 4.5C UC/C IR # 25
Cooler(s)/Kit(s): 44990
Date/Time sample(s) sent to storage: 07/12/2019 16:00
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336  
Everett, WA  
+1 425 356 2600

Fort Collins, CO  
+1 970 490 1511  
Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page 1 of 1

COC ID: 196133

HS19070609

Golder Associates Inc.  
Houston TX-Wood Preserving Works



WV

Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/F	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TB0_-201801			Water	1	2		X									
2	WG-1620-MW40B-20190711	7-11-19	1200	W		6	X		X	X							
3	WG-1620-MW42B-20190711		1300	W		6	X		X	X							
4	WG-1620-P11-20190711		1355	W		6	X		X	X							
5	WG-1620-MW05-20190711		1450	W		6	X		X	X							
6	WG-1620-MW64A-20190711		1550	W		6	X		X	X							
7	WG-1620-FB02-20190711		1605	W		6	X		X	X							
8	WG-1620-TB02-20190711			W		2	X		X	X							
9																	
10																	

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>	Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Dctys <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour	Results Due Date:
Relinquished by <b>John Br</b>	Date: <b>7-12-19</b>	Time: <b>12:15</b>	Received by:
Relinquished by <b>John Br</b>	Date: <b>7/12/19</b>	Time: <b>12:15</b>	Received by (Laboratory): <b>J. MAURIN</b>
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):
Notes: <b>UPRR Houston MW/PW</b>		Cooler ID: <b>44990</b>	Cooler Temp: <b>4.5</b>
QC Package: (Check One Box Below)		Level II Std QC <input checked="" type="checkbox"/> TERP Checklist	
		Level III Std QC/Raw Data <input type="checkbox"/> TERP Level IV	
		Level IV SW846/CLP <input type="checkbox"/>	
		Other: <b>CFO-2</b>	

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

July 23, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19070735**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 6 sample(s) on Jul 15, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/23/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070735			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143051,143254,R342562			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			2
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			3
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/23/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070735			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143051,143254,R342562			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 07/23/2019
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS19070735
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 143051,143254,R342562
ER# <sup>5</sup>	Description	
1	Semivolatile Organics Method Sw8270, sample WG-1620-TW41B-20190712, the surrogate recoveries could not be determined due to dilution below the calibration range.	
2	Batch 143051, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch	
3	Batch R342562, Volatile Organics Method SW8260, sample HS19070681-04, MSD was performed on unrelated sample.	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.</p> <p>O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);          NA = Not Applicable;          NR = Not Reviewed;          R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070735

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19070735-01	WG-1620-MW09-20190712	Water		12-Jul-2019 07:35	15-Jul-2019 16:30	<input type="checkbox"/>
HS19070735-02	WG-1620-MW03-20190712	Water		12-Jul-2019 08:35	15-Jul-2019 16:30	<input type="checkbox"/>
HS19070735-03	WG-1620-MW04-20190712	Water		12-Jul-2019 09:25	15-Jul-2019 16:30	<input type="checkbox"/>
HS19070735-04	WG-1620-TW41B-20190712	Water		12-Jul-2019 10:40	15-Jul-2019 16:30	<input type="checkbox"/>
HS19070735-05	WG-1620-FB03-20190712	Water		12-Jul-2019 11:00	15-Jul-2019 16:30	<input type="checkbox"/>
HS19070735-06	WG-1620-TB03-20190712	Water	CG 062119 -79	12-Jul-2019 00:00	15-Jul-2019 16:30	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW09-20190712  
 Collection Date: 12-Jul-2019 07:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jul-2019 19:17
Benzene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 19:17
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 19:17
Ethylbenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 19:17
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jul-2019 19:17
Toluene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 19:17
Xylenes, Total	U		0.00030	0.0010	mg/L	1	16-Jul-2019 19:17
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:17</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:17</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:17</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:17</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW09-20190712  
 Collection Date: 12-Jul-2019 07:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 16:18
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 16:18
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 16:18
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 16:18
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 16:18
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	18-Jul-2019 16:18
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 16:18
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 16:18
Acenaphthene	U		0.000027	0.00010	mg/L	1	18-Jul-2019 16:18
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 16:18
<b>Anthracene</b>	<b>0.000061</b>	<b>J</b>	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:18</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 16:18
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 16:18
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 16:18
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 16:18
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 16:18
Dibenzofuran	U		0.000020	0.00010	mg/L	1	18-Jul-2019 16:18
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 16:18
Fluoranthene	U		0.000010	0.00010	mg/L	1	18-Jul-2019 16:18
Fluorene	U		0.000030	0.00010	mg/L	1	18-Jul-2019 16:18
<b>Naphthalene</b>	<b>0.00015</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:18</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 16:18
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 16:18
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 16:18
Phenanthrene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 16:18
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 16:18
Pyrene	U		0.000019	0.00010	mg/L	1	18-Jul-2019 16:18
<i>Surr: 2,4,6-Tribromophenol</i>	<i>91.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>68.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>69.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:18</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>83.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:18</i>
<i>Surr: Nitrobenzene-d5</i>	<i>74.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:18</i>
<i>Surr: Phenol-d6</i>	<i>79.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:18</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000901</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 22:25</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW03-20190712  
 Collection Date: 12-Jul-2019 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jul-2019 19:42
Benzene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 19:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 19:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 19:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jul-2019 19:42
Toluene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 19:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	16-Jul-2019 19:42
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:42</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:42</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:42</i>
<i>Surr: Toluene-d8</i>	<i>98.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 19:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW03-20190712  
 Collection Date: 12-Jul-2019 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 16:37
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 16:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 16:37
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 16:37
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 16:37
<b>2-Methylnaphthalene</b>	<b>0.000035</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:37</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 16:37
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 16:37
Acenaphthene	U		0.000027	0.00010	mg/L	1	18-Jul-2019 16:37
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 16:37
<b>Anthracene</b>	<b>0.000055</b>	<b>J</b>	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:37</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 16:37
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 16:37
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 16:37
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 16:37
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 16:37
Dibenzofuran	U		0.000020	0.00010	mg/L	1	18-Jul-2019 16:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 16:37
Fluoranthene	U		0.000010	0.00010	mg/L	1	18-Jul-2019 16:37
Fluorene	U		0.000030	0.00010	mg/L	1	18-Jul-2019 16:37
<b>Naphthalene</b>	<b>0.00029</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:37</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 16:37
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 16:37
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 16:37
Phenanthrene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 16:37
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 16:37
<b>Pyrene</b>	<b>0.000020</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:37</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>62.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:37</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:37</i>
<i>Surr: Nitrobenzene-d5</i>	<i>67.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:37</i>
<i>Surr: Phenol-d6</i>	<i>74.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:37</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000582</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 23:04</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW04-20190712  
 Collection Date: 12-Jul-2019 09:25

**ANALYTICAL REPORT**

WorkOrder:HS19070735  
 Lab ID:HS19070735-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jul-2019 20:07
Benzene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 20:07
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 20:07
Ethylbenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 20:07
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jul-2019 20:07
Toluene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 20:07
Xylenes, Total	U		0.00030	0.0010	mg/L	1	16-Jul-2019 20:07
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:07</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:07</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:07</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:07</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW04-20190712  
 Collection Date: 12-Jul-2019 09:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 16:57
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 16:57
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 16:57
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 16:57
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 16:57
<b>2-Methylnaphthalene</b>	<b>0.000059</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:57</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 16:57
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 16:57
<b>Acenaphthene</b>	<b>0.000052</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:57</b>
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 16:57
Anthracene	U		0.000014	0.00010	mg/L	1	18-Jul-2019 16:57
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 16:57
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 16:57
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 16:57
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 16:57
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 16:57
<b>Dibenzofuran</b>	<b>0.000058</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:57</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 16:57
Fluoranthene	U		0.000010	0.00010	mg/L	1	18-Jul-2019 16:57
<b>Fluorene</b>	<b>0.000053</b>	<b>J</b>	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:57</b>
<b>Naphthalene</b>	<b>0.000053</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:57</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 16:57
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 16:57
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 16:57
Phenanthrene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 16:57
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 16:57
<b>Pyrene</b>	<b>0.000016</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 16:57</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>83.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:57</i>
<i>Surr: Nitrobenzene-d5</i>	<i>80.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:57</i>
<i>Surr: Phenol-d6</i>	<i>76.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 16:57</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0127</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 23:06</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TW41B-20190712  
 Collection Date: 12-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jul-2019 20:32
<b>Benzene</b>	<b>0.0013</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 20:32
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 20:32
<b>Ethylbenzene</b>	<b>0.0011</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 20:32
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jul-2019 20:32
<b>Toluene</b>	<b>0.0011</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 20:32
<b>Xylenes, Total</b>	<b>0.019</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 20:32
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:32</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>107</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:32</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:32</i>
<i>Surr: Toluene-d8</i>	<i>96.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 20:32</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TW41B-20190712  
 Collection Date: 12-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 17:16
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 17:16
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 17:16
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 17:16
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 17:16
<b>2-Methylnaphthalene</b>	<b>0.067</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	19-Jul-2019 17:34
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 17:16
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 17:16
<b>Acenaphthene</b>	<b>0.15</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	19-Jul-2019 17:54
<b>Acenaphthylene</b>	<b>0.0014</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 17:16
<b>Anthracene</b>	<b>0.0055</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 17:16
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 17:16
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 17:16
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 17:16
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 17:16
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 17:16
<b>Dibenzofuran</b>	<b>0.074</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	19-Jul-2019 17:34
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 17:16
<b>Fluoranthene</b>	<b>0.0034</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 17:16
<b>Fluorene</b>	<b>0.085</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	19-Jul-2019 17:34
<b>Naphthalene</b>	<b>0.69</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	19-Jul-2019 17:54
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 17:16
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 17:16
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 17:16
<b>Phenanthrene</b>	<b>0.027</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	19-Jul-2019 17:34
Phenol	U		0.000035	0.00020	mg/L	1	18-Jul-2019 17:16
<b>Pyrene</b>	<b>0.0014</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Jul-2019 17:16
Surr: 2,4,6-Tribromophenol	118			34-129	%REC	10	19-Jul-2019 17:34
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	19-Jul-2019 17:54
Surr: 2,4,6-Tribromophenol	69.5			34-129	%REC	1	18-Jul-2019 17:16
Surr: 2-Fluorobiphenyl	54.2			40-125	%REC	1	18-Jul-2019 17:16
Surr: 2-Fluorobiphenyl	81.3			40-125	%REC	10	19-Jul-2019 17:34
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	19-Jul-2019 17:54
Surr: 2-Fluorophenol	60.1			20-120	%REC	10	19-Jul-2019 17:34
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	19-Jul-2019 17:54
Surr: 2-Fluorophenol	57.7			20-120	%REC	1	18-Jul-2019 17:16
Surr: 4-Terphenyl-d14	84.7			40-135	%REC	1	18-Jul-2019 17:16
Surr: 4-Terphenyl-d14	83.2			40-135	%REC	10	19-Jul-2019 17:34
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	19-Jul-2019 17:54

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TW41B-20190712  
 Collection Date: 12-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	67.6			41-120	%REC	10	19-Jul-2019 17:34
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	19-Jul-2019 17:54
Surr: Nitrobenzene-d5	65.5			41-120	%REC	1	18-Jul-2019 17:16
Surr: Phenol-d6	63.7			20-120	%REC	1	18-Jul-2019 17:16
Surr: Phenol-d6	70.5			20-120	%REC	10	19-Jul-2019 17:34
Surr: Phenol-d6	0	JS		20-120	%REC	100	19-Jul-2019 17:54
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jul-2019		Analyst: JHD	
Arsenic	0.113		0.000400	0.00200	mg/L	1	22-Jul-2019 23:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB03-20190712  
 Collection Date: 12-Jul-2019 11:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jul-2019 18:26
<b>Benzene</b>	<b>0.00052</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 18:26
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 18:26
<b>Ethylbenzene</b>	<b>0.00032</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 18:26
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jul-2019 18:26
<b>Toluene</b>	<b>0.00037</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	16-Jul-2019 18:26
Xylenes, Total	U		0.00030	0.0010	mg/L	1	16-Jul-2019 18:26
Surr: 1,2-Dichloroethane-d4	97.5			70-126	%REC	1	16-Jul-2019 18:26
Surr: 4-Bromofluorobenzene	97.1			81-113	%REC	1	16-Jul-2019 18:26
Surr: Dibromofluoromethane	100			77-123	%REC	1	16-Jul-2019 18:26
Surr: Toluene-d8	101			82-127	%REC	1	16-Jul-2019 18:26

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB03-20190712  
 Collection Date: 12-Jul-2019 11:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	18-Jul-2019 17:35
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	18-Jul-2019 17:35
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	18-Jul-2019 17:35
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	18-Jul-2019 17:35
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	18-Jul-2019 17:35
<b>2-Methylnaphthalene</b>	<b>0.000052</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 17:35</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	18-Jul-2019 17:35
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	18-Jul-2019 17:35
Acenaphthene	U		0.000027	0.00010	mg/L	1	18-Jul-2019 17:35
Acenaphthylene	U		0.000015	0.00010	mg/L	1	18-Jul-2019 17:35
Anthracene	U		0.000014	0.00010	mg/L	1	18-Jul-2019 17:35
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	18-Jul-2019 17:35
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	18-Jul-2019 17:35
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	18-Jul-2019 17:35
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	18-Jul-2019 17:35
Chrysene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 17:35
<b>Dibenzofuran</b>	<b>0.000026</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 17:35</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	18-Jul-2019 17:35
Fluoranthene	U		0.000010	0.00010	mg/L	1	18-Jul-2019 17:35
Fluorene	U		0.000030	0.00010	mg/L	1	18-Jul-2019 17:35
<b>Naphthalene</b>	<b>0.00028</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 17:35</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	18-Jul-2019 17:35
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	18-Jul-2019 17:35
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	18-Jul-2019 17:35
Phenanthrene	U		0.000021	0.00010	mg/L	1	18-Jul-2019 17:35
<b>Phenol</b>	<b>0.000052</b>	<b>J</b>	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>18-Jul-2019 17:35</b>
Pyrene	U		0.000019	0.00010	mg/L	1	18-Jul-2019 17:35
<i>Surr: 2,4,6-Tribromophenol</i>	<i>83.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 17:35</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>98.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 17:35</i>
<i>Surr: 2-Fluorophenol</i>	<i>77.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 17:35</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>95.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 17:35</i>
<i>Surr: Nitrobenzene-d5</i>	<i>100</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 17:35</i>
<i>Surr: Phenol-d6</i>	<i>91.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 17:35</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	22-Jul-2019 23:11

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB03-20190712  
 Collection Date: 12-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070735  
 Lab ID:HS19070735-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jul-2019 18:52
Benzene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 18:52
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 18:52
Ethylbenzene	U		0.00030	0.0010	mg/L	1	16-Jul-2019 18:52
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jul-2019 18:52
Toluene	U		0.00020	0.0010	mg/L	1	16-Jul-2019 18:52
Xylenes, Total	U		0.00030	0.0010	mg/L	1	16-Jul-2019 18:52
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 18:52</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 18:52</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 18:52</i>
<i>Surr: Toluene-d8</i>	<i>96.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>16-Jul-2019 18:52</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**Batch ID:** 143051      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070735-01	1	1000	1 (mL)	0.001
HS19070735-02	1	1000	1 (mL)	0.001
HS19070735-03	1	1000	1 (mL)	0.001
HS19070735-04	1	1000	1 (mL)	0.001
HS19070735-05	1	1000	1 (mL)	0.001

**Batch ID:** 143254      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070735-01	1	10	10 (mL)	1
HS19070735-02	1	10	10 (mL)	1
HS19070735-03	1	10	10 (mL)	1
HS19070735-04	1	10	10 (mL)	1
HS19070735-05	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID:</b> 143051 ( 0 )		<b>Test Name :</b> LOW-LEVEL SEMIVOLATILES BY 8270D			<b>Matrix:</b> Water	
HS19070735-01	WG-1620-MW09-20190712	12 Jul 2019 07:35		16 Jul 2019 09:08	18 Jul 2019 16:18	1
HS19070735-02	WG-1620-MW03-20190712	12 Jul 2019 08:35		16 Jul 2019 09:08	18 Jul 2019 16:37	1
HS19070735-03	WG-1620-MW04-20190712	12 Jul 2019 09:25		16 Jul 2019 09:08	18 Jul 2019 16:57	1
HS19070735-04	WG-1620-TW41B-20190712	12 Jul 2019 10:40		16 Jul 2019 09:08	19 Jul 2019 17:54	100
HS19070735-04	WG-1620-TW41B-20190712	12 Jul 2019 10:40		16 Jul 2019 09:08	19 Jul 2019 17:34	10
HS19070735-04	WG-1620-TW41B-20190712	12 Jul 2019 10:40		16 Jul 2019 09:08	18 Jul 2019 17:16	1
HS19070735-05	WG-1620-FB03-20190712	12 Jul 2019 11:00		16 Jul 2019 09:08	18 Jul 2019 17:35	1
<b>Batch ID:</b> 143254 ( 0 )		<b>Test Name :</b> ICP-MS METALS BY SW6020A			<b>Matrix:</b> Water	
HS19070735-01	WG-1620-MW09-20190712	12 Jul 2019 07:35		22 Jul 2019 12:00	22 Jul 2019 22:25	1
HS19070735-02	WG-1620-MW03-20190712	12 Jul 2019 08:35		22 Jul 2019 12:00	22 Jul 2019 23:04	1
HS19070735-03	WG-1620-MW04-20190712	12 Jul 2019 09:25		22 Jul 2019 12:00	22 Jul 2019 23:06	1
HS19070735-04	WG-1620-TW41B-20190712	12 Jul 2019 10:40		22 Jul 2019 12:00	22 Jul 2019 23:09	1
HS19070735-05	WG-1620-FB03-20190712	12 Jul 2019 11:00		22 Jul 2019 12:00	22 Jul 2019 23:11	1
<b>Batch ID:</b> R342562 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Water	
HS19070735-01	WG-1620-MW09-20190712	12 Jul 2019 07:35			16 Jul 2019 19:17	1
HS19070735-02	WG-1620-MW03-20190712	12 Jul 2019 08:35			16 Jul 2019 19:42	1
HS19070735-03	WG-1620-MW04-20190712	12 Jul 2019 09:25			16 Jul 2019 20:07	1
HS19070735-04	WG-1620-TW41B-20190712	12 Jul 2019 10:40			16 Jul 2019 20:32	1
HS19070735-05	WG-1620-FB03-20190712	12 Jul 2019 11:00			16 Jul 2019 18:26	1
HS19070735-06	WG-1620-TB03-20190712	12 Jul 2019 00:00			16 Jul 2019 18:52	1

WorkOrder: HS19070735  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200



WorkOrder: HS19070735  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

Matrix: Aqueous

Units: mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00010	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000096	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00012	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00011	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.00010	0.00012	0.000019	0.00010
A	2-Methylnaphthalene	91-57-6	0.000050	0.000049	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000013	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00013	0.000047	0.0010
A	Acenaphthene	83-32-9	0.00010	0.00011	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Acenaphthylene	208-96-8	0.00010	0.00010	0.000015	0.00010
A	Anthracene	120-12-7	0.00010	0.00011	0.000014	0.00010
A	Anthracene	120-12-7	0.000050	0.000055	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000068	0.000050	0.00010
A	Benz(a)anthracene	56-55-3	0.00010	0.000099	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.00010	0.000076	0.000020	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000053	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000086	0.000037	0.00020
A	Chrysene	218-01-9	0.00010	0.00011	0.000021	0.00010
A	Chrysene	218-01-9	0.000050	0.000066	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000052	0.000020	0.00010
A	Dibenzofuran	132-64-9	0.00010	0.000094	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.000099	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluoranthene	206-44-0	0.00010	0.00012	0.000010	0.00010
A	Fluorene	86-73-7	0.00010	0.00012	0.000030	0.00010
A	Fluorene	86-73-7	0.000050	0.000052	0.000030	0.00010
A	Naphthalene	91-20-3	0.00010	0.00010	0.000020	0.00010
A	Naphthalene	91-20-3	0.000050	0.000061	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00013	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000085	0.000079	0.00020
A	Phenanthrene	85-01-8	0.00010	0.00011	0.000021	0.00010
A	Phenanthrene	85-01-8	0.000050	0.000055	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000094	0.000035	0.00020
A	Pyrene	129-00-0	0.00010	0.00011	0.000019	0.00010
A	Pyrene	129-00-0	0.000050	0.000061	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020

WorkOrder: HS19070735

InstrumentID: SV-7

Test Code: 8270\_LOW\_W

Test Number: SW8270

Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19070735  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00064	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00060	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00061	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00077	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00058	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00062	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QC BATCH REPORT**

Batch ID: 143254 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-143254</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jul-2019 22:21</b>					
Client ID:		Run ID: <b>ICPMS05_342799</b>	SeqNo: <b>5176129</b>	PrepDate: <b>22-Jul-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-143254</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jul-2019 22:23</b>					
Client ID:		Run ID: <b>ICPMS05_342799</b>	SeqNo: <b>5176130</b>	PrepDate: <b>22-Jul-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05183	0.00200	0.05	0	104	80 - 120			
<b>MS</b>	Sample ID: <b>HS19070735-01MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jul-2019 22:30</b>					
Client ID: <b>WG-1620-MW09-20190712</b>		Run ID: <b>ICPMS05_342799</b>	SeqNo: <b>5176133</b>	PrepDate: <b>22-Jul-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05366	0.00200	0.05	0.000901	106	80 - 120			
<b>MSD</b>	Sample ID: <b>HS19070735-01MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jul-2019 22:32</b>					
Client ID: <b>WG-1620-MW09-20190712</b>		Run ID: <b>ICPMS05_342799</b>	SeqNo: <b>5176134</b>	PrepDate: <b>22-Jul-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.0522	0.00200	0.05	0.000901	103	80 - 120	0.05366	2.76	20
<b>PDS</b>	Sample ID: <b>HS19070735-01PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jul-2019 22:35</b>					
Client ID: <b>WG-1620-MW09-20190712</b>		Run ID: <b>ICPMS05_342799</b>	SeqNo: <b>5176135</b>	PrepDate: <b>22-Jul-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.0984	0.00200	0.1	0.000901	97.5	75 - 125			
<b>SD</b>	Sample ID: <b>HS19070735-01SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jul-2019 22:28</b>					
Client ID: <b>WG-1620-MW09-20190712</b>		Run ID: <b>ICPMS05_342799</b>	SeqNo: <b>5176132</b>	PrepDate: <b>22-Jul-2019</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	U	0.0100					0.000901	0	10
The following samples were analyzed in this batch:									
	HS19070735-01	HS19070735-02	HS19070735-03	HS19070735-04	HS19070735-05				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QC BATCH REPORT**

Batch ID: 143051 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143051	Units: ug/L			Analysis Date: 17-Jul-2019 11:25					
Client ID:	Run ID: SV-7_342594	SeqNo: 5170297	PrepDate: 16-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.164</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.71</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.02</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>100</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.296</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.9</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.224</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.169</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QC BATCH REPORT**

Batch ID: 143051 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143051	Units: ug/L			Analysis Date: 17-Jul-2019 11:44					
Client ID:	Run ID: SV-7_342594	SeqNo: 5170298		PrepDate: 16-Jul-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.891	0.20	5	0	97.8	39 - 127				
2,4-Dimethylphenol	4.456	0.20	5	0	89.1	35 - 120				
2,4-Dinitrotoluene	5.665	0.20	5	0	113	50 - 122				
2,6-Dinitrotoluene	5.12	0.20	5	0	102	50 - 120				
2-Chloronaphthalene	4.568	0.20	5	0	91.4	50 - 120				
2-Methylnaphthalene	4.785	0.10	5	0	95.7	50 - 120				
4,6-Dinitro-2-methylphenol	4.433	0.20	5	0	88.7	25 - 121				
4-Nitrophenol	5.291	1.0	5	0	106	30 - 130				
Acenaphthene	4.568	0.10	5	0	91.4	45 - 120				
Acenaphthylene	4.495	0.10	5	0	89.9	47 - 120				
Anthracene	4.502	0.10	5	0	90.0	45 - 120				
Benz(a)anthracene	5.224	0.10	5	0	104	40 - 120				
Benzo(a)pyrene	4.184	0.10	5	0	83.7	45 - 120				
Bis(2-chloroethoxy)methane	4.62	0.20	5	0	92.4	45 - 120				
Bis(2-ethylhexyl)phthalate	5.383	0.20	5	0	108	40 - 139				
Chrysene	5.78	0.10	5	0	116	43 - 120				
Dibenzofuran	4.573	0.10	5	0	91.5	50 - 120				
Di-n-butyl phthalate	4.964	0.20	5	0	99.3	45 - 123				
Fluoranthene	5.235	0.10	5	0	105	45 - 125				
Fluorene	4.312	0.10	5	0	86.2	49 - 120				
Naphthalene	4.422	0.10	5	0	88.4	45 - 120				
Nitrobenzene	5.307	0.20	5	0	106	44 - 120				
N-Nitrosodiphenylamine	4.846	0.20	5	0	96.9	40 - 125				
Pentachlorophenol	3.998	0.20	5	0	80.0	19 - 121				
Phenanthrene	4.511	0.10	5	0	90.2	45 - 121				
Phenol	4.734	0.20	5	0	94.7	20 - 124				
Pyrene	4.84	0.10	5	0	96.8	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.277</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>106</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.715</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.401</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.682</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.128</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.86</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.2</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QC BATCH REPORT**

**Batch ID:** 143051 ( 0 )      **Instrument:** SV-7      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

LCSD		Sample ID: LCSD-143051			Units: ug/L		Analysis Date: 17-Jul-2019 12:04			
Client ID:		Run ID: SV-7_342594			SeqNo: 5170299		PrepDate: 16-Jul-2019		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.027	0.20	5	0	101	39 - 127	4.891	2.75	20	
2,4-Dimethylphenol	4.508	0.20	5	0	90.2	35 - 120	4.456	1.17	20	
2,4-Dinitrotoluene	6.009	0.20	5	0	120	50 - 122	5.665	5.91	20	
2,6-Dinitrotoluene	5.413	0.20	5	0	108	50 - 120	5.12	5.56	20	
2-Chloronaphthalene	4.685	0.20	5	0	93.7	50 - 120	4.568	2.52	20	
2-Methylnaphthalene	4.646	0.10	5	0	92.9	50 - 120	4.785	2.96	20	
4,6-Dinitro-2-methylphenol	4.386	0.20	5	0	87.7	25 - 121	4.433	1.06	30	
4-Nitrophenol	5.432	1.0	5	0	109	30 - 130	5.291	2.62	20	
Acenaphthene	4.65	0.10	5	0	93.0	45 - 120	4.568	1.77	20	
Acenaphthylene	4.597	0.10	5	0	91.9	47 - 120	4.495	2.24	20	
Anthracene	4.626	0.10	5	0	92.5	45 - 120	4.502	2.72	20	
Benz(a)anthracene	5.314	0.10	5	0	106	40 - 120	5.224	1.7	20	
Benzo(a)pyrene	4.37	0.10	5	0	87.4	45 - 120	4.184	4.35	20	
Bis(2-chloroethoxy)methane	4.471	0.20	5	0	89.4	45 - 120	4.62	3.27	20	
Bis(2-ethylhexyl)phthalate	5.015	0.20	5	0	100	40 - 139	5.383	7.09	20	
Chrysene	5.4	0.10	5	0	108	43 - 120	5.78	6.79	20	
Dibenzofuran	4.789	0.10	5	0	95.8	50 - 120	4.573	4.61	20	
Di-n-butyl phthalate	4.9	0.20	5	0	98.0	45 - 123	4.964	1.29	20	
Fluoranthene	5.005	0.10	5	0	100	45 - 125	5.235	4.5	20	
Fluorene	4.585	0.10	5	0	91.7	49 - 120	4.312	6.13	20	
Naphthalene	4.38	0.10	5	0	87.6	45 - 120	4.422	0.956	20	
Nitrobenzene	5.058	0.20	5	0	101	44 - 120	5.307	4.82	20	
N-Nitrosodiphenylamine	4.98	0.20	5	0	99.6	40 - 125	4.846	2.72	20	
Pentachlorophenol	3.749	0.20	5	0	75.0	19 - 121	3.998	6.45	20	
Phenanthrene	4.452	0.10	5	0	89.0	45 - 121	4.511	1.31	20	
Phenol	4.892	0.20	5	0	97.8	20 - 124	4.734	3.28	20	
Pyrene	4.992	0.10	5	0	99.8	40 - 130	4.84	3.09	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.252</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>34 - 129</i>	<i>5.277</i>	<i>0.469</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.793</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.9</i>	<i>40 - 125</i>	<i>4.715</i>	<i>1.64</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.505</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.1</i>	<i>20 - 120</i>	<i>4.401</i>	<i>2.33</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.494</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.9</i>	<i>40 - 135</i>	<i>4.682</i>	<i>4.08</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>5.029</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>41 - 120</i>	<i>5.128</i>	<i>1.95</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.973</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.5</i>	<i>20 - 120</i>	<i>4.86</i>	<i>2.3</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070735-01    HS19070735-02    HS19070735-03    HS19070735-04  
 HS19070735-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QC BATCH REPORT**

**Batch ID:** R342562 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190716</b>			Units: <b>ug/L</b>		Analysis Date: <b>16-Jul-2019 14:10</b>			
Client ID:		Run ID: <b>VOA4_342562</b>			SeqNo: <b>5169824</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.6</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.44</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-190716</b>			Units: <b>ug/L</b>		Analysis Date: <b>16-Jul-2019 13:18</b>			
Client ID:		Run ID: <b>VOA4_342562</b>			SeqNo: <b>5169823</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.58	1.0	20	0	87.9	70 - 124				
Benzene	18.78	1.0	20	0	93.9	74 - 120				
Chlorobenzene	19.65	1.0	20	0	98.3	76 - 113				
Ethylbenzene	19.35	1.0	20	0	96.7	77 - 117				
Methylene chloride	19.06	2.0	20	0	95.3	70 - 127				
Toluene	19.08	1.0	20	0	95.4	77 - 118				
Xylenes, Total	64.06	1.0	60	0	107	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.1</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>48.58</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>81 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QC BATCH REPORT**

Batch ID: R342562 ( 0 )		Instrument: VOA4		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>	Sample ID: <b>HS19070681-04MS</b>	Units: <b>ug/L</b>			Analysis Date: <b>16-Jul-2019 16:45</b>					
Client ID:	Run ID: <b>VOA4_342562</b>	SeqNo: <b>5169830</b>		PrepDate:		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.82	1.0	20	0	74.1	70 - 127				
Benzene	124.2	1.0	20	109.8	71.6	70 - 127				O
Chlorobenzene	16.67	1.0	20	0	83.4	70 - 114				
Ethylbenzene	37.48	1.0	20	17.94	97.7	70 - 124				
Methylene chloride	17	2.0	20	0	85.0	70 - 128				
Toluene	21.24	1.0	20	2.685	92.8	70 - 123				
Xylenes, Total	62.26	1.0	60	4.367	96.5	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.06</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.1</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>52.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.33</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>				

<b>MSD</b>	Sample ID: <b>HS19070681-04MSD</b>	Units: <b>ug/L</b>			Analysis Date: <b>16-Jul-2019 17:11</b>					
Client ID:	Run ID: <b>VOA4_342562</b>	SeqNo: <b>5169831</b>		PrepDate:		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.45	1.0	20	0	72.2	70 - 127	14.82	2.55	20	
Benzene	115.9	1.0	20	109.8	30.4	70 - 127	124.2	6.87	20	SO
Chlorobenzene	15.87	1.0	20	0	79.4	70 - 114	16.67	4.94	20	
Ethylbenzene	33.81	1.0	20	17.94	79.3	70 - 124	37.48	10.3	20	
Methylene chloride	14.72	2.0	20	0	73.6	70 - 128	17	14.3	20	
Toluene	19.3	1.0	20	2.685	83.1	70 - 123	21.24	9.56	20	
Xylenes, Total	55.27	1.0	60	4.367	84.8	70 - 130	62.26	11.9	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.42</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>70 - 126</i>	<i>47.06</i>	<i>0.76</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.53</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>81 - 113</i>	<i>52.4</i>	<i>5.63</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>77 - 123</i>	<i>48.33</i>	<i>0.983</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>	<i>50.02</i>	<i>0.425</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070735-01 HS19070735-02 HS19070735-03 HS19070735-04  
 HS19070735-05 HS19070735-06

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070735

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kansas	E-10352 2018-2019	31-Jul-2019
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070735

**SAMPLE TRACKING**

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Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS19070735-01	WG-1620-MW09-20190712	Login	15/07/2019 18:21:52	JRM	MET044
HS19070735-01	WG-1620-MW09-20190712	Login	15/07/2019 18:21:52	JRM	EXT108
HS19070735-01	WG-1620-MW09-20190712	Login	15/07/2019 18:21:52	JRM	VOA074
HS19070735-02	WG-1620-MW03-20190712	Login	15/07/2019 18:21:52	JRM	MET044
HS19070735-02	WG-1620-MW03-20190712	Login	15/07/2019 18:21:52	JRM	EXT108
HS19070735-02	WG-1620-MW03-20190712	Login	15/07/2019 18:21:52	JRM	VOA074
HS19070735-03	WG-1620-MW04-20190712	Login	15/07/2019 18:21:52	JRM	MET044
HS19070735-03	WG-1620-MW04-20190712	Login	15/07/2019 18:21:52	JRM	EXT108
HS19070735-03	WG-1620-MW04-20190712	Login	15/07/2019 18:21:52	JRM	VOA074
HS19070735-04	WG-1620-TW41B-20190712	Login	15/07/2019 18:21:52	JRM	MET044
HS19070735-04	WG-1620-TW41B-20190712	Login	15/07/2019 18:21:52	JRM	EXT135
HS19070735-04	WG-1620-TW41B-20190712	Login	15/07/2019 18:21:52	JRM	VOA074
HS19070735-05	WG-1620-FB03-20190712	Login	15/07/2019 18:21:52	JRM	MET044
HS19070735-05	WG-1620-FB03-20190712	Login	15/07/2019 18:21:52	JRM	EXT135
HS19070735-05	WG-1620-FB03-20190712	Login	15/07/2019 18:21:52	JRM	VOA074
HS19070735-06	WG-1620-TB03-20190712	Login	15/07/2019 18:21:52	JRM	VOA074

---

Sample Receipt Checklist

Client Name: PBW
Work Order: HS19070735

Date/Time Received: 15-Jul-2019 16:30
Received by: JRM

Checklist completed by: Jared R. Makan
eSignature
Date: 15-Jul-2019

Reviewed by: Dane J. Wacasey
eSignature
Date: 18-Jul-2019

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:196142

Temperature(s)/Thermometer(s): 1.0c/1.0c UC/C IR25
Cooler(s)/Kit(s): 45005
Date/Time sample(s) sent to storage: 07/15/2019 18:40
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

## HS19070735

Golder Associates Inc.  
Houston TX-Wood Preserving Works

Page 1 of 1

COC ID: 196142



ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A 8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B 8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C 8270_LOW_W (5632532 Semi/Volatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D ICP_TW (5636002 5652546 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E
	Suite 4004		Stop 0750	F
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G
Phone	(512) 671-3434	Phone		H
Fax	(512) 671-3446	Fax		I
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TB01-20190712</del>			<del>Water</del>	<del>+</del>	<del>2</del>		*									
2	WG-1620-MW09-20190712	7-12-19	0735	W		6	X		X	X							
3	WG-1620-MW03-20190712		0835	W		6	X		X	X							
4	WG-1620-MW04-20190712		0925	W		6	X		X	X							
5	WG-1620-TW41B-20190712		1040	W		6	X		X	X							
6	WG-1620-FB03-20190712		1100	W		6	X		X	X							
7	WG-1620-TB03-20190712			W		2	X										
8																	
9																	
10																	

Sampler(s) Please Print & Sign <b>John Beayton</b>		Shipment Method <b>Hand Delivered</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STB 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour		Results Due Date:
Relinquished by: <b>John Beayton</b>	Date: <b>7-15-19</b>	Time: <b>16:30</b>	Received by:	Notes: <b>UPRR Houston MWPW</b>	
Relinquished by: <b>John Beayton</b>	Date: <b>7/15/19</b>	Time: <b>16:30</b>	Received by (Laboratory): <b>S. M. ...</b>	Cooler ID: <b>45005</b>	Cooler Temp.: <b>10</b>
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	QC Package: (Check One Box Below) <input type="checkbox"/> Level II Site QC <input checked="" type="checkbox"/> TRRP Checklist <input type="checkbox"/> Level III Site QC/Raw Data <input type="checkbox"/> TRRP Level IV <input type="checkbox"/> Level IV SW/813/CLP <input type="checkbox"/> Other:	
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035					

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

July 25, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19070864**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 13 sample(s) on Jul 17, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.



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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/25/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19070864			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143194,143304,143325,R342655,R342706			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Were MS/MSD RPDs within laboratory QC limits?		X			3
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?	X				
		Were analytical duplicates analyzed at the appropriate frequency?	X				
		Were RPDs or relative standard deviations within the laboratory QC limits?	X				
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				4
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group			LRC Date: 07/25/2019				
Project Name: Houston TX-Wood Preserving Works			Laboratory Job Number: HS19070864				
Reviewer Name: Dane Wacasey			Prep Batch Number(s): 143194,143304,143325,R342655,R342706				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 07/25/2019
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS19070864
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 143194,143304,143325,R342655,R342706
ER# <sup>5</sup>	Description	
1	Semivolatile Organics Method SW8270, samples WG-1620-MW25C-20190716, WG-1620-MW63B-20190716: the surrogate recoveries could not be determined due to dilution below the calibration range	
2	Batch 143194, Semivolatile Organics Method SW8270, sample WG-1620-MW28C-20190716, MSD recovered outside the control limit for Phenol due to possible matrix effect.	
3	Batch 143194, Semivolatile Organics Method SW8270, sample WG-1620-MW28C-20190716, MS/MSD RPD recovered above the RPD limit for 2-Chloronaphthalene due to possible matrix effect.	
4	Batch 143194, Semivolatile Organics Method SW8270, samples WG-1620-MW25C-20190716, the GCMS semi-volatile extract of this sample was run at a dilution due to a high level of matrix interference.	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.                      O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);                      NA = Not Applicable;                      NR = Not Reviewed;                      R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070864

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19070864-01	WG-1620-MW21C-20190716	Water		16-Jul-2019 07:35	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-02	WG-1620-FD01-20190716	Water		16-Jul-2019 07:35	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-03	WG-1620-MW62B-20190716	Water		16-Jul-2019 08:25	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-04	WG-1620-MW53C-20190716	Water		16-Jul-2019 10:15	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-05	WG-1620-MW54C-20190716	Water		16-Jul-2019 11:15	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-06	WG-1620-MW25C-20190716	Water		16-Jul-2019 12:15	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-07	WG-1620-MW25A-20190716	Water		16-Jul-2019 13:10	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-08	WG-1620-MW36B-20190716	Water		16-Jul-2019 14:20	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-09	WG-1620-MW36A-20190716	Water		16-Jul-2019 15:20	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-10	WG-1620-MW28C-20190716	Water		16-Jul-2019 16:15	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-11	WG-1620-MW28A-20190716	Water		16-Jul-2019 17:00	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-12	WG-1620-MW63B-20190716	Water		16-Jul-2019 17:55	17-Jul-2019 14:24	<input type="checkbox"/>
HS19070864-13	WG-1620-TB04-20190716	Water		16-Jul-2019 00:00	17-Jul-2019 14:24	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW21C-20190716  
 Collection Date: 16-Jul-2019 07:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jul-2019 14:48
Benzene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 14:48
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 14:48
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 14:48
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jul-2019 14:48
Toluene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 14:48
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jul-2019 14:48
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:48</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:48</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:48</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 14:48</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW21C-20190716  
 Collection Date: 16-Jul-2019 07:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jul-2019 01:12
<b>2,4-Dimethylphenol</b>	<b>0.00010</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jul-2019 01:12
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jul-2019 01:12
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jul-2019 01:12
<b>2-Methylnaphthalene</b>	<b>0.00040</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jul-2019 01:12
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jul-2019 01:12
<b>Acenaphthene</b>	<b>0.0057</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<b>Acenaphthylene</b>	<b>0.000074</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<b>Anthracene</b>	<b>0.00044</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jul-2019 01:12
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jul-2019 01:12
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jul-2019 01:12
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000040</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jul-2019 01:12
<b>Dibenzofuran</b>	<b>0.0032</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jul-2019 01:12
<b>Fluoranthene</b>	<b>0.00081</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<b>Fluorene</b>	<b>0.0019</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<b>Naphthalene</b>	<b>0.0044</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jul-2019 01:12
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jul-2019 01:12
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jul-2019 01:12
<b>Phenanthrene</b>	<b>0.00087</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<b>Phenol</b>	<b>0.00018</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<b>Pyrene</b>	<b>0.00039</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:12
<i>Surr: 2,4,6-Tribromophenol</i>	77.5			34-129	%REC	1	20-Jul-2019 01:12
<i>Surr: 2-Fluorobiphenyl</i>	61.6			40-125	%REC	1	20-Jul-2019 01:12
<i>Surr: 2-Fluorophenol</i>	53.9			20-120	%REC	1	20-Jul-2019 01:12
<i>Surr: 4-Terphenyl-d14</i>	72.9			40-135	%REC	1	20-Jul-2019 01:12
<i>Surr: Nitrobenzene-d5</i>	67.9			41-120	%REC	1	20-Jul-2019 01:12
<i>Surr: Phenol-d6</i>	67.1			20-120	%REC	1	20-Jul-2019 01:12
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00130</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	23-Jul-2019 22:28

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20190716  
 Collection Date: 16-Jul-2019 07:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 07:08
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 07:08
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 07:08
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 07:08
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 07:08
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 07:08
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 07:08
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>88.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 07:08</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 07:08</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 07:08</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 07:08</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20190716  
 Collection Date: 16-Jul-2019 07:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jul-2019 01:32
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	20-Jul-2019 01:32
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jul-2019 01:32
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jul-2019 01:32
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jul-2019 01:32
<b>2-Methylnaphthalene</b>	<b>0.000050</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jul-2019 01:32
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jul-2019 01:32
<b>Acenaphthene</b>	<b>0.00081</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
Acenaphthylene	U		0.000015	0.00010	mg/L	1	20-Jul-2019 01:32
<b>Anthracene</b>	<b>0.000050</b>	<b>J</b>	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jul-2019 01:32
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jul-2019 01:32
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jul-2019 01:32
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jul-2019 01:32
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jul-2019 01:32
<b>Dibenzofuran</b>	<b>0.00040</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<b>Di-n-butyl phthalate</b>	<b>0.000020</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<b>Fluoranthene</b>	<b>0.000087</b>	<b>J</b>	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<b>Fluorene</b>	<b>0.00025</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<b>Naphthalene</b>	<b>0.00080</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jul-2019 01:32
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jul-2019 01:32
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jul-2019 01:32
<b>Phenanthrene</b>	<b>0.00011</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<b>Phenol</b>	<b>0.00081</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<b>Pyrene</b>	<b>0.000033</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jul-2019 01:32</b>
<i>Surr: 2,4,6-Tribromophenol</i>	76.9			34-129	%REC	1	20-Jul-2019 01:32
<i>Surr: 2-Fluorobiphenyl</i>	65.2			40-125	%REC	1	20-Jul-2019 01:32
<i>Surr: 2-Fluorophenol</i>	58.3			20-120	%REC	1	20-Jul-2019 01:32
<i>Surr: 4-Terphenyl-d14</i>	68.4			40-135	%REC	1	20-Jul-2019 01:32
<i>Surr: Nitrobenzene-d5</i>	68.2			41-120	%REC	1	20-Jul-2019 01:32
<i>Surr: Phenol-d6</i>	64.6			20-120	%REC	1	20-Jul-2019 01:32
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00326</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 22:31</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW62B-20190716  
 Collection Date: 16-Jul-2019 08:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 04:42
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 04:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 04:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 04:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 04:42
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 04:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 04:42
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:42</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:42</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:42</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW62B-20190716  
 Collection Date: 16-Jul-2019 08:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jul-2019 01:51
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	20-Jul-2019 01:51
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jul-2019 01:51
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jul-2019 01:51
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jul-2019 01:51
<b>2-Methylnaphthalene</b>	<b>0.000096</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jul-2019 01:51
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jul-2019 01:51
<b>Acenaphthene</b>	<b>0.000099</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
Acenaphthylene	U		0.000015	0.00010	mg/L	1	20-Jul-2019 01:51
Anthracene	U		0.000014	0.00010	mg/L	1	20-Jul-2019 01:51
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jul-2019 01:51
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jul-2019 01:51
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jul-2019 01:51
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jul-2019 01:51
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jul-2019 01:51
<b>Dibenzofuran</b>	<b>0.000048</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jul-2019 01:51
<b>Fluoranthene</b>	<b>0.000017</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
<b>Fluorene</b>	<b>0.000036</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
<b>Naphthalene</b>	<b>0.00028</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jul-2019 01:51
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jul-2019 01:51
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jul-2019 01:51
<b>Phenanthrene</b>	<b>0.000021</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
<b>Phenol</b>	<b>0.00056</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 01:51
Pyrene	U		0.000019	0.00010	mg/L	1	20-Jul-2019 01:51
<i>Surr: 2,4,6-Tribromophenol</i>	85.2			34-129	%REC	1	20-Jul-2019 01:51
<i>Surr: 2-Fluorobiphenyl</i>	73.5			40-125	%REC	1	20-Jul-2019 01:51
<i>Surr: 2-Fluorophenol</i>	60.4			20-120	%REC	1	20-Jul-2019 01:51
<i>Surr: 4-Terphenyl-d14</i>	73.3			40-135	%REC	1	20-Jul-2019 01:51
<i>Surr: Nitrobenzene-d5</i>	70.2			41-120	%REC	1	20-Jul-2019 01:51
<i>Surr: Phenol-d6</i>	70.9			20-120	%REC	1	20-Jul-2019 01:51
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00194</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	23-Jul-2019 22:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW53C-20190716  
 Collection Date: 16-Jul-2019 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jul-2019 20:06
Benzene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 20:06
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 20:06
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 20:06
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jul-2019 20:06
Toluene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 20:06
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jul-2019 20:06
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>89.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:06</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:06</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:06</i>
<i>Surr: Toluene-d8</i>	<i>99.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:06</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW53C-20190716  
 Collection Date: 16-Jul-2019 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jul-2019 02:11
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	20-Jul-2019 02:11
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jul-2019 02:11
<b>2,6-Dinitrotoluene</b>	<b>0.0028</b>		<b>0.000042</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 02:11
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jul-2019 02:11
<b>2-Methylnaphthalene</b>	<b>0.000071</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:11
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jul-2019 02:11
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jul-2019 02:11
Acenaphthene	U		0.000027	0.00010	mg/L	1	20-Jul-2019 02:11
Acenaphthylene	U		0.000015	0.00010	mg/L	1	20-Jul-2019 02:11
Anthracene	U		0.000014	0.00010	mg/L	1	20-Jul-2019 02:11
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jul-2019 02:11
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jul-2019 02:11
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jul-2019 02:11
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jul-2019 02:11
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jul-2019 02:11
Dibenzofuran	U		0.000020	0.00010	mg/L	1	20-Jul-2019 02:11
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jul-2019 02:11
Fluoranthene	U		0.000010	0.00010	mg/L	1	20-Jul-2019 02:11
Fluorene	U		0.000030	0.00010	mg/L	1	20-Jul-2019 02:11
<b>Naphthalene</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:11
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jul-2019 02:11
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jul-2019 02:11
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jul-2019 02:11
Phenanthrene	U		0.000021	0.00010	mg/L	1	20-Jul-2019 02:11
<b>Phenol</b>	<b>0.00016</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 02:11
Pyrene	U		0.000019	0.00010	mg/L	1	20-Jul-2019 02:11
<i>Surr: 2,4,6-Tribromophenol</i>	<i>77.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:11</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>70.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:11</i>
<i>Surr: 2-Fluorophenol</i>	<i>63.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:11</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>73.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:11</i>
<i>Surr: Nitrobenzene-d5</i>	<i>69.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:11</i>
<i>Surr: Phenol-d6</i>	<i>70.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:11</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000569</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	23-Jul-2019 22:40

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54C-20190716  
 Collection Date: 16-Jul-2019 11:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jul-2019 23:48
Benzene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 23:48
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 23:48
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 23:48
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jul-2019 23:48
Toluene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 23:48
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jul-2019 23:48
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 23:48</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 23:48</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 23:48</i>
<i>Surr: Toluene-d8</i>	<i>103</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 23:48</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54C-20190716  
 Collection Date: 16-Jul-2019 11:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jul-2019 02:31
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	20-Jul-2019 02:31
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jul-2019 02:31
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jul-2019 02:31
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jul-2019 02:31
<b>2-Methylnaphthalene</b>	<b>0.0021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jul-2019 02:31
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jul-2019 02:31
<b>Acenaphthene</b>	<b>0.041</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:38
<b>Acenaphthylene</b>	<b>0.00037</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
<b>Anthracene</b>	<b>0.0023</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jul-2019 02:31
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jul-2019 02:31
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jul-2019 02:31
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jul-2019 02:31
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jul-2019 02:31
<b>Dibenzofuran</b>	<b>0.032</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:38
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jul-2019 02:31
<b>Fluoranthene</b>	<b>0.0033</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
<b>Fluorene</b>	<b>0.018</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:38
<b>Naphthalene</b>	<b>0.032</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:38
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jul-2019 02:31
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jul-2019 02:31
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jul-2019 02:31
<b>Phenanthrene</b>	<b>0.0065</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
<b>Phenol</b>	<b>0.00044</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
<b>Pyrene</b>	<b>0.0016</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jul-2019 02:31
<i>Surr: 2,4,6-Tribromophenol</i>	<i>81.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:31</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>95.1</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.3</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>58.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:31</i>
<i>Surr: 2-Fluorophenol</i>	<i>58.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:31</i>
<i>Surr: 2-Fluorophenol</i>	<i>69.2</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>93.8</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>67.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:31</i>
<i>Surr: Nitrobenzene-d5</i>	<i>57.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:31</i>
<i>Surr: Nitrobenzene-d5</i>	<i>86.8</i>			<i>41-120</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:38</i>
<i>Surr: Phenol-d6</i>	<i>64.3</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:38</i>
<i>Surr: Phenol-d6</i>	<i>62.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jul-2019 02:31</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54C-20190716  
 Collection Date: 16-Jul-2019 11:15

**ANALYTICAL REPORT**

WorkOrder:HS19070864  
 Lab ID:HS19070864-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jul-2019		Analyst: JHD	
Arsenic	0.00103	J	0.000400	0.00200	mg/L	1	23-Jul-2019 22:42

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25C-20190716  
 Collection Date: 16-Jul-2019 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	18-Jul-2019 20:30
<b>Benzene</b>	<b>0.0013</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	18-Jul-2019 20:30
Chlorobenzene		U	0.00030	0.0010	mg/L	1	18-Jul-2019 20:30
<b>Ethylbenzene</b>	<b>0.035</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	18-Jul-2019 20:30
Methylene chloride		U	0.0010	0.0020	mg/L	1	18-Jul-2019 20:30
<b>Toluene</b>	<b>0.011</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	18-Jul-2019 20:30
Vinyl chloride		U	0.00020	0.0010	mg/L	1	18-Jul-2019 20:30
<b>Xylenes, Total</b>	<b>0.25</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	18-Jul-2019 20:30
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:30</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:30</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:30</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 20:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25C-20190716  
 Collection Date: 16-Jul-2019 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.00021	0.0020	mg/L	10	22-Jul-2019 14:58
<b>2,4-Dimethylphenol</b>	<b>0.0075</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
2,4-Dinitrotoluene		U	0.00058	0.0020	mg/L	10	22-Jul-2019 14:58
2,6-Dinitrotoluene		U	0.00042	0.0020	mg/L	10	22-Jul-2019 14:58
2-Chloronaphthalene		U	0.00021	0.0020	mg/L	10	22-Jul-2019 14:58
<b>2-Methylnaphthalene</b>	<b>0.54</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	22-Jul-2019 15:18
4,6-Dinitro-2-methylphenol		U	0.00020	0.0020	mg/L	10	22-Jul-2019 14:58
4-Nitrophenol		U	0.00047	0.010	mg/L	10	22-Jul-2019 14:58
<b>Acenaphthene</b>	<b>0.20</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	22-Jul-2019 15:18
<b>Acenaphthylene</b>	<b>0.0020</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<b>Anthracene</b>	<b>0.016</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<b>Benz(a)anthracene</b>	<b>0.0013</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<b>Benzo(a)pyrene</b>	<b>0.00037</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
Bis(2-chloroethoxy)methane		U	0.00030	0.0020	mg/L	10	22-Jul-2019 14:58
Bis(2-ethylhexyl)phthalate		U	0.00037	0.0020	mg/L	10	22-Jul-2019 14:58
<b>Chrysene</b>	<b>0.0014</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<b>Dibenzofuran</b>	<b>0.19</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	22-Jul-2019 15:18
Di-n-butyl phthalate		U	0.00020	0.0020	mg/L	10	22-Jul-2019 14:58
<b>Fluoranthene</b>	<b>0.014</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<b>Fluorene</b>	<b>0.081</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<b>Naphthalene</b>	<b>4.2</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	22-Jul-2019 15:38
Nitrobenzene		U	0.00024	0.0020	mg/L	10	22-Jul-2019 14:58
N-Nitrosodiphenylamine		U	0.00025	0.0020	mg/L	10	22-Jul-2019 14:58
Pentachlorophenol		U	0.00079	0.0020	mg/L	10	22-Jul-2019 14:58
<b>Phenanthrene</b>	<b>0.12</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	22-Jul-2019 15:18
Phenol		U	0.00035	0.0020	mg/L	10	22-Jul-2019 14:58
<b>Pyrene</b>	<b>0.0085</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 14:58
<i>Surr: 2,4,6-Tribromophenol</i>	<i>119</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 14:58</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>22-Jul-2019 15:18</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 15:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>78.7</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 14:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>22-Jul-2019 15:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 15:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>74.1</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 14:58</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>22-Jul-2019 15:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 15:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>96.6</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 14:58</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>22-Jul-2019 15:18</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 15:38</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25C-20190716  
 Collection Date: 16-Jul-2019 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	73.3			41-120	%REC	10	22-Jul-2019 14:58
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	22-Jul-2019 15:18
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	22-Jul-2019 15:38
Surr: Phenol-d6	0	JS		20-120	%REC	1000	22-Jul-2019 15:38
Surr: Phenol-d6	84.5			20-120	%REC	10	22-Jul-2019 14:58
Surr: Phenol-d6	0	JS		20-120	%REC	100	22-Jul-2019 15:18
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jul-2019		Analyst: JHD	
Arsenic	0.00487		0.000400	0.00200	mg/L	1	23-Jul-2019 22:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25A-20190716  
 Collection Date: 16-Jul-2019 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:28
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:28
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 03:28
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 03:28
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 03:28
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:28
Vinyl chloride	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:28
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 03:28
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>89.1</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:28</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.1</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:28</i>
<i>Surr: Dibromofluoromethane</i>		<i>98.5</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:28</i>
<i>Surr: Toluene-d8</i>		<i>101</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:28</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25A-20190716  
 Collection Date: 16-Jul-2019 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 12:00
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 12:00
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 12:00
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 12:00
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 12:00
<b>2-Methylnaphthalene</b>	<b>0.000057</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 12:00</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 12:00
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 12:00
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 12:00
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 12:00
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 12:00
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 12:00
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 12:00
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 12:00
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	22-Jul-2019 12:00
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 12:00
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 12:00
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 12:00
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 12:00
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 12:00
<b>Naphthalene</b>	<b>0.00023</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 12:00</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 12:00
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 12:00
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 12:00
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 12:00
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 12:00
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 12:00
<i>Surr: 2,4,6-Tribromophenol</i>	<i>92.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:00</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:00</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:00</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>82.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:00</i>
<i>Surr: Nitrobenzene-d5</i>	<i>74.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:00</i>
<i>Surr: Phenol-d6</i>	<i>68.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:00</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00285</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jul-2019 15:44</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36B-20190716  
 Collection Date: 16-Jul-2019 14:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:37
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 00:37
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 00:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 00:37
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:37
Vinyl chloride	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:37
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 00:37
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>87.9</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:37</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>96.2</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:37</i>
<i>Surr: Dibromofluoromethane</i>		<i>96.5</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:37</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:37</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36B-20190716  
 Collection Date: 16-Jul-2019 14:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 12:20
<b>2,4-Dimethylphenol</b>	<b>0.000078</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 12:20
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 12:20
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 12:20
<b>2-Methylnaphthalene</b>	<b>0.000058</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 12:20
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 12:20
<b>Acenaphthene</b>	<b>0.00014</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 12:20
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 12:20
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 12:20
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 12:20
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 12:20
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000088</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 12:20
<b>Dibenzofuran</b>	<b>0.000087</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 12:20
<b>Fluoranthene</b>	<b>0.000017</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
<b>Fluorene</b>	<b>0.000047</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
<b>Naphthalene</b>	<b>0.0060</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 12:20
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 12:20
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 12:20
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 12:20
<b>Phenol</b>	<b>0.00017</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 12:20
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 12:20
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:20</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>72.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>74.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:20</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>82.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:20</i>
<i>Surr: Nitrobenzene-d5</i>	<i>78.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:20</i>
<i>Surr: Phenol-d6</i>	<i>75.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 12:20</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00127</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jul-2019 15:47

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36A-20190716  
 Collection Date: 16-Jul-2019 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:53
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:53
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 03:53
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 03:53
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 03:53
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:53
Vinyl chloride	U		0.00020	0.0010	mg/L	1	19-Jul-2019 03:53
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 03:53
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>87.0</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:53</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>96.0</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:53</i>
<i>Surr: Dibromofluoromethane</i>		<i>97.3</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:53</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 03:53</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36A-20190716  
 Collection Date: 16-Jul-2019 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 15:58
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 15:58
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 15:58
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 15:58
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 15:58
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 15:58
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 15:58
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 15:58
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 15:58
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 15:58
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 15:58
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 15:58
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 15:58
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 15:58
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	22-Jul-2019 15:58
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 15:58
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 15:58
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 15:58
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 15:58
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 15:58
<b>Naphthalene</b>	<b>0.00015</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 15:58
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 15:58
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 15:58
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 15:58
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 15:58
<b>Phenol</b>	<b>0.00014</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 15:58
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 15:58
<i>Surr: 2,4,6-Tribromophenol</i>	<i>84.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 15:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 15:58</i>
<i>Surr: 2-Fluorophenol</i>	<i>63.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 15:58</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>82.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 15:58</i>
<i>Surr: Nitrobenzene-d5</i>	<i>75.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 15:58</i>
<i>Surr: Phenol-d6</i>	<i>63.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 15:58</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00244</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jul-2019 15:49

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28C-20190716  
 Collection Date: 16-Jul-2019 16:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jul-2019 13:34
Benzene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 13:34
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 13:34
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jul-2019 13:34
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jul-2019 13:34
Toluene	U		0.00020	0.0010	mg/L	1	18-Jul-2019 13:34
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jul-2019 13:34
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>88.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 13:34</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 13:34</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 13:34</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jul-2019 13:34</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28C-20190716  
 Collection Date: 16-Jul-2019 16:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 16:18
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 16:18
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 16:18
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 16:18
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 16:18
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 16:18
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 16:18
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 16:18
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 16:18
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 16:18
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 16:18
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 16:18
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 16:18
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 16:18
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00013</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 16:18
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 16:18
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 16:18
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 16:18
<b>Fluoranthene</b>	<b>0.000015</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 16:18
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 16:18
<b>Naphthalene</b>	<b>0.000095</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 16:18
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 16:18
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 16:18
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 16:18
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 16:18
<b>Phenol</b>	<b>0.0025</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 16:18
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 16:18
<i>Surr: 2,4,6-Tribromophenol</i>	<i>109</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 16:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 16:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>70.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 16:18</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 16:18</i>
<i>Surr: Nitrobenzene-d5</i>	<i>79.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 16:18</i>
<i>Surr: Phenol-d6</i>	<i>69.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 16:18</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000456</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jul-2019 15:51

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28A-20190716  
 Collection Date: 16-Jul-2019 17:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 04:17
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 04:17
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 04:17
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 04:17
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 04:17
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 04:17
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 04:17
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>88.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:17</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:17</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:17</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 04:17</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28A-20190716  
 Collection Date: 16-Jul-2019 17:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 18:37
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 18:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 18:37
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 18:37
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 18:37
<b>2-Methylnaphthalene</b>	<b>0.00020</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 18:37
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 18:37
<b>Acenaphthene</b>	<b>0.000082</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 18:37
<b>Anthracene</b>	<b>0.000036</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Benz(a)anthracene</b>	<b>0.000055</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Benzo(a)pyrene</b>	<b>0.000095</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 18:37
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00023</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Chrysene</b>	<b>0.00010</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Dibenzofuran</b>	<b>0.000084</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 18:37
<b>Fluoranthene</b>	<b>0.00012</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Fluorene</b>	<b>0.000053</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Naphthalene</b>	<b>0.0013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 18:37
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 18:37
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 18:37
<b>Phenanthrene</b>	<b>0.000049</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Phenol</b>	<b>0.000073</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<b>Pyrene</b>	<b>0.00016</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 18:37
<i>Surr: 2,4,6-Tribromophenol</i>	111			34-129	%REC	1	22-Jul-2019 18:37
<i>Surr: 2-Fluorobiphenyl</i>	76.9			40-125	%REC	1	22-Jul-2019 18:37
<i>Surr: 2-Fluorophenol</i>	65.8			20-120	%REC	1	22-Jul-2019 18:37
<i>Surr: 4-Terphenyl-d14</i>	82.4			40-135	%REC	1	22-Jul-2019 18:37
<i>Surr: Nitrobenzene-d5</i>	72.0			41-120	%REC	1	22-Jul-2019 18:37
<i>Surr: Phenol-d6</i>	74.5			20-120	%REC	1	22-Jul-2019 18:37
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00920</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jul-2019 16:19

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW63B-20190716  
 Collection Date: 16-Jul-2019 17:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 08:00
<b>Benzene</b>	<b>0.11</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	19-Jul-2019 08:00
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 08:00
<b>Ethylbenzene</b>	<b>0.14</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	19-Jul-2019 08:00
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 08:00
<b>Toluene</b>	<b>0.00071</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	19-Jul-2019 08:00
<b>Xylenes, Total</b>	<b>0.041</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	19-Jul-2019 08:00
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>85.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 08:00</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 08:00</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 08:00</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 08:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW63B-20190716  
 Collection Date: 16-Jul-2019 17:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 13:59
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 13:59
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 13:59
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 13:59
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 13:59
<b>2-Methylnaphthalene</b>	<b>0.037</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:58
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 13:59
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 13:59
<b>Acenaphthene</b>	<b>0.012</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:58
<b>Acenaphthylene</b>	<b>0.00016</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 13:59
<b>Anthracene</b>	<b>0.00034</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 13:59
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 13:59
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 13:59
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 13:59
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	22-Jul-2019 13:59
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 13:59
<b>Dibenzofuran</b>	<b>0.011</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	22-Jul-2019 16:58
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 13:59
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 13:59
<b>Fluorene</b>	<b>0.0037</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 13:59
<b>Naphthalene</b>	<b>1.5</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	22-Jul-2019 17:57
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 13:59
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 13:59
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 13:59
<b>Phenanthrene</b>	<b>0.0013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 13:59
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 13:59
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 13:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>107</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 13:59</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>106</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:58</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 17:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>71.6</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 17:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 13:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 13:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>83.3</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:58</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 17:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>84.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 13:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.7</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>22-Jul-2019 16:58</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>1000</i>	<i>22-Jul-2019 17:57</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW63B-20190716  
 Collection Date: 16-Jul-2019 17:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	85.7			41-120	%REC	1	22-Jul-2019 13:59
Surr: Nitrobenzene-d5	75.0			41-120	%REC	10	22-Jul-2019 16:58
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	22-Jul-2019 17:57
Surr: Phenol-d6	0	JS		20-120	%REC	1000	22-Jul-2019 17:57
Surr: Phenol-d6	68.4			20-120	%REC	1	22-Jul-2019 13:59
Surr: Phenol-d6	66.8			20-120	%REC	10	22-Jul-2019 16:58
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jul-2019		Analyst: JHD	
Arsenic	0.00156	J	0.000400	0.00200	mg/L	1	24-Jul-2019 16:21

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB04-20190716  
 Collection Date: 16-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070864  
 Lab ID:HS19070864-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:12
Benzene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:12
Chlorobenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 00:12
Ethylbenzene	U		0.00030	0.0010	mg/L	1	19-Jul-2019 00:12
Methylene chloride	U		0.0010	0.0020	mg/L	1	19-Jul-2019 00:12
Toluene	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:12
Vinyl chloride	U		0.00020	0.0010	mg/L	1	19-Jul-2019 00:12
Xylenes, Total	U		0.00030	0.0010	mg/L	1	19-Jul-2019 00:12
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:12</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:12</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:12</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>19-Jul-2019 00:12</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**Batch ID:** 143194      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070864-01	1	1000	1 (mL)	0.001
HS19070864-02	1	1000	1 (mL)	0.001
HS19070864-03	1	1000	1 (mL)	0.001
HS19070864-04	1	1000	1 (mL)	0.001
HS19070864-05	1	1000	1 (mL)	0.001
HS19070864-06	1	1000	1 (mL)	0.001
HS19070864-07	1	1000	1 (mL)	0.001
HS19070864-08	1	1000	1 (mL)	0.001
HS19070864-09	1	1000	1 (mL)	0.001
HS19070864-10	1	1000	1 (mL)	0.001
HS19070864-11	1	1000	1 (mL)	0.001
HS19070864-12	1	1000	1 (mL)	0.001

**Batch ID:** 143304      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070864-01	1	10	10 (mL)	1
HS19070864-02	1	10	10 (mL)	1
HS19070864-03	1	10	10 (mL)	1
HS19070864-04	1	10	10 (mL)	1
HS19070864-05	1	10	10 (mL)	1
HS19070864-06	1	10	10 (mL)	1

**Batch ID:** 143325      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070864-07	1	10	10 (mL)	1
HS19070864-08	1	10	10 (mL)	1
HS19070864-09	1	10	10 (mL)	1
HS19070864-10	1	10	10 (mL)	1
HS19070864-11	1	10	10 (mL)	1
HS19070864-12	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143194 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS19070864-01	WG-1620-MW21C-20190716	16 Jul 2019 07:35		19 Jul 2019 11:41	20 Jul 2019 01:12	1
HS19070864-02	WG-1620-FD01-20190716	16 Jul 2019 07:35		19 Jul 2019 11:41	20 Jul 2019 01:32	1
HS19070864-03	WG-1620-MW62B-20190716	16 Jul 2019 08:25		19 Jul 2019 11:41	20 Jul 2019 01:51	1
HS19070864-04	WG-1620-MW53C-20190716	16 Jul 2019 10:15		19 Jul 2019 11:41	20 Jul 2019 02:11	1
HS19070864-05	WG-1620-MW54C-20190716	16 Jul 2019 11:15		19 Jul 2019 11:41	22 Jul 2019 16:38	10
HS19070864-05	WG-1620-MW54C-20190716	16 Jul 2019 11:15		19 Jul 2019 11:41	20 Jul 2019 02:31	1
HS19070864-06	WG-1620-MW25C-20190716	16 Jul 2019 12:15		19 Jul 2019 11:41	22 Jul 2019 15:38	1000
HS19070864-06	WG-1620-MW25C-20190716	16 Jul 2019 12:15		19 Jul 2019 11:41	22 Jul 2019 15:18	100
HS19070864-06	WG-1620-MW25C-20190716	16 Jul 2019 12:15		19 Jul 2019 11:41	22 Jul 2019 14:58	10
HS19070864-07	WG-1620-MW25A-20190716	16 Jul 2019 13:10		19 Jul 2019 11:41	22 Jul 2019 12:00	1
HS19070864-08	WG-1620-MW36B-20190716	16 Jul 2019 14:20		19 Jul 2019 11:41	22 Jul 2019 12:20	1
HS19070864-09	WG-1620-MW36A-20190716	16 Jul 2019 15:20		19 Jul 2019 11:41	22 Jul 2019 15:58	1
HS19070864-10	WG-1620-MW28C-20190716	16 Jul 2019 16:15		19 Jul 2019 11:41	22 Jul 2019 16:18	1
HS19070864-11	WG-1620-MW28A-20190716	16 Jul 2019 17:00		19 Jul 2019 11:41	22 Jul 2019 18:37	1
HS19070864-12	WG-1620-MW63B-20190716	16 Jul 2019 17:55		19 Jul 2019 11:41	22 Jul 2019 17:57	1000
HS19070864-12	WG-1620-MW63B-20190716	16 Jul 2019 17:55		19 Jul 2019 11:41	22 Jul 2019 16:58	10
HS19070864-12	WG-1620-MW63B-20190716	16 Jul 2019 17:55		19 Jul 2019 11:41	22 Jul 2019 13:59	1
<b>Batch ID: 143304 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19070864-01	WG-1620-MW21C-20190716	16 Jul 2019 07:35		23 Jul 2019 12:00	23 Jul 2019 22:28	1
HS19070864-02	WG-1620-FD01-20190716	16 Jul 2019 07:35		23 Jul 2019 12:00	23 Jul 2019 22:31	1
HS19070864-03	WG-1620-MW62B-20190716	16 Jul 2019 08:25		23 Jul 2019 12:00	23 Jul 2019 22:37	1
HS19070864-04	WG-1620-MW53C-20190716	16 Jul 2019 10:15		23 Jul 2019 12:00	23 Jul 2019 22:40	1
HS19070864-05	WG-1620-MW54C-20190716	16 Jul 2019 11:15		23 Jul 2019 12:00	23 Jul 2019 22:42	1
HS19070864-06	WG-1620-MW25C-20190716	16 Jul 2019 12:15		23 Jul 2019 12:00	23 Jul 2019 22:44	1
<b>Batch ID: 143325 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19070864-07	WG-1620-MW25A-20190716	16 Jul 2019 13:10		24 Jul 2019 09:30	24 Jul 2019 15:44	1
HS19070864-08	WG-1620-MW36B-20190716	16 Jul 2019 14:20		24 Jul 2019 09:30	24 Jul 2019 15:47	1
HS19070864-09	WG-1620-MW36A-20190716	16 Jul 2019 15:20		24 Jul 2019 09:30	24 Jul 2019 15:49	1
HS19070864-10	WG-1620-MW28C-20190716	16 Jul 2019 16:15		24 Jul 2019 09:30	24 Jul 2019 15:51	1
HS19070864-11	WG-1620-MW28A-20190716	16 Jul 2019 17:00		24 Jul 2019 09:30	24 Jul 2019 16:19	1
HS19070864-12	WG-1620-MW63B-20190716	16 Jul 2019 17:55		24 Jul 2019 09:30	24 Jul 2019 16:21	1
<b>Batch ID: R342655 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19070864-01	WG-1620-MW21C-20190716	16 Jul 2019 07:35			18 Jul 2019 14:48	1
HS19070864-04	WG-1620-MW53C-20190716	16 Jul 2019 10:15			18 Jul 2019 20:06	1
HS19070864-06	WG-1620-MW25C-20190716	16 Jul 2019 12:15			18 Jul 2019 20:30	1
HS19070864-10	WG-1620-MW28C-20190716	16 Jul 2019 16:15			18 Jul 2019 13:34	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: R342706 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19070864-02	WG-1620-FD01-20190716	16 Jul 2019 07:35			19 Jul 2019 07:08	1
HS19070864-03	WG-1620-MW62B-20190716	16 Jul 2019 08:25			19 Jul 2019 04:42	1
HS19070864-05	WG-1620-MW54C-20190716	16 Jul 2019 11:15			18 Jul 2019 23:48	1
HS19070864-07	WG-1620-MW25A-20190716	16 Jul 2019 13:10			19 Jul 2019 03:28	1
HS19070864-08	WG-1620-MW36B-20190716	16 Jul 2019 14:20			19 Jul 2019 00:37	1
HS19070864-09	WG-1620-MW36A-20190716	16 Jul 2019 15:20			19 Jul 2019 03:53	1
HS19070864-11	WG-1620-MW28A-20190716	16 Jul 2019 17:00			19 Jul 2019 04:17	1
HS19070864-12	WG-1620-MW63B-20190716	16 Jul 2019 17:55			19 Jul 2019 08:00	1
HS19070864-13	WG-1620-TB04-20190716	16 Jul 2019 00:00			19 Jul 2019 00:12	1

WorkOrder: HS19070864  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200

WorkOrder: HS19070864  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00010	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000096	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00012	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00011	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000049	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000013	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00013	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000051	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000055	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000068	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000053	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000086	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000066	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000052	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.000099	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000052	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000061	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00013	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000085	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000055	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000094	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000061	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19070864  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00067	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00056	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00059	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00051	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00054	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00056	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00046	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00052	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

<b>Batch ID:</b> 143304 ( 0 )	<b>Instrument:</b> ICPMS05	<b>Method:</b> ICP-MS METALS BY SW6020A
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<b>MBLK</b>	Sample ID: <b>MBLK-143304</b>	Units: <b>mg/L</b>	Analysis Date: <b>23-Jul-2019 21:27</b>							
Client ID:	Run ID: <b>ICPMS05_342915</b>	SeqNo: <b>5177608</b>	PrepDate: <b>23-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic U 0.00200

<b>LCS</b>	Sample ID: <b>LCS-143304</b>	Units: <b>mg/L</b>	Analysis Date: <b>23-Jul-2019 21:29</b>							
Client ID:	Run ID: <b>ICPMS05_342915</b>	SeqNo: <b>5177609</b>	PrepDate: <b>23-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.04867 0.00200 0.05 0 97.3 80 - 120

<b>MS</b>	Sample ID: <b>HS19070566-06MS</b>	Units: <b>mg/L</b>	Analysis Date: <b>23-Jul-2019 21:36</b>							
Client ID:	Run ID: <b>ICPMS05_342915</b>	SeqNo: <b>5177612</b>	PrepDate: <b>23-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.04953 0.00200 0.05 0.001142 96.8 80 - 120

<b>MSD</b>	Sample ID: <b>HS19070566-06MSD</b>	Units: <b>mg/L</b>	Analysis Date: <b>23-Jul-2019 21:38</b>							
Client ID:	Run ID: <b>ICPMS05_342915</b>	SeqNo: <b>5177613</b>	PrepDate: <b>23-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.04774 0.00200 0.05 0.001142 93.2 80 - 120 0.04953 3.68 20

<b>SD</b>	Sample ID: <b>HS19070566-06SD</b>	Units: <b>mg/L</b>	Analysis Date: <b>23-Jul-2019 21:33</b>							
Client ID:	Run ID: <b>ICPMS05_342915</b>	SeqNo: <b>5177611</b>	PrepDate: <b>23-Jul-2019</b> DF: <b>5</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual

Arsenic U 0.0100 0.001142 0 10

<b>The following samples were analyzed in this batch:</b>	HS19070864-01	HS19070864-02	HS19070864-03	HS19070864-04
	HS19070864-05	HS19070864-06		



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

Batch ID: 143325 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A						
<b>MBLK</b>	Sample ID: <b>MBLK-143325</b>	Units: <b>mg/L</b>		Analysis Date: <b>25-Jul-2019 12:47</b>						
Client ID:		Run ID: <b>ICPMS05_343031</b>	SeqNo: <b>5180640</b>	PrepDate: <b>24-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-143325</b>	Units: <b>mg/L</b>		Analysis Date: <b>25-Jul-2019 12:49</b>						
Client ID:		Run ID: <b>ICPMS05_343031</b>	SeqNo: <b>5180641</b>	PrepDate: <b>24-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.04998	0.00200	0.05	0	100.0	80 - 120				
<b>MS</b>	Sample ID: <b>HS19070864-10MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>25-Jul-2019 13:01</b>						
Client ID: <b>WG-1620-MW28C-20190716</b>		Run ID: <b>ICPMS05_343031</b>	SeqNo: <b>5180628</b>	PrepDate: <b>24-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.05185	0.00200	0.05	0.000456	103	80 - 120				
<b>MSD</b>	Sample ID: <b>HS19070864-10MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jul-2019 15:58</b>						
Client ID: <b>WG-1620-MW28C-20190716</b>		Run ID: <b>ICPMS05_342951</b>	SeqNo: <b>5179479</b>	PrepDate: <b>24-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.05134	0.00200	0.05	0.000456	102	80 - 120	0.05087	0.902	20	
<b>PDS</b>	Sample ID: <b>HS19070864-10PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jul-2019 16:00</b>						
Client ID: <b>WG-1620-MW28C-20190716</b>		Run ID: <b>ICPMS05_342951</b>	SeqNo: <b>5179480</b>	PrepDate: <b>24-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.1061	0.00200	0.1	0.000456	106	75 - 125				
<b>SD</b>	Sample ID: <b>HS19070864-10SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jul-2019 15:53</b>						
Client ID: <b>WG-1620-MW28C-20190716</b>		Run ID: <b>ICPMS05_342951</b>	SeqNo: <b>5179477</b>	PrepDate: <b>24-Jul-2019</b>	DF: <b>5</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual	
Arsenic	U	0.0100					0.000456	0	10	

The following samples were analyzed in this batch: HS19070864-07 HS19070864-08 HS19070864-09 HS19070864-10  
 HS19070864-11 HS19070864-12

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

Batch ID: 143194 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143194	Units: ug/L			Analysis Date: 22-Jul-2019 11:00					
Client ID:	Run ID: SV-7_342875	SeqNo: 5176674		PrepDate: 19-Jul-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.914</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.077</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>81.5</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.22</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>84.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.338</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.8</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.575</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.5</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.538</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

Batch ID: 143194 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143194	Units: ug/L			Analysis Date: 22-Jul-2019 11:20					
Client ID:	Run ID: SV-7_342875	SeqNo: 5176675		PrepDate: 19-Jul-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.448	0.20	5	0	89.0	39 - 127				
2,4-Dimethylphenol	4.065	0.20	5	0	81.3	35 - 120				
2,4-Dinitrotoluene	5.282	0.20	5	0	106	50 - 122				
2,6-Dinitrotoluene	4.965	0.20	5	0	99.3	50 - 120				
2-Chloronaphthalene	4.463	0.20	5	0	89.3	50 - 120				
2-Methylnaphthalene	4.445	0.10	5	0	88.9	50 - 120				
4,6-Dinitro-2-methylphenol	4.939	0.20	5	0	98.8	25 - 121				
4-Nitrophenol	5.208	1.0	5	0	104	30 - 130				
Acenaphthene	4.419	0.10	5	0	88.4	45 - 120				
Acenaphthylene	4.131	0.10	5	0	82.6	47 - 120				
Anthracene	4.283	0.10	5	0	85.7	45 - 120				
Benz(a)anthracene	4.742	0.10	5	0	94.8	40 - 120				
Benzo(a)pyrene	5.119	0.10	5	0	102	45 - 120				
Bis(2-chloroethoxy)methane	4.08	0.20	5	0	81.6	45 - 120				
Bis(2-ethylhexyl)phthalate	4.229	0.20	5	0	84.6	40 - 139				
Chrysene	4.691	0.10	5	0	93.8	43 - 120				
Dibenzofuran	4.419	0.10	5	0	88.4	50 - 120				
Di-n-butyl phthalate	4.404	0.20	5	0	88.1	45 - 123				
Fluoranthene	4.651	0.10	5	0	93.0	45 - 125				
Fluorene	4.206	0.10	5	0	84.1	49 - 120				
Naphthalene	4.123	0.10	5	0	82.5	45 - 120				
Nitrobenzene	4.585	0.20	5	0	91.7	44 - 120				
N-Nitrosodiphenylamine	4.889	0.20	5	0	97.8	40 - 125				
Pentachlorophenol	4.714	0.20	5	0	94.3	19 - 121				
Phenanthrene	4.063	0.10	5	0	81.3	45 - 121				
Phenol	3.959	0.20	5	0	79.2	20 - 124				
Pyrene	4.472	0.10	5	0	89.4	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.604</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>112</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.713</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.535</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.7</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.418</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.4</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.682</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.612</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.2</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

Batch ID: 143194 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS19070864-10MS		Units: ug/L		Analysis Date: 22-Jul-2019 14:19				
Client ID: WG-1620-MW28C-20190716		Run ID: SV-7_342875		SeqNo: 5176679		PrepDate: 19-Jul-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.693	0.20	5	0	73.9	39 - 127				
2,4-Dimethylphenol	3.397	0.20	5	0	67.9	35 - 120				
2,4-Dinitrotoluene	4.804	0.20	5	0	96.1	50 - 122				
2,6-Dinitrotoluene	4.2	0.20	5	0	84.0	50 - 120				
2-Chloronaphthalene	3.834	0.20	5	0	76.7	50 - 120				
2-Methylnaphthalene	3.758	0.10	5	0	75.2	50 - 120				
4,6-Dinitro-2-methylphenol	4.56	0.20	5	0	91.2	25 - 121				
4-Nitrophenol	4.615	1.0	5	0	92.3	30 - 130				
Acenaphthene	3.72	0.10	5	0	74.4	45 - 120				
Acenaphthylene	3.54	0.10	5	0	70.8	47 - 120				
Anthracene	3.902	0.10	5	0	78.0	45 - 120				
Benz(a)anthracene	4.637	0.10	5	0	92.7	40 - 120				
Benzo(a)pyrene	5.199	0.10	5	0	104	45 - 120				
Bis(2-chloroethoxy)methane	3.411	0.20	5	0	68.2	45 - 120				
Bis(2-ethylhexyl)phthalate	4.294	0.20	5	0.1317	83.2	40 - 139				
Chrysene	4.71	0.10	5	0	94.2	43 - 120				
Dibenzofuran	3.788	0.10	5	0	75.8	50 - 120				
Di-n-butyl phthalate	4.444	0.20	5	0	88.9	45 - 123				
Fluoranthene	4.84	0.10	5	0.01464	96.5	45 - 125				
Fluorene	3.733	0.10	5	0	74.7	49 - 120				
Naphthalene	3.498	0.10	5	0.09505	68.1	45 - 120				
Nitrobenzene	4.018	0.20	5	0	80.4	44 - 120				
N-Nitrosodiphenylamine	4.429	0.20	5	0	88.6	40 - 125				
Pentachlorophenol	4.657	0.20	5	0	93.1	19 - 121				
Phenanthrene	3.943	0.10	5	0	78.9	45 - 121				
Phenol	4.038	0.20	5	2.478	31.2	20 - 124				
Pyrene	4.3	0.10	5	0.01185	85.8	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.507</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>110</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.746</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.9</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.67</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.25</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.0</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.838</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.8</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.934</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.7</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

Batch ID: 143194 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS19070864-10MSD	Units: ug/L			Analysis Date: 22-Jul-2019 14:39					
Client ID: WG-1620-MW28C-20190716	Run ID: SV-7_342875	SeqNo: 5176680	PrepDate: 19-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.672	0.20	5	0	73.4	39 - 127	3.693	0.547	20	
2,4-Dimethylphenol	3.328	0.20	5	0	66.6	35 - 120	3.397	2.04	20	
2,4-Dinitrotoluene	5.13	0.20	5	0	103	50 - 122	4.804	6.56	20	
2,6-Dinitrotoluene	4.241	0.20	5	0	84.8	50 - 120	4.2	0.976	20	
2-Chloronaphthalene	3.063	0.20	5	0	61.3	50 - 120	3.834	22.4	20	R
2-Methylnaphthalene	3.431	0.10	5	0	68.6	50 - 120	3.758	9.09	20	
4,6-Dinitro-2-methylphenol	4.711	0.20	5	0	94.2	25 - 121	4.56	3.25	30	
4-Nitrophenol	4.266	1.0	5	0	85.3	30 - 130	4.615	7.87	20	
Acenaphthene	3.557	0.10	5	0	71.1	45 - 120	3.72	4.47	20	
Acenaphthylene	3.316	0.10	5	0	66.3	47 - 120	3.54	6.53	20	
Anthracene	3.912	0.10	5	0	78.2	45 - 120	3.902	0.239	20	
Benz(a)anthracene	4.648	0.10	5	0	93.0	40 - 120	4.637	0.248	20	
Benzo(a)pyrene	5.69	0.10	5	0	114	45 - 120	5.199	9.02	20	
Bis(2-chloroethoxy)methane	3.137	0.20	5	0	62.7	45 - 120	3.411	8.37	20	
Bis(2-ethylhexyl)phthalate	4.518	0.20	5	0.1317	87.7	40 - 139	4.294	5.09	20	
Chrysene	5.022	0.10	5	0	100	43 - 120	4.71	6.41	20	
Dibenzofuran	3.658	0.10	5	0	73.2	50 - 120	3.788	3.49	20	
Di-n-butyl phthalate	4.461	0.20	5	0	89.2	45 - 123	4.444	0.39	20	
Fluoranthene	4.861	0.10	5	0.01464	96.9	45 - 125	4.84	0.439	20	
Fluorene	3.634	0.10	5	0	72.7	49 - 120	3.733	2.7	20	
Naphthalene	3.326	0.10	5	0.09505	64.6	45 - 120	3.498	5.04	20	
Nitrobenzene	3.644	0.20	5	0	72.9	44 - 120	4.018	9.78	20	
N-Nitrosodiphenylamine	4.775	0.20	5	0	95.5	40 - 125	4.429	7.52	20	
Pentachlorophenol	4.782	0.20	5	0	95.6	19 - 121	4.657	2.66	20	
Phenanthrene	4.08	0.10	5	0	81.6	45 - 121	3.943	3.41	20	
Phenol	3.46	0.20	5	2.478	19.6	20 - 124	4.038	15.4	20	S
Pyrene	4.476	0.10	5	0.01185	89.3	40 - 130	4.3	4	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.092</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>34 - 129</i>	<i>5.507</i>	<i>7.83</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.506</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>70.1</i>	<i>40 - 125</i>	<i>3.746</i>	<i>6.6</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.369</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>67.4</i>	<i>20 - 120</i>	<i>3.67</i>	<i>8.54</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.3</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.0</i>	<i>40 - 135</i>	<i>4.25</i>	<i>1.16</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.541</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>70.8</i>	<i>41 - 120</i>	<i>3.838</i>	<i>8.05</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.444</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>68.9</i>	<i>20 - 120</i>	<i>3.934</i>	<i>13.3</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

**Batch ID:** 143194 ( 0 )      **Instrument:** SV-7      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

The following samples were analyzed in this batch:

HS19070864-01	HS19070864-02	HS19070864-03	HS19070864-04
HS19070864-05	HS19070864-06	HS19070864-07	HS19070864-08
HS19070864-09	HS19070864-10	HS19070864-11	HS19070864-12

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

**Batch ID:** R342655 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190718</b>			Units: <b>ug/L</b>		Analysis Date: <b>18-Jul-2019 12:21</b>			
Client ID:		Run ID: <b>VOA2_342655</b>			SeqNo: <b>5171721</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Vinyl chloride	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>43.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>87.1</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.0</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.0</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>51.34</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-190718</b>			Units: <b>ug/L</b>		Analysis Date: <b>18-Jul-2019 11:57</b>			
Client ID:		Run ID: <b>VOA2_342655</b>			SeqNo: <b>5171720</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	19.53	1.0	20	0	97.7	70 - 124				
Benzene	19.25	1.0	20	0	96.2	74 - 120				
Chlorobenzene	19.1	1.0	20	0	95.5	76 - 113				
Ethylbenzene	19.59	1.0	20	0	97.9	77 - 117				
Methylene chloride	21.04	2.0	20	0	105	70 - 127				
Toluene	21.75	1.0	20	0	109	77 - 118				
Vinyl chloride	21.23	1.0	20	0	106	70 - 130				
Xylenes, Total	63.51	1.0	60	0	106	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.6</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>46.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.8</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

**Batch ID:** R342655 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS19070864-10MS			Units: ug/L		Analysis Date: 18-Jul-2019 13:59			
Client ID: WG-1620-MW28C-20190716		Run ID: VOA2_342655			SeqNo: 5171839		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.04	1.0	20	0	80.2	70 - 127				
Benzene	17.2	1.0	20	0	86.0	70 - 127				
Chlorobenzene	17.15	1.0	20	0	85.7	70 - 114				
Ethylbenzene	17.86	1.0	20	0	89.3	70 - 124				
Methylene chloride	17.39	2.0	20	0	86.9	70 - 128				
Toluene	19.38	1.0	20	0	96.9	70 - 123				
Vinyl chloride	17.32	1.0	20	0	86.6	70 - 130				
Xylenes, Total	56.86	1.0	60	0	94.8	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.22</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>90.4</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>46.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.8</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.33</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

MSD		Sample ID: HS19070864-10MSD			Units: ug/L		Analysis Date: 18-Jul-2019 14:23			
Client ID: WG-1620-MW28C-20190716		Run ID: VOA2_342655			SeqNo: 5171840		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.85	1.0	20	0	84.3	70 - 127	16.04	4.93	20	
Benzene	16.52	1.0	20	0	82.6	70 - 127	17.2	3.98	20	
Chlorobenzene	16.89	1.0	20	0	84.5	70 - 114	17.15	1.5	20	
Ethylbenzene	17.1	1.0	20	0	85.5	70 - 124	17.86	4.3	20	
Methylene chloride	17	2.0	20	0	85.0	70 - 128	17.39	2.26	20	
Toluene	18.9	1.0	20	0	94.5	70 - 123	19.38	2.49	20	
Vinyl chloride	17.85	1.0	20	0	89.3	70 - 130	17.32	3.02	20	
Xylenes, Total	56.52	1.0	60	0	94.2	70 - 130	56.86	0.595	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.82</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.6</i>	<i>70 - 126</i>	<i>45.22</i>	<i>1.33</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>81 - 113</i>	<i>47.86</i>	<i>1.62</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>77 - 123</i>	<i>46.88</i>	<i>1.95</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>50.33</i>	<i>0.435</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070864-01    HS19070864-04    HS19070864-06    HS19070864-10



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

<b>Batch ID:</b> R342706 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-190718</b>	Units: <b>ug/L</b>			Analysis Date: <b>18-Jul-2019 23:24</b>				
Client ID:	Run ID: <b>VOA2_342706</b>	SeqNo: <b>5172917</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>44.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>89.6</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.2</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.99</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-190718</b>	Units: <b>ug/L</b>			Analysis Date: <b>18-Jul-2019 22:59</b>				
Client ID:	Run ID: <b>VOA2_342706</b>	SeqNo: <b>5172916</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	19.09	1.0	20	0	95.5	70 - 124			
Benzene	18.84	1.0	20	0	94.2	74 - 120			
Chlorobenzene	18.95	1.0	20	0	94.8	76 - 113			
Ethylbenzene	18.88	1.0	20	0	94.4	77 - 117			
Methylene chloride	20.56	2.0	20	0	103	70 - 127			
Toluene	21.32	1.0	20	0	107	77 - 118			
Vinyl chloride	20.66	1.0	20	0	103	70 - 130			
Xylenes, Total	61.87	1.0	60	0	103	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.4</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.3</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>47.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QC BATCH REPORT**

**Batch ID:** R342706 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS19070864-08MS			Units: ug/L		Analysis Date: 19-Jul-2019 02:39			
Client ID: WG-1620-MW36B-20190716		Run ID: VOA2_342706			SeqNo: 5172921		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.7	1.0	20	0	78.5	70 - 127				
Benzene	16.86	1.0	20	0	84.3	70 - 127				
Chlorobenzene	16.83	1.0	20	0	84.2	70 - 114				
Ethylbenzene	17.18	1.0	20	0	85.9	70 - 124				
Methylene chloride	15.53	2.0	20	0	77.6	70 - 128				
Toluene	19.17	1.0	20	0	95.9	70 - 123				
Vinyl chloride	18.84	1.0	20	0	94.2	70 - 130				
Xylenes, Total	55.11	1.0	60	0	91.8	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.7</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.0</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.11</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.2</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 127</i>				

MSD		Sample ID: HS19070864-08MSD			Units: ug/L		Analysis Date: 19-Jul-2019 03:04			
Client ID: WG-1620-MW36B-20190716		Run ID: VOA2_342706			SeqNo: 5172922		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.2	1.0	20	0	81.0	70 - 127	15.7	3.11	20	
Benzene	16.39	1.0	20	0	82.0	70 - 127	16.86	2.81	20	
Chlorobenzene	16.02	1.0	20	0	80.1	70 - 114	16.83	4.92	20	
Ethylbenzene	16.88	1.0	20	0	84.4	70 - 124	17.18	1.79	20	
Methylene chloride	16.85	2.0	20	0	84.3	70 - 128	15.53	8.19	20	
Toluene	18.14	1.0	20	0	90.7	70 - 123	19.17	5.55	20	
Vinyl chloride	18.13	1.0	20	0	90.6	70 - 130	18.84	3.87	20	
Xylenes, Total	53.8	1.0	60	0	89.7	70 - 130	55.11	2.4	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.6</i>	<i>70 - 126</i>	<i>45.83</i>	<i>1.04</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>81 - 113</i>	<i>48</i>	<i>1.18</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>77 - 123</i>	<i>48.11</i>	<i>1.13</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 127</i>	<i>49.8</i>	<i>0.0359</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS19070864-02	HS19070864-03	HS19070864-05	HS19070864-07
HS19070864-08	HS19070864-09	HS19070864-11	HS19070864-12
HS19070864-13			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070864

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kansas	E-10352 2018-2019	31-Jul-2019
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070864

**SAMPLE TRACKING**

Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS19070864-01	WG-1620-MW21C-20190716	Login	17/07/2019 18:07:39	RPG	MET074
HS19070864-01	WG-1620-MW21C-20190716	Login	17/07/2019 18:07:39	RPG	EXT011
HS19070864-01	WG-1620-MW21C-20190716	Login	17/07/2019 18:07:39	RPG	VOA153
HS19070864-02	WG-1620-FD01-20190716	Login	17/07/2019 18:07:39	RPG	MET074
HS19070864-02	WG-1620-FD01-20190716	Login	17/07/2019 18:07:39	RPG	EXT011
HS19070864-02	WG-1620-FD01-20190716	Login	17/07/2019 18:07:39	RPG	VOA153
HS19070864-03	WG-1620-MW62B-20190716	Login	17/07/2019 18:07:39	RPG	MET074
HS19070864-03	WG-1620-MW62B-20190716	Login	17/07/2019 18:07:39	RPG	EXT011
HS19070864-03	WG-1620-MW62B-20190716	Login	17/07/2019 18:07:39	RPG	VOA153
HS19070864-04	WG-1620-MW53C-20190716	Login	17/07/2019 18:07:39	RPG	MET074
HS19070864-04	WG-1620-MW53C-20190716	Login	17/07/2019 18:07:39	RPG	EXT012
HS19070864-04	WG-1620-MW53C-20190716	Login	17/07/2019 18:07:39	RPG	VOA153
HS19070864-05	WG-1620-MW54C-20190716	Login	17/07/2019 18:07:39	RPG	MET074
HS19070864-05	WG-1620-MW54C-20190716	Login	17/07/2019 18:07:39	RPG	EXT012
HS19070864-05	WG-1620-MW54C-20190716	Login	17/07/2019 18:07:39	RPG	VOA153
HS19070864-13	WG-1620-TB04-20190716	Login	17/07/2019 18:07:39	RPG	VOA153

**Sample Receipt Checklist**

Client Name: PBW  
 Work Order: HS19070864

Date/Time Received: **17-Jul-2019 14:24**  
 Received by: **AC**

Checklist completed by: Raegen Giga 17-Jul-2019  
 eSignature Date

Reviewed by: Dane J. Wacasey 18-Jul-2019  
 eSignature Date

Matrices: **Water**

Carrier name: **Client**

- Shipping container/cooler in good condition? Yes  No  Not Present
- Custody seals intact on shipping container/cooler? Yes  No  Not Present
- Custody seals intact on sample bottles? Yes  No  Not Present
- VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes  No  Not Present
- Chain of custody present? Yes  No  1 Page(s)
- Chain of custody signed when relinquished and received? Yes  No  COC IDs:196141/196139
- Samplers name present on COC? Yes  No
- Chain of custody agrees with sample labels? Yes  No
- Samples in proper container/bottle? Yes  No
- Sample containers intact? Yes  No
- Sufficient sample volume for indicated test? Yes  No
- All samples received within holding time? Yes  No
- Container/Temp Blank temperature in compliance? Yes  No
- Temperature(s)/Thermometer(s): 0.2c , 0.7c , 2.2c UC/C IR25
- Cooler(s)/Kit(s): 44161/44877/45083
- Date/Time sample(s) sent to storage: 07/17/2019 18:40
- Water - VOA vials have zero headspace? Yes  No  No VOA vials submitted
- Water - pH acceptable upon receipt? Yes  No  N/A
- pH adjusted? Yes  No  N/A
- pH adjusted by:

Login Notes: Vial # 3 ID Differs: COC = WG-1620-MW36A-20190716; Label = WG-1620-MW36B-20190716; Logged per COC using vial #1 and vial #2

Client Contacted: \_\_\_\_\_ Date Contacted: \_\_\_\_\_ Person Contacted: \_\_\_\_\_  
 Contacted By: \_\_\_\_\_ Regarding: \_\_\_\_\_  
 Comments:   
 Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page 1 of 2

COC ID: 196141

Houston, TX  
+1 281 530 5656

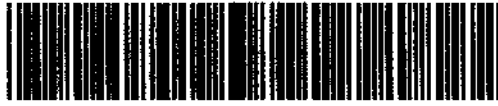
Spring City, PA  
+1 610 948 4903

South Charleston, WV  
+1 304 356 3168

Middletown, PA  
+1 717 944 5541

Salt Lake City, UT  
+1 801 266 7700

York, PA  
+1 717 505 5280

Customer Information		ALS Project Manager:		ALS Work Order #:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	<p><b>HS19070864</b></p> <p>Golder Associates Inc.</p> <p>Houston TX-Wood Preserving Works</p> 
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-T00-20190716</del>			Water	←	6		*									
2	WG-1620-MW21C-20190716	7-16-19	0735	W		6	X		X	X							
3	WG-1620-FD01-20190716		0735	W		6	X		X	X							
4	WG-1620-MW62B-20190716		0825	W		6	X		X	X							
5	WG-1620-MW53C-20190716		1015	W		6	X		X	X							
6	WG-1620-MW54C-20190716		1115	W		6	X		X	X							
7	WG-1620-MW25C-20190716		1215	W		6		X	X	X							
8	WG-1620-MW25A-20190716		1310	W		6		X	X	X							
9	WG-1620-MW36B-20190716		1420	W		6		X	X	X							
10	WG-1620-MW36A-20190716		1520	W		6		X	X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>	Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hrs	Results Due Date:
Relinquished by: <b>John Br</b> Date: <b>7-17-19</b> Time: <b>14:24</b>	Received by: <b>AC</b>	Notes: UPRR Houston MWPW	
Relinquished by: <b>John Br</b> Date: <b>7-17-19</b> Time: <b>14:24</b>	Received by (Laboratory): <b>AC</b>	QC Package: (Check One Box Below)	
Logged by (Laboratory):	Checked by (Laboratory):	Cooler ID: <b>44161</b> Cooler Temp: <b>0.2</b>	<input type="checkbox"/> Level II Std OC <input checked="" type="checkbox"/> TRRP Checklist
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035		Cooler ID: <b>44877</b> Cooler Temp: <b>0.7</b>	<input type="checkbox"/> Level III Std OC/Raw Data <input type="checkbox"/> TRRP Level IV
		Cooler ID: <b>45083</b> Cooler Temp: <b>0.2</b>	<input type="checkbox"/> Level IV SW/MS/CLP <input type="checkbox"/> Other

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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# Chain of Custody Form

Page 2 of 2

COC ID: 196139

HS19070864

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A 8260_LL_W (5632528 Volatile Organics Site specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B 8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C 8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D ICP_TW (5636002 5652646 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E MS/MSD
	Suite 4004		Stop 0750	F
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G
Phone	(512) 671-3434	Phone		H
Fax	(512) 671-3446	Fax		I
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TB0-20190716</del>			Water	-	2		*									
2	WG-1620-MW28C-20190716	7-16-19	1615	W		12	X		X	X	X						
3	WG-1620-MW28A-20190716		1700	W		6	X		X	X							
4	WG-1620-MW63B-20190716		1755	W		6	X		X	X							
5	WG-1620-TB04-20190716		-	W		2											
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <b>JOHN PRAY</b>	Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour	Results Due Date:
Relinquished by: <b>John B</b>	Date: <b>7-17-19</b> Time: <b>14:24</b>	Received by: <b>AC</b>	Notes: <b>UPRR Houston MWPW</b>
Relinquished by: <b>John B</b>	Date: <b>7-17-19</b> Time: <b>14:24</b>	Received by (Laboratory): <b>AC</b>	QC Package: (Check One Box Below) <input type="checkbox"/> Level II Std QC <input checked="" type="checkbox"/> TRRP Checklist <input type="checkbox"/> TRRP Level IV <input type="checkbox"/> Level III Std QCRaw Data <input type="checkbox"/> Level IV SM/MS/CLP <input type="checkbox"/> Other
Logged by (Laboratory):	Date: Time:	Checked by (Laboratory):	
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035			

- Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
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July 31, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19070994**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 34 sample(s) on Jul 19, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL

Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

**Laboratory Review Checklist: Reportable Data**

Laboratory Name: ALS Laboratory Group		LRC Date: 07/31/2019					
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS19071233					
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 143239,143295,143526,143558,R342868,R342877,R342879					
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			3
		Were MS/MSD RPDs within laboratory QC limits?		X			4
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?	X				
		Were analytical duplicates analyzed at the appropriate frequency?	X				
		Were RPDs or relative standard deviations within the laboratory QC limits?	X				
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				5
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 07/31/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19071233			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143239,143295,143526,143558,R342868,R342877,R342879			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 07/31/2019
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS19071233
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 143239,143295,143526,143558,R342868,R342877,R342879

ER# <sup>5</sup>	Description
1	Volatile Organics Method SW8260, sample WG-1620-MW71B-20190717, WG-1620-MW68B-20190718, WG-1620-FD03-20190718, WG-1620-MW83B-20190718, WG-1620-MW35B-20190718, the surrogate recoveries could not be determined due to dilution below the calibration range.
2	Batch 143295, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.
3	Batch 143239, Semivolatile Organics Method SW8270, sample WG-1620-MW61A-20190717, MSD recovered outside the control limit for Benzo(a)pyrene due to possible matrix effect.  Batch R342877, Volatile Organics Method SW8260, sample WG-1620-MW83C-20190718, MS and MSD recovered outside the control limit for surrogate 4-Bromofluorobenzene due to possible matrix effect.  Batch R342879, Volatile Organics Method SW8260, sample WG-1620-MW61A-20190717, MSD recovered below the control limit for 1,2-Dichloroethane due to possible matrix effect.
4	Batch 143239, Semivolatile Organics Method SW8270, sample WG-1620-MW61A-20190717, MS/MSD RPD recovered above the RPD limits for 2,6-Dinitrotoluene, Benzo(a)pyrene and surrogates 2,4,6-Tribromophenol, 2-Fluorobiphenyl due to possible matrix effect.
5	Batch 143239, Semivolatile Organics Method SW8270, samples WG-1620-MW68B-20190718, WG-1620-FD03-20190718, the GCMS semi-volatile extract of the samples were run at a dilution due to a high level of matrix interference.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070994

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19070994-01	WG-1620-FB04-20190716	Water		16-Jul-2019 18:20	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-02	WG-1620-MW61A-20190717	Water		17-Jul-2019 07:50	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-03	WG-1620-MW60A-20190717	Water		17-Jul-2019 08:35	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-04	WG-1620-MW69A-20190717	Water		17-Jul-2019 09:35	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-05	WG-1620-MW47C-20190717	Water		17-Jul-2019 10:25	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-06	WG-1620-MW48C-20190717	Water		17-Jul-2019 11:20	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-07	WG-1620-MW59A-20190717	Water		17-Jul-2019 12:20	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-08	WG-1620-MW59B-20190717	Water		17-Jul-2019 13:10	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-09	WG-1620-MW44A-20190717	Water		17-Jul-2019 14:00	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-10	WG-1620-MW87C-20190717	Water		17-Jul-2019 15:40	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-11	WG-1620-MW71B-20190717	Water		17-Jul-2019 16:40	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-12	WG-1620-MW33A-20190717	Water		17-Jul-2019 17:50	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-13	WG-1620-FD02-20190717	Water		17-Jul-2019 17:50	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-14	WG-1620-MW26A-20190717	Water		17-Jul-2019 18:45	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-15	WG-1620-FB05-20190717	Water		17-Jul-2019 19:05	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-16	WG-1620-MW68A-20190718	Water		18-Jul-2019 07:55	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-17	WG-1620-MW68B-20190718	Water		18-Jul-2019 08:55	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-18	WG-1620-FD03-20190718	Water		18-Jul-2019 08:55	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-19	WG-1620-MW68C-20190718	Water		18-Jul-2019 09:45	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-20	WG-1620-MW83B-20190718	Water		18-Jul-2019 10:40	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-21	WG-1620-MW83C-20190718	Water		18-Jul-2019 11:30	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-22	WG-1620-MW35A-20190718	Water		18-Jul-2019 12:40	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-23	WG-1620-MW35B-20190718	Water		18-Jul-2019 13:30	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-24	WG-1620-MW90B-20190718	Water		18-Jul-2019 14:25	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-25	WG-1620-MW89B-20190718	Water		18-Jul-2019 15:15	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-26	WG-1620-MW38B-20190718	Water		18-Jul-2019 16:10	19-Jul-2019 13:15	<input type="checkbox"/>

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19070994

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19070994-27	WG-1620-MW27A-20190718	Water		18-Jul-2019 17:15	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-28	WG-1620-MW27C-20190718	Water		18-Jul-2019 18:05	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-29	WG-1620-MW51A-20190719	Water		19-Jul-2019 07:25	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-30	WG-1620-MW51C-20190719	Water		19-Jul-2019 08:55	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-31	WG-1620-MW81B-20190719	Water		19-Jul-2019 09:45	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-32	WG-1620-MW50A-20190719	Water		19-Jul-2019 10:40	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-33	WG-1620-TB05-20190719	Water	C&G-062119-82	19-Jul-2019 00:00	19-Jul-2019 13:15	<input type="checkbox"/>
HS19070994-34	WG-1620-TB06-20190719	Water	C&G-062119-82	19-Jul-2019 00:00	19-Jul-2019 13:15	<input type="checkbox"/>



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB04-20190716  
 Collection Date: 16-Jul-2019 18:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 00:42
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 00:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 00:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 00:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 00:42
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 00:42
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 00:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 00:42
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>102</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 00:42</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.6</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 00:42</i>
<i>Surr: Dibromofluoromethane</i>		<i>104</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 00:42</i>
<i>Surr: Toluene-d8</i>		<i>98.6</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 00:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB04-20190716  
 Collection Date: 16-Jul-2019 18:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 11:41
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 11:41
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 11:41
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 11:41
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 11:41
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 11:41
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 11:41
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 11:41
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 11:41
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 11:41
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 11:41
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 11:41
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 11:41
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 11:41
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00045</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 11:41
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 11:41
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 11:41
<b>Di-n-butyl phthalate</b>	<b>0.000043</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 11:41
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 11:41
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 11:41
Naphthalene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 11:41
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 11:41
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 11:41
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 11:41
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 11:41
<b>Phenol</b>	<b>0.000084</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 11:41
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 11:41
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 11:41</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>72.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 11:41</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 11:41</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 11:41</i>
<i>Surr: Nitrobenzene-d5</i>	<i>70.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 11:41</i>
<i>Surr: Phenol-d6</i>	<i>70.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 11:41</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 21:42

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW61A-20190717  
 Collection Date: 17-Jul-2019 07:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:32
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:32
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 01:32
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 01:32
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 01:32
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:32
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:32
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 01:32
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>102</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:32</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>95.9</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:32</i>
<i>Surr: Dibromofluoromethane</i>		<i>107</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:32</i>
<i>Surr: Toluene-d8</i>		<i>99.0</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:32</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW61A-20190717  
 Collection Date: 17-Jul-2019 07:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 12:21
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 12:21
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 12:21
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 12:21
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 12:21
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 12:21
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 12:21
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 12:21
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 12:21
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 12:21
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 12:21
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 12:21
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 12:21
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 12:21
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00037</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 12:21
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 12:21
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 12:21
<b>Di-n-butyl phthalate</b>	<b>0.000056</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 12:21
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 12:21
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 12:21
Naphthalene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 12:21
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 12:21
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 12:21
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 12:21
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 12:21
<b>Phenol</b>	<b>0.00011</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 12:21
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 12:21
<i>Surr: 2,4,6-Tribromophenol</i>	<i>109</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 12:21</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>83.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 12:21</i>
<i>Surr: 2-Fluorophenol</i>	<i>74.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 12:21</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>99.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 12:21</i>
<i>Surr: Nitrobenzene-d5</i>	<i>84.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 12:21</i>
<i>Surr: Phenol-d6</i>	<i>83.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 12:21</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00117</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 21:26

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW60A-20190717  
 Collection Date: 17-Jul-2019 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:10
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:10
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:10
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:10
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 03:10
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:10
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:10
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:10
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>102</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:10</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>96.1</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:10</i>
<i>Surr: Dibromofluoromethane</i>		<i>104</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:10</i>
<i>Surr: Toluene-d8</i>		<i>101</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:10</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW60A-20190717  
 Collection Date: 17-Jul-2019 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 18:57
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 18:57
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 18:57
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 18:57
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 18:57
<b>2-Methylnaphthalene</b>	<b>0.000072</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 18:57</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 18:57
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 18:57
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 18:57
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 18:57
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 18:57
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 18:57
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 18:57
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 18:57
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00055</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 18:57</b>
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 18:57
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 18:57
<b>Di-n-butyl phthalate</b>	<b>0.000075</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 18:57</b>
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 18:57
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 18:57
<b>Naphthalene</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 18:57</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 18:57
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 18:57
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 18:57
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 18:57
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 18:57
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 18:57
<i>Surr: 2,4,6-Tribromophenol</i>	<i>96.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 18:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>71.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 18:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 18:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>75.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 18:57</i>
<i>Surr: Nitrobenzene-d5</i>	<i>79.7</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 18:57</i>
<i>Surr: Phenol-d6</i>	<i>70.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 18:57</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000440</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>30-Jul-2019 21:44</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW69A-20190717  
 Collection Date: 17-Jul-2019 09:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:35
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:35
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:35
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:35
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 03:35
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:35
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:35
<i>Surr: 1,2-Dichloroethane-d4</i>		102		70-126	%REC	1	23-Jul-2019 03:35
<i>Surr: 4-Bromofluorobenzene</i>		102		81-113	%REC	1	23-Jul-2019 03:35
<i>Surr: Dibromofluoromethane</i>		106		77-123	%REC	1	23-Jul-2019 03:35
<i>Surr: Toluene-d8</i>		97.9		82-127	%REC	1	23-Jul-2019 03:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW69A-20190717  
 Collection Date: 17-Jul-2019 09:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 19:17
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 19:17
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 19:17
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 19:17
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 19:17
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 19:17
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 19:17
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 19:17
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 19:17
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 19:17
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 19:17
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 19:17
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:17
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 19:17
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00034</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 19:17
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 19:17
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:17
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 19:17
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 19:17
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 19:17
<b>Naphthalene</b>	<b>0.000078</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 19:17
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 19:17
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 19:17
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 19:17
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 19:17
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 19:17
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 19:17
<i>Surr: 2,4,6-Tribromophenol</i>	<i>102</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>71.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:17</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>77.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:17</i>
<i>Surr: Nitrobenzene-d5</i>	<i>73.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:17</i>
<i>Surr: Phenol-d6</i>	<i>70.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:17</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000642</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 21:47

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW47C-20190717  
 Collection Date: 17-Jul-2019 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:59
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:59
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:59
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:59
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 03:59
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:59
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:59
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:59</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>106</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:59</i>
<i>Surr: Toluene-d8</i>	<i>97.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW47C-20190717  
 Collection Date: 17-Jul-2019 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 19:36
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 19:36
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 19:36
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 19:36
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 19:36
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 19:36
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 19:36
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 19:36
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 19:36
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 19:36
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 19:36
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 19:36
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:36
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 19:36
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00027</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 19:36
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 19:36
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:36
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 19:36
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 19:36
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 19:36
Naphthalene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:36
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 19:36
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 19:36
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 19:36
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 19:36
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 19:36
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 19:36
<i>Surr: 2,4,6-Tribromophenol</i>	<i>101</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:36</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>82.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:36</i>
<i>Surr: 2-Fluorophenol</i>	<i>75.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:36</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>80.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:36</i>
<i>Surr: Nitrobenzene-d5</i>	<i>91.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:36</i>
<i>Surr: Phenol-d6</i>	<i>75.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 19:36</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000440</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>30-Jul-2019 21:49</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW48C-20190717  
 Collection Date: 17-Jul-2019 11:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:24
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:24
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:24
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:24
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 04:24
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:24
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:24
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:24</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:24</i>
<i>Surr: Dibromofluoromethane</i>	<i>106</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:24</i>
<i>Surr: Toluene-d8</i>	<i>99.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:24</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW48C-20190717  
 Collection Date: 17-Jul-2019 11:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 19:56
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 19:56
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 19:56
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 19:56
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 19:56
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 19:56
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 19:56
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 19:56
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 19:56
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 19:56
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 19:56
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 19:56
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:56
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 19:56
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00030</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 19:56
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 19:56
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:56
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 19:56
<b>Fluoranthene</b>	<b>0.000035</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 19:56
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 19:56
Naphthalene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 19:56
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 19:56
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 19:56
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 19:56
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 19:56
<b>Phenol</b>	<b>0.000074</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 19:56
<b>Pyrene</b>	<b>0.000036</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 19:56
<i>Surr: 2,4,6-Tribromophenol</i>	93.3			34-129	%REC	1	22-Jul-2019 19:56
<i>Surr: 2-Fluorobiphenyl</i>	71.4			40-125	%REC	1	22-Jul-2019 19:56
<i>Surr: 2-Fluorophenol</i>	68.3			20-120	%REC	1	22-Jul-2019 19:56
<i>Surr: 4-Terphenyl-d14</i>	77.0			40-135	%REC	1	22-Jul-2019 19:56
<i>Surr: Nitrobenzene-d5</i>	74.2			41-120	%REC	1	22-Jul-2019 19:56
<i>Surr: Phenol-d6</i>	65.4			20-120	%REC	1	22-Jul-2019 19:56
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00167</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 21:51

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59A-20190717  
 Collection Date: 17-Jul-2019 12:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:48
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:48
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:48
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:48
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 04:48
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:48
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:48
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:48
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>101</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:48</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>101</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:48</i>
<i>Surr: Dibromofluoromethane</i>		<i>105</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:48</i>
<i>Surr: Toluene-d8</i>		<i>97.3</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:48</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59A-20190717  
 Collection Date: 17-Jul-2019 12:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 20:16
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 20:16
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 20:16
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 20:16
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 20:16
<b>2-Methylnaphthalene</b>	<b>0.000036</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 20:16</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 20:16
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 20:16
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 20:16
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 20:16
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 20:16
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 20:16
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 20:16
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 20:16
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00050</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 20:16</b>
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 20:16
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 20:16
<b>Di-n-butyl phthalate</b>	<b>0.000064</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 20:16</b>
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 20:16
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 20:16
<b>Naphthalene</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 20:16</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 20:16
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 20:16
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 20:16
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 20:16
<b>Phenol</b>	<b>0.000058</b>	<b>J</b>	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>22-Jul-2019 20:16</b>
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 20:16
<i>Surr: 2,4,6-Tribromophenol</i>	<i>102</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:16</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>83.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>83.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:16</i>
<i>Surr: Phenol-d6</i>	<i>75.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:16</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00455</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>30-Jul-2019 21:53</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59B-20190717  
 Collection Date: 17-Jul-2019 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:13
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:13
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 05:13
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:13
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>102</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>95.1</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>
<i>Surr: Dibromofluoromethane</i>		<i>104</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>
<i>Surr: Toluene-d8</i>		<i>99.1</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59B-20190717  
 Collection Date: 17-Jul-2019 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 20:36
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 20:36
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 20:36
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 20:36
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 20:36
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 20:36
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 20:36
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 20:36
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 20:36
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 20:36
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 20:36
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 20:36
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 20:36
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 20:36
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00039</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 20:36
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 20:36
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 20:36
<b>Di-n-butyl phthalate</b>	<b>0.000055</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 20:36
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 20:36
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 20:36
<b>Naphthalene</b>	<b>0.00015</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:36
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 20:36
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 20:36
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 20:36
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 20:36
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 20:36
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 20:36
<i>Surr: 2,4,6-Tribromophenol</i>	<i>109</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:36</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>70.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:36</i>
<i>Surr: 2-Fluorophenol</i>	<i>60.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:36</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>77.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:36</i>
<i>Surr: Nitrobenzene-d5</i>	<i>68.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:36</i>
<i>Surr: Phenol-d6</i>	<i>60.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:36</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000542</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 21:56

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44A-20190717  
 Collection Date: 17-Jul-2019 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:37
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 05:37
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:37
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>100</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>98.9</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>
<i>Surr: Dibromofluoromethane</i>		<i>104</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>
<i>Surr: Toluene-d8</i>		<i>96.0</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44A-20190717  
 Collection Date: 17-Jul-2019 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 20:56
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 20:56
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 20:56
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 20:56
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 20:56
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 20:56
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 20:56
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 20:56
<b>Acenaphthene</b>	<b>0.056</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 15:59
<b>Acenaphthylene</b>	<b>0.00028</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
<b>Anthracene</b>	<b>0.00089</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
<b>Benz(a)anthracene</b>	<b>0.000085</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 20:56
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 20:56
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00025</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
<b>Chrysene</b>	<b>0.000052</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 20:56
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 20:56
<b>Fluoranthene</b>	<b>0.0091</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
<b>Fluorene</b>	<b>0.014</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 15:59
<b>Naphthalene</b>	<b>0.00019</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 20:56
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 20:56
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 20:56
<b>Phenanthrene</b>	<b>0.00012</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
Phenol	U		0.000035	0.00020	mg/L	1	22-Jul-2019 20:56
<b>Pyrene</b>	<b>0.0057</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 20:56
<i>Surr: 2,4,6-Tribromophenol</i>	<i>103</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 15:59</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>80.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>48.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>61.1</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 15:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.0</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 15:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>55.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:56</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>72.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:56</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>87.2</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 15:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>64.4</i>			<i>41-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 15:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>59.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:56</i>
<i>Surr: Phenol-d6</i>	<i>53.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 20:56</i>
<i>Surr: Phenol-d6</i>	<i>62.1</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 15:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44A-20190717  
 Collection Date: 17-Jul-2019 14:00

**ANALYTICAL REPORT**

WorkOrder:HS19070994  
 Lab ID:HS19070994-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>	<b>Method:SW6020</b>			Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	0.0303		0.000400	0.00200	mg/L	1	30-Jul-2019 21:58

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW87C-20190717  
 Collection Date: 17-Jul-2019 15:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:02
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:02
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:02
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:02
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:02
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:02
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:02
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:02</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:02</i>
<i>Surr: Dibromofluoromethane</i>	<i>104</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:02</i>
<i>Surr: Toluene-d8</i>	<i>98.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:02</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW87C-20190717  
 Collection Date: 17-Jul-2019 15:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 21:16
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 21:16
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 21:16
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 21:16
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 21:16
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 21:16
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 21:16
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 21:16
Acenaphthene	U		0.000027	0.00010	mg/L	1	22-Jul-2019 21:16
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 21:16
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 21:16
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 21:16
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 21:16
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 21:16
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00018</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:16
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 21:16
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 21:16
<b>Di-n-butyl phthalate</b>	<b>0.00017</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:16
Fluoranthene	U		0.000010	0.00010	mg/L	1	22-Jul-2019 21:16
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 21:16
<b>Naphthalene</b>	<b>0.00017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:16
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 21:16
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 21:16
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 21:16
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 21:16
<b>Phenol</b>	<b>0.00013</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:16
Pyrene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 21:16
<i>Surr: 2,4,6-Tribromophenol</i>	<i>106</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>73.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>61.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:16</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>78.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>73.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:16</i>
<i>Surr: Phenol-d6</i>	<i>67.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:16</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 22:00

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW71B-20190717  
 Collection Date: 17-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:26
<b>Benzene</b>	<b>0.13</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:26
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:26
<b>Ethylbenzene</b>	<b>0.031</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:26
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:26
<b>Toluene</b>	<b>0.023</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:26
<b>Xylenes, Total</b>	<b>0.054</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:26
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:26</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:26</i>
<i>Surr: Dibromofluoromethane</i>	<i>104</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:26</i>
<i>Surr: Toluene-d8</i>	<i>98.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:26</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW71B-20190717  
 Collection Date: 17-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 21:36
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 21:36
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 21:36
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 21:36
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 21:36
<b>2-Methylnaphthalene</b>	<b>0.024</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 16:19
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 21:36
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 21:36
<b>Acenaphthene</b>	<b>0.016</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 16:19
<b>Acenaphthylene</b>	<b>0.00033</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Anthracene</b>	<b>0.0066</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Benz(a)anthracene</b>	<b>0.0026</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Benzo(a)pyrene</b>	<b>0.00083</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 21:36
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00014</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Chrysene</b>	<b>0.0024</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Dibenzofuran</b>	<b>0.016</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 16:19
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 21:36
<b>Fluoranthene</b>	<b>0.017</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 16:19
<b>Fluorene</b>	<b>0.0093</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Naphthalene</b>	<b>0.44</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	23-Jul-2019 16:39
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 21:36
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 21:36
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 21:36
<b>Phenanthrene</b>	<b>0.030</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 16:19
<b>Phenol</b>	<b>0.000051</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<b>Pyrene</b>	<b>0.0088</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:36
<i>Surr: 2,4,6-Tribromophenol</i>	96.3			34-129	%REC	1	22-Jul-2019 21:36
<i>Surr: 2,4,6-Tribromophenol</i>	113			34-129	%REC	10	23-Jul-2019 16:19
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	100	23-Jul-2019 16:39
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	100	23-Jul-2019 16:39
<i>Surr: 2-Fluorobiphenyl</i>	77.7			40-125	%REC	10	23-Jul-2019 16:19
<i>Surr: 2-Fluorobiphenyl</i>	58.7			40-125	%REC	1	22-Jul-2019 21:36
<i>Surr: 2-Fluorophenol</i>	64.3			20-120	%REC	1	22-Jul-2019 21:36
<i>Surr: 2-Fluorophenol</i>	90.1			20-120	%REC	10	23-Jul-2019 16:19
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	100	23-Jul-2019 16:39
<i>Surr: 4-Terphenyl-d14</i>	0	JS		40-135	%REC	100	23-Jul-2019 16:39
<i>Surr: 4-Terphenyl-d14</i>	88.9			40-135	%REC	10	23-Jul-2019 16:19
<i>Surr: 4-Terphenyl-d14</i>	74.0			40-135	%REC	1	22-Jul-2019 21:36

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW71B-20190717  
 Collection Date: 17-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	67.5			41-120	%REC	1	22-Jul-2019 21:36
Surr: Nitrobenzene-d5	72.8			41-120	%REC	10	23-Jul-2019 16:19
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	23-Jul-2019 16:39
Surr: Phenol-d6	71.5			20-120	%REC	10	23-Jul-2019 16:19
Surr: Phenol-d6	0	JS		20-120	%REC	100	23-Jul-2019 16:39
Surr: Phenol-d6	63.8			20-120	%REC	1	22-Jul-2019 21:36
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	0.000626	J	0.000400	0.00200	mg/L	1	30-Jul-2019 22:02

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33A-20190717  
 Collection Date: 17-Jul-2019 17:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:51
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:51
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:51
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:51
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:51
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:51
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:51
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:51</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:51</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:51</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33A-20190717  
 Collection Date: 17-Jul-2019 17:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jul-2019 21:55
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	22-Jul-2019 21:55
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jul-2019 21:55
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jul-2019 21:55
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jul-2019 21:55
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	22-Jul-2019 21:55
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jul-2019 21:55
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jul-2019 21:55
<b>Acenaphthene</b>	<b>0.000098</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:55
Acenaphthylene	U		0.000015	0.00010	mg/L	1	22-Jul-2019 21:55
Anthracene	U		0.000014	0.00010	mg/L	1	22-Jul-2019 21:55
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	22-Jul-2019 21:55
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	22-Jul-2019 21:55
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jul-2019 21:55
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00022</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:55
Chrysene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 21:55
Dibenzofuran	U		0.000020	0.00010	mg/L	1	22-Jul-2019 21:55
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	22-Jul-2019 21:55
<b>Fluoranthene</b>	<b>0.00016</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:55
Fluorene	U		0.000030	0.00010	mg/L	1	22-Jul-2019 21:55
<b>Naphthalene</b>	<b>0.000076</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:55
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jul-2019 21:55
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jul-2019 21:55
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	22-Jul-2019 21:55
Phenanthrene	U		0.000021	0.00010	mg/L	1	22-Jul-2019 21:55
<b>Phenol</b>	<b>0.00011</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jul-2019 21:55
<b>Pyrene</b>	<b>0.00028</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jul-2019 21:55
<i>Surr: 2,4,6-Tribromophenol</i>	<i>103</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:55</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>81.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:55</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:55</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>78.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:55</i>
<i>Surr: Nitrobenzene-d5</i>	<i>75.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:55</i>
<i>Surr: Phenol-d6</i>	<i>76.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jul-2019 21:55</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0155</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 22:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD02-20190717  
 Collection Date: 17-Jul-2019 17:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:16
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:16
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:16
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:16
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 07:16
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:16
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:16
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:16</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:16</i>
<i>Surr: Dibromofluoromethane</i>	<i>104</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:16</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:16</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD02-20190717  
 Collection Date: 17-Jul-2019 17:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 13:20
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 13:20
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 13:20
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 13:20
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 13:20
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 13:20
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 13:20
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 13:20
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 13:20
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 13:20
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 13:20
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 13:20
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 13:20
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 13:20
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000084</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 13:20
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 13:20
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 13:20
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 13:20
<b>Fluoranthene</b>	<b>0.00017</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 13:20
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 13:20
<b>Naphthalene</b>	<b>0.000096</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 13:20
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 13:20
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 13:20
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 13:20
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 13:20
<b>Phenol</b>	<b>0.000053</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 13:20
<b>Pyrene</b>	<b>0.00069</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 13:20
<i>Surr: 2,4,6-Tribromophenol</i>	<i>105</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:20</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:20</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>73.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:20</i>
<i>Surr: Nitrobenzene-d5</i>	<i>69.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:20</i>
<i>Surr: Phenol-d6</i>	<i>67.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:20</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0147</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 22:11

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW26A-20190717  
 Collection Date: 17-Jul-2019 18:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:40
<b>Benzene</b>	<b>0.00036</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 07:40
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:40
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:40
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 07:40
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:40
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:40
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:40</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:40</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:40</i>
<i>Surr: Toluene-d8</i>	<i>98.3</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW26A-20190717  
 Collection Date: 17-Jul-2019 18:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 13:40
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 13:40
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 13:40
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 13:40
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 13:40
<b>2-Methylnaphthalene</b>	<b>0.000059</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 13:40
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 13:40
<b>Acenaphthene</b>	<b>0.028</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	<b>10</b>	<b>23-Jul-2019 16:59</b>
<b>Acenaphthylene</b>	<b>0.00013</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
<b>Anthracene</b>	<b>0.00047</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 13:40
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 13:40
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 13:40
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00019</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 13:40
<b>Dibenzofuran</b>	<b>0.00021</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 13:40
<b>Fluoranthene</b>	<b>0.0029</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
<b>Fluorene</b>	<b>0.0017</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
<b>Naphthalene</b>	<b>0.00032</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 13:40
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 13:40
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 13:40
<b>Phenanthrene</b>	<b>0.000083</b>	<b>J</b>	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 13:40
<b>Pyrene</b>	<b>0.0015</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 13:40</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>95.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:40</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>83.1</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 16:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>104</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 16:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>81.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>72.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>84.3</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 16:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>79.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:40</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>84.8</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 16:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>82.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:40</i>
<i>Surr: Nitrobenzene-d5</i>	<i>109</i>			<i>41-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 16:59</i>
<i>Surr: Phenol-d6</i>	<i>82.9</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 16:59</i>
<i>Surr: Phenol-d6</i>	<i>75.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 13:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW26A-20190717  
 Collection Date: 17-Jul-2019 18:45

**ANALYTICAL REPORT**

WorkOrder:HS19070994  
 Lab ID:HS19070994-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 29-Jul-2019		Analyst: JHD
Arsenic	0.0933		0.000400	0.00200	mg/L	1	30-Jul-2019 22:14

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB05-20190717  
 Collection Date: 17-Jul-2019 19:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:07
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:07
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 01:07
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 01:07
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 01:07
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 01:07
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 01:07
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:07</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:07</i>
<i>Surr: Dibromofluoromethane</i>	<i>105</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:07</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 01:07</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB05-20190717  
 Collection Date: 17-Jul-2019 19:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 14:00
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 14:00
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 14:00
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 14:00
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 14:00
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 14:00
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 14:00
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 14:00
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 14:00
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 14:00
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 14:00
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 14:00
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 14:00
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 14:00
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00023</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 14:00
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 14:00
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 14:00
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 14:00
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 14:00
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 14:00
<b>Naphthalene</b>	<b>0.00023</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:00
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 14:00
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 14:00
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 14:00
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 14:00
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 14:00
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 14:00
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:00</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>104</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:00</i>
<i>Surr: 2-Fluorophenol</i>	<i>97.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:00</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:00</i>
<i>Surr: Nitrobenzene-d5</i>	<i>107</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:00</i>
<i>Surr: Phenol-d6</i>	<i>98.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:00</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 22:16

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68A-20190718  
 Collection Date: 18-Jul-2019 07:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:05
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:05
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:05
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:05
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 08:05
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:05
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:05
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:05</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:05</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:05</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:05</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68A-20190718  
 Collection Date: 18-Jul-2019 07:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 14:20
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 14:20
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 14:20
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 14:20
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 14:20
<b>2-Methylnaphthalene</b>	<b>0.00012</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 14:20
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 14:20
<b>Acenaphthene</b>	<b>0.0019</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 14:20
<b>Anthracene</b>	<b>0.000028</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 14:20
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 14:20
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 14:20
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00024</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 14:20
<b>Dibenzofuran</b>	<b>0.000082</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
<b>Di-n-butyl phthalate</b>	<b>0.00015</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 14:20
<b>Fluorene</b>	<b>0.00012</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
<b>Naphthalene</b>	<b>0.00035</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 14:20
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 14:20
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 14:20
<b>Phenanthrene</b>	<b>0.000062</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 14:20
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 14:20
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 14:20
<i>Surr: 2,4,6-Tribromophenol</i>	<i>110</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:20</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>82.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>92.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:20</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>95.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:20</i>
<i>Surr: Nitrobenzene-d5</i>	<i>80.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:20</i>
<i>Surr: Phenol-d6</i>	<i>85.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 14:20</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0353</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 22:18

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68B-20190718  
 Collection Date: 18-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0020	0.010	mg/L	10	23-Jul-2019 08:32
<b>Benzene</b>	<b>1.4</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	100	23-Jul-2019 08:59
Chlorobenzene	U		0.0030	0.010	mg/L	10	23-Jul-2019 08:32
<b>Ethylbenzene</b>	<b>0.52</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	23-Jul-2019 08:32
Methylene chloride	U		0.010	0.020	mg/L	10	23-Jul-2019 08:32
<b>Toluene</b>	<b>0.37</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	23-Jul-2019 08:32
<b>Xylenes, Total</b>	<b>1.5</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	23-Jul-2019 08:32
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.0</i>			<i>70-126</i>	<i>%REC</i>	<i>100</i>	<i>23-Jul-2019 08:59</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.2</i>			<i>70-126</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 08:32</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>104</i>			<i>81-113</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 08:32</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.3</i>			<i>81-113</i>	<i>%REC</i>	<i>100</i>	<i>23-Jul-2019 08:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 08:32</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>100</i>	<i>23-Jul-2019 08:59</i>
<i>Surr: Toluene-d8</i>	<i>99.6</i>			<i>82-127</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 08:32</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>100</i>	<i>23-Jul-2019 08:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68B-20190718  
 Collection Date: 18-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	23-Jul-2019 21:16
<b>2,4-Dimethylphenol</b>	<b>0.076</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	23-Jul-2019 21:16
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	23-Jul-2019 21:16
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	23-Jul-2019 21:16
<b>2-Methylnaphthalene</b>	<b>0.72</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	25-Jul-2019 18:57
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	23-Jul-2019 21:16
4-Nitrophenol	U		0.00047	0.010	mg/L	10	23-Jul-2019 21:16
<b>Acenaphthene</b>	<b>0.20</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	25-Jul-2019 18:57
<b>Acenaphthylene</b>	<b>0.0027</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
<b>Anthracene</b>	<b>0.030</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
<b>Benz(a)anthracene</b>	<b>0.0051</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
<b>Benzo(a)pyrene</b>	<b>0.0018</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	23-Jul-2019 21:16
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	23-Jul-2019 21:16
<b>Chrysene</b>	<b>0.0050</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
<b>Dibenzofuran</b>	<b>0.21</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	25-Jul-2019 18:57
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	23-Jul-2019 21:16
<b>Fluoranthene</b>	<b>0.043</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
<b>Fluorene</b>	<b>0.11</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	25-Jul-2019 18:57
<b>Naphthalene</b>	<b>9.6</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	24-Jul-2019 15:22
Nitrobenzene	U		0.00024	0.0020	mg/L	10	23-Jul-2019 21:16
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	23-Jul-2019 21:16
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	23-Jul-2019 21:16
<b>Phenanthrene</b>	<b>0.21</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	25-Jul-2019 18:57
Phenol	U		0.00035	0.0020	mg/L	10	23-Jul-2019 21:16
<b>Pyrene</b>	<b>0.025</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:16
<i>Surr: 2,4,6-Tribromophenol</i>	<i>102</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:16</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 15:22</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>25-Jul-2019 18:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>25-Jul-2019 18:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 15:22</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>80.5</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 15:22</i>
<i>Surr: 2-Fluorophenol</i>	<i>116</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>25-Jul-2019 18:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>25-Jul-2019 18:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>94.2</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:16</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 15:22</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68B-20190718  
 Collection Date: 18-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	24-Jul-2019 15:22
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	25-Jul-2019 18:57
Surr: Nitrobenzene-d5	99.5			41-120	%REC	10	23-Jul-2019 21:16
Surr: Phenol-d6	99.7			20-120	%REC	10	23-Jul-2019 21:16
Surr: Phenol-d6	0	JS		20-120	%REC	100	25-Jul-2019 18:57
Surr: Phenol-d6	0	JS		20-120	%REC	1000	24-Jul-2019 15:22
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	0.0134		0.000400	0.00200	mg/L	1	30-Jul-2019 22:21

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD03-20190718  
 Collection Date: 18-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0020	0.010	mg/L	10	23-Jul-2019 09:26
<b>Benzene</b>	<b>1.4</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	100	23-Jul-2019 09:53
Chlorobenzene	U		0.0030	0.010	mg/L	10	23-Jul-2019 09:26
<b>Ethylbenzene</b>	<b>0.51</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	23-Jul-2019 09:26
Methylene chloride	U		0.010	0.020	mg/L	10	23-Jul-2019 09:26
<b>Toluene</b>	<b>0.37</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	23-Jul-2019 09:26
<b>Xylenes, Total</b>	<b>1.4</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	23-Jul-2019 09:26
<i>Surr: 1,2-Dichloroethane-d4</i>	98.5			70-126	%REC	10	23-Jul-2019 09:26
<i>Surr: 1,2-Dichloroethane-d4</i>	97.1			70-126	%REC	100	23-Jul-2019 09:53
<i>Surr: 4-Bromofluorobenzene</i>	102			81-113	%REC	10	23-Jul-2019 09:26
<i>Surr: 4-Bromofluorobenzene</i>	102			81-113	%REC	100	23-Jul-2019 09:53
<i>Surr: Dibromofluoromethane</i>	103			77-123	%REC	10	23-Jul-2019 09:26
<i>Surr: Dibromofluoromethane</i>	99.5			77-123	%REC	100	23-Jul-2019 09:53
<i>Surr: Toluene-d8</i>	98.3			82-127	%REC	10	23-Jul-2019 09:26
<i>Surr: Toluene-d8</i>	97.2			82-127	%REC	100	23-Jul-2019 09:53

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD03-20190718  
 Collection Date: 18-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.00021	0.0020	mg/L	10	23-Jul-2019 21:36
<b>2,4-Dimethylphenol</b>	<b>0.076</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
2,4-Dinitrotoluene		U	0.00058	0.0020	mg/L	10	23-Jul-2019 21:36
2,6-Dinitrotoluene		U	0.00042	0.0020	mg/L	10	23-Jul-2019 21:36
2-Chloronaphthalene		U	0.00021	0.0020	mg/L	10	23-Jul-2019 21:36
<b>2-Methylnaphthalene</b>	<b>0.66</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	24-Jul-2019 14:23
4,6-Dinitro-2-methylphenol		U	0.00020	0.0020	mg/L	10	23-Jul-2019 21:36
4-Nitrophenol		U	0.00047	0.010	mg/L	10	23-Jul-2019 21:36
<b>Acenaphthene</b>	<b>0.17</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	24-Jul-2019 14:23
<b>Acenaphthylene</b>	<b>0.0022</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<b>Anthracene</b>	<b>0.022</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<b>Benz(a)anthracene</b>	<b>0.0030</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<b>Benzo(a)pyrene</b>	<b>0.0012</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
Bis(2-chloroethoxy)methane		U	0.00030	0.0020	mg/L	10	23-Jul-2019 21:36
Bis(2-ethylhexyl)phthalate		U	0.00037	0.0020	mg/L	10	23-Jul-2019 21:36
<b>Chrysene</b>	<b>0.0029</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<b>Dibenzofuran</b>	<b>0.18</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	24-Jul-2019 14:23
Di-n-butyl phthalate		U	0.00020	0.0020	mg/L	10	23-Jul-2019 21:36
<b>Fluoranthene</b>	<b>0.027</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<b>Fluorene</b>	<b>0.083</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<b>Naphthalene</b>	<b>9.9</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	24-Jul-2019 14:43
Nitrobenzene		U	0.00024	0.0020	mg/L	10	23-Jul-2019 21:36
N-Nitrosodiphenylamine		U	0.00025	0.0020	mg/L	10	23-Jul-2019 21:36
Pentachlorophenol		U	0.00079	0.0020	mg/L	10	23-Jul-2019 21:36
<b>Phenanthrene</b>	<b>0.16</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	24-Jul-2019 14:23
Phenol		U	0.00035	0.0020	mg/L	10	23-Jul-2019 21:36
<b>Pyrene</b>	<b>0.015</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:36
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 14:43</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>24-Jul-2019 14:23</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>91.9</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:36</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>77.8</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:36</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>24-Jul-2019 14:23</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 14:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>24-Jul-2019 14:23</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 14:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>105</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:36</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>76.3</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>23-Jul-2019 21:36</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>24-Jul-2019 14:23</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>1000</i>	<i>24-Jul-2019 14:43</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD03-20190718  
 Collection Date: 18-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	24-Jul-2019 14:23
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	24-Jul-2019 14:43
Surr: Nitrobenzene-d5	106			41-120	%REC	10	23-Jul-2019 21:36
Surr: Phenol-d6	76.0			20-120	%REC	10	23-Jul-2019 21:36
Surr: Phenol-d6	0	JS		20-120	%REC	100	24-Jul-2019 14:23
Surr: Phenol-d6	0	JS		20-120	%REC	1000	24-Jul-2019 14:43
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	0.0129		0.000400	0.00200	mg/L	1	30-Jul-2019 22:23

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68C-20190718  
 Collection Date: 18-Jul-2019 09:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:01
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:01
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:01
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:01
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 02:01
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:01
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:01
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>88.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:01</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:01</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:01</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:01</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68C-20190718  
 Collection Date: 18-Jul-2019 09:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 15:19
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 15:19
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 15:19
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 15:19
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 15:19
<b>2-Methylnaphthalene</b>	<b>0.000072</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 15:19
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 15:19
<b>Acenaphthene</b>	<b>0.00012</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
<b>Acenaphthylene</b>	<b>0.000053</b>	<b>J</b>	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 15:19
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 15:19
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 15:19
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 15:19
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00032</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 15:19
<b>Dibenzofuran</b>	<b>0.000064</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
<b>Di-n-butyl phthalate</b>	<b>0.000067</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 15:19
<b>Fluorene</b>	<b>0.000051</b>	<b>J</b>	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
<b>Naphthalene</b>	<b>0.0011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 15:19
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 15:19
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 15:19
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 15:19
<b>Phenol</b>	<b>0.00026</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 15:19</b>
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 15:19
<i>Surr: 2,4,6-Tribromophenol</i>	<i>106</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 15:19</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>84.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 15:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>79.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 15:19</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>84.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 15:19</i>
<i>Surr: Nitrobenzene-d5</i>	<i>80.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 15:19</i>
<i>Surr: Phenol-d6</i>	<i>80.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 15:19</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 22:25

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83B-20190718  
 Collection Date: 18-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:30
<b>Benzene</b>	<b>0.021</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:30
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:30
<b>Ethylbenzene</b>	<b>0.055</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:30
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:30
<b>Toluene</b>	<b>0.0042</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:30
<b>Xylenes, Total</b>	<b>0.061</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 06:30
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>87.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:30</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:30</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:30</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83B-20190718  
 Collection Date: 18-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 15:39
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 15:39
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 15:39
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 15:39
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 15:39
<b>2-Methylnaphthalene</b>	<b>0.050</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:56
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 15:39
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 15:39
<b>Acenaphthene</b>	<b>0.021</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:56
<b>Acenaphthylene</b>	<b>0.00025</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
<b>Anthracene</b>	<b>0.00099</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 15:39
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 15:39
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 15:39
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00036</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 15:39
<b>Dibenzofuran</b>	<b>0.014</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	23-Jul-2019 21:56
<b>Di-n-butyl phthalate</b>	<b>0.00012</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
<b>Fluoranthene</b>	<b>0.00038</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
<b>Fluorene</b>	<b>0.0076</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
<b>Naphthalene</b>	<b>0.77</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	24-Jul-2019 15:02
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 15:39
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 15:39
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 15:39
<b>Phenanthrene</b>	<b>0.0059</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
<b>Phenol</b>	<b>0.00051</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
<b>Pyrene</b>	<b>0.00020</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 15:39
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	24-Jul-2019 15:02
Surr: 2,4,6-Tribromophenol	94.6			34-129	%REC	10	23-Jul-2019 21:56
Surr: 2,4,6-Tribromophenol	101			34-129	%REC	1	23-Jul-2019 15:39
Surr: 2-Fluorobiphenyl	58.6			40-125	%REC	1	23-Jul-2019 15:39
Surr: 2-Fluorobiphenyl	55.7			40-125	%REC	10	23-Jul-2019 21:56
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	24-Jul-2019 15:02
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	24-Jul-2019 15:02
Surr: 2-Fluorophenol	64.1			20-120	%REC	10	23-Jul-2019 21:56
Surr: 2-Fluorophenol	61.1			20-120	%REC	1	23-Jul-2019 15:39
Surr: 4-Terphenyl-d14	91.3			40-135	%REC	1	23-Jul-2019 15:39
Surr: 4-Terphenyl-d14	86.5			40-135	%REC	10	23-Jul-2019 21:56
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	24-Jul-2019 15:02

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83B-20190718  
 Collection Date: 18-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	24-Jul-2019 15:02
Surr: Nitrobenzene-d5	51.9			41-120	%REC	1	23-Jul-2019 15:39
Surr: Nitrobenzene-d5	56.9			41-120	%REC	10	23-Jul-2019 21:56
Surr: Phenol-d6	59.6			20-120	%REC	1	23-Jul-2019 15:39
Surr: Phenol-d6	47.6			20-120	%REC	10	23-Jul-2019 21:56
Surr: Phenol-d6	0	JS		20-120	%REC	100	24-Jul-2019 15:02
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 29-Jul-2019		Analyst: JHD	
Arsenic	0.0648		0.000400	0.00200	mg/L	1	30-Jul-2019 22:27

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83C-20190718  
 Collection Date: 18-Jul-2019 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:13
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:13
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:13
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:13
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 03:13
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 03:13
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 03:13
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:13</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:13</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:13</i>
<i>Surr: Toluene-d8</i>	<i>105</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 03:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83C-20190718  
 Collection Date: 18-Jul-2019 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 17:19
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 17:19
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 17:19
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 17:19
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 17:19
<b>2-Methylnaphthalene</b>	<b>0.00012</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 17:19
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 17:19
<b>Acenaphthene</b>	<b>0.000094</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 17:19
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 17:19
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 17:19
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 17:19
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 17:19
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00023</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 17:19
<b>Dibenzofuran</b>	<b>0.000070</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
<b>Di-n-butyl phthalate</b>	<b>0.000040</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 17:19
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 17:19
<b>Naphthalene</b>	<b>0.0014</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 17:19
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 17:19
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 17:19
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 17:19
<b>Phenol</b>	<b>0.000068</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 17:19
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 17:19
<i>Surr: 2,4,6-Tribromophenol</i>	<i>78.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:19</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>59.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>54.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:19</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:19</i>
<i>Surr: Nitrobenzene-d5</i>	<i>71.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:19</i>
<i>Surr: Phenol-d6</i>	<i>56.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:19</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00617</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 22:41

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35A-20190718  
 Collection Date: 18-Jul-2019 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:25
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:25
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:25
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:25
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 04:25
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 04:25
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 04:25
<i>Surr: 1,2-Dichloroethane-d4</i>		104		70-126	%REC	1	23-Jul-2019 04:25
<i>Surr: 4-Bromofluorobenzene</i>		103		81-113	%REC	1	23-Jul-2019 04:25
<i>Surr: Dibromofluoromethane</i>		102		77-123	%REC	1	23-Jul-2019 04:25
<i>Surr: Toluene-d8</i>		107		82-127	%REC	1	23-Jul-2019 04:25

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35A-20190718  
 Collection Date: 18-Jul-2019 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 17:38
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 17:38
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 17:38
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 17:38
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 17:38
<b>2-Methylnaphthalene</b>	<b>0.00012</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 17:38
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 17:38
<b>Acenaphthene</b>	<b>0.0028</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
<b>Acenaphthylene</b>	<b>0.000071</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
<b>Anthracene</b>	<b>0.00011</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 17:38
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 17:38
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 17:38
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000057</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 17:38
<b>Dibenzofuran</b>	<b>0.00039</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 17:38
<b>Fluoranthene</b>	<b>0.00015</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
<b>Fluorene</b>	<b>0.00036</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
<b>Naphthalene</b>	<b>0.00083</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 17:38
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 17:38
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 17:38
<b>Phenanthrene</b>	<b>0.000042</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 17:38
<b>Pyrene</b>	<b>0.000096</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:38
<i>Surr: 2,4,6-Tribromophenol</i>	<i>103</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>70.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>69.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:38</i>
<i>Surr: Nitrobenzene-d5</i>	<i>83.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:38</i>
<i>Surr: Phenol-d6</i>	<i>73.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:38</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0548</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 22:57

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35B-20190718  
 Collection Date: 18-Jul-2019 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	23-Jul-2019 04:49
<b>Benzene</b>	<b>0.0045</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 04:49
Chlorobenzene		U	0.00030	0.0010	mg/L	1	23-Jul-2019 04:49
<b>Ethylbenzene</b>	<b>0.014</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 04:49
Methylene chloride		U	0.0010	0.0020	mg/L	1	23-Jul-2019 04:49
<b>Toluene</b>	<b>0.00054</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 04:49
<b>Xylenes, Total</b>	<b>0.0062</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	23-Jul-2019 04:49
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:49</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:49</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:49</i>
<i>Surr: Toluene-d8</i>	<i>104</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 04:49</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35B-20190718  
 Collection Date: 18-Jul-2019 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 17:58
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 17:58
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 17:58
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 17:58
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 17:58
<b>2-Methylnaphthalene</b>	<b>0.041</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	25-Jul-2019 19:17
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 17:58
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 17:58
<b>Acenaphthene</b>	<b>0.029</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	25-Jul-2019 19:17
<b>Acenaphthylene</b>	<b>0.00022</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Anthracene</b>	<b>0.0022</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Benz(a)anthracene</b>	<b>0.000083</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Benzo(a)pyrene</b>	<b>0.000053</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 17:58
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00012</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Chrysene</b>	<b>0.000091</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Dibenzofuran</b>	<b>0.029</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	25-Jul-2019 19:17
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 17:58
<b>Fluoranthene</b>	<b>0.0017</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Fluorene</b>	<b>0.014</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	25-Jul-2019 19:17
<b>Naphthalene</b>	<b>1.1</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	500	26-Jul-2019 19:55
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 17:58
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 17:58
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 17:58
<b>Phenanthrene</b>	<b>0.019</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	25-Jul-2019 19:17
<b>Phenol</b>	<b>0.00017</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<b>Pyrene</b>	<b>0.00079</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 17:58
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>500</i>	<i>26-Jul-2019 19:55</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>112</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:58</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.6</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>25-Jul-2019 19:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>97.8</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>25-Jul-2019 19:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>88.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>500</i>	<i>26-Jul-2019 19:55</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>500</i>	<i>26-Jul-2019 19:55</i>
<i>Surr: 2-Fluorophenol</i>	<i>86.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:58</i>
<i>Surr: 2-Fluorophenol</i>	<i>90.8</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>25-Jul-2019 19:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>105</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>25-Jul-2019 19:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 17:58</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>500</i>	<i>26-Jul-2019 19:55</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35B-20190718  
 Collection Date: 18-Jul-2019 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
Surr: Nitrobenzene-d5	89.4			41-120	%REC	1	23-Jul-2019 17:58
Surr: Nitrobenzene-d5	97.7			41-120	%REC	10	25-Jul-2019 19:17
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	500	26-Jul-2019 19:55
Surr: Phenol-d6	0	JS		20-120	%REC	500	26-Jul-2019 19:55
Surr: Phenol-d6	89.7			20-120	%REC	10	25-Jul-2019 19:17
Surr: Phenol-d6	93.6			20-120	%REC	1	23-Jul-2019 17:58
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
Arsenic	0.00120	J	0.000400	0.00200	mg/L	1	30-Jul-2019 22:59

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW90B-20190718  
 Collection Date: 18-Jul-2019 14:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:13
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:13
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 05:13
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:13
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:13
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>109</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>107</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>
<i>Surr: Dibromofluoromethane</i>	<i>109</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>
<i>Surr: Toluene-d8</i>	<i>111</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW90B-20190718  
 Collection Date: 18-Jul-2019 14:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 18:18
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 18:18
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 18:18
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 18:18
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 18:18
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 18:18
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 18:18
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 18:18
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 18:18
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 18:18
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 18:18
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 18:18
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 18:18
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 18:18
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 18:18
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 18:18
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 18:18
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 18:18
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 18:18
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 18:18
<b>Naphthalene</b>	<b>0.000091</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:18
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 18:18
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 18:18
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 18:18
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 18:18
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 18:18
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 18:18
<i>Surr: 2,4,6-Tribromophenol</i>	<i>101</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>84.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:18</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:18</i>
<i>Surr: Nitrobenzene-d5</i>	<i>87.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:18</i>
<i>Surr: Phenol-d6</i>	<i>85.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:18</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0135</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 23:01

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW89B-20190718  
 Collection Date: 18-Jul-2019 15:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:37
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 05:37
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 05:37
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 05:37
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>105</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>105</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>
<i>Surr: Dibromofluoromethane</i>	<i>104</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>
<i>Surr: Toluene-d8</i>	<i>112</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 05:37</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW89B-20190718  
 Collection Date: 18-Jul-2019 15:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 18:38
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 18:38
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 18:38
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 18:38
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 18:38
<b>2-Methylnaphthalene</b>	<b>0.000050</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 18:38</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 18:38
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 18:38
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 18:38
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 18:38
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 18:38
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 18:38
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 18:38
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 18:38
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00015</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 18:38</b>
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 18:38
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 18:38
<b>Di-n-butyl phthalate</b>	<b>0.000038</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 18:38</b>
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 18:38
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 18:38
<b>Naphthalene</b>	<b>0.00025</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 18:38</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 18:38
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 18:38
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 18:38
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 18:38
<b>Phenol</b>	<b>0.00042</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 18:38</b>
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 18:38
<i>Surr: 2,4,6-Tribromophenol</i>	<i>105</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>76.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:38</i>
<i>Surr: Nitrobenzene-d5</i>	<i>80.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:38</i>
<i>Surr: Phenol-d6</i>	<i>76.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:38</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 23:04

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38B-20190718  
 Collection Date: 18-Jul-2019 16:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-26  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:01
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:01
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:01
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:01
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:01
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:01
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:01
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>114</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:01</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>111</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:01</i>
<i>Surr: Dibromofluoromethane</i>	<i>113</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:01</i>
<i>Surr: Toluene-d8</i>	<i>115</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:01</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38B-20190718  
 Collection Date: 18-Jul-2019 16:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-26  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 18:58
<b>2,4-Dimethylphenol</b>	<b>0.000056</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 18:58
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 18:58
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 18:58
<b>2-Methylnaphthalene</b>	<b>0.00030</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 18:58
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 18:58
<b>Acenaphthene</b>	<b>0.00012</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 18:58
<b>Anthracene</b>	<b>0.000037</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 18:58
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 18:58
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 18:58
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00031</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 18:58
<b>Dibenzofuran</b>	<b>0.00014</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 18:58
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 18:58
<b>Fluorene</b>	<b>0.000060</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
<b>Naphthalene</b>	<b>0.0026</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 18:58
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 18:58
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 18:58
<b>Phenanthrene</b>	<b>0.000079</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
<b>Phenol</b>	<b>0.000056</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 18:58
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 18:58
<i>Surr: 2,4,6-Tribromophenol</i>	<i>101</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:58</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>96.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:58</i>
<i>Surr: 2-Fluorophenol</i>	<i>83.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:58</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:58</i>
<i>Surr: Nitrobenzene-d5</i>	<i>90.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:58</i>
<i>Surr: Phenol-d6</i>	<i>90.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 18:58</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000553</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 23:06

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27A-20190718  
 Collection Date: 18-Jul-2019 17:15

**ANALYTICAL REPORT**

WorkOrder:HS19070994  
 Lab ID:HS19070994-27  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:25
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:25
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:25
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:25
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:25
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:25
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:25
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:25</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>103</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:25</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:25</i>
<i>Surr: Toluene-d8</i>	<i>108</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 06:25</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27A-20190718  
 Collection Date: 18-Jul-2019 17:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-27  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 19:18
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 19:18
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 19:18
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 19:18
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 19:18
<b>2-Methylnaphthalene</b>	<b>0.00034</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 19:18
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 19:18
<b>Acenaphthene</b>	<b>0.00011</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 19:18
<b>Anthracene</b>	<b>0.000036</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 19:18
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 19:18
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 19:18
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00020</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 19:18
<b>Dibenzofuran</b>	<b>0.00014</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
<b>Di-n-butyl phthalate</b>	<b>0.000034</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
<b>Fluoranthene</b>	<b>0.000024</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
<b>Fluorene</b>	<b>0.000074</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
<b>Naphthalene</b>	<b>0.0026</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 19:18
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 19:18
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 19:18
<b>Phenanthrene</b>	<b>0.00013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:18
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 19:18
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 19:18
<i>Surr: 2,4,6-Tribromophenol</i>	<i>106</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>86.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:18</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>93.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:18</i>
<i>Surr: Nitrobenzene-d5</i>	<i>90.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:18</i>
<i>Surr: Phenol-d6</i>	<i>83.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:18</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000498</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 23:08

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27C-20190718  
 Collection Date: 18-Jul-2019 18:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-28  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:48
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:48
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:48
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:48
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 06:48
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 06:48
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 06:48
<i>Surr: 1,2-Dichloroethane-d4</i>		110		70-126	%REC	1	23-Jul-2019 06:48
<i>Surr: 4-Bromofluorobenzene</i>		107		81-113	%REC	1	23-Jul-2019 06:48
<i>Surr: Dibromofluoromethane</i>		108		77-123	%REC	1	23-Jul-2019 06:48
<i>Surr: Toluene-d8</i>		113		82-127	%REC	1	23-Jul-2019 06:48

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27C-20190718  
 Collection Date: 18-Jul-2019 18:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-28  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 19:37
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 19:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 19:37
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 19:37
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 19:37
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 19:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 19:37
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 19:37
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 19:37
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 19:37
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 19:37
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 19:37
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 19:37
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 19:37
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000092</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 19:37
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 19:37
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 19:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 19:37
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 19:37
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 19:37
<b>Naphthalene</b>	<b>0.000066</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:37
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 19:37
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 19:37
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 19:37
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 19:37
<b>Phenol</b>	<b>0.00023</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 19:37
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 19:37
<i>Surr: 2,4,6-Tribromophenol</i>	<i>106</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>85.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>81.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:37</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:37</i>
<i>Surr: Nitrobenzene-d5</i>	<i>91.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:37</i>
<i>Surr: Phenol-d6</i>	<i>87.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:37</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000428</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 23:10

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51A-20190719  
 Collection Date: 19-Jul-2019 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-29  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:12
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:12
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:12
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:12
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 07:12
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:12
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:12
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>115</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:12</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>109</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:12</i>
<i>Surr: Dibromofluoromethane</i>	<i>114</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:12</i>
<i>Surr: Toluene-d8</i>	<i>116</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:12</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51A-20190719  
 Collection Date: 19-Jul-2019 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-29  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 19:57
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 19:57
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 19:57
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 19:57
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 19:57
<b>2-Methylnaphthalene</b>	<b>0.000067</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:57
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 19:57
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 19:57
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 19:57
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 19:57
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 19:57
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 19:57
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 19:57
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 19:57
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00013</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 19:57
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 19:57
<b>Dibenzofuran</b>	<b>0.000030</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:57
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 19:57
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 19:57
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 19:57
<b>Naphthalene</b>	<b>0.00055</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:57
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 19:57
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 19:57
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 19:57
<b>Phenanthrene</b>	<b>0.000026</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 19:57
<b>Phenol</b>	<b>0.00019</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 19:57
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 19:57
<i>Surr: 2,4,6-Tribromophenol</i>	<i>112</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>95.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>92.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:57</i>
<i>Surr: Nitrobenzene-d5</i>	<i>99.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:57</i>
<i>Surr: Phenol-d6</i>	<i>99.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 19:57</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 23:13

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51C-20190719  
 Collection Date: 19-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-30  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:36
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:36
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:36
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:36
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 07:36
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 07:36
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 07:36
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>111</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:36</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>109</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:36</i>
<i>Surr: Dibromofluoromethane</i>	<i>110</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:36</i>
<i>Surr: Toluene-d8</i>	<i>116</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 07:36</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51C-20190719  
 Collection Date: 19-Jul-2019 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-30  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 20:17
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 20:17
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 20:17
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 20:17
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 20:17
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 20:17
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 20:17
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 20:17
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 20:17
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 20:17
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 20:17
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 20:17
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 20:17
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 20:17
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000079</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 20:17
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 20:17
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 20:17
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 20:17
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 20:17
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 20:17
<b>Naphthalene</b>	<b>0.00023</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jul-2019 20:17
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 20:17
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 20:17
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 20:17
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 20:17
<b>Phenol</b>	<b>0.00027</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jul-2019 20:17
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 20:17
<i>Surr: 2,4,6-Tribromophenol</i>	<i>106</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>87.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:17</i>
<i>Surr: 2-Fluorophenol</i>	<i>83.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:17</i>
<i>Surr: Nitrobenzene-d5</i>	<i>92.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:17</i>
<i>Surr: Phenol-d6</i>	<i>89.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:17</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	30-Jul-2019 23:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW81B-20190719  
 Collection Date: 19-Jul-2019 09:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-31  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:00
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:00
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:00
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:00
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 08:00
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:00
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:00
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>110</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:00</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>109</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:00</i>
<i>Surr: Dibromofluoromethane</i>		<i>109</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:00</i>
<i>Surr: Toluene-d8</i>		<i>114</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW81B-20190719  
 Collection Date: 19-Jul-2019 09:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-31  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jul-2019 20:37
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jul-2019 20:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jul-2019 20:37
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jul-2019 20:37
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jul-2019 20:37
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 20:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jul-2019 20:37
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jul-2019 20:37
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jul-2019 20:37
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jul-2019 20:37
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jul-2019 20:37
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jul-2019 20:37
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jul-2019 20:37
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jul-2019 20:37
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	23-Jul-2019 20:37
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 20:37
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jul-2019 20:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jul-2019 20:37
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jul-2019 20:37
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jul-2019 20:37
<b>Naphthalene</b>	<b>0.000099</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jul-2019 20:37</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jul-2019 20:37
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jul-2019 20:37
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jul-2019 20:37
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jul-2019 20:37
Phenol	U		0.000035	0.00020	mg/L	1	23-Jul-2019 20:37
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jul-2019 20:37
<i>Surr: 2,4,6-Tribromophenol</i>	<i>85.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:37</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>80.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:37</i>
<i>Surr: Nitrobenzene-d5</i>	<i>80.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:37</i>
<i>Surr: Phenol-d6</i>	<i>79.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 20:37</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000563</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>30-Jul-2019 23:17</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50A-20190719  
 Collection Date: 19-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-32  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:24
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:24
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:24
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:24
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 08:24
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 08:24
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 08:24
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>105</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:24</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>103</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:24</i>
<i>Surr: Dibromofluoromethane</i>	<i>106</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:24</i>
<i>Surr: Toluene-d8</i>	<i>110</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 08:24</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50A-20190719  
 Collection Date: 19-Jul-2019 10:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-32  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Jul-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	24-Jul-2019 15:41
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	24-Jul-2019 15:41
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	24-Jul-2019 15:41
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	24-Jul-2019 15:41
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	24-Jul-2019 15:41
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	24-Jul-2019 15:41
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	24-Jul-2019 15:41
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	24-Jul-2019 15:41
Acenaphthene	U		0.000027	0.00010	mg/L	1	24-Jul-2019 15:41
Acenaphthylene	U		0.000015	0.00010	mg/L	1	24-Jul-2019 15:41
Anthracene	U		0.000014	0.00010	mg/L	1	24-Jul-2019 15:41
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	24-Jul-2019 15:41
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	24-Jul-2019 15:41
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	24-Jul-2019 15:41
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	24-Jul-2019 15:41
Chrysene	U		0.000021	0.00010	mg/L	1	24-Jul-2019 15:41
Dibenzofuran	U		0.000020	0.00010	mg/L	1	24-Jul-2019 15:41
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	24-Jul-2019 15:41
Fluoranthene	U		0.000010	0.00010	mg/L	1	24-Jul-2019 15:41
Fluorene	U		0.000030	0.00010	mg/L	1	24-Jul-2019 15:41
<b>Naphthalene</b>	<b>0.00027</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	24-Jul-2019 15:41
Nitrobenzene	U		0.000024	0.00020	mg/L	1	24-Jul-2019 15:41
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	24-Jul-2019 15:41
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	24-Jul-2019 15:41
Phenanthrene	U		0.000021	0.00010	mg/L	1	24-Jul-2019 15:41
<b>Phenol</b>	<b>0.00016</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	24-Jul-2019 15:41
Pyrene	U		0.000019	0.00010	mg/L	1	24-Jul-2019 15:41
<i>Surr: 2,4,6-Tribromophenol</i>	<i>87.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>24-Jul-2019 15:41</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>24-Jul-2019 15:41</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>24-Jul-2019 15:41</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>24-Jul-2019 15:41</i>
<i>Surr: Nitrobenzene-d5</i>	<i>73.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>24-Jul-2019 15:41</i>
<i>Surr: Phenol-d6</i>	<i>83.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>24-Jul-2019 15:41</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jul-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000642</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	30-Jul-2019 23:24

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB05-20190719  
 Collection Date: 19-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-33  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:25
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:25
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:25
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:25
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 02:25
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:25
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:25
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:25
<i>Surr: 1,2-Dichloroethane-d4</i>		93.4		70-126	%REC	1	23-Jul-2019 02:25
<i>Surr: 4-Bromofluorobenzene</i>		96.4		81-113	%REC	1	23-Jul-2019 02:25
<i>Surr: Dibromofluoromethane</i>		92.9		77-123	%REC	1	23-Jul-2019 02:25
<i>Surr: Toluene-d8</i>		103		82-127	%REC	1	23-Jul-2019 02:25

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB06-20190719  
 Collection Date: 19-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19070994  
 Lab ID:HS19070994-34  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:49
Benzene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:49
Chlorobenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:49
Ethylbenzene	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:49
Methylene chloride	U		0.0010	0.0020	mg/L	1	23-Jul-2019 02:49
Toluene	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:49
Vinyl chloride	U		0.00020	0.0010	mg/L	1	23-Jul-2019 02:49
Xylenes, Total	U		0.00030	0.0010	mg/L	1	23-Jul-2019 02:49
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:49</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:49</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:49</i>
<i>Surr: Toluene-d8</i>	<i>105</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Jul-2019 02:49</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**Batch ID:** 143239      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070994-01	1	1000	1 (mL)	0.001
HS19070994-02	1	1000	1 (mL)	0.001
HS19070994-03	1	1000	1 (mL)	0.001
HS19070994-04	1	1000	1 (mL)	0.001
HS19070994-05	1	1000	1 (mL)	0.001
HS19070994-06	1	1000	1 (mL)	0.001
HS19070994-07	1	1000	1 (mL)	0.001
HS19070994-08	1	1000	1 (mL)	0.001
HS19070994-09	1	1000	1 (mL)	0.001
HS19070994-10	1	1000	1 (mL)	0.001
HS19070994-11	1	1000	1 (mL)	0.001
HS19070994-12	1	1000	1 (mL)	0.001
HS19070994-13	1	1000	1 (mL)	0.001
HS19070994-14	1	1000	1 (mL)	0.001
HS19070994-15	1	1000	1 (mL)	0.001
HS19070994-16	1	1000	1 (mL)	0.001
HS19070994-17	1	1000	1 (mL)	0.001
HS19070994-18	1	1000	1 (mL)	0.001
HS19070994-19	1	1000	1 (mL)	0.001
HS19070994-20	1	1000	1 (mL)	0.001

**Batch ID:** 143295      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070994-21	1	1000	1 (mL)	0.001
HS19070994-22	1	1000	1 (mL)	0.001
HS19070994-23	1	1000	1 (mL)	0.001
HS19070994-24	1	1000	1 (mL)	0.001
HS19070994-25	1	1000	1 (mL)	0.001
HS19070994-26	1	1000	1 (mL)	0.001
HS19070994-27	1	1000	1 (mL)	0.001
HS19070994-28	1	1000	1 (mL)	0.001
HS19070994-29	1	1000	1 (mL)	0.001
HS19070994-30	1	1000	1 (mL)	0.001
HS19070994-31	1	1000	1 (mL)	0.001
HS19070994-32	1	1000	1 (mL)	0.001

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**Batch ID:** 143526      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070994-01	1	10	10 (mL)	1
HS19070994-02	1	10	10 (mL)	1
HS19070994-03	1	10	10 (mL)	1
HS19070994-04	1	10	10 (mL)	1
HS19070994-05	1	10	10 (mL)	1
HS19070994-06	1	10	10 (mL)	1
HS19070994-07	1	10	10 (mL)	1
HS19070994-08	1	10	10 (mL)	1
HS19070994-09	1	10	10 (mL)	1
HS19070994-10	1	10	10 (mL)	1
HS19070994-11	1	10	10 (mL)	1
HS19070994-12	1	10	10 (mL)	1
HS19070994-13	1	10	10 (mL)	1
HS19070994-14	1	10	10 (mL)	1
HS19070994-15	1	10	10 (mL)	1
HS19070994-16	1	10	10 (mL)	1
HS19070994-17	1	10	10 (mL)	1
HS19070994-18	1	10	10 (mL)	1
HS19070994-19	1	10	10 (mL)	1
HS19070994-20	1	10	10 (mL)	1

**Batch ID:** 143558      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19070994-21	1	10	10 (mL)	1
HS19070994-22	1	10	10 (mL)	1
HS19070994-23	1	10	10 (mL)	1
HS19070994-24	1	10	10 (mL)	1
HS19070994-25	1	10	10 (mL)	1
HS19070994-26	1	10	10 (mL)	1
HS19070994-27	1	10	10 (mL)	1
HS19070994-28	1	10	10 (mL)	1
HS19070994-29	1	10	10 (mL)	1
HS19070994-30	1	10	10 (mL)	1
HS19070994-31	1	10	10 (mL)	1
HS19070994-32	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID:</b> 143239 ( 0 )		<b>Test Name :</b> LOW-LEVEL SEMIVOLATILES BY 8270D			<b>Matrix:</b> Water	
HS19070994-01	WG-1620-FB04-20190716	16 Jul 2019 18:20		22 Jul 2019 11:13	23 Jul 2019 11:41	1
HS19070994-02	WG-1620-MW61A-20190717	17 Jul 2019 07:50		22 Jul 2019 11:13	23 Jul 2019 12:21	1
HS19070994-03	WG-1620-MW60A-20190717	17 Jul 2019 08:35		22 Jul 2019 11:13	22 Jul 2019 18:57	1
HS19070994-04	WG-1620-MW69A-20190717	17 Jul 2019 09:35		22 Jul 2019 11:13	22 Jul 2019 19:17	1
HS19070994-05	WG-1620-MW47C-20190717	17 Jul 2019 10:25		22 Jul 2019 11:13	22 Jul 2019 19:36	1
HS19070994-06	WG-1620-MW48C-20190717	17 Jul 2019 11:20		22 Jul 2019 11:13	22 Jul 2019 19:56	1
HS19070994-07	WG-1620-MW59A-20190717	17 Jul 2019 12:20		22 Jul 2019 11:13	22 Jul 2019 20:16	1
HS19070994-08	WG-1620-MW59B-20190717	17 Jul 2019 13:10		22 Jul 2019 11:13	22 Jul 2019 20:36	1
HS19070994-09	WG-1620-MW44A-20190717	17 Jul 2019 14:00		22 Jul 2019 11:13	23 Jul 2019 15:59	10
HS19070994-09	WG-1620-MW44A-20190717	17 Jul 2019 14:00		22 Jul 2019 11:13	22 Jul 2019 20:56	1
HS19070994-10	WG-1620-MW87C-20190717	17 Jul 2019 15:40		22 Jul 2019 11:13	22 Jul 2019 21:16	1
HS19070994-11	WG-1620-MW71B-20190717	17 Jul 2019 16:40		22 Jul 2019 11:13	23 Jul 2019 16:39	100
HS19070994-11	WG-1620-MW71B-20190717	17 Jul 2019 16:40		22 Jul 2019 11:13	23 Jul 2019 16:19	10
HS19070994-11	WG-1620-MW71B-20190717	17 Jul 2019 16:40		22 Jul 2019 11:13	22 Jul 2019 21:36	1
HS19070994-12	WG-1620-MW33A-20190717	17 Jul 2019 17:50		22 Jul 2019 11:13	22 Jul 2019 21:55	1
HS19070994-13	WG-1620-FD02-20190717	17 Jul 2019 17:50		22 Jul 2019 11:13	23 Jul 2019 13:20	1
HS19070994-14	WG-1620-MW26A-20190717	17 Jul 2019 18:45		22 Jul 2019 11:13	23 Jul 2019 16:59	10
HS19070994-14	WG-1620-MW26A-20190717	17 Jul 2019 18:45		22 Jul 2019 11:13	23 Jul 2019 13:40	1
HS19070994-15	WG-1620-FB05-20190717	17 Jul 2019 19:05		22 Jul 2019 11:13	23 Jul 2019 14:00	1
HS19070994-16	WG-1620-MW68A-20190718	18 Jul 2019 07:55		22 Jul 2019 11:13	23 Jul 2019 14:20	1
HS19070994-17	WG-1620-MW68B-20190718	18 Jul 2019 08:55		22 Jul 2019 11:13	25 Jul 2019 18:57	100
HS19070994-17	WG-1620-MW68B-20190718	18 Jul 2019 08:55		22 Jul 2019 11:13	24 Jul 2019 15:22	1000
HS19070994-17	WG-1620-MW68B-20190718	18 Jul 2019 08:55		22 Jul 2019 11:13	23 Jul 2019 21:16	10
HS19070994-18	WG-1620-FD03-20190718	18 Jul 2019 08:55		22 Jul 2019 11:13	24 Jul 2019 14:43	1000
HS19070994-18	WG-1620-FD03-20190718	18 Jul 2019 08:55		22 Jul 2019 11:13	24 Jul 2019 14:23	100
HS19070994-18	WG-1620-FD03-20190718	18 Jul 2019 08:55		22 Jul 2019 11:13	23 Jul 2019 21:36	10
HS19070994-19	WG-1620-MW68C-20190718	18 Jul 2019 09:45		22 Jul 2019 11:13	23 Jul 2019 15:19	1
HS19070994-20	WG-1620-MW83B-20190718	18 Jul 2019 10:40		22 Jul 2019 11:13	24 Jul 2019 15:02	100
HS19070994-20	WG-1620-MW83B-20190718	18 Jul 2019 10:40		22 Jul 2019 11:13	23 Jul 2019 21:56	10
HS19070994-20	WG-1620-MW83B-20190718	18 Jul 2019 10:40		22 Jul 2019 11:13	23 Jul 2019 15:39	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143295 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS19070994-21	WG-1620-MW83C-20190718	18 Jul 2019 11:30		23 Jul 2019 13:00	23 Jul 2019 17:19	1
HS19070994-22	WG-1620-MW35A-20190718	18 Jul 2019 12:40		23 Jul 2019 13:00	23 Jul 2019 17:38	1
HS19070994-23	WG-1620-MW35B-20190718	18 Jul 2019 13:30		23 Jul 2019 13:00	26 Jul 2019 19:55	500
HS19070994-23	WG-1620-MW35B-20190718	18 Jul 2019 13:30		23 Jul 2019 13:00	25 Jul 2019 19:17	10
HS19070994-23	WG-1620-MW35B-20190718	18 Jul 2019 13:30		23 Jul 2019 13:00	23 Jul 2019 17:58	1
HS19070994-24	WG-1620-MW90B-20190718	18 Jul 2019 14:25		23 Jul 2019 13:00	23 Jul 2019 18:18	1
HS19070994-25	WG-1620-MW89B-20190718	18 Jul 2019 15:15		23 Jul 2019 13:00	23 Jul 2019 18:38	1
HS19070994-26	WG-1620-MW38B-20190718	18 Jul 2019 16:10		23 Jul 2019 13:00	23 Jul 2019 18:58	1
HS19070994-27	WG-1620-MW27A-20190718	18 Jul 2019 17:15		23 Jul 2019 13:00	23 Jul 2019 19:18	1
HS19070994-28	WG-1620-MW27C-20190718	18 Jul 2019 18:05		23 Jul 2019 13:00	23 Jul 2019 19:37	1
HS19070994-29	WG-1620-MW51A-20190719	19 Jul 2019 07:25		23 Jul 2019 13:00	23 Jul 2019 19:57	1
HS19070994-30	WG-1620-MW51C-20190719	19 Jul 2019 08:55		23 Jul 2019 13:00	23 Jul 2019 20:17	1
HS19070994-31	WG-1620-MW81B-20190719	19 Jul 2019 09:45		23 Jul 2019 13:00	23 Jul 2019 20:37	1
HS19070994-32	WG-1620-MW50A-20190719	19 Jul 2019 10:40		23 Jul 2019 13:00	24 Jul 2019 15:41	1
<b>Batch ID: 143526 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19070994-01	WG-1620-FB04-20190716	16 Jul 2019 18:20		29 Jul 2019 14:00	30 Jul 2019 21:42	1
HS19070994-02	WG-1620-MW61A-20190717	17 Jul 2019 07:50		29 Jul 2019 14:00	30 Jul 2019 21:26	1
HS19070994-03	WG-1620-MW60A-20190717	17 Jul 2019 08:35		29 Jul 2019 14:00	30 Jul 2019 21:44	1
HS19070994-04	WG-1620-MW69A-20190717	17 Jul 2019 09:35		29 Jul 2019 14:00	30 Jul 2019 21:47	1
HS19070994-05	WG-1620-MW47C-20190717	17 Jul 2019 10:25		29 Jul 2019 14:00	30 Jul 2019 21:49	1
HS19070994-06	WG-1620-MW48C-20190717	17 Jul 2019 11:20		29 Jul 2019 14:00	30 Jul 2019 21:51	1
HS19070994-07	WG-1620-MW59A-20190717	17 Jul 2019 12:20		29 Jul 2019 14:00	30 Jul 2019 21:53	1
HS19070994-08	WG-1620-MW59B-20190717	17 Jul 2019 13:10		29 Jul 2019 14:00	30 Jul 2019 21:56	1
HS19070994-09	WG-1620-MW44A-20190717	17 Jul 2019 14:00		29 Jul 2019 14:00	30 Jul 2019 21:58	1
HS19070994-10	WG-1620-MW87C-20190717	17 Jul 2019 15:40		29 Jul 2019 14:00	30 Jul 2019 22:00	1
HS19070994-11	WG-1620-MW71B-20190717	17 Jul 2019 16:40		29 Jul 2019 14:00	30 Jul 2019 22:02	1
HS19070994-12	WG-1620-MW33A-20190717	17 Jul 2019 17:50		29 Jul 2019 14:00	30 Jul 2019 22:09	1
HS19070994-13	WG-1620-FD02-20190717	17 Jul 2019 17:50		29 Jul 2019 14:00	30 Jul 2019 22:11	1
HS19070994-14	WG-1620-MW26A-20190717	17 Jul 2019 18:45		29 Jul 2019 14:00	30 Jul 2019 22:14	1
HS19070994-15	WG-1620-FB05-20190717	17 Jul 2019 19:05		29 Jul 2019 14:00	30 Jul 2019 22:16	1
HS19070994-16	WG-1620-MW68A-20190718	18 Jul 2019 07:55		29 Jul 2019 14:00	30 Jul 2019 22:18	1
HS19070994-17	WG-1620-MW68B-20190718	18 Jul 2019 08:55		29 Jul 2019 14:00	30 Jul 2019 22:21	1
HS19070994-18	WG-1620-FD03-20190718	18 Jul 2019 08:55		29 Jul 2019 14:00	30 Jul 2019 22:23	1
HS19070994-19	WG-1620-MW68C-20190718	18 Jul 2019 09:45		29 Jul 2019 14:00	30 Jul 2019 22:25	1
HS19070994-20	WG-1620-MW83B-20190718	18 Jul 2019 10:40		29 Jul 2019 14:00	30 Jul 2019 22:27	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143558 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19070994-21	WG-1620-MW83C-20190718	18 Jul 2019 11:30		30 Jul 2019 12:00	30 Jul 2019 22:41	1
HS19070994-22	WG-1620-MW35A-20190718	18 Jul 2019 12:40		30 Jul 2019 12:00	30 Jul 2019 22:57	1
HS19070994-23	WG-1620-MW35B-20190718	18 Jul 2019 13:30		30 Jul 2019 12:00	30 Jul 2019 22:59	1
HS19070994-24	WG-1620-MW90B-20190718	18 Jul 2019 14:25		30 Jul 2019 12:00	30 Jul 2019 23:01	1
HS19070994-25	WG-1620-MW89B-20190718	18 Jul 2019 15:15		30 Jul 2019 12:00	30 Jul 2019 23:04	1
HS19070994-26	WG-1620-MW38B-20190718	18 Jul 2019 16:10		30 Jul 2019 12:00	30 Jul 2019 23:06	1
HS19070994-27	WG-1620-MW27A-20190718	18 Jul 2019 17:15		30 Jul 2019 12:00	30 Jul 2019 23:08	1
HS19070994-28	WG-1620-MW27C-20190718	18 Jul 2019 18:05		30 Jul 2019 12:00	30 Jul 2019 23:10	1
HS19070994-29	WG-1620-MW51A-20190719	19 Jul 2019 07:25		30 Jul 2019 12:00	30 Jul 2019 23:13	1
HS19070994-30	WG-1620-MW51C-20190719	19 Jul 2019 08:55		30 Jul 2019 12:00	30 Jul 2019 23:15	1
HS19070994-31	WG-1620-MW81B-20190719	19 Jul 2019 09:45		30 Jul 2019 12:00	30 Jul 2019 23:17	1
HS19070994-32	WG-1620-MW50A-20190719	19 Jul 2019 10:40		30 Jul 2019 12:00	30 Jul 2019 23:24	1
<b>Batch ID: R342868 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19070994-19	WG-1620-MW68C-20190718	18 Jul 2019 09:45			23 Jul 2019 02:01	1
HS19070994-20	WG-1620-MW83B-20190718	18 Jul 2019 10:40			23 Jul 2019 06:30	1
<b>Batch ID: R342877 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19070994-21	WG-1620-MW83C-20190718	18 Jul 2019 11:30			23 Jul 2019 03:13	1
HS19070994-22	WG-1620-MW35A-20190718	18 Jul 2019 12:40			23 Jul 2019 04:25	1
HS19070994-23	WG-1620-MW35B-20190718	18 Jul 2019 13:30			23 Jul 2019 04:49	1
HS19070994-24	WG-1620-MW90B-20190718	18 Jul 2019 14:25			23 Jul 2019 05:13	1
HS19070994-25	WG-1620-MW89B-20190718	18 Jul 2019 15:15			23 Jul 2019 05:37	1
HS19070994-26	WG-1620-MW38B-20190718	18 Jul 2019 16:10			23 Jul 2019 06:01	1
HS19070994-27	WG-1620-MW27A-20190718	18 Jul 2019 17:15			23 Jul 2019 06:25	1
HS19070994-28	WG-1620-MW27C-20190718	18 Jul 2019 18:05			23 Jul 2019 06:48	1
HS19070994-29	WG-1620-MW51A-20190719	19 Jul 2019 07:25			23 Jul 2019 07:12	1
HS19070994-30	WG-1620-MW51C-20190719	19 Jul 2019 08:55			23 Jul 2019 07:36	1
HS19070994-31	WG-1620-MW81B-20190719	19 Jul 2019 09:45			23 Jul 2019 08:00	1
HS19070994-32	WG-1620-MW50A-20190719	19 Jul 2019 10:40			23 Jul 2019 08:24	1
HS19070994-33	WG-1620-TB05-20190719	19 Jul 2019 00:00			23 Jul 2019 02:25	1
HS19070994-34	WG-1620-TB06-20190719	19 Jul 2019 00:00			23 Jul 2019 02:49	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: R342879 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19070994-01	WG-1620-FB04-20190716	16 Jul 2019 18:20			23 Jul 2019 00:42	1
HS19070994-02	WG-1620-MW61A-20190717	17 Jul 2019 07:50			23 Jul 2019 01:32	1
HS19070994-03	WG-1620-MW60A-20190717	17 Jul 2019 08:35			23 Jul 2019 03:10	1
HS19070994-04	WG-1620-MW69A-20190717	17 Jul 2019 09:35			23 Jul 2019 03:35	1
HS19070994-05	WG-1620-MW47C-20190717	17 Jul 2019 10:25			23 Jul 2019 03:59	1
HS19070994-06	WG-1620-MW48C-20190717	17 Jul 2019 11:20			23 Jul 2019 04:24	1
HS19070994-07	WG-1620-MW59A-20190717	17 Jul 2019 12:20			23 Jul 2019 04:48	1
HS19070994-08	WG-1620-MW59B-20190717	17 Jul 2019 13:10			23 Jul 2019 05:13	1
HS19070994-09	WG-1620-MW44A-20190717	17 Jul 2019 14:00			23 Jul 2019 05:37	1
HS19070994-10	WG-1620-MW87C-20190717	17 Jul 2019 15:40			23 Jul 2019 06:02	1
HS19070994-11	WG-1620-MW71B-20190717	17 Jul 2019 16:40			23 Jul 2019 06:26	1
HS19070994-12	WG-1620-MW33A-20190717	17 Jul 2019 17:50			23 Jul 2019 06:51	1
HS19070994-13	WG-1620-FD02-20190717	17 Jul 2019 17:50			23 Jul 2019 07:16	1
HS19070994-14	WG-1620-MW26A-20190717	17 Jul 2019 18:45			23 Jul 2019 07:40	1
HS19070994-15	WG-1620-FB05-20190717	17 Jul 2019 19:05			23 Jul 2019 01:07	1
HS19070994-16	WG-1620-MW68A-20190718	18 Jul 2019 07:55			23 Jul 2019 08:05	1
HS19070994-17	WG-1620-MW68B-20190718	18 Jul 2019 08:55			23 Jul 2019 08:59	100
HS19070994-17	WG-1620-MW68B-20190718	18 Jul 2019 08:55			23 Jul 2019 08:32	10
HS19070994-18	WG-1620-FD03-20190718	18 Jul 2019 08:55			23 Jul 2019 09:53	100
HS19070994-18	WG-1620-FD03-20190718	18 Jul 2019 08:55			23 Jul 2019 09:26	10

WorkOrder: HS19070994  
InstrumentID: ICPMS05  
Test Code: ICP\_TW  
Test Number: SW6020  
Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200



WorkOrder: HS19070994  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

Matrix: Aqueous

Units: mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00010	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000096	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00012	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00011	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.00010	0.00012	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000013	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00013	0.000047	0.0010
A	Acenaphthene	83-32-9	0.00010	0.00011	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.00010	0.00010	0.000015	0.00010
A	Anthracene	120-12-7	0.00010	0.00011	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.00010	0.000099	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.00010	0.000076	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000086	0.000037	0.00020
A	Chrysene	218-01-9	0.00010	0.00011	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.00010	0.000094	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.000099	0.000020	0.00020
A	Fluoranthene	206-44-0	0.00010	0.00012	0.000010	0.00010
A	Fluorene	86-73-7	0.00010	0.00012	0.000030	0.00010
A	Naphthalene	91-20-3	0.00010	0.00010	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00013	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000085	0.000079	0.00020
A	Phenanthrene	85-01-8	0.00010	0.00011	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000094	0.000035	0.00020
A	Pyrene	129-00-0	0.00010	0.00011	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19070994  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00064	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00060	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00061	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00077	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00058	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00070	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00062	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS19070994  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00067	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00056	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00059	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00051	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00054	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00056	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00052	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS19070994  
 InstrumentID: VOA6  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00065	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00062	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00061	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00062	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00067	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00060	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00049	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.0019	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

<b>Batch ID:</b> 143526 ( 0 )	<b>Instrument:</b> ICPMS05	<b>Method:</b> ICP-MS METALS BY SW6020A
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<b>MBLK</b>	Sample ID: <b>MBLK-143526</b>	Units: <b>mg/L</b>	Analysis Date: <b>30-Jul-2019 21:22</b>							
Client ID:	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187971</b>	PrepDate: <b>29-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic U 0.00200

<b>LCS</b>	Sample ID: <b>LCS-143526</b>	Units: <b>mg/L</b>	Analysis Date: <b>30-Jul-2019 21:24</b>							
Client ID:	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187972</b>	PrepDate: <b>29-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.04941 0.00200 0.05 0 98.8 80 - 120

<b>MS</b>	Sample ID: <b>HS19070994-02MS</b>	Units: <b>mg/L</b>	Analysis Date: <b>30-Jul-2019 21:31</b>							
Client ID: <b>WG-1620-MW61A-20190717</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187975</b>	PrepDate: <b>29-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.0527 0.00200 0.05 0.00117 103 80 - 120

<b>MSD</b>	Sample ID: <b>HS19070994-02MSD</b>	Units: <b>mg/L</b>	Analysis Date: <b>30-Jul-2019 21:33</b>							
Client ID: <b>WG-1620-MW61A-20190717</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187976</b>	PrepDate: <b>29-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.05075 0.00200 0.05 0.00117 99.2 80 - 120 0.0527 3.77 20

<b>PDS</b>	Sample ID: <b>HS19070994-02PDS</b>	Units: <b>mg/L</b>	Analysis Date: <b>30-Jul-2019 21:35</b>							
Client ID: <b>WG-1620-MW61A-20190717</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187977</b>	PrepDate: <b>29-Jul-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.1026 0.00200 0.1 0.00117 101 75 - 125

<b>SD</b>	Sample ID: <b>HS19070994-02SD</b>	Units: <b>mg/L</b>	Analysis Date: <b>30-Jul-2019 21:29</b>							
Client ID: <b>WG-1620-MW61A-20190717</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187974</b>	PrepDate: <b>29-Jul-2019</b> DF: <b>5</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual

Arsenic U 0.0100 0.00117 0 10

<b>The following samples were analyzed in this batch:</b>	HS19070994-01	HS19070994-02	HS19070994-03	HS19070994-04
	HS19070994-05	HS19070994-06	HS19070994-07	HS19070994-08
	HS19070994-09	HS19070994-10	HS19070994-11	HS19070994-12
	HS19070994-13	HS19070994-14	HS19070994-15	HS19070994-16
	HS19070994-17	HS19070994-18	HS19070994-19	HS19070994-20

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143558 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-143558</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Jul-2019 22:36</b>					
Client ID:	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187939</b>	PrepDate: <b>30-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-143558</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Jul-2019 22:39</b>					
Client ID:	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187940</b>	PrepDate: <b>30-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05011	0.00200	0.05	0	100	80 - 120			
<b>MS</b>	Sample ID: <b>HS19070994-21MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Jul-2019 22:45</b>					
Client ID: <b>WG-1620-MW83C-20190718</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187943</b>	PrepDate: <b>30-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05667	0.00200	0.05	0.006174	101	80 - 120			
<b>MSD</b>	Sample ID: <b>HS19070994-21MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Jul-2019 22:48</b>					
Client ID: <b>WG-1620-MW83C-20190718</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187944</b>	PrepDate: <b>30-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05623	0.00200	0.05	0.006174	100	80 - 120	0.05667	0.771	20
<b>PDS</b>	Sample ID: <b>HS19070994-21PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Jul-2019 22:50</b>					
Client ID: <b>WG-1620-MW83C-20190718</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187945</b>	PrepDate: <b>30-Jul-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.1089	0.00200	0.1	0.006174	103	75 - 125			
<b>SD</b>	Sample ID: <b>HS19070994-21SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Jul-2019 22:43</b>					
Client ID: <b>WG-1620-MW83C-20190718</b>	Run ID: <b>ICPMS05_343276</b>	SeqNo: <b>5187942</b>	PrepDate: <b>30-Jul-2019</b>	DF: <b>5</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	0.006418	0.0100					0.006174	0 10	J

The following samples were analyzed in this batch:

HS19070994-21	HS19070994-22	HS19070994-23	HS19070994-24
HS19070994-25	HS19070994-26	HS19070994-27	HS19070994-28
HS19070994-29	HS19070994-30	HS19070994-31	HS19070994-32

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143239 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143239	Units: ug/L			Analysis Date: 23-Jul-2019 11:01					
Client ID:	Run ID: SV-7_342905	SeqNo: 5177235	PrepDate: 22-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.901</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.092</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.088</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.374</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.5</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.314</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>106</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.31</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>106</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143239 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143239	Units: ug/L			Analysis Date: 23-Jul-2019 12:01					
Client ID:	Run ID: SV-7_342905	SeqNo: 5177237		PrepDate: 22-Jul-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.199	0.20	5	0	84.0	39 - 127				
2,4-Dimethylphenol	4.42	0.20	5	0	88.4	35 - 120				
2,4-Dinitrotoluene	5.269	0.20	5	0	105	50 - 122				
2,6-Dinitrotoluene	4.941	0.20	5	0	98.8	50 - 120				
2-Chloronaphthalene	4.227	0.20	5	0	84.5	50 - 120				
2-Methylnaphthalene	4.244	0.10	5	0	84.9	50 - 120				
4,6-Dinitro-2-methylphenol	4.356	0.20	5	0	87.1	25 - 121				
4-Nitrophenol	5.086	1.0	5	0	102	30 - 130				
Acenaphthene	4.208	0.10	5	0	84.2	45 - 120				
Acenaphthylene	4.309	0.10	5	0	86.2	47 - 120				
Anthracene	4.111	0.10	5	0	82.2	45 - 120				
Benz(a)anthracene	4.208	0.10	5	0	84.2	40 - 120				
Benzo(a)pyrene	4.855	0.10	5	0	97.1	45 - 120				
Bis(2-chloroethoxy)methane	4.151	0.20	5	0	83.0	45 - 120				
Bis(2-ethylhexyl)phthalate	3.673	0.20	5	0	73.5	40 - 139				
Chrysene	4.083	0.10	5	0	81.7	43 - 120				
Dibenzofuran	4.51	0.10	5	0	90.2	50 - 120				
Di-n-butyl phthalate	3.819	0.20	5	0	76.4	45 - 123				
Fluoranthene	4.501	0.10	5	0	90.0	45 - 125				
Fluorene	4.41	0.10	5	0	88.2	49 - 120				
Naphthalene	4.048	0.10	5	0	81.0	45 - 120				
Nitrobenzene	4.645	0.20	5	0	92.9	44 - 120				
N-Nitrosodiphenylamine	4.632	0.20	5	0	92.6	40 - 125				
Pentachlorophenol	3.799	0.20	5	0	76.0	19 - 121				
Phenanthrene	3.815	0.10	5	0	76.3	45 - 121				
Phenol	4.879	0.20	5	0	97.6	20 - 124				
Pyrene	4.227	0.10	5	0	84.5	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.4</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>108</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.924</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.5</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.63</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.6</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.258</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.2</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.988</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.8</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.053</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>20 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143239 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS19070994-02MS		Units: ug/L		Analysis Date: 23-Jul-2019 12:41				
Client ID: WG-1620-MW61A-20190717		Run ID: SV-7_342905		SeqNo: 5177239		PrepDate: 22-Jul-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.415	0.20	5	0	88.3	39 - 127				
2,4-Dimethylphenol	4.037	0.20	5	0	80.7	35 - 120				
2,4-Dinitrotoluene	5.451	0.20	5	0	109	50 - 122				
2,6-Dinitrotoluene	4.473	0.20	5	0	89.5	50 - 120				
2-Chloronaphthalene	3.33	0.20	5	0	66.6	50 - 120				
2-Methylnaphthalene	3.896	0.10	5	0	77.9	50 - 120				
4,6-Dinitro-2-methylphenol	5.429	0.20	5	0	109	25 - 121				
4-Nitrophenol	5.9	1.0	5	0	118	30 - 130				
Acenaphthene	3.872	0.10	5	0	77.4	45 - 120				
Acenaphthylene	3.749	0.10	5	0	75.0	47 - 120				
Anthracene	4.161	0.10	5	0	83.2	45 - 120				
Benz(a)anthracene	5.184	0.10	5	0	104	40 - 120				
Benzo(a)pyrene	5.7	0.10	5	0	114	45 - 120				
Bis(2-chloroethoxy)methane	3.56	0.20	5	0	71.2	45 - 120				
Bis(2-ethylhexyl)phthalate	4.733	0.20	5	0.3733	87.2	40 - 139				
Chrysene	5.104	0.10	5	0	102	43 - 120				
Dibenzofuran	4.1	0.10	5	0	82.0	50 - 120				
Di-n-butyl phthalate	4.681	0.20	5	0.0555	92.5	45 - 123				
Fluoranthene	5.584	0.10	5	0	112	45 - 125				
Fluorene	3.523	0.10	5	0	70.5	49 - 120				
Naphthalene	3.717	0.10	5	0	74.3	45 - 120				
Nitrobenzene	4.017	0.20	5	0	80.3	44 - 120				
N-Nitrosodiphenylamine	5.187	0.20	5	0	104	40 - 125				
Pentachlorophenol	5.322	0.20	5	0	106	19 - 121				
Phenanthrene	4.06	0.10	5	0	81.2	45 - 121				
Phenol	3.917	0.20	5	0.1053	76.2	20 - 124				
Pyrene	4.987	0.10	5	0	99.7	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.013</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>100</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.557</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>71.1</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.895</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.9</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.264</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.008</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>80.2</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.273</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.5</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143239 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS19070994-02MSD		Units: ug/L		Analysis Date: 23-Jul-2019 13:00				
Client ID: WG-1620-MW61A-20190717		Run ID: SV-7_342905		SeqNo: 5177240		PrepDate: 22-Jul-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	3.835	0.20	5	0	76.7	39 - 127	4.415	14.1	20	
2,4-Dimethylphenol	4.214	0.20	5	0	84.3	35 - 120	4.037	4.29	20	
2,4-Dinitrotoluene	5.713	0.20	5	0	114	50 - 122	5.451	4.68	20	
2,6-Dinitrotoluene	5.485	0.20	5	0	110	50 - 120	4.473	20.3	20 R	
2-Chloronaphthalene	3.935	0.20	5	0	78.7	50 - 120	3.33	16.7	20	
2-Methylnaphthalene	4.33	0.10	5	0	86.6	50 - 120	3.896	10.6	20	
4,6-Dinitro-2-methylphenol	4.985	0.20	5	0	99.7	25 - 121	5.429	8.53	30	
4-Nitrophenol	5.681	1.0	5	0	114	30 - 130	5.9	3.79	20	
Acenaphthene	3.837	0.10	5	0	76.7	45 - 120	3.872	0.909	20	
Acenaphthylene	4.093	0.10	5	0	81.9	47 - 120	3.749	8.78	20	
Anthracene	4.614	0.10	5	0	92.3	45 - 120	4.161	10.3	20	
Benz(a)anthracene	5.589	0.10	5	0	112	40 - 120	5.184	7.52	20	
Benzo(a)pyrene	7.055	0.10	5	0	141	45 - 120	5.7	21.2	20 SR	
Bis(2-chloroethoxy)methane	3.926	0.20	5	0	78.5	45 - 120	3.56	9.79	20	
Bis(2-ethylhexyl)phthalate	5.527	0.20	5	0.3733	103	40 - 139	4.733	15.5	20	
Chrysene	5.555	0.10	5	0	111	43 - 120	5.104	8.47	20	
Dibenzofuran	4.216	0.10	5	0	84.3	50 - 120	4.1	2.79	20	
Di-n-butyl phthalate	5.227	0.20	5	0.0555	103	45 - 123	4.681	11	20	
Fluoranthene	5.526	0.10	5	0	111	45 - 125	5.584	1.05	20	
Fluorene	4.025	0.10	5	0	80.5	49 - 120	3.523	13.3	20	
Naphthalene	4.11	0.10	5	0	82.2	45 - 120	3.717	10	20	
Nitrobenzene	4.765	0.20	5	0	95.3	44 - 120	4.017	17	20	
N-Nitrosodiphenylamine	4.526	0.20	5	0	90.5	40 - 125	5.187	13.6	20	
Pentachlorophenol	5.029	0.20	5	0	101	19 - 121	5.322	5.66	20	
Phenanthrene	4.459	0.10	5	0	89.2	45 - 121	4.06	9.37	20	
Phenol	4.499	0.20	5	0.1053	87.9	20 - 124	3.917	13.8	20	
Pyrene	4.913	0.10	5	0	98.3	40 - 130	4.987	1.49	20	
Surr: 2,4,6-Tribromophenol	6.449	0.20	5	0	129	34 - 129	5.013	25.1	20 R	
Surr: 2-Fluorobiphenyl	4.764	0.20	5	0	95.3	40 - 125	3.557	29	20 R	
Surr: 2-Fluorophenol	4.185	0.20	5	0	83.7	20 - 120	3.895	7.17	20	
Surr: 4-Terphenyl-d14	4.986	0.20	5	0	99.7	40 - 135	5.264	5.43	20	
Surr: Nitrobenzene-d5	4.82	0.20	5	0	96.4	41 - 120	4.008	18.4	20	
Surr: Phenol-d6	4.591	0.20	5	0	91.8	20 - 120	4.273	7.17	20	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

**Batch ID:** 143239 ( 0 )      **Instrument:** SV-7      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

The following samples were analyzed in this batch:

HS19070994-01	HS19070994-02	HS19070994-03	HS19070994-04
HS19070994-05	HS19070994-06	HS19070994-07	HS19070994-08
HS19070994-09	HS19070994-10	HS19070994-11	HS19070994-12
HS19070994-13	HS19070994-14	HS19070994-15	HS19070994-16
HS19070994-17	HS19070994-18	HS19070994-19	HS19070994-20

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143295 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143295	Units: ug/L			Analysis Date: 24-Jul-2019 11:06					
Client ID:	Run ID: SV-7_342956	SeqNo: 5178273	PrepDate: 23-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	4.93	0.20	5	0	98.6	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.783	0.20	5	0	95.7	40 - 125				
<i>Surr: 2-Fluorophenol</i>	4.587	0.20	5	0	91.7	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.364	0.20	5	0	87.3	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	5.101	0.20	5	0	102	41 - 120				
<i>Surr: Phenol-d6</i>	4.662	0.20	5	0	93.2	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143295 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143295	Units: ug/L			Analysis Date: 24-Jul-2019 11:26					
Client ID:	Run ID: SV-7_342956	SeqNo: 5178274	PrepDate: 23-Jul-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.27	0.20	5	0	85.4	39 - 127				
2,4-Dimethylphenol	4.441	0.20	5	0	88.8	35 - 120				
2,4-Dinitrotoluene	5.322	0.20	5	0	106	50 - 122				
2,6-Dinitrotoluene	5.12	0.20	5	0	102	50 - 120				
2-Chloronaphthalene	4.199	0.20	5	0	84.0	50 - 120				
2-Methylnaphthalene	4.573	0.10	5	0	91.5	50 - 120				
4,6-Dinitro-2-methylphenol	4.011	0.20	5	0	80.2	25 - 121				
4-Nitrophenol	4.862	1.0	5	0	97.2	30 - 130				
Acenaphthene	4.17	0.10	5	0	83.4	45 - 120				
Acenaphthylene	4.231	0.10	5	0	84.6	47 - 120				
Anthracene	4.089	0.10	5	0	81.8	45 - 120				
Benz(a)anthracene	4.331	0.10	5	0	86.6	40 - 120				
Benzo(a)pyrene	4.809	0.10	5	0	96.2	45 - 120				
Bis(2-chloroethoxy)methane	4.11	0.20	5	0	82.2	45 - 120				
Bis(2-ethylhexyl)phthalate	3.802	0.20	5	0	76.0	40 - 139				
Chrysene	4.35	0.10	5	0	87.0	43 - 120				
Dibenzofuran	4.37	0.10	5	0	87.4	50 - 120				
Di-n-butyl phthalate	4.066	0.20	5	0	81.3	45 - 123				
Fluoranthene	4.558	0.10	5	0	91.2	45 - 125				
Fluorene	4.23	0.10	5	0	84.6	49 - 120				
Naphthalene	4.296	0.10	5	0	85.9	45 - 120				
Nitrobenzene	4.503	0.20	5	0	90.1	44 - 120				
N-Nitrosodiphenylamine	4.726	0.20	5	0	94.5	40 - 125				
Pentachlorophenol	3.475	0.20	5	0	69.5	19 - 121				
Phenanthrene	4.047	0.10	5	0	80.9	45 - 121				
Phenol	4.067	0.20	5	0	81.3	20 - 124				
Pyrene	4.074	0.10	5	0	81.5	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.763</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>115</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.062</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.062</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.484</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.7</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.287</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>106</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.182</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: 143295 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-143295		Units: ug/L		Analysis Date: 24-Jul-2019 11:45				
Client ID:		Run ID: SV-7_342956		SeqNo: 5178275		PrepDate: 23-Jul-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.086	0.20	5	0	81.7	39 - 127	4.27	4.39	20	
2,4-Dimethylphenol	3.98	0.20	5	0	79.6	35 - 120	4.441	10.9	20	
2,4-Dinitrotoluene	4.984	0.20	5	0	99.7	50 - 122	5.322	6.57	20	
2,6-Dinitrotoluene	4.803	0.20	5	0	96.1	50 - 120	5.12	6.38	20	
2-Chloronaphthalene	4.174	0.20	5	0	83.5	50 - 120	4.199	0.592	20	
2-Methylnaphthalene	4.27	0.10	5	0	85.4	50 - 120	4.573	6.86	20	
4,6-Dinitro-2-methylphenol	4.202	0.20	5	0	84.0	25 - 121	4.011	4.64	30	
4-Nitrophenol	4.668	1.0	5	0	93.4	30 - 130	4.862	4.08	20	
Acenaphthene	4.187	0.10	5	0	83.7	45 - 120	4.17	0.401	20	
Acenaphthylene	4.194	0.10	5	0	83.9	47 - 120	4.231	0.87	20	
Anthracene	3.968	0.10	5	0	79.4	45 - 120	4.089	2.99	20	
Benz(a)anthracene	4.302	0.10	5	0	86.0	40 - 120	4.331	0.675	20	
Benzo(a)pyrene	4.814	0.10	5	0	96.3	45 - 120	4.809	0.104	20	
Bis(2-chloroethoxy)methane	3.975	0.20	5	0	79.5	45 - 120	4.11	3.33	20	
Bis(2-ethylhexyl)phthalate	4.142	0.20	5	0	82.8	40 - 139	3.802	8.55	20	
Chrysene	4.582	0.10	5	0	91.6	43 - 120	4.35	5.2	20	
Dibenzofuran	4.408	0.10	5	0	88.2	50 - 120	4.37	0.854	20	
Di-n-butyl phthalate	3.886	0.20	5	0	77.7	45 - 123	4.066	4.51	20	
Fluoranthene	4.17	0.10	5	0	83.4	45 - 125	4.558	8.9	20	
Fluorene	4.027	0.10	5	0	80.5	49 - 120	4.23	4.91	20	
Naphthalene	4.055	0.10	5	0	81.1	45 - 120	4.296	5.76	20	
Nitrobenzene	4.324	0.20	5	0	86.5	44 - 120	4.503	4.06	20	
N-Nitrosodiphenylamine	4.328	0.20	5	0	86.6	40 - 125	4.726	8.8	20	
Pentachlorophenol	3.469	0.20	5	0	69.4	19 - 121	3.475	0.178	20	
Phenanthrene	3.88	0.10	5	0	77.6	45 - 121	4.047	4.21	20	
Phenol	3.851	0.20	5	0	77.0	20 - 124	4.067	5.47	20	
Pyrene	4.178	0.10	5	0	83.6	40 - 130	4.074	2.54	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.553</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>34 - 129</i>	<i>5.763</i>	<i>3.72</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.19</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>40 - 125</i>	<i>5.062</i>	<i>2.49</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.776</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.5</i>	<i>20 - 120</i>	<i>5.062</i>	<i>5.8</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.505</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.1</i>	<i>40 - 135</i>	<i>4.484</i>	<i>0.476</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>4.96</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.2</i>	<i>41 - 120</i>	<i>5.287</i>	<i>6.39</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.968</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.4</i>	<i>20 - 120</i>	<i>5.182</i>	<i>4.22</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

<b>Batch ID:</b> 143295 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS19070994-21	HS19070994-22	HS19070994-23	HS19070994-24
HS19070994-25	HS19070994-26	HS19070994-27	HS19070994-28
HS19070994-29	HS19070994-30	HS19070994-31	HS19070994-32

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

**Batch ID:** R342868 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190722</b>			Units: <b>ug/L</b>		Analysis Date: <b>22-Jul-2019 23:59</b>			
Client ID:		Run ID: <b>VOA2_342868</b>			SeqNo: <b>5176537</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>44.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>89.3</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.5</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.12</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.2</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-190722</b>			Units: <b>ug/L</b>		Analysis Date: <b>22-Jul-2019 23:34</b>			
Client ID:		Run ID: <b>VOA2_342868</b>			SeqNo: <b>5176536</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	19.73	1.0	20	0	98.6	70 - 124				
Benzene	19.51	1.0	20	0	97.6	74 - 120				
Chlorobenzene	18.74	1.0	20	0	93.7	76 - 113				
Ethylbenzene	19.69	1.0	20	0	98.5	77 - 117				
Methylene chloride	22.08	2.0	20	0	110	70 - 127				
Toluene	21.86	1.0	20	0	109	77 - 118				
Xylenes, Total	62.78	1.0	60	0	105	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.76</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.9</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.22</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.4</i>	<i>81 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

**Batch ID:** R342868 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS19070949-09MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>23-Jul-2019 01:12</b>			
Client ID:		Run ID: <b>VOA2_342868</b>			SeqNo: <b>5176540</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.08	1.0	20	0	80.4	70 - 127				
Benzene	16.82	1.0	20	0	84.1	70 - 127				
Chlorobenzene	16.71	1.0	20	0	83.5	70 - 114				
Ethylbenzene	17.05	1.0	20	0	85.2	70 - 124				
Methylene chloride	16.4	2.0	20	0	82.0	70 - 128				
Toluene	18.99	1.0	20	0	94.9	70 - 123				
Xylenes, Total	55.12	1.0	60	0	91.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.06</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.1</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.34</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.7</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS19070949-09MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>23-Jul-2019 01:36</b>			
Client ID:		Run ID: <b>VOA2_342868</b>			SeqNo: <b>5176541</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.97	1.0	20	0	79.9	70 - 127	16.08	0.65	20	
Benzene	16.22	1.0	20	0	81.1	70 - 127	16.82	3.64	20	
Chlorobenzene	16.12	1.0	20	0	80.6	70 - 114	16.71	3.54	20	
Ethylbenzene	16.48	1.0	20	0	82.4	70 - 124	17.05	3.37	20	
Methylene chloride	17.05	2.0	20	0	85.2	70 - 128	16.4	3.88	20	
Toluene	18.03	1.0	20	0	90.2	70 - 123	18.99	5.16	20	
Xylenes, Total	53.8	1.0	60	0	89.7	70 - 130	55.12	2.42	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.15</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.3</i>	<i>70 - 126</i>	<i>46.06</i>	<i>0.195</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>81 - 113</i>	<i>47.94</i>	<i>0.0444</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.5</i>	<i>77 - 123</i>	<i>47.34</i>	<i>0.827</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>82 - 127</i>	<i>49.78</i>	<i>0.281</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19070994-19      HS19070994-20

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

<b>Batch ID:</b> R342877 ( 0 )		<b>Instrument:</b> VOA6		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-190722</b>	Units: <b>ug/L</b>			Analysis Date: <b>23-Jul-2019 02:01</b>				
Client ID:	Run ID: <b>VOA6_342877</b>	SeqNo: <b>5176718</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.06</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.41</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>46.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.8</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>52.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-190722</b>	Units: <b>ug/L</b>			Analysis Date: <b>23-Jul-2019 01:13</b>				
Client ID:	Run ID: <b>VOA6_342877</b>	SeqNo: <b>5176717</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	20.99	1.0	20	0	105	70 - 124			
Benzene	20.69	1.0	20	0	103	74 - 120			
Chlorobenzene	21.43	1.0	20	0	107	76 - 113			
Ethylbenzene	22.09	1.0	20	0	110	77 - 117			
Methylene chloride	20.58	2.0	20	0	103	70 - 127			
Toluene	21.9	1.0	20	0	110	77 - 118			
Vinyl chloride	19	1.0	20	0	95.0	70 - 130			
Xylenes, Total	65.67	1.0	60	0	109	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.16</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.3</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>45.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.4</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>46.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.6</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

Batch ID: R342877 ( 0 )		Instrument: VOA6		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>		Sample ID: <b>HS19070994-21MS</b>		Units: <b>ug/L</b>		Analysis Date: <b>23-Jul-2019 03:37</b>				
Client ID: <b>WG-1620-MW83C-20190718</b>		Run ID: <b>VOA6_342877</b>		SeqNo: <b>5176722</b>		PrepDate:		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	20.44	1.0	20	0	102	70 - 127				
Benzene	20.94	1.0	20	0	105	70 - 127				
Chlorobenzene	21.37	1.0	20	0	107	70 - 114				
Ethylbenzene	22.48	1.0	20	0	112	70 - 124				
Methylene chloride	20.11	2.0	20	0	101	70 - 128				
Toluene	22.22	1.0	20	0	111	70 - 123				
Vinyl chloride	17.71	1.0	20	0	88.6	70 - 130				
Xylenes, Total	65.87	1.0	60	0	110	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>55.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>112</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>57.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>114</i>	<i>81 - 113</i>			<b>S</b>	
<i>Surr: Dibromofluoromethane</i>	<i>56.54</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>113</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>58.54</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>117</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS19070994-21MSD</b>		Units: <b>ug/L</b>		Analysis Date: <b>23-Jul-2019 04:01</b>			
Client ID: <b>WG-1620-MW83C-20190718</b>		Run ID: <b>VOA6_342877</b>		SeqNo: <b>5176723</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	20.9	1.0	20	0	104	70 - 127	20.44	2.23	20
Benzene	20.94	1.0	20	0	105	70 - 127	20.94	0.0341	20
Chlorobenzene	21.3	1.0	20	0	107	70 - 114	21.37	0.321	20
Ethylbenzene	22.27	1.0	20	0	111	70 - 124	22.48	0.963	20
Methylene chloride	19.56	2.0	20	0	97.8	70 - 128	20.11	2.8	20
Toluene	22.05	1.0	20	0	110	70 - 123	22.22	0.788	20
Vinyl chloride	17.32	1.0	20	0	86.6	70 - 130	17.71	2.25	20
Xylenes, Total	66.41	1.0	60	0	111	70 - 130	65.87	0.815	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>57.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>115</i>	<i>70 - 126</i>	<i>55.84</i>	<i>2.72</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>57.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>115</i>	<i>81 - 113</i>	<i>57.14</i>	<i>0.606</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>56.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>114</i>	<i>77 - 123</i>	<i>56.54</i>	<i>0.628</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>58.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>116</i>	<i>82 - 127</i>	<i>58.54</i>	<i>0.502</i>	<i>20</i>

The following samples were analyzed in this batch:

HS19070994-21	HS19070994-22	HS19070994-23	HS19070994-24
HS19070994-25	HS19070994-26	HS19070994-27	HS19070994-28
HS19070994-29	HS19070994-30	HS19070994-31	HS19070994-32
HS19070994-33	HS19070994-34		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

<b>Batch ID:</b> R342879 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-190722</b>	Units: <b>ug/L</b>			Analysis Date: <b>23-Jul-2019 00:18</b>				
Client ID:	Run ID: <b>VOA4_342879</b>	SeqNo: <b>5176775</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.5</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>53.07</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>106</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.85</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-190722</b>	Units: <b>ug/L</b>			Analysis Date: <b>22-Jul-2019 23:29</b>				
Client ID:	Run ID: <b>VOA4_342879</b>	SeqNo: <b>5176774</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	17.24	1.0	20	0	86.2	70 - 124			
Benzene	18.53	1.0	20	0	92.7	74 - 120			
Chlorobenzene	18.8	1.0	20	0	94.0	76 - 113			
Ethylbenzene	18.76	1.0	20	0	93.8	77 - 117			
Methylene chloride	21.91	2.0	20	0	110	70 - 127			
Toluene	18.63	1.0	20	0	93.2	77 - 118			
Vinyl chloride	17.76	1.0	20	0	88.8	70 - 130			
Xylenes, Total	58.77	1.0	60	0	97.9	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.6</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.3</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>52.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QC BATCH REPORT**

**Batch ID:** R342879 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS19070994-02MS			Units: ug/L		Analysis Date: 23-Jul-2019 01:56			
Client ID: WG-1620-MW61A-20190717		Run ID: VOA4_342879			SeqNo: 5176779		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.52	1.0	20	0	72.6	70 - 127				
Benzene	16.38	1.0	20	0	81.9	70 - 127				
Chlorobenzene	16.17	1.0	20	0	80.8	70 - 114				
Ethylbenzene	16.41	1.0	20	0	82.0	70 - 124				
Methylene chloride	16.8	2.0	20	0	84.0	70 - 128				
Toluene	16.66	1.0	20	0	83.3	70 - 123				
Vinyl chloride	16.09	1.0	20	0	80.4	70 - 130				
Xylenes, Total	52.27	1.0	60	0	87.1	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>52.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.27</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

MSD		Sample ID: HS19070994-02MSD			Units: ug/L		Analysis Date: 23-Jul-2019 02:21			
Client ID: WG-1620-MW61A-20190717		Run ID: VOA4_342879			SeqNo: 5176780		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	13.98	1.0	20	0	69.9	70 - 127	14.52	3.83	20	S
Benzene	15.56	1.0	20	0	77.8	70 - 127	16.38	5.12	20	
Chlorobenzene	15.68	1.0	20	0	78.4	70 - 114	16.17	3.03	20	
Ethylbenzene	15.79	1.0	20	0	79.0	70 - 124	16.41	3.81	20	
Methylene chloride	15.84	2.0	20	0	79.2	70 - 128	16.8	5.86	20	
Toluene	15.71	1.0	20	0	78.5	70 - 123	16.66	5.9	20	
Vinyl chloride	15.34	1.0	20	0	76.7	70 - 130	16.09	4.78	20	
Xylenes, Total	51.57	1.0	60	0	86.0	70 - 130	52.27	1.34	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.3</i>	<i>70 - 126</i>	<i>48.95</i>	<i>1.61</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>50.73</i>	<i>1.61</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>50.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>77 - 123</i>	<i>52.02</i>	<i>2.39</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.0</i>	<i>82 - 127</i>	<i>50.27</i>	<i>2.57</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS19070994-01	HS19070994-02	HS19070994-03	HS19070994-04
HS19070994-05	HS19070994-06	HS19070994-07	HS19070994-08
HS19070994-09	HS19070994-10	HS19070994-11	HS19070994-12
HS19070994-13	HS19070994-14	HS19070994-15	HS19070994-16
HS19070994-17	HS19070994-18		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19070994

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020

**Sample Receipt Checklist**

Client Name: PBW  
 Work Order: HS19070994

Date/Time Received: **19-Jul-2019 13:15**  
 Received by: **AC**

Checklist completed by: Asad Chaudhry 19-Jul-2019  
 eSignature Date

Reviewed by: Dane J. Wacasey 20-Jul-2019  
 eSignature Date

Matrices: **Water**

Carrier name: **Client**

- Shipping container/cooler in good condition? Yes  No  Not Present
- Custody seals intact on shipping container/cooler? Yes  No  Not Present
- Custody seals intact on sample bottles? Yes  No  Not Present
- VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes  No  Not Present
- Chain of custody present? Yes  No  4 Page(s)
- Chain of custody signed when relinquished and received? Yes  No  COC IDs:196129, 196108, 196107, 196109
- Samplers name present on COC? Yes  No
- Chain of custody agrees with sample labels? Yes  No
- Samples in proper container/bottle? Yes  No
- Sample containers intact? Yes  No
- Sufficient sample volume for indicated test? Yes  No
- All samples received within holding time? Yes  No
- Container/Temp Blank temperature in compliance? Yes  No

Temperature(s)/Thermometer(s): 3.5c, 0.4c, 0.3c, 03c, 0.4c, 0.5c, 0.4c C/UC IR 25  
 Cooler(s)/Kit(s): 43620, 44883, 43052, 44142, 44164, 45160, 45144  
 Date/Time sample(s) sent to storage: 07/19/2019 19:00

- Water - VOA vials have zero headspace? Yes  No  No VOA vials submitted
- Water - pH acceptable upon receipt? Yes  No  N/A
- pH adjusted? Yes  No  N/A

pH adjusted by:

Login Notes:

Client Contacted: \_\_\_\_\_ Date Contacted: \_\_\_\_\_ Person Contacted: \_\_\_\_\_  
 Contacted By: \_\_\_\_\_ Regarding: \_\_\_\_\_

Comments:

Corrective Action:





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# Chain of Custody Form

Page 1 of 4

COC ID: 196129

## HS19070994

Golder Associates Inc.  
Houston TX-Wood Preserving Works

n, WV



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UFRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8230_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8230_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Maltzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5632646 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	ms/msD
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.maltzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TRD-201801</del>			<del>Water</del>		2		*									
2	WG-1620-FB04-20190716	7-16-19	1820			6		X	X	X							
3	WG-1620-MW61A-20190717	7-17-19	0750			12		X	X	X	X						
4	WG-1620-MW60A-20190717		0835			6		X	X	X							
5	WG-1620-MW69A-20190717		0935			6		X	X	X							
6	WG-1620-MW47C-20190717		1025			6	X		X	X							
7	WG-1620-MW48C-20190717		1120			6	X		X	X							
8	WG-1620-MW59A-20190717		1220			6		X	X	X							
9	WG-1620-MW59B-20190717		1310			6		X	X	X							
10	WG-1620-MW44A-20190717		1400			6		X	X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>	Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour	Other: _____	Results Due Date: _____
Relinquished by: <b>John Br</b>	Date: <b>7-19-19</b> Time: <b>13:15</b>	Received by: _____	Notes: <b>LIPRR Houston MWPW</b>	
Relinquished by: _____	Date: <b>7-19-19</b> Time: <b>13:15</b>	Received by (Laboratory): <b>AC</b>	Cooler ID	Cooler Temp.
Logged by (Laboratory): _____	Date: _____ Time: _____	Checked by (Laboratory): _____	<b>43620</b>	<b>3.5</b>
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035			<b>44883</b>	<b>0.4</b>
			<b>43052</b>	<b>0.3</b>
			<b>44142</b>	<b>0.3</b>
			<b>44164</b>	<b>0.4</b>
			<b>45160</b>	<b>0.5</b>
			<b>45111</b>	<b>0.5</b>

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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# Chain of Custody Form

Page 2 of 4

COC ID: 196108

HS19070994

Golder Associates Inc.  
Houston TX-Wood Preserving Works

n. WV  
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ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A
Work Order		Project Number	1620-07-Rev0 SR 92688	B
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E
	Suite 4004		Stop 0750	F
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G
Phone	(512) 671-3434	Phone		H
Fax	(512) 671-3446	Fax		I
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TB0-201901</del>			Water	T	2		X									
2	WG-1620-MW87C-20190717	7-17-19	1540			6	X		X	X							
3	WG-1620-MW71B-20190717		1640			6	X		X	X							
4	WG-1620-MW33A-20190717		1750			6	X		X	X							
5	WG-1626-FD02-20190717		1750			6	X		X	X							
6	WG-1620-MW26A-20190717		1845			6	X		X	X							
7	WG-1620-FB05-20190717		1905			6	X		X	X							
8	WG-1620-MW68A-20190718	7-18-19	0755			6	X		X	X							
9	WG-1620-MW68B-20190718		0855			6	X		X	X							
10	WG-1620-FD03-20190718		0855			6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> 1-3 Wk. Days <input type="checkbox"/> 5 Wk. Days <input type="checkbox"/> 2 Wk. Days <input type="checkbox"/> 24-hour				Results Due Date:			
Relinquished by: <i>John</i>	Date: 7-19-19	Time: 13:15	Received by:	Notes: UPRR Houston MWPW							
Relinquished by:	Date:	Time:	Received by (Laboratory): AC	Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)					
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):			<input type="checkbox"/> Level II Std. QC	<input checked="" type="checkbox"/> TRRP Check/Std				
						<input type="checkbox"/> Level III Std. QC/Run/Date	<input type="checkbox"/> TRRP Level IV				
						<input type="checkbox"/> Level IV SW-846/CLP					

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
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# Chain of Custody Form

Page 3 of 4

COC ID: 196107

HS19070994

Golder Associates Inc.  
Houston TX-Wood Preserving Works

88, WV  
8

0



ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Paterburs	Project Name	Houston TX-Wood Preserving Works	A
Work Order		Project Number	1620-07-Rev0 SR 92688	B
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- AVF	C
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E
	Suite 4004		Stop 0750	F
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G
Phone	(512) 671-3434	Phone		H
Fax	(512) 671-3446	Fax		I
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-T80-201801</del>			Water	-	2		*									
2	WG-1620-MW68C-20190718	7-18-19	0945			6	X		X	X							
3	WG-1620-MW83B-20190718		1040			6	X		X	X							
4	WG-1620-MW83C-20190718		1130			6	X		X	X							
5	WG-1620-MW35A-20190718		1240			6	X		X	X							
6	WG-1620-MW35B-20190718		1330			6	X		X	X							
7	WG-1620-MW90B-20190718		1425			6	X		X	X							
8	WG-1620-MW89B-20190718		1515			6	X		X	X							
9	WG-1620-MW38B-20190718		1610			6	X		X	X							
10	WG-1620-MW27A-20190718		1715			6	X		X	X							

Sampler(s) Please Print & Sign  
**JOHN BRAYTON** *John Br*  
 Relinquished by: *John Br* Date: **7-19-19** Time: **13:15**  
 Relinquished by: *John Br* Date: **7-19-19** Time: **13:15**  
 Logged by (Laboratory): \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 Shipment Method: **HAND DELIVERED**  
 Required Turnaround Time: (Check Box)  STD 10 Wk Days  5 Wk Days  2 Wk Days  24 Hour  
 Results Due Date: \_\_\_\_\_  
 Notes: **UPRR Houston MWPVW**  
 Cooler ID: \_\_\_\_\_ Cooler Temp: \_\_\_\_\_  
 QC Package: (Check One Box Below)  
 Level II Std QC  TRRP Check at  
 Level III Std. COC Raw Data  TRRP Level I /  
 Level IV SW-846/CLP

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

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# Chain of Custody For

Page 4 of 4

COC ID: 196109

## HS19070994

Golder Associates Inc.  
Houston TX-Wood Preserving Works

30, WV  
18

10



ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A
Work Order		Project Number	1620-07-Rev0 SR 92688	B
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E
				F
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G
Phone	(512) 671-3434	Phone		H
Fax	(512) 671-3446	Fax		I
e-Mail Address	eric.matzner@obwllc.com	e-Mail Address		J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>W6-1620-TB0-20190719</del>			<del>Water</del>	<del>+</del>	<del>2</del>		X									
2	W6-1620-MW27C-20190718	7-18-19	1805			6	X		X	X							
3	W6-1620-MW51A-20190719	7-19-19	0725			6	X		X	X							
4	W6-1620-MW51C-20190719		0855			6	X		X	X							
5	W6-1620-MW81B-20190719		0945			6	X		X	X							
6	W6-1620-MW50A-20190719		1040			6	X		X	X							
7	W6-1620-TB05-20190719					2		X									
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>John Bergstrom</i>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box)				Results Due Date:	
Relinquished by <i>John Bergstrom</i>		Date: 7-19-19 Time: 1315		<input checked="" type="checkbox"/> STD 10 Wk. Days		<input type="checkbox"/> 5 Wk. Days		<input type="checkbox"/> 2 Wk. Days	
Relinquished by <i>John Bergstrom</i>		Date: 7-19-19 Time: 1315		Received by: <b>AC</b>		Notes: UPRR Houston MW/PV			
Logged by (Laboratory):		Date: Time:		Checked by (Laboratory):		Cooler ID		Cooler Temp.	
								QC Package: (Check One Box Below)	
								<input checked="" type="checkbox"/> Level II Std. QC	
								<input type="checkbox"/> Level III Std. QC/Run Date	
								<input type="checkbox"/> Level IV SW846/CLP	
								<input type="checkbox"/> TRRP Check 1	
								<input type="checkbox"/> TRRP Level 1	

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>4</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

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10450 Stancliff Rd. Suite 210  
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F: +1 281 530 5887

August 12, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19080069**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 13 sample(s) on Aug 01, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: DAYNA.FISHER  
Dane J. Wacasey

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 08/12/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19080069			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143727, 143772, R343615, R343708			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Were MS/MSD RPDs within laboratory QC limits?		X			3
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				4
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				



<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 08/12/2019			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19080069			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143727, 143772, R343615, R343708			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?			X		
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section)					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?			X		
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?			X		
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 08/12/2019
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS19080069
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 143727, 143772, R343615, R343708
ER# <sup>5</sup>	Description	
1	<p>Batch 143727, Semivolatiles by Method SW8270, Samples WG-1620-MW33BR-20190730, WG-1620-MW77A-20190730, WG-1620-MW74B-20190730, WG-1620-MW79A-20190730: Surrogate recoveries could not be determined due to dilution below the calibration range.</p> <p>Batch R343708, Volatiles by Method SW8260, Sample WG-1620-MW79A-20190730, surrogate 4-Bromofluorobenzene recovered above upper control limits due to possible matrix effect.</p>	
2	Batch R343708, Volatiles by Method SW8260, Sample HS19080060-04, MSD was performed on an unrelated sample	
3	Batch 143727, Semivolatiles by Method SW8270, Sample WG-1620-MW82B-20190730, MS/MSD RPD recovered above control limits for Naphthalene.	
4	Batch R343708, Volatiles by Method SW8260, Sample WG-1620-MW74B-20190730, lowest practical dilution performed due to high concentration of non-target analyte(s).	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.</p> <p>O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);          NA = Not Applicable;          NR = Not Reviewed;          R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19080069

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19080069-01	WG-1620-MW32AR-20190730	Water		30-Jul-2019 07:25	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-02	WG-1620-MW33BR-20190730	Water		30-Jul-2019 08:20	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-03	WG-1620-MW84B-20190730	Water		30-Jul-2019 09:15	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-04	WG-1620-MW82B-20190730	Water		30-Jul-2019 10:10	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-05	WG-1620-MW80B-20190730	Water		30-Jul-2019 11:10	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-06	WG-1620-MW77A-20190730	Water		30-Jul-2019 11:55	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-07	WG-1620-MW85C-20190730	Water		30-Jul-2019 15:35	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-08	WG-1620-MW86C-20190730	Water		30-Jul-2019 16:40	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-09	WG-1620-FD04-20190730	Water		30-Jul-2019 16:40	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-10	WG-1620-MW76C-20190730	Water		30-Jul-2019 17:30	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-11	WG-1620-MW74B-20190730	Water		30-Jul-2019 18:20	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-12	WG-1620-MW79A-20190730	Water		30-Jul-2019 19:10	01-Aug-2019 14:40	<input type="checkbox"/>
HS19080069-13	WG-1620-TB05-20190730	Water	CG 062119 -77	30-Jul-2019 00:00	01-Aug-2019 14:40	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32AR-20190730  
 Collection Date: 30-Jul-2019 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:33
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:33
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:33
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:33
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 06:33
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:33
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:33
<i>Surr: 1,2-Dichloroethane-d4</i>		96.7		70-126	%REC	1	03-Aug-2019 06:33
<i>Surr: 4-Bromofluorobenzene</i>		95.2		81-113	%REC	1	03-Aug-2019 06:33
<i>Surr: Dibromofluoromethane</i>		99.3		77-123	%REC	1	03-Aug-2019 06:33
<i>Surr: Toluene-d8</i>		100		82-127	%REC	1	03-Aug-2019 06:33

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32AR-20190730  
 Collection Date: 30-Jul-2019 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 18:21
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	02-Aug-2019 18:21
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 18:21
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 18:21
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 18:21
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 18:21
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 18:21
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 18:21
Acenaphthene	U		0.000027	0.00010	mg/L	1	02-Aug-2019 18:21
Acenaphthylene	U		0.000015	0.00010	mg/L	1	02-Aug-2019 18:21
Anthracene	U		0.000014	0.00010	mg/L	1	02-Aug-2019 18:21
<b>Benz(a)anthracene</b>	<b>0.000054</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:21
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 18:21
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 18:21
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000074</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:21
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 18:21
Dibenzofuran	U		0.000020	0.00010	mg/L	1	02-Aug-2019 18:21
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	02-Aug-2019 18:21
Fluoranthene	U		0.000010	0.00010	mg/L	1	02-Aug-2019 18:21
Fluorene	U		0.000030	0.00010	mg/L	1	02-Aug-2019 18:21
Naphthalene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 18:21
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 18:21
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 18:21
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 18:21
Phenanthrene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 18:21
Phenol	U		0.000035	0.00020	mg/L	1	02-Aug-2019 18:21
<b>Pyrene</b>	<b>0.00045</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:21
<i>Surr: 2,4,6-Tribromophenol</i>	63.8			34-129	%REC	1	02-Aug-2019 18:21
<i>Surr: 2-Fluorobiphenyl</i>	74.7			40-125	%REC	1	02-Aug-2019 18:21
<i>Surr: 2-Fluorophenol</i>	59.9			20-120	%REC	1	02-Aug-2019 18:21
<i>Surr: 4-Terphenyl-d14</i>	80.7			40-135	%REC	1	02-Aug-2019 18:21
<i>Surr: Nitrobenzene-d5</i>	62.7			41-120	%REC	1	02-Aug-2019 18:21
<i>Surr: Phenol-d6</i>	63.3			20-120	%REC	1	02-Aug-2019 18:21
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.0628</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	08-Aug-2019 22:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33BR-20190730  
 Collection Date: 30-Jul-2019 08:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 00:29
<b>Benzene</b>	<b>0.25</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	5	03-Aug-2019 01:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 00:29
<b>Ethylbenzene</b>	<b>0.065</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 00:29
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 00:29
<b>Toluene</b>	<b>0.0031</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 00:29
Vinyl chloride	U		0.00020	0.0010	mg/L	1	03-Aug-2019 00:29
<b>Xylenes, Total</b>	<b>0.016</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 00:29
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.3</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:29</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.5</i>			<i>70-126</i>	<i>%REC</i>	<i>5</i>	<i>03-Aug-2019 01:42</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:29</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.7</i>			<i>81-113</i>	<i>%REC</i>	<i>5</i>	<i>03-Aug-2019 01:42</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:29</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.8</i>			<i>77-123</i>	<i>%REC</i>	<i>5</i>	<i>03-Aug-2019 01:42</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:29</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>5</i>	<i>03-Aug-2019 01:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33BR-20190730  
 Collection Date: 30-Jul-2019 08:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 18:40
<b>2,4-Dimethylphenol</b>	<b>0.00028</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 18:40
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 18:40
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 18:40
<b>2-Methylnaphthalene</b>	<b>0.014</b>		<b>0.000095</b>	<b>0.00050</b>	<b>mg/L</b>	5	06-Aug-2019 21:12
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 18:40
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 18:40
<b>Acenaphthene</b>	<b>0.0078</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
Acenaphthylene	U		0.000015	0.00010	mg/L	1	02-Aug-2019 18:40
<b>Anthracene</b>	<b>0.00055</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	02-Aug-2019 18:40
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 18:40
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 18:40
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000070</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 18:40
<b>Dibenzofuran</b>	<b>0.012</b>		<b>0.00010</b>	<b>0.00050</b>	<b>mg/L</b>	5	06-Aug-2019 21:12
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	02-Aug-2019 18:40
<b>Fluoranthene</b>	<b>0.00039</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
<b>Fluorene</b>	<b>0.0042</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
<b>Naphthalene</b>	<b>0.53</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	06-Aug-2019 21:31
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 18:40
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 18:40
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 18:40
<b>Phenanthrene</b>	<b>0.0048</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
<b>Phenol</b>	<b>0.00021</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
<b>Pyrene</b>	<b>0.00026</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:40
Surr: 2,4,6-Tribromophenol	79.8			34-129	%REC	5	06-Aug-2019 21:12
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	06-Aug-2019 21:31
Surr: 2,4,6-Tribromophenol	67.0			34-129	%REC	1	02-Aug-2019 18:40
Surr: 2-Fluorobiphenyl	71.9			40-125	%REC	1	02-Aug-2019 18:40
Surr: 2-Fluorobiphenyl	78.3			40-125	%REC	5	06-Aug-2019 21:12
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	06-Aug-2019 21:31
Surr: 2-Fluorophenol	88.8			20-120	%REC	5	06-Aug-2019 21:12
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	06-Aug-2019 21:31
Surr: 2-Fluorophenol	80.7			20-120	%REC	1	02-Aug-2019 18:40
Surr: 4-Terphenyl-d14	84.8			40-135	%REC	1	02-Aug-2019 18:40
Surr: 4-Terphenyl-d14	91.4			40-135	%REC	5	06-Aug-2019 21:12
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	06-Aug-2019 21:31

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33BR-20190730  
 Collection Date: 30-Jul-2019 08:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
Surr: Nitrobenzene-d5	71.1			41-120	%REC	1	02-Aug-2019 18:40
Surr: Nitrobenzene-d5	81.6			41-120	%REC	5	06-Aug-2019 21:12
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	06-Aug-2019 21:31
Surr: Phenol-d6	68.0			20-120	%REC	5	06-Aug-2019 21:12
Surr: Phenol-d6	0	JS		20-120	%REC	100	06-Aug-2019 21:31
Surr: Phenol-d6	67.9			20-120	%REC	1	02-Aug-2019 18:40
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
Arsenic	0.00117	J	0.000400	0.00200	mg/L	1	08-Aug-2019 22:53

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84B-20190730  
 Collection Date: 30-Jul-2019 09:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:58
<b>Benzene</b>	<b>0.010</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 06:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:58
<b>Ethylbenzene</b>	<b>0.013</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 06:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 06:58
<b>Toluene</b>	<b>0.0015</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 06:58
<b>Xylenes, Total</b>	<b>0.0041</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 06:58
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:58</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:58</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:58</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84B-20190730  
 Collection Date: 30-Jul-2019 09:15

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	02-Aug-2019 18:59
<b>2,4-Dimethylphenol</b>	<b>0.00081</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	02-Aug-2019 18:59
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	02-Aug-2019 18:59
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	02-Aug-2019 18:59
<b>2-Methylnaphthalene</b>	<b>0.0027</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	02-Aug-2019 18:59
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	02-Aug-2019 18:59
<b>Acenaphthene</b>	<b>0.0080</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<b>Acenaphthylene</b>	<b>0.00030</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<b>Anthracene</b>	<b>0.00042</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	02-Aug-2019 18:59
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	02-Aug-2019 18:59
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	02-Aug-2019 18:59
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00026</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
Chrysene		U	0.000021	0.00010	mg/L	1	02-Aug-2019 18:59
<b>Dibenzofuran</b>	<b>0.0054</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<b>Di-n-butyl phthalate</b>	<b>0.000055</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<b>Fluoranthene</b>	<b>0.00015</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<b>Fluorene</b>	<b>0.0026</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<b>Naphthalene</b>	<b>0.060</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	06-Aug-2019 18:21
Nitrobenzene		U	0.000024	0.00020	mg/L	1	02-Aug-2019 18:59
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	02-Aug-2019 18:59
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	02-Aug-2019 18:59
<b>Phenanthrene</b>	<b>0.0021</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
Phenol		U	0.000035	0.00020	mg/L	1	02-Aug-2019 18:59
<b>Pyrene</b>	<b>0.00011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 18:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>74.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 18:59</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>66.0</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 18:21</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.0</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 18:21</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>66.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 18:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 18:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.6</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 18:21</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>113</i>			<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 18:21</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 18:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>53.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 18:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>74.0</i>			<i>41-120</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 18:21</i>
<i>Surr: Phenol-d6</i>	<i>77.3</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 18:21</i>
<i>Surr: Phenol-d6</i>	<i>53.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 18:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84B-20190730  
 Collection Date: 30-Jul-2019 09:15

**ANALYTICAL REPORT**

WorkOrder:HS19080069  
 Lab ID:HS19080069-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>	<b>Method:SW6020</b>			Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
Arsenic	0.00838		0.000400	0.00200	mg/L	1	08-Aug-2019 22:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW82B-20190730  
 Collection Date: 30-Jul-2019 10:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 00:05
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 00:05
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 00:05
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 00:05
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 00:05
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 00:05
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 00:05
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:05</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:05</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:05</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 00:05</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW82B-20190730  
 Collection Date: 30-Jul-2019 10:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 19:18
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	02-Aug-2019 19:18
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 19:18
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 19:18
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 19:18
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 19:18
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 19:18
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 19:18
Acenaphthene	U		0.000027	0.00010	mg/L	1	02-Aug-2019 19:18
Acenaphthylene	U		0.000015	0.00010	mg/L	1	02-Aug-2019 19:18
Anthracene	U		0.000014	0.00010	mg/L	1	02-Aug-2019 19:18
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	02-Aug-2019 19:18
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 19:18
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 19:18
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	02-Aug-2019 19:18
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 19:18
Dibenzofuran	U		0.000020	0.00010	mg/L	1	02-Aug-2019 19:18
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	02-Aug-2019 19:18
Fluoranthene	U		0.000010	0.00010	mg/L	1	02-Aug-2019 19:18
Fluorene	U		0.000030	0.00010	mg/L	1	02-Aug-2019 19:18
Naphthalene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 19:18
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 19:18
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 19:18
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 19:18
Phenanthrene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 19:18
Phenol	U		0.000035	0.00020	mg/L	1	02-Aug-2019 19:18
Pyrene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 19:18
<i>Surr: 2,4,6-Tribromophenol</i>		78.5		34-129	%REC	1	02-Aug-2019 19:18
<i>Surr: 2-Fluorobiphenyl</i>		75.6		40-125	%REC	1	02-Aug-2019 19:18
<i>Surr: 2-Fluorophenol</i>		59.2		20-120	%REC	1	02-Aug-2019 19:18
<i>Surr: 4-Terphenyl-d14</i>		95.7		40-135	%REC	1	02-Aug-2019 19:18
<i>Surr: Nitrobenzene-d5</i>		68.6		41-120	%REC	1	02-Aug-2019 19:18
<i>Surr: Phenol-d6</i>		64.6		20-120	%REC	1	02-Aug-2019 19:18
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>		<b>0.00873</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	08-Aug-2019 22:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW80B-20190730  
 Collection Date: 30-Jul-2019 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 07:22
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 07:22
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 07:22
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 07:22
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 07:22
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 07:22
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 07:22
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.3</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:22</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:22</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:22</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:22</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW80B-20190730  
 Collection Date: 30-Jul-2019 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 20:16
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	02-Aug-2019 20:16
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 20:16
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 20:16
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 20:16
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 20:16
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 20:16
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 20:16
Acenaphthene	U		0.000027	0.00010	mg/L	1	02-Aug-2019 20:16
Acenaphthylene	U		0.000015	0.00010	mg/L	1	02-Aug-2019 20:16
Anthracene	U		0.000014	0.00010	mg/L	1	02-Aug-2019 20:16
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	02-Aug-2019 20:16
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 20:16
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 20:16
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000070</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 20:16
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 20:16
Dibenzofuran	U		0.000020	0.00010	mg/L	1	02-Aug-2019 20:16
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	02-Aug-2019 20:16
Fluoranthene	U		0.000010	0.00010	mg/L	1	02-Aug-2019 20:16
Fluorene	U		0.000030	0.00010	mg/L	1	02-Aug-2019 20:16
Naphthalene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 20:16
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 20:16
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 20:16
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 20:16
Phenanthrene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 20:16
Phenol	U		0.000035	0.00020	mg/L	1	02-Aug-2019 20:16
Pyrene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 20:16
<i>Surr: 2,4,6-Tribromophenol</i>	<i>85.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>81.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>55.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:16</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>59.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:16</i>
<i>Surr: Phenol-d6</i>	<i>65.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:16</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00162</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	08-Aug-2019 22:57

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW77A-20190730  
 Collection Date: 30-Jul-2019 11:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 09:23
<b>Benzene</b>	<b>0.063</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 09:23
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 09:23
<b>Ethylbenzene</b>	<b>0.072</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 09:23
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 09:23
<b>Toluene</b>	<b>0.012</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 09:23
<b>Xylenes, Total</b>	<b>0.097</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 09:23
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 09:23</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 09:23</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 09:23</i>
<i>Surr: Toluene-d8</i>	<i>99.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 09:23</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW77A-20190730  
 Collection Date: 30-Jul-2019 11:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 20:35
<b>2,4-Dimethylphenol</b>	<b>0.062</b>		<b>0.00080</b>	<b>0.0040</b>	<b>mg/L</b>	20	06-Aug-2019 18:40
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 20:35
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 20:35
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 20:35
<b>2-Methylnaphthalene</b>	<b>0.20</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	06-Aug-2019 18:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 20:35
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 20:35
<b>Acenaphthene</b>	<b>0.16</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	06-Aug-2019 18:40
<b>Acenaphthylene</b>	<b>0.0013</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 20:35
<b>Anthracene</b>	<b>0.0028</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 20:35
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	02-Aug-2019 20:35
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 20:35
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 20:35
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	02-Aug-2019 20:35
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 20:35
<b>Dibenzofuran</b>	<b>0.096</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	06-Aug-2019 18:40
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	02-Aug-2019 20:35
<b>Fluoranthene</b>	<b>0.00052</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 20:35
<b>Fluorene</b>	<b>0.078</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	06-Aug-2019 18:40
<b>Naphthalene</b>	<b>4.9</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	06-Aug-2019 18:59
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 20:35
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 20:35
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 20:35
<b>Phenanthrene</b>	<b>0.032</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	06-Aug-2019 18:40
Phenol	U		0.000035	0.00020	mg/L	1	02-Aug-2019 20:35
<b>Pyrene</b>	<b>0.00041</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 20:35
<i>Surr: 2,4,6-Tribromophenol</i>	78.6			34-129	%REC	1	02-Aug-2019 20:35
<i>Surr: 2,4,6-Tribromophenol</i>	125			34-129	%REC	20	06-Aug-2019 18:40
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	2000	06-Aug-2019 18:59
<i>Surr: 2-Fluorobiphenyl</i>	90.5			40-125	%REC	20	06-Aug-2019 18:40
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	2000	06-Aug-2019 18:59
<i>Surr: 2-Fluorobiphenyl</i>	69.1			40-125	%REC	1	02-Aug-2019 20:35
<i>Surr: 2-Fluorophenol</i>	50.6			20-120	%REC	1	02-Aug-2019 20:35
<i>Surr: 2-Fluorophenol</i>	109			20-120	%REC	20	06-Aug-2019 18:40
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	2000	06-Aug-2019 18:59
<i>Surr: 4-Terphenyl-d14</i>	121			40-135	%REC	20	06-Aug-2019 18:40
<i>Surr: 4-Terphenyl-d14</i>	0	JS		40-135	%REC	2000	06-Aug-2019 18:59
<i>Surr: 4-Terphenyl-d14</i>	84.8			40-135	%REC	1	02-Aug-2019 20:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW77A-20190730  
 Collection Date: 30-Jul-2019 11:55

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
Surr: Nitrobenzene-d5	99.9			41-120	%REC	1	02-Aug-2019 20:35
Surr: Nitrobenzene-d5	63.6	J		41-120	%REC	20	06-Aug-2019 18:40
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	06-Aug-2019 18:59
Surr: Phenol-d6	51.9	J		20-120	%REC	20	06-Aug-2019 18:40
Surr: Phenol-d6	0	JS		20-120	%REC	2000	06-Aug-2019 18:59
Surr: Phenol-d6	61.3			20-120	%REC	1	02-Aug-2019 20:35
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
Arsenic	0.0231		0.000400	0.00200	mg/L	1	08-Aug-2019 22:59

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW85C-20190730  
 Collection Date: 30-Jul-2019 15:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 07:46
<b>Benzene</b>	<b>0.00093</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 07:46
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 07:46
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 07:46
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 07:46
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 07:46
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 07:46
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:46</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:46</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:46</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 07:46</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW85C-20190730  
 Collection Date: 30-Jul-2019 15:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 20:54
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	02-Aug-2019 20:54
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 20:54
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 20:54
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 20:54
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 20:54
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 20:54
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 20:54
Acenaphthene	U		0.000027	0.00010	mg/L	1	02-Aug-2019 20:54
Acenaphthylene	U		0.000015	0.00010	mg/L	1	02-Aug-2019 20:54
Anthracene	U		0.000014	0.00010	mg/L	1	02-Aug-2019 20:54
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	02-Aug-2019 20:54
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 20:54
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 20:54
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	02-Aug-2019 20:54
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 20:54
Dibenzofuran	U		0.000020	0.00010	mg/L	1	02-Aug-2019 20:54
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	02-Aug-2019 20:54
Fluoranthene	U		0.000010	0.00010	mg/L	1	02-Aug-2019 20:54
Fluorene	U		0.000030	0.00010	mg/L	1	02-Aug-2019 20:54
<b>Naphthalene</b>	<b>0.0012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 20:54
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 20:54
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 20:54
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 20:54
Phenanthrene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 20:54
Phenol	U		0.000035	0.00020	mg/L	1	02-Aug-2019 20:54
Pyrene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 20:54
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:54</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>67.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:54</i>
<i>Surr: 2-Fluorophenol</i>	<i>55.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:54</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>87.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:54</i>
<i>Surr: Nitrobenzene-d5</i>	<i>57.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:54</i>
<i>Surr: Phenol-d6</i>	<i>56.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 20:54</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.000633</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>08-Aug-2019 23:01</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW86C-20190730  
 Collection Date: 30-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:10
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:10
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:10
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:10
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 08:10
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:10
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:10
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.3</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:10</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:10</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:10</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:10</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW86C-20190730  
 Collection Date: 30-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	02-Aug-2019 21:13
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	02-Aug-2019 21:13
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	02-Aug-2019 21:13
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	02-Aug-2019 21:13
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	02-Aug-2019 21:13
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 21:13
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	02-Aug-2019 21:13
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	02-Aug-2019 21:13
Acenaphthene	U		0.000027	0.00010	mg/L	1	02-Aug-2019 21:13
Acenaphthylene	U		0.000015	0.00010	mg/L	1	02-Aug-2019 21:13
Anthracene	U		0.000014	0.00010	mg/L	1	02-Aug-2019 21:13
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	02-Aug-2019 21:13
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	02-Aug-2019 21:13
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	02-Aug-2019 21:13
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000060</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 21:13
Chrysene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 21:13
Dibenzofuran	U		0.000020	0.00010	mg/L	1	02-Aug-2019 21:13
<b>Di-n-butyl phthalate</b>	<b>0.00011</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	02-Aug-2019 21:13
Fluoranthene	U		0.000010	0.00010	mg/L	1	02-Aug-2019 21:13
Fluorene	U		0.000030	0.00010	mg/L	1	02-Aug-2019 21:13
<b>Naphthalene</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	02-Aug-2019 21:13
Nitrobenzene	U		0.000024	0.00020	mg/L	1	02-Aug-2019 21:13
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	02-Aug-2019 21:13
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	02-Aug-2019 21:13
Phenanthrene	U		0.000021	0.00010	mg/L	1	02-Aug-2019 21:13
Phenol	U		0.000035	0.00020	mg/L	1	02-Aug-2019 21:13
Pyrene	U		0.000019	0.00010	mg/L	1	02-Aug-2019 21:13
<i>Surr: 2,4,6-Tribromophenol</i>	<i>52.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 21:13</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>53.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 21:13</i>
<i>Surr: 2-Fluorophenol</i>	<i>52.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 21:13</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>80.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 21:13</i>
<i>Surr: Nitrobenzene-d5</i>	<i>49.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 21:13</i>
<i>Surr: Phenol-d6</i>	<i>49.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>02-Aug-2019 21:13</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00236</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	08-Aug-2019 23:04

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD04-20190730  
 Collection Date: 30-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:35
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:35
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:35
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:35
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 08:35
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:35
Vinyl chloride	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:35
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>99.2</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:35</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>97.4</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:35</i>
<i>Surr: Dibromofluoromethane</i>		<i>102</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:35</i>
<i>Surr: Toluene-d8</i>		<i>100</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:35</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD04-20190730  
 Collection Date: 30-Jul-2019 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 17:05
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 17:05
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 17:05
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 17:05
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 17:05
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 17:05
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 17:05
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 17:05
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 17:05
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 17:05
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 17:05
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 17:05
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 17:05
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 17:05
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000050</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 17:05
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 17:05
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 17:05
<b>Di-n-butyl phthalate</b>	<b>0.000073</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 17:05
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 17:05
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 17:05
<b>Naphthalene</b>	<b>0.00031</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 17:05
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 17:05
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 17:05
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 17:05
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 17:05
Phenol	U		0.000035	0.00020	mg/L	1	06-Aug-2019 17:05
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 17:05
<i>Surr: 2,4,6-Tribromophenol</i>	<i>56.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:05</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>63.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:05</i>
<i>Surr: 2-Fluorophenol</i>	<i>59.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:05</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:05</i>
<i>Surr: Nitrobenzene-d5</i>	<i>56.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:05</i>
<i>Surr: Phenol-d6</i>	<i>51.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:05</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00190</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	08-Aug-2019 23:06

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76C-20190730  
 Collection Date: 30-Jul-2019 17:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:59
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:59
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:59
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:59
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 08:59
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 08:59
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 08:59
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:59</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:59</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 08:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76C-20190730  
 Collection Date: 30-Jul-2019 17:30

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 17:24
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 17:24
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 17:24
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 17:24
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 17:24
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 17:24
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 17:24
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 17:24
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 17:24
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 17:24
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 17:24
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 17:24
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 17:24
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 17:24
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 17:24
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 17:24
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 17:24
<b>Di-n-butyl phthalate</b>	<b>0.000054</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>06-Aug-2019 17:24</b>
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 17:24
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 17:24
<b>Naphthalene</b>	<b>0.00018</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>06-Aug-2019 17:24</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 17:24
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 17:24
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 17:24
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 17:24
Phenol	U		0.000035	0.00020	mg/L	1	06-Aug-2019 17:24
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 17:24
<i>Surr: 2,4,6-Tribromophenol</i>	<i>47.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>55.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:24</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:24</i>
<i>Surr: Nitrobenzene-d5</i>	<i>63.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:24</i>
<i>Surr: Phenol-d6</i>	<i>57.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 17:24</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00216</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>08-Aug-2019 23:08</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW74B-20190730  
 Collection Date: 30-Jul-2019 18:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0010	0.0050	mg/L	5	05-Aug-2019 18:44
<b>Benzene</b>	<b>0.59</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	5	05-Aug-2019 18:44
Chlorobenzene	U		0.0015	0.0050	mg/L	5	05-Aug-2019 18:44
<b>Ethylbenzene</b>	<b>0.15</b>		<b>0.0015</b>	<b>0.0050</b>	<b>mg/L</b>	5	05-Aug-2019 18:44
Methylene chloride	U		0.0050	0.010	mg/L	5	05-Aug-2019 18:44
<b>Toluene</b>	<b>0.52</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	5	05-Aug-2019 18:44
<b>Xylenes, Total</b>	<b>0.42</b>		<b>0.0015</b>	<b>0.0050</b>	<b>mg/L</b>	5	05-Aug-2019 18:44
Surr: 1,2-Dichloroethane-d4	97.0			70-126	%REC	5	05-Aug-2019 18:44
Surr: 4-Bromofluorobenzene	99.2			81-113	%REC	5	05-Aug-2019 18:44
Surr: Dibromofluoromethane	106			77-123	%REC	5	05-Aug-2019 18:44
Surr: Toluene-d8	101			82-127	%REC	5	05-Aug-2019 18:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW74B-20190730  
 Collection Date: 30-Jul-2019 18:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 15:43
<b>2,4-Dimethylphenol</b>	<b>37</b>		<b>0.40</b>	<b>2.0</b>	<b>mg/L</b>	10000	06-Aug-2019 19:56
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 15:43
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 15:43
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 15:43
<b>2-Methylnaphthalene</b>	<b>0.48</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	06-Aug-2019 19:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 15:43
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 15:43
<b>Acenaphthene</b>	<b>0.24</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	06-Aug-2019 19:37
<b>Acenaphthylene</b>	<b>0.0037</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 15:43
<b>Anthracene</b>	<b>0.0071</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 15:43
<b>Benz(a)anthracene</b>	<b>0.00022</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 15:43
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 15:43
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 15:43
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 15:43
<b>Chrysene</b>	<b>0.00016</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 15:43
<b>Dibenzofuran</b>	<b>0.19</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	06-Aug-2019 19:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 15:43
<b>Fluoranthene</b>	<b>0.0035</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 15:43
<b>Fluorene</b>	<b>0.14</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	06-Aug-2019 19:37
<b>Naphthalene</b>	<b>13</b>		<b>0.20</b>	<b>1.0</b>	<b>mg/L</b>	10000	06-Aug-2019 19:56
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 15:43
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 15:43
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 15:43
<b>Phenanthrene</b>	<b>0.087</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	06-Aug-2019 19:18
<b>Phenol</b>	<b>33</b>		<b>0.35</b>	<b>2.0</b>	<b>mg/L</b>	10000	06-Aug-2019 19:56
<b>Pyrene</b>	<b>0.0023</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 15:43
<i>Surr: 2,4,6-Tribromophenol</i>	<i>40.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 15:43</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>85.9</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 19:18</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>117</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>06-Aug-2019 19:37</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>10000</i>	<i>06-Aug-2019 19:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>10000</i>	<i>06-Aug-2019 19:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>89.7</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 19:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>104</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>06-Aug-2019 19:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>49.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 15:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>102</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 15:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>87.1</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>06-Aug-2019 19:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>117</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>06-Aug-2019 19:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>10000</i>	<i>06-Aug-2019 19:56</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW74B-20190730  
 Collection Date: 30-Jul-2019 18:20

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	10000	06-Aug-2019 19:56
Surr: 4-Terphenyl-d14	95.1			40-135	%REC	10	06-Aug-2019 19:18
Surr: 4-Terphenyl-d14	120	J		40-135	%REC	100	06-Aug-2019 19:37
Surr: 4-Terphenyl-d14	58.4			40-135	%REC	1	06-Aug-2019 15:43
Surr: Nitrobenzene-d5	73.3			41-120	%REC	1	06-Aug-2019 15:43
Surr: Nitrobenzene-d5	61.6			41-120	%REC	10	06-Aug-2019 19:18
Surr: Nitrobenzene-d5	106	J		41-120	%REC	100	06-Aug-2019 19:37
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	10000	06-Aug-2019 19:56
Surr: Phenol-d6	72.9			20-120	%REC	1	06-Aug-2019 15:43
Surr: Phenol-d6	88.6			20-120	%REC	10	06-Aug-2019 19:18
Surr: Phenol-d6	117	J		20-120	%REC	100	06-Aug-2019 19:37
Surr: Phenol-d6	0	JS		20-120	%REC	10000	06-Aug-2019 19:56
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
Arsenic	0.00128	J	0.000400	0.00200	mg/L	1	09-Aug-2019 16:08

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW79A-20190730  
 Collection Date: 30-Jul-2019 19:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	05-Aug-2019 18:17
<b>Benzene</b>	<b>0.013</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	05-Aug-2019 18:17
Chlorobenzene	U		0.00030	0.0010	mg/L	1	05-Aug-2019 18:17
<b>Ethylbenzene</b>	<b>0.0067</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	05-Aug-2019 18:17
Methylene chloride	U		0.0010	0.0020	mg/L	1	05-Aug-2019 18:17
<b>Toluene</b>	<b>0.018</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	05-Aug-2019 18:17
<b>Xylenes, Total</b>	<b>0.023</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	05-Aug-2019 18:17
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 18:17</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>121</i>	<i>S</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 18:17</i>
<i>Surr: Dibromofluoromethane</i>	<i>104</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 18:17</i>
<i>Surr: Toluene-d8</i>	<i>96.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 18:17</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW79A-20190730  
 Collection Date: 30-Jul-2019 19:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 16:02
<b>2,4-Dimethylphenol</b>	<b>0.33</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	100	06-Aug-2019 20:34
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 16:02
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 16:02
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 16:02
<b>2-Methylnaphthalene</b>	<b>0.051</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	06-Aug-2019 20:15
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 16:02
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 16:02
<b>Acenaphthene</b>	<b>0.034</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	06-Aug-2019 20:15
<b>Acenaphthylene</b>	<b>0.00076</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
<b>Anthracene</b>	<b>0.0020</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
<b>Benz(a)anthracene</b>	<b>0.00023</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 16:02
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 16:02
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00055</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
<b>Chrysene</b>	<b>0.00027</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
<b>Dibenzofuran</b>	<b>0.026</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	06-Aug-2019 20:15
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 16:02
<b>Fluoranthene</b>	<b>0.0012</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
<b>Fluorene</b>	<b>0.017</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	06-Aug-2019 20:15
<b>Naphthalene</b>	<b>0.74</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	06-Aug-2019 20:34
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 16:02
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 16:02
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 16:02
<b>Phenanthrene</b>	<b>0.0082</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
<b>Phenol</b>	<b>0.063</b>		<b>0.00035</b>	<b>0.0020</b>	<b>mg/L</b>	10	06-Aug-2019 20:15
<b>Pyrene</b>	<b>0.0011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	06-Aug-2019 16:02
Surr: 2,4,6-Tribromophenol	124			34-129	%REC	10	06-Aug-2019 20:15
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	06-Aug-2019 20:34
Surr: 2,4,6-Tribromophenol	72.9			34-129	%REC	1	06-Aug-2019 16:02
Surr: 2-Fluorobiphenyl	95.2			40-125	%REC	1	06-Aug-2019 16:02
Surr: 2-Fluorobiphenyl	114			40-125	%REC	10	06-Aug-2019 20:15
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	06-Aug-2019 20:34
Surr: 2-Fluorophenol	84.7			20-120	%REC	10	06-Aug-2019 20:15
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	06-Aug-2019 20:34
Surr: 2-Fluorophenol	93.9			20-120	%REC	1	06-Aug-2019 16:02
Surr: 4-Terphenyl-d14	104			40-135	%REC	1	06-Aug-2019 16:02
Surr: 4-Terphenyl-d14	103			40-135	%REC	10	06-Aug-2019 20:15
Surr: 4-Terphenyl-d14	161	JS		40-135	%REC	100	06-Aug-2019 20:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW79A-20190730  
 Collection Date: 30-Jul-2019 19:10

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 02-Aug-2019		Analyst: LG	
Surr: Nitrobenzene-d5	68.7			41-120	%REC	1	06-Aug-2019 16:02
Surr: Nitrobenzene-d5	103			41-120	%REC	10	06-Aug-2019 20:15
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	06-Aug-2019 20:34
Surr: Phenol-d6	92.9			20-120	%REC	10	06-Aug-2019 20:15
Surr: Phenol-d6	0	JS		20-120	%REC	100	06-Aug-2019 20:34
Surr: Phenol-d6	85.4			20-120	%REC	1	06-Aug-2019 16:02
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 05-Aug-2019		Analyst: JHD	
Arsenic	0.00991		0.000400	0.00200	mg/L	1	08-Aug-2019 23:17

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB05-20190730  
 Collection Date: 30-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19080069  
 Lab ID:HS19080069-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: PC			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:09
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:09
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:09
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:09
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 06:09
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:09
Vinyl chloride	U		0.00020	0.0010	mg/L	1	03-Aug-2019 06:09
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 06:09
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:09</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:09</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:09</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 06:09</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**Batch ID:** 143727      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19080069-01	1	1000	1 (mL)	0.001
HS19080069-02	1	1000	1 (mL)	0.001
HS19080069-03	1	1000	1 (mL)	0.001
HS19080069-04	1	1000	1 (mL)	0.001
HS19080069-05	1	1000	1 (mL)	0.001
HS19080069-06	1	1000	1 (mL)	0.001
HS19080069-07	1	1000	1 (mL)	0.001
HS19080069-08	1	1000	1 (mL)	0.001
HS19080069-09	1	1000	1 (mL)	0.001
HS19080069-10	1	1000	1 (mL)	0.001
HS19080069-11	1	1000	1 (mL)	0.001
HS19080069-12	1	1000	1 (mL)	0.001

**Batch ID:** 143772      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19080069-01	1	10	10 (mL)	1
HS19080069-02	1	10	10 (mL)	1
HS19080069-03	1	10	10 (mL)	1
HS19080069-04	1	10	10 (mL)	1
HS19080069-05	1	10	10 (mL)	1
HS19080069-06	1	10	10 (mL)	1
HS19080069-07	1	10	10 (mL)	1
HS19080069-08	1	10	10 (mL)	1
HS19080069-09	1	10	10 (mL)	1
HS19080069-10	1	10	10 (mL)	1
HS19080069-11	1	10	10 (mL)	1
HS19080069-12	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143727 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS19080069-01	WG-1620-MW32AR-20190730	30 Jul 2019 07:25		02 Aug 2019 07:00	02 Aug 2019 18:21	1
HS19080069-02	WG-1620-MW33BR-20190730	30 Jul 2019 08:20		02 Aug 2019 07:00	06 Aug 2019 21:31	100
HS19080069-02	WG-1620-MW33BR-20190730	30 Jul 2019 08:20		02 Aug 2019 07:00	06 Aug 2019 21:12	5
HS19080069-02	WG-1620-MW33BR-20190730	30 Jul 2019 08:20		02 Aug 2019 07:00	02 Aug 2019 18:40	1
HS19080069-03	WG-1620-MW84B-20190730	30 Jul 2019 09:15		02 Aug 2019 07:00	06 Aug 2019 18:21	10
HS19080069-03	WG-1620-MW84B-20190730	30 Jul 2019 09:15		02 Aug 2019 07:00	02 Aug 2019 18:59	1
HS19080069-04	WG-1620-MW82B-20190730	30 Jul 2019 10:10		02 Aug 2019 07:00	02 Aug 2019 19:18	1
HS19080069-05	WG-1620-MW80B-20190730	30 Jul 2019 11:10		02 Aug 2019 07:00	02 Aug 2019 20:16	1
HS19080069-06	WG-1620-MW77A-20190730	30 Jul 2019 11:55		02 Aug 2019 07:00	06 Aug 2019 18:59	2000
HS19080069-06	WG-1620-MW77A-20190730	30 Jul 2019 11:55		02 Aug 2019 07:00	06 Aug 2019 18:40	20
HS19080069-06	WG-1620-MW77A-20190730	30 Jul 2019 11:55		02 Aug 2019 07:00	02 Aug 2019 20:35	1
HS19080069-07	WG-1620-MW85C-20190730	30 Jul 2019 15:35		02 Aug 2019 07:00	02 Aug 2019 20:54	1
HS19080069-08	WG-1620-MW86C-20190730	30 Jul 2019 16:40		02 Aug 2019 07:00	02 Aug 2019 21:13	1
HS19080069-09	WG-1620-FD04-20190730	30 Jul 2019 16:40		02 Aug 2019 07:00	06 Aug 2019 17:05	1
HS19080069-10	WG-1620-MW76C-20190730	30 Jul 2019 17:30		02 Aug 2019 07:00	06 Aug 2019 17:24	1
HS19080069-11	WG-1620-MW74B-20190730	30 Jul 2019 18:20		02 Aug 2019 07:00	06 Aug 2019 19:56	1000 0
HS19080069-11	WG-1620-MW74B-20190730	30 Jul 2019 18:20		02 Aug 2019 07:00	06 Aug 2019 19:37	100
HS19080069-11	WG-1620-MW74B-20190730	30 Jul 2019 18:20		02 Aug 2019 07:00	06 Aug 2019 19:18	10
HS19080069-11	WG-1620-MW74B-20190730	30 Jul 2019 18:20		02 Aug 2019 07:00	06 Aug 2019 15:43	1
HS19080069-12	WG-1620-MW79A-20190730	30 Jul 2019 19:10		02 Aug 2019 07:00	06 Aug 2019 20:34	100
HS19080069-12	WG-1620-MW79A-20190730	30 Jul 2019 19:10		02 Aug 2019 07:00	06 Aug 2019 20:15	10
HS19080069-12	WG-1620-MW79A-20190730	30 Jul 2019 19:10		02 Aug 2019 07:00	06 Aug 2019 16:02	1
<b>Batch ID: 143772 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19080069-01	WG-1620-MW32AR-20190730	30 Jul 2019 07:25		05 Aug 2019 09:00	08 Aug 2019 22:50	1
HS19080069-02	WG-1620-MW33BR-20190730	30 Jul 2019 08:20		05 Aug 2019 09:00	08 Aug 2019 22:53	1
HS19080069-03	WG-1620-MW84B-20190730	30 Jul 2019 09:15		05 Aug 2019 09:00	08 Aug 2019 22:55	1
HS19080069-04	WG-1620-MW82B-20190730	30 Jul 2019 10:10		05 Aug 2019 09:00	08 Aug 2019 22:34	1
HS19080069-05	WG-1620-MW80B-20190730	30 Jul 2019 11:10		05 Aug 2019 09:00	08 Aug 2019 22:57	1
HS19080069-06	WG-1620-MW77A-20190730	30 Jul 2019 11:55		05 Aug 2019 09:00	08 Aug 2019 22:59	1
HS19080069-07	WG-1620-MW85C-20190730	30 Jul 2019 15:35		05 Aug 2019 09:00	08 Aug 2019 23:01	1
HS19080069-08	WG-1620-MW86C-20190730	30 Jul 2019 16:40		05 Aug 2019 09:00	08 Aug 2019 23:04	1
HS19080069-09	WG-1620-FD04-20190730	30 Jul 2019 16:40		05 Aug 2019 09:00	08 Aug 2019 23:06	1
HS19080069-10	WG-1620-MW76C-20190730	30 Jul 2019 17:30		05 Aug 2019 09:00	08 Aug 2019 23:08	1
HS19080069-11	WG-1620-MW74B-20190730	30 Jul 2019 18:20		05 Aug 2019 09:00	09 Aug 2019 16:08	1
HS19080069-12	WG-1620-MW79A-20190730	30 Jul 2019 19:10		05 Aug 2019 09:00	08 Aug 2019 23:17	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: R343615 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS19080069-01	WG-1620-MW32AR-20190730	30 Jul 2019 07:25			03 Aug 2019 06:33	1
HS19080069-02	WG-1620-MW33BR-20190730	30 Jul 2019 08:20			03 Aug 2019 01:42	5
HS19080069-02	WG-1620-MW33BR-20190730	30 Jul 2019 08:20			03 Aug 2019 00:29	1
HS19080069-03	WG-1620-MW84B-20190730	30 Jul 2019 09:15			03 Aug 2019 06:58	1
HS19080069-04	WG-1620-MW82B-20190730	30 Jul 2019 10:10			03 Aug 2019 00:05	1
HS19080069-05	WG-1620-MW80B-20190730	30 Jul 2019 11:10			03 Aug 2019 07:22	1
HS19080069-06	WG-1620-MW77A-20190730	30 Jul 2019 11:55			03 Aug 2019 09:23	1
HS19080069-07	WG-1620-MW85C-20190730	30 Jul 2019 15:35			03 Aug 2019 07:46	1
HS19080069-08	WG-1620-MW86C-20190730	30 Jul 2019 16:40			03 Aug 2019 08:10	1
HS19080069-09	WG-1620-FD04-20190730	30 Jul 2019 16:40			03 Aug 2019 08:35	1
HS19080069-10	WG-1620-MW76C-20190730	30 Jul 2019 17:30			03 Aug 2019 08:59	1
HS19080069-13	WG-1620-TB05-20190730	30 Jul 2019 00:00			03 Aug 2019 06:09	1
<b>Batch ID: R343708 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS19080069-11	WG-1620-MW74B-20190730	30 Jul 2019 18:20			05 Aug 2019 18:44	5
HS19080069-12	WG-1620-MW79A-20190730	30 Jul 2019 19:10			05 Aug 2019 18:17	1

WorkOrder: HS19080069  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200

WorkOrder: HS19080069  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00013	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00012	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000053	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00014	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.00010	0.00010	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.00016	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00018	0.000047	0.0010
A	Acenaphthene	83-32-9	0.00010	0.00011	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.00010	0.00010	0.000015	0.00010
A	Anthracene	120-12-7	0.00010	0.00012	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.00010	0.00012	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.00010	0.00013	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000091	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000084	0.000037	0.00020
A	Chrysene	218-01-9	0.00010	0.00010	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.00010	0.00011	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.00010	0.00011	0.000010	0.00010
A	Fluorene	86-73-7	0.00010	0.00010	0.000030	0.00010
A	Naphthalene	91-20-3	0.00010	0.00012	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00017	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00015	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.00015	0.000079	0.00020
A	Phenanthrene	85-01-8	0.00010	0.00011	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00013	0.000035	0.00020
A	Pyrene	129-00-0	0.00010	0.000099	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19080069  
 InstrumentID: VOA9  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00058	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00060	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00061	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00058	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00052	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00063	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS19080069  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00067	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00056	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00059	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00051	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00054	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00056	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00052	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

Batch ID: 143772 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A						
<b>MBLK</b>	Sample ID: <b>MBLK-143772</b>	Units: <b>mg/L</b>		Analysis Date: <b>08-Aug-2019 22:30</b>						
Client ID:		Run ID: <b>ICPMS05_343891</b>	SeqNo: <b>5202657</b>	PrepDate: <b>05-Aug-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-143772</b>	Units: <b>mg/L</b>		Analysis Date: <b>08-Aug-2019 22:32</b>						
Client ID:		Run ID: <b>ICPMS05_343891</b>	SeqNo: <b>5202658</b>	PrepDate: <b>05-Aug-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.0475	0.00200	0.05	0	95.0	80 - 120				
<b>MS</b>	Sample ID: <b>HS19080069-04MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>08-Aug-2019 22:39</b>						
Client ID: <b>WG-1620-MW82B-20190730</b>		Run ID: <b>ICPMS05_343891</b>	SeqNo: <b>5202661</b>	PrepDate: <b>05-Aug-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.06004	0.00200	0.05	0.008727	103	80 - 120				
<b>MSD</b>	Sample ID: <b>HS19080069-04MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>08-Aug-2019 22:41</b>						
Client ID: <b>WG-1620-MW82B-20190730</b>		Run ID: <b>ICPMS05_343891</b>	SeqNo: <b>5202662</b>	PrepDate: <b>05-Aug-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.05823	0.00200	0.05	0.008727	99.0	80 - 120	0.06004	3.06	20	
<b>PDS</b>	Sample ID: <b>HS19080069-04PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>08-Aug-2019 22:44</b>						
Client ID: <b>WG-1620-MW82B-20190730</b>		Run ID: <b>ICPMS05_343891</b>	SeqNo: <b>5202663</b>	PrepDate: <b>05-Aug-2019</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.1112	0.00200	0.1	0.008727	103	75 - 125				
<b>SD</b>	Sample ID: <b>HS19080069-04SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>08-Aug-2019 22:37</b>						
Client ID: <b>WG-1620-MW82B-20190730</b>		Run ID: <b>ICPMS05_343891</b>	SeqNo: <b>5202660</b>	PrepDate: <b>05-Aug-2019</b>	DF: <b>5</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual	
Arsenic	0.008077	0.0100					0.008727	0 10	J	

The following samples were analyzed in this batch:

HS19080069-01	HS19080069-02	HS19080069-03	HS19080069-04
HS19080069-05	HS19080069-06	HS19080069-07	HS19080069-08
HS19080069-09	HS19080069-10	HS19080069-11	HS19080069-12

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

Batch ID: 143727 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143727	Units: ug/L			Analysis Date: 02-Aug-2019 13:29					
Client ID:	Run ID: SV-6_343627	SeqNo: 5196073		PrepDate: 02-Aug-2019		DF: 1				
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	3.503	0.20	5	0	70.1	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.588	0.20	5	0	91.8	40 - 125				
<i>Surr: 2-Fluorophenol</i>	4.162	0.20	5	0	83.2	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.367	0.20	5	0	87.3	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	4.55	0.20	5	0	91.0	41 - 120				
<i>Surr: Phenol-d6</i>	4.055	0.20	5	0	81.1	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

Batch ID: 143727 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143727	Units: ug/L			Analysis Date: 02-Aug-2019 13:48					
Client ID:	Run ID: SV-6_343627	SeqNo: 5196074		PrepDate: 02-Aug-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.518	0.20	5	0	90.4	39 - 127				
2,4-Dimethylphenol	4.308	0.20	5	0	86.2	35 - 120				
2,4-Dinitrotoluene	4.312	0.20	5	0	86.2	50 - 122				
2,6-Dinitrotoluene	4.261	0.20	5	0	85.2	50 - 120				
2-Chloronaphthalene	4.56	0.20	5	0	91.2	50 - 120				
2-Methylnaphthalene	4.221	0.10	5	0	84.4	50 - 120				
4,6-Dinitro-2-methylphenol	4.02	0.20	5	0	80.4	25 - 121				
4-Nitrophenol	4.598	1.0	5	0	92.0	30 - 130				
Acenaphthene	4.112	0.10	5	0	82.2	45 - 120				
Acenaphthylene	4.516	0.10	5	0	90.3	47 - 120				
Anthracene	4.121	0.10	5	0	82.4	45 - 120				
Benz(a)anthracene	4.275	0.10	5	0	85.5	40 - 120				
Benzo(a)pyrene	4.39	0.10	5	0	87.8	45 - 120				
Bis(2-chloroethoxy)methane	4.166	0.20	5	0	83.3	45 - 120				
Bis(2-ethylhexyl)phthalate	4.912	0.20	5	0	98.2	40 - 139				
Chrysene	4.374	0.10	5	0	87.5	43 - 120				
Dibenzofuran	4.495	0.10	5	0	89.9	50 - 120				
Di-n-butyl phthalate	4.752	0.20	5	0	95.0	45 - 123				
Fluoranthene	4.286	0.10	5	0	85.7	45 - 125				
Fluorene	4.419	0.10	5	0	88.4	49 - 120				
Naphthalene	4.236	0.10	5	0	84.7	45 - 120				
Nitrobenzene	3.994	0.20	5	0	79.9	44 - 120				
N-Nitrosodiphenylamine	4.38	0.20	5	0	87.6	40 - 125				
Pentachlorophenol	1.98	0.20	5	0	39.6	19 - 121				
Phenanthrene	4.208	0.10	5	0	84.2	45 - 121				
Phenol	4.791	0.20	5	0	95.8	20 - 124				
Pyrene	4.756	0.10	5	0	95.1	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.335</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.7</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.065</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.627</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.5</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.919</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.4</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.524</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.5</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.738</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

Batch ID: 143727 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS19080069-04MS		Units: ug/L		Analysis Date: 02-Aug-2019 19:37				
Client ID: WG-1620-MW82B-20190730		Run ID: SV-6_343627		SeqNo: 5196079		PrepDate: 02-Aug-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.089	0.20	5	0	81.8	39 - 127				
2,4-Dimethylphenol	3.156	0.20	5	0	63.1	35 - 120				
2,4-Dinitrotoluene	4.03	0.20	5	0	80.6	50 - 122				
2,6-Dinitrotoluene	3.94	0.20	5	0	78.8	50 - 120				
2-Chloronaphthalene	3.639	0.20	5	0	72.8	50 - 120				
2-Methylnaphthalene	3.38	0.10	5	0	67.6	50 - 120				
4,6-Dinitro-2-methylphenol	3.401	0.20	5	0	68.0	25 - 121				
4-Nitrophenol	4.39	1.0	5	0	87.8	30 - 130				
Acenaphthene	3.739	0.10	5	0	74.8	45 - 120				
Acenaphthylene	3.891	0.10	5	0	77.8	47 - 120				
Anthracene	3.992	0.10	5	0	79.8	45 - 120				
Benz(a)anthracene	4.248	0.10	5	0	85.0	40 - 120				
Benzo(a)pyrene	4.45	0.10	5	0	89.0	45 - 120				
Bis(2-chloroethoxy)methane	3.177	0.20	5	0	63.5	45 - 120				
Bis(2-ethylhexyl)phthalate	4.984	0.20	5	0	99.7	40 - 139				
Chrysene	4.216	0.10	5	0	84.3	43 - 120				
Dibenzofuran	3.933	0.10	5	0	78.7	50 - 120				
Di-n-butyl phthalate	4.711	0.20	5	0	94.2	45 - 123				
Fluoranthene	4.273	0.10	5	0	85.5	45 - 125				
Fluorene	4.049	0.10	5	0	81.0	49 - 120				
Naphthalene	3.253	0.10	5	0	65.1	45 - 120				
Nitrobenzene	3.184	0.20	5	0	63.7	44 - 120				
N-Nitrosodiphenylamine	4.228	0.20	5	0	84.6	40 - 125				
Pentachlorophenol	1.635	0.20	5	0	32.7	19 - 121				
Phenanthrene	4.215	0.10	5	0	84.3	45 - 121				
Phenol	3.315	0.20	5	0	66.3	20 - 124				
Pyrene	4.551	0.10	5	0	91.0	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.145</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>82.9</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.013</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>80.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.282</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>65.6</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.438</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.8</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.521</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>70.4</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.445</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>68.9</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

Batch ID: 143727 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS19080069-04MSD	Units: ug/L			Analysis Date: 02-Aug-2019 19:57					
Client ID: WG-1620-MW82B-20190730	Run ID: SV-6_343627	SeqNo: 5196080	PrepDate: 02-Aug-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.066	0.20	5	0	81.3	39 - 127	4.089	0.576	20	
2,4-Dimethylphenol	3.175	0.20	5	0	63.5	35 - 120	3.156	0.601	20	
2,4-Dinitrotoluene	4.251	0.20	5	0	85.0	50 - 122	4.03	5.36	20	
2,6-Dinitrotoluene	3.931	0.20	5	0	78.6	50 - 120	3.94	0.236	20	
2-Chloronaphthalene	3.31	0.20	5	0	66.2	50 - 120	3.639	9.48	20	
2-Methylnaphthalene	3.116	0.10	5	0	62.3	50 - 120	3.38	8.15	20	
4,6-Dinitro-2-methylphenol	3.77	0.20	5	0	75.4	25 - 121	3.401	10.3	30	
4-Nitrophenol	4.608	1.0	5	0	92.2	30 - 130	4.39	4.86	20	
Acenaphthene	3.471	0.10	5	0	69.4	45 - 120	3.739	7.43	20	
Acenaphthylene	3.67	0.10	5	0	73.4	47 - 120	3.891	5.84	20	
Anthracene	3.982	0.10	5	0	79.6	45 - 120	3.992	0.252	20	
Benz(a)anthracene	4.498	0.10	5	0	90.0	40 - 120	4.248	5.71	20	
Benzo(a)pyrene	4.831	0.10	5	0	96.6	45 - 120	4.45	8.21	20	
Bis(2-chloroethoxy)methane	3.181	0.20	5	0	63.6	45 - 120	3.177	0.126	20	
Bis(2-ethylhexyl)phthalate	5.362	0.20	5	0	107	40 - 139	4.984	7.3	20	
Chrysene	4.526	0.10	5	0	90.5	43 - 120	4.216	7.1	20	
Dibenzofuran	3.769	0.10	5	0	75.4	50 - 120	3.933	4.26	20	
Di-n-butyl phthalate	4.984	0.20	5	0	99.7	45 - 123	4.711	5.63	20	
Fluoranthene	4.559	0.10	5	0	91.2	45 - 125	4.273	6.47	20	
Fluorene	4.064	0.10	5	0	81.3	49 - 120	4.049	0.359	20	
Naphthalene	5.202	0.10	5	0	104	45 - 120	3.253	46.1	20	R
Nitrobenzene	3.038	0.20	5	0	60.8	44 - 120	3.184	4.7	20	
N-Nitrosodiphenylamine	4.292	0.20	5	0	85.8	40 - 125	4.228	1.5	20	
Pentachlorophenol	1.991	0.20	5	0	39.8	19 - 121	1.635	19.6	20	
Phenanthrene	4.176	0.10	5	0	83.5	45 - 121	4.215	0.934	20	
Phenol	3.678	0.20	5	0	73.6	20 - 124	3.315	10.4	20	
Pyrene	4.807	0.10	5	0	96.1	40 - 130	4.551	5.49	20	
Surr: 2,4,6-Tribromophenol	4.112	0.20	5	0	82.2	34 - 129	4.145	0.793	20	
Surr: 2-Fluorobiphenyl	3.568	0.20	5	0	71.4	40 - 125	4.013	11.7	20	
Surr: 2-Fluorophenol	3.662	0.20	5	0	73.2	20 - 120	3.282	11	20	
Surr: 4-Terphenyl-d14	4.548	0.20	5	0	91.0	40 - 135	4.438	2.46	20	
Surr: Nitrobenzene-d5	3.299	0.20	5	0	66.0	41 - 120	3.521	6.51	20	
Surr: Phenol-d6	3.547	0.20	5	0	70.9	20 - 120	3.445	2.91	20	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

<b>Batch ID:</b> 143727 ( 0 )	<b>Instrument:</b> SV-6	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS19080069-01	HS19080069-02	HS19080069-03	HS19080069-04
HS19080069-05	HS19080069-06	HS19080069-07	HS19080069-08
HS19080069-09	HS19080069-10	HS19080069-11	HS19080069-12

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

<b>Batch ID:</b> R343615 ( 0 )		<b>Instrument:</b> VOA9		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-190802</b>	Units: <b>ug/L</b>			Analysis Date: <b>02-Aug-2019 23:41</b>				
Client ID:	Run ID: <b>VOA9_343615</b>	SeqNo: <b>5195198</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.44</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.9</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.54</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.58</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-190802</b>	Units: <b>ug/L</b>			Analysis Date: <b>02-Aug-2019 23:17</b>				
Client ID:	Run ID: <b>VOA9_343615</b>	SeqNo: <b>5195197</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	17.43	1.0	20	0	87.1	70 - 124			
Benzene	19.45	1.0	20	0	97.2	74 - 120			
Chlorobenzene	18.2	1.0	20	0	91.0	76 - 113			
Ethylbenzene	18.86	1.0	20	0	94.3	77 - 117			
Methylene chloride	19.47	2.0	20	0	97.3	70 - 127			
Toluene	18.57	1.0	20	0	92.8	77 - 118			
Vinyl chloride	19.04	1.0	20	0	95.2	70 - 130			
Xylenes, Total	56.17	1.0	60	0	93.6	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.0</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

**Batch ID:** R343615 ( 0 )      **Instrument:** VOA9      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS19080069-04MS			Units: ug/L		Analysis Date: 03-Aug-2019 00:54			
Client ID: WG-1620-MW82B-20190730		Run ID: VOA9_343615			SeqNo: 5195201		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.05	1.0	20	0	90.2	70 - 127				
Benzene	20.65	1.0	20	0	103	70 - 127				
Chlorobenzene	19.12	1.0	20	0	95.6	70 - 114				
Ethylbenzene	20.29	1.0	20	0	101	70 - 124				
Methylene chloride	19.82	2.0	20	0	99.1	70 - 128				
Toluene	19.74	1.0	20	0	98.7	70 - 123				
Vinyl chloride	16.73	1.0	20	0	83.7	70 - 130				
Xylenes, Total	60.26	1.0	60	0	100	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.6</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 127</i>				

MSD		Sample ID: HS19080069-04MSD			Units: ug/L		Analysis Date: 03-Aug-2019 01:18			
Client ID: WG-1620-MW82B-20190730		Run ID: VOA9_343615			SeqNo: 5195202		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.6	1.0	20	0	83.0	70 - 127	18.05	8.33	20	
Benzene	18.87	1.0	20	0	94.3	70 - 127	20.65	9.02	20	
Chlorobenzene	18.21	1.0	20	0	91.1	70 - 114	19.12	4.87	20	
Ethylbenzene	18.92	1.0	20	0	94.6	70 - 124	20.29	6.95	20	
Methylene chloride	18.04	2.0	20	0	90.2	70 - 128	19.82	9.39	20	
Toluene	18.52	1.0	20	0	92.6	70 - 123	19.74	6.39	20	
Vinyl chloride	15.02	1.0	20	0	75.1	70 - 130	16.73	10.8	20	
Xylenes, Total	57.07	1.0	60	0	95.1	70 - 130	60.26	5.44	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.93</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.9</i>	<i>70 - 126</i>	<i>47.31</i>	<i>2.97</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.08</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>	<i>50.26</i>	<i>1.62</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.7</i>	<i>77 - 123</i>	<i>50.29</i>	<i>1.87</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>51.66</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>82 - 127</i>	<i>50.78</i>	<i>1.71</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS19080069-01	HS19080069-02	HS19080069-03	HS19080069-04
HS19080069-05	HS19080069-06	HS19080069-07	HS19080069-08
HS19080069-09	HS19080069-10	HS19080069-13	



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

**Batch ID:** R343708 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190805</b>			Units: <b>ug/L</b>		Analysis Date: <b>05-Aug-2019 13:43</b>			
Client ID:		Run ID: <b>VOA2_343708</b>			SeqNo: <b>5197254</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>51.41</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>51.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-190805</b>			Units: <b>ug/L</b>		Analysis Date: <b>05-Aug-2019 13:19</b>			
Client ID:		Run ID: <b>VOA2_343708</b>			SeqNo: <b>5197253</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	19.63	1.0	20	0	98.2	70 - 124				
Benzene	19.95	1.0	20	0	99.7	74 - 120				
Chlorobenzene	20.62	1.0	20	0	103	76 - 113				
Ethylbenzene	20.77	1.0	20	0	104	77 - 117				
Methylene chloride	21.15	2.0	20	0	106	70 - 127				
Toluene	20.27	1.0	20	0	101	77 - 118				
Xylenes, Total	69.61	1.0	60	0	116	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>52.59</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>57.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>115</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.38</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>48.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QC BATCH REPORT**

**Batch ID:** R343708 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS19080060-04MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>05-Aug-2019 17:01</b>			
Client ID:		Run ID: <b>VOA2_343708</b>			SeqNo: <b>5197262</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.34	1.0	20	0	76.7	70 - 127				
Benzene	17.05	1.0	20	0	85.3	70 - 127				
Chlorobenzene	17.94	1.0	20	0	89.7	70 - 114				
Ethylbenzene	18.08	1.0	20	0	90.4	70 - 124				
Methylene chloride	15.91	2.0	20	0	79.5	70 - 128				
Toluene	18.72	1.0	20	0	93.6	70 - 123				
Xylenes, Total	56.34	1.0	60	0	93.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.1</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.2</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS19080060-04MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>05-Aug-2019 17:26</b>			
Client ID:		Run ID: <b>VOA2_343708</b>			SeqNo: <b>5197263</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.1	1.0	20	0	90.5	70 - 127	15.34		16.5	20
Benzene	20.16	1.0	20	0	101	70 - 127	17.05		16.7	20
Chlorobenzene	17.72	1.0	20	0	88.6	70 - 114	17.94		1.22	20
Ethylbenzene	19.47	1.0	20	0	97.3	70 - 124	18.08		7.39	20
Methylene chloride	14.72	2.0	20	0	73.6	70 - 128	15.91		7.76	20
Toluene	17.49	1.0	20	0	87.4	70 - 123	18.72		6.81	20
Xylenes, Total	61.55	1.0	60	0	103	70 - 130	56.34		8.84	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 126</i>	<i>47.1</i>		<i>7.44</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>57.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>116</i>	<i>81 - 113</i>	<i>48.39</i>		<i>18</i>	<i>20 S</i>
<i>Surr: Dibromofluoromethane</i>	<i>49.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>77 - 123</i>	<i>50.39</i>		<i>1.64</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>48.82</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>82 - 127</i>	<i>49.52</i>		<i>1.43</i>	<i>20</i>

The following samples were analyzed in this batch: HS19080069-11      HS19080069-12

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080069

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

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**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19080069

**SAMPLE TRACKING**

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Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS19080069-01	WG-1620-MW32AR-20190730	Login	01/08/2019 17:01:24	JRM	MET074
HS19080069-01	WG-1620-MW32AR-20190730	Login	01/08/2019 17:01:24	JRM	EXT011
HS19080069-01	WG-1620-MW32AR-20190730	Login	01/08/2019 17:01:24	JRM	VOA153

Sample Receipt Checklist

Client Name: PBW
Work Order: HS19080069

Date/Time Received: 01-Aug-2019 14:40
Received by: AC

Checklist completed by: Jared R. Makan
eSignature
Date: 1-Aug-2019

Reviewed by: Dane J. Wacasey
eSignature
Date: 6-Aug-2019

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition?
Custody seals intact on shipping container/cooler?
Custody seals intact on sample bottles?
VOA/TX1005/TX1006 Solids in hermetically sealed vials?
Chain of custody present?
Chain of custody signed when relinquished and received?
Samplers name present on COC?
Chain of custody agrees with sample labels?
Samples in proper container/bottle?
Sample containers intact?
Sufficient sample volume for indicated test?
All samples received within holding time?
Container/Temp Blank temperature in compliance?

- Yes/No/Not Present checkboxes for each item in the list above.

2 Page(s)
COC IDs:196137, 196130

Temperature(s)/Thermometer(s): 1.9c/1.9c, 1.6c/1.6c, 1.7c/1.7c UC/C IR25
Cooler(s)/Kit(s): 45149, 44874, 45101
Date/Time sample(s) sent to storage: 08/01/2019 17:20

- Water - VOA vials have zero headspace?
Water - pH acceptable upon receipt?
pH adjusted?
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



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# Chain of Custody Form

Page 1 of 2

COC ID: 196137

## HS19080069

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Malzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive	Address	1400 Douglas Street
	Suite 400-1		Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.malzner@pbwllc.com	e-Mail Address	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TB0_-201801			Water	1	2		X									
2	WG-1620-MW32AR-20190730	7-30-19	0725			6	X		X	X							
3	WG-1620-MW33BR-20190730		0820			6		X	X	X							
4	WG-1620-MW84B-20190730		0915			6	X		X	X							
5	WG-1620-MW82B-20190730		1010			12	X		X	X	X						
6	WG-1620-MW80B-20190730		1110			6	X		X	X							
7	WG-1620-MW77A-20190730		1155			6	X		X	X							
8	WG-1620-MW85C-20190730		1535			6	X		X	X							
9	WG-1620-MW86C-20190730		1640			6	X		X	X							
10	WG-1620-PD04-20190730		1640			6		X	X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>	Shipment Method <b>Hand Delivered</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour	Results Due Date:
Relinquished by: <b>John</b>	Date: <b>8-1-19</b>	Time: <b>14:40</b>	Received by: <b>AC</b>
Relinquished by:	Date:	Time:	Received by (Laboratory):
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):
Notes: UPRR Houston MWPW		QC Package: (Check One Box Below)	
Cooler ID	Cooler Temp.	<input type="checkbox"/> Level II Std QC <input checked="" type="checkbox"/> TRRF Checklist <input type="checkbox"/> Level III Std QC/Raw Data <input type="checkbox"/> TRRF Level IV <input type="checkbox"/> Level IV SWB/CLP <input type="checkbox"/> Other	
45149	1.9		
44874	1.6		
45101	1.7		

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.



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+1 616 399 6070

# Chain of Custody Form

Page 2 of 2

COC ID: 196130

HS19080069

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information		ALS Project Manager:											
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatiles Organics Site Specific)										
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)										
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 Semi/Volatiles Site specific)										
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5838002 5652646 Metals - As, Pb)										
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E											
	Suite 4004		Stop 0750	F											
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G											
Phone	(512) 671-3434	Phone		H											
Fax	(512) 671-3446	Fax		I											
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J											

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TBD -201801			Water	1	2		X									
2	WG-1620-MW76C-20190730	7-30-19	1730			6	X		X	X							
3	WG-1620-MW74B-20190730		1820			6	X		X	X							
4	WG-1620-MW79A-20190730		1910			6	X		X	X							
5	WG-1620-TBOS-20190730					2		X									
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>John Beaman</i>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour				Results Due Date:			
Relinquished by: <i>John Beaman</i>		Date: 8-1-19	Time: 14:40	Received by:		Notes: UPRR Houston MWPW					
Relinquished by: <i>John Beaman</i>		Date: 8-1-19	Time: 14:40	Received by (Laboratory): AC		Cooler ID:	Cooler Temp.:	QC Package: (Check One Box Below)			
Logged by (Laboratory):		Date:	Time:	Checked by (Laboratory):		<input type="checkbox"/> Level II Std QC	<input checked="" type="checkbox"/> TRRF Checklist				
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035						<input type="checkbox"/> Level III Std QCR as Data	<input type="checkbox"/> TRRF Level IV				
						<input type="checkbox"/> Level IV SW946/CLP					

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

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August 12, 2019

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS19080113**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 12 sample(s) on Aug 02, 2019 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: DAYNA.FISHER  
Dane J. Wacasey

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

**Laboratory Review Checklist: Reportable Data**

Laboratory Name: ALS Laboratory Group			LRC Date: 08/12/2019				
Project Name: Houston TX-Wood Preserving Works			Laboratory Job Number: HS19080113				
Reviewer Name: Dane Wacasey			Prep Batch Number(s): 143759, 143819, 143865, R343626, R343632, R343643				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?	X				
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			1
		Were MS/MSD RPDs within laboratory QC limits?		X			2
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data								
Laboratory Name: ALS Laboratory Group				LRC Date: 08/12/2019				
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS19080113				
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 143759, 143819, 143865, R343626, R343632, R343643				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER <sup>5</sup>	
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>						
		Were response factors and/or relative response factors for each analyte within QC limits?	X					
		Were percent RSDs or correlation coefficient criteria met?	X					
		Was the number of standards recommended in the method used for all analytes?	X					
		Were all points generated between the lowest and highest standard used to calculate the curve?	X					
		Are ICAL data available for all instruments used?	X					
		Has the initial calibration curve been verified using an appropriate second source standard?	X					
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>						
		Was the CCV analyzed at the method-required frequency?	X					
		Were percent differences for each analyte within the method-required QC limits?	X					
		Was the ICAL curve verified for each analyte?	X					
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X					
<b>S3</b>	O	<b>Mass spectral tuning:</b>						
		Was the appropriate compound for the method used for tuning?	X					
		Were ion abundance data within the method-required QC limits?	X					
<b>S4</b>	O	<b>Internal standards (IS):</b>						
		Were IS area counts and retention times within the method-required QC limits?	X					
<b>S5</b>	OI	<b>Raw data (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section</b>						
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X					
		Were data associated with manual integrations flagged on the raw data?	X					
<b>S6</b>	O	<b>Dual column confirmation</b>						
		Did dual column confirmation results meet the method-required QC?			X			
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>						
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X			
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>						
		Were percent recoveries within method QC limits?	X					
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>						
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X					
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>						
		Was a MDL study performed for each reported analyte?	X					
		Is the MDL either adjusted or supported by the analysis of DCSs?	X					
<b>S11</b>	OI	<b>Proficiency test reports:</b>						
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X					
<b>S12</b>	OI	<b>Standards documentation</b>						
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X					
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>						
		Are the procedures for compound/analyte identification documented?	X					
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>						
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X					
		Is documentation of the analyst's competency up-to-date and on file?	X					
<b>S15</b>	OI	<b>Verification/validation documentation for methods (NELAC Chap 5 or ISO/IEC 17025 Section 5)</b>						
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X					
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>						
		Are laboratory SOPs current and on file for each method performed?	X					

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 08/12/2019
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS19080113
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 143759, 143819, 143865, R343626, R343632, R343643

<b>ER#<sup>5</sup></b>	<b>Description</b>
1	Batch 143759, Semivolatiles by Method SW8270, Sample WG-1620-MW67B-20190731 and WG-1620-MW65D-20190731, MS and/or MSD recovered outside control limits for multiple compounds due to sample matrix interference.
2	Batch 143759, Semivolatiles by Method SW8270, samples WG-1620-MW67B-20190731 and WG-1620-MW65D-20190731, MS/MSD RPD recovered above control limits for 4-Nitrophenol.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19080113

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS19080113-01	WG-1620-MW36D-20190731	Water		31-Jul-2019 07:40	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-02	WG-1620-MW38A-20190731	Water		31-Jul-2019 08:35	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-03	WG-1620-MW65D-20190731	Water		31-Jul-2019 09:45	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-04	WG-1620-MW66D-20190731	Water		31-Jul-2019 10:45	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-05	WG-1620-MW59D-20190731	Water		31-Jul-2019 11:50	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-06	WG-1620-FD05-20190731	Water		31-Jul-2019 11:50	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-07	WG-1620-MW22AR-20190731	Water		31-Jul-2019 13:05	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-08	WG-1620-MW22BR-20190731	Water		31-Jul-2019 13:50	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-09	WG-1620-MW88C-20190731	Water		31-Jul-2019 15:45	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-10	WG-1620-MW67B-20190731	Water		31-Jul-2019 16:45	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-11	WG-1620-MW49A-20190731	Water		31-Jul-2019 17:45	02-Aug-2019 12:00	<input type="checkbox"/>
HS19080113-12	WG-1620-TB06-20190731	Water		31-Jul-2019 00:00	02-Aug-2019 12:00	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36D-20190731  
 Collection Date: 31-Jul-2019 07:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 19:20
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 19:20
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 19:20
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 19:20
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 19:20
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 19:20
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 19:20
<i>Surr: 1,2-Dichloroethane-d4</i>	98.9			70-126	%REC	1	03-Aug-2019 19:20
<i>Surr: 4-Bromofluorobenzene</i>	99.5			81-113	%REC	1	03-Aug-2019 19:20
<i>Surr: Dibromofluoromethane</i>	96.9			77-123	%REC	1	03-Aug-2019 19:20
<i>Surr: Toluene-d8</i>	99.7			82-127	%REC	1	03-Aug-2019 19:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36D-20190731  
 Collection Date: 31-Jul-2019 07:40

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	05-Aug-2019 19:43
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	05-Aug-2019 19:43
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	05-Aug-2019 19:43
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	05-Aug-2019 19:43
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	05-Aug-2019 19:43
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	05-Aug-2019 19:43
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	05-Aug-2019 19:43
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	05-Aug-2019 19:43
Acenaphthene	U		0.000027	0.00010	mg/L	1	05-Aug-2019 19:43
Acenaphthylene	U		0.000015	0.00010	mg/L	1	05-Aug-2019 19:43
<b>Anthracene</b>	<b>0.000017</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
<b>Benz(a)anthracene</b>	<b>0.000053</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
<b>Benzo(a)pyrene</b>	<b>0.000087</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	05-Aug-2019 19:43
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000080</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
<b>Chrysene</b>	<b>0.000058</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
Dibenzofuran	U		0.000020	0.00010	mg/L	1	05-Aug-2019 19:43
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	05-Aug-2019 19:43
<b>Fluoranthene</b>	<b>0.00011</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
Fluorene	U		0.000030	0.00010	mg/L	1	05-Aug-2019 19:43
Naphthalene	U		0.000020	0.00010	mg/L	1	05-Aug-2019 19:43
Nitrobenzene	U		0.000024	0.00020	mg/L	1	05-Aug-2019 19:43
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	05-Aug-2019 19:43
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	05-Aug-2019 19:43
<b>Phenanthrene</b>	<b>0.000066</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
Phenol	U		0.000035	0.00020	mg/L	1	05-Aug-2019 19:43
<b>Pyrene</b>	<b>0.000099</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 19:43
<i>Surr: 2,4,6-Tribromophenol</i>	<i>83.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 19:43</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>90.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 19:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 19:43</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 19:43</i>
<i>Surr: Nitrobenzene-d5</i>	<i>94.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 19:43</i>
<i>Surr: Phenol-d6</i>	<i>91.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 19:43</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 06-Aug-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	08-Aug-2019 00:06

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38A-20190731  
 Collection Date: 31-Jul-2019 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
<b>1,2-Dichloroethane</b>	<b>0.00031</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	03-Aug-2019 19:44
Benzene		U	0.00020	0.0010	mg/L	1	03-Aug-2019 19:44
Chlorobenzene		U	0.00030	0.0010	mg/L	1	03-Aug-2019 19:44
Ethylbenzene		U	0.00030	0.0010	mg/L	1	03-Aug-2019 19:44
Methylene chloride		U	0.0010	0.0020	mg/L	1	03-Aug-2019 19:44
Toluene		U	0.00020	0.0010	mg/L	1	03-Aug-2019 19:44
Xylenes, Total		U	0.00030	0.0010	mg/L	1	03-Aug-2019 19:44
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 19:44</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>92.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 19:44</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 19:44</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 19:44</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38A-20190731  
 Collection Date: 31-Jul-2019 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	05-Aug-2019 20:02
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	05-Aug-2019 20:02
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	05-Aug-2019 20:02
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	05-Aug-2019 20:02
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	05-Aug-2019 20:02
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	05-Aug-2019 20:02
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	05-Aug-2019 20:02
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	05-Aug-2019 20:02
Acenaphthene	U		0.000027	0.00010	mg/L	1	05-Aug-2019 20:02
Acenaphthylene	U		0.000015	0.00010	mg/L	1	05-Aug-2019 20:02
<b>Anthracene</b>	<b>0.000022</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
<b>Benz(a)anthracene</b>	<b>0.000087</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
<b>Benzo(a)pyrene</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	05-Aug-2019 20:02
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000084</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
<b>Chrysene</b>	<b>0.000092</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
Dibenzofuran	U		0.000020	0.00010	mg/L	1	05-Aug-2019 20:02
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	05-Aug-2019 20:02
<b>Fluoranthene</b>	<b>0.00018</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
Fluorene	U		0.000030	0.00010	mg/L	1	05-Aug-2019 20:02
<b>Naphthalene</b>	<b>0.000068</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
Nitrobenzene	U		0.000024	0.00020	mg/L	1	05-Aug-2019 20:02
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	05-Aug-2019 20:02
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	05-Aug-2019 20:02
<b>Phenanthrene</b>	<b>0.00010</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
Phenol	U		0.000035	0.00020	mg/L	1	05-Aug-2019 20:02
<b>Pyrene</b>	<b>0.00014</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	05-Aug-2019 20:02
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:02</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>91.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:02</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:02</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>80.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:02</i>
<i>Surr: Nitrobenzene-d5</i>	<i>90.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:02</i>
<i>Surr: Phenol-d6</i>	<i>84.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:02</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 06-Aug-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	08-Aug-2019 00:08

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW65D-20190731  
 Collection Date: 31-Jul-2019 09:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 14:42
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 14:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 14:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 14:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 14:42
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 14:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 14:42
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 14:42</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 14:42</i>
<i>Surr: Dibromofluoromethane</i>	<i>103</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 14:42</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 14:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW65D-20190731  
 Collection Date: 31-Jul-2019 09:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	05-Aug-2019 20:21
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	05-Aug-2019 20:21
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	05-Aug-2019 20:21
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	05-Aug-2019 20:21
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	05-Aug-2019 20:21
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	05-Aug-2019 20:21
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	05-Aug-2019 20:21
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	05-Aug-2019 20:21
Acenaphthene	U		0.000027	0.00010	mg/L	1	05-Aug-2019 20:21
Acenaphthylene	U		0.000015	0.00010	mg/L	1	05-Aug-2019 20:21
Anthracene	U		0.000014	0.00010	mg/L	1	05-Aug-2019 20:21
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	05-Aug-2019 20:21
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	05-Aug-2019 20:21
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	05-Aug-2019 20:21
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00016</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	05-Aug-2019 20:21
Chrysene	U		0.000021	0.00010	mg/L	1	05-Aug-2019 20:21
Dibenzofuran	U		0.000020	0.00010	mg/L	1	05-Aug-2019 20:21
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	05-Aug-2019 20:21
Fluoranthene	U		0.000010	0.00010	mg/L	1	05-Aug-2019 20:21
Fluorene	U		0.000030	0.00010	mg/L	1	05-Aug-2019 20:21
Naphthalene	U		0.000020	0.00010	mg/L	1	05-Aug-2019 20:21
Nitrobenzene	U		0.000024	0.00020	mg/L	1	05-Aug-2019 20:21
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	05-Aug-2019 20:21
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	05-Aug-2019 20:21
Phenanthrene	U		0.000021	0.00010	mg/L	1	05-Aug-2019 20:21
Phenol	U		0.000035	0.00020	mg/L	1	05-Aug-2019 20:21
Pyrene	U		0.000019	0.00010	mg/L	1	05-Aug-2019 20:21
<i>Surr: 2,4,6-Tribromophenol</i>	<i>91.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:21</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>95.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:21</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:21</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:21</i>
<i>Surr: Nitrobenzene-d5</i>	<i>102</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:21</i>
<i>Surr: Phenol-d6</i>	<i>89.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>05-Aug-2019 20:21</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 06-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00135</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	07-Aug-2019 23:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW66D-20190731  
 Collection Date: 31-Jul-2019 10:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:09
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:09
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:09
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:09
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 20:09
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:09
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:09
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:09</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>93.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:09</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:09</i>
<i>Surr: Toluene-d8</i>	<i>104</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:09</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW66D-20190731  
 Collection Date: 31-Jul-2019 10:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 13:29
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 13:29
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 13:29
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 13:29
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 13:29
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 13:29
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 13:29
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 13:29
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 13:29
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 13:29
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 13:29
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 13:29
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 13:29
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 13:29
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00022</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 13:29
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 13:29
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 13:29
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 13:29
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 13:29
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 13:29
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 13:29
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 13:29
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 13:29
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 13:29
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 13:29
Phenol	U		0.000035	0.00020	mg/L	1	06-Aug-2019 13:29
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 13:29
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:29</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>89.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:29</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:29</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>79.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:29</i>
<i>Surr: Nitrobenzene-d5</i>	<i>91.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:29</i>
<i>Surr: Phenol-d6</i>	<i>84.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:29</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 06-Aug-2019		Analyst: JHD	
<b>Arsenic</b>	<b>0.00124</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>08-Aug-2019 00:10</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59D-20190731  
 Collection Date: 31-Jul-2019 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:34
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:34
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:34
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:34
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 20:34
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:34
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:34
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:34</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:34</i>
<i>Surr: Dibromofluoromethane</i>	<i>95.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:34</i>
<i>Surr: Toluene-d8</i>	<i>105</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:34</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59D-20190731  
 Collection Date: 31-Jul-2019 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 13:48
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 13:48
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 13:48
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 13:48
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 13:48
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 13:48
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 13:48
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 13:48
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 13:48
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 13:48
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 13:48
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 13:48
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 13:48
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 13:48
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 13:48
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 13:48
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 13:48
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 13:48
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 13:48
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 13:48
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 13:48
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 13:48
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 13:48
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 13:48
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 13:48
<b>Phenol</b>	<b>0.00011</b>	<b>J</b>	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>06-Aug-2019 13:48</b>
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 13:48
<i>Surr: 2,4,6-Tribromophenol</i>	<i>87.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:48</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>78.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:48</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:48</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>83.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:48</i>
<i>Surr: Nitrobenzene-d5</i>	<i>81.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:48</i>
<i>Surr: Phenol-d6</i>	<i>74.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 13:48</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 06-Aug-2019		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	08-Aug-2019 00:13

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD05-20190731  
 Collection Date: 31-Jul-2019 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:58
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:58
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 20:58
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 20:58
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 20:58
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:58</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:58</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:58</i>
<i>Surr: Toluene-d8</i>	<i>103</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>03-Aug-2019 20:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD05-20190731  
 Collection Date: 31-Jul-2019 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 14:07
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 14:07
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 14:07
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 14:07
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 14:07
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 14:07
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 14:07
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 14:07
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 14:07
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 14:07
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 14:07
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 14:07
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 14:07
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 14:07
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 14:07
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 14:07
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 14:07
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 14:07
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 14:07
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 14:07
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 14:07
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 14:07
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 14:07
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 14:07
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 14:07
Phenol	U		0.000035	0.00020	mg/L	1	06-Aug-2019 14:07
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 14:07
<i>Surr: 2,4,6-Tribromophenol</i>		82.7		34-129	%REC	1	06-Aug-2019 14:07
<i>Surr: 2-Fluorobiphenyl</i>		79.6		40-125	%REC	1	06-Aug-2019 14:07
<i>Surr: 2-Fluorophenol</i>		72.8		20-120	%REC	1	06-Aug-2019 14:07
<i>Surr: 4-Terphenyl-d14</i>		83.9		40-135	%REC	1	06-Aug-2019 14:07
<i>Surr: Nitrobenzene-d5</i>		83.3		41-120	%REC	1	06-Aug-2019 14:07
<i>Surr: Phenol-d6</i>		77.9		20-120	%REC	1	06-Aug-2019 14:07
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 07-Aug-2019		Analyst: JC	
<b>Arsenic</b>	<b>0.000502</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>07-Aug-2019 18:26</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22AR-20190731  
 Collection Date: 31-Jul-2019 13:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	04-Aug-2019 01:54
Benzene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 01:54
Chlorobenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 01:54
Ethylbenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 01:54
Methylene chloride	U		0.0010	0.0020	mg/L	1	04-Aug-2019 01:54
Toluene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 01:54
Xylenes, Total	U		0.00030	0.0010	mg/L	1	04-Aug-2019 01:54
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 01:54</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 01:54</i>
<i>Surr: Dibromofluoromethane</i>	<i>95.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 01:54</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 01:54</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22AR-20190731  
 Collection Date: 31-Jul-2019 13:05

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 14:26
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 14:26
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 14:26
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 14:26
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 14:26
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 14:26
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 14:26
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 14:26
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 14:26
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 14:26
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 14:26
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 14:26
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 14:26
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 14:26
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 14:26
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 14:26
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 14:26
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 14:26
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 14:26
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 14:26
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 14:26
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 14:26
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 14:26
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 14:26
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 14:26
Phenol	U		0.000035	0.00020	mg/L	1	06-Aug-2019 14:26
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 14:26
<i>Surr: 2,4,6-Tribromophenol</i>	<i>85.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 14:26</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>90.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 14:26</i>
<i>Surr: 2-Fluorophenol</i>	<i>82.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 14:26</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>83.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 14:26</i>
<i>Surr: Nitrobenzene-d5</i>	<i>93.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 14:26</i>
<i>Surr: Phenol-d6</i>	<i>85.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 14:26</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 07-Aug-2019		Analyst: JC	
<b>Arsenic</b>	<b>0.000481</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>07-Aug-2019 18:28</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22BR-20190731  
 Collection Date: 31-Jul-2019 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:02
Benzene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:02
Chlorobenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:02
Ethylbenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:02
Methylene chloride	U		0.0010	0.0020	mg/L	1	04-Aug-2019 04:02
Toluene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:02
Xylenes, Total	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:02
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:02</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:02</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:02</i>
<i>Surr: Toluene-d8</i>	<i>105</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:02</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22BR-20190731  
 Collection Date: 31-Jul-2019 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 18:33
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 18:33
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 18:33
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 18:33
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 18:33
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 18:33
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 18:33
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 18:33
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 18:33
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 18:33
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 18:33
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 18:33
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 18:33
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 18:33
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 18:33
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 18:33
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 18:33
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 18:33
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 18:33
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 18:33
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 18:33
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 18:33
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 18:33
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 18:33
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 18:33
<b>Phenol</b>	<b>0.00069</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 18:33
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 18:33
<i>Surr: 2,4,6-Tribromophenol</i>	<i>59.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:33</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>82.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:33</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:33</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:33</i>
<i>Surr: Nitrobenzene-d5</i>	<i>65.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:33</i>
<i>Surr: Phenol-d6</i>	<i>72.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:33</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 07-Aug-2019		Analyst: JC	
<b>Arsenic</b>	<b>0.000559</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	07-Aug-2019 18:30

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88C-20190731  
 Collection Date: 31-Jul-2019 15:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:27
Benzene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:27
Chlorobenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:27
Ethylbenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:27
Methylene chloride	U		0.0010	0.0020	mg/L	1	04-Aug-2019 04:27
Toluene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:27
Xylenes, Total	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:27
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>93.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:27</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>100</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:27</i>
<i>Surr: Dibromofluoromethane</i>	<i>94.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:27</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:27</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88C-20190731  
 Collection Date: 31-Jul-2019 15:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 18:52
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 18:52
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 18:52
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 18:52
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 18:52
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 18:52
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 18:52
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 18:52
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 18:52
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 18:52
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 18:52
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 18:52
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 18:52
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 18:52
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 18:52
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 18:52
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 18:52
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 18:52
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 18:52
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 18:52
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 18:52
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 18:52
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 18:52
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 18:52
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 18:52
<b>Phenol</b>	<b>0.00055</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 18:52
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 18:52
<i>Surr: 2,4,6-Tribromophenol</i>	<i>46.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:52</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>67.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:52</i>
<i>Surr: 2-Fluorophenol</i>	<i>65.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:52</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:52</i>
<i>Surr: Nitrobenzene-d5</i>	<i>58.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:52</i>
<i>Surr: Phenol-d6</i>	<i>66.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 18:52</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 07-Aug-2019		Analyst: JC	
Arsenic	U		0.000400	0.00200	mg/L	1	07-Aug-2019 18:32

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW67B-20190731  
 Collection Date: 31-Jul-2019 16:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	03-Aug-2019 14:49
Benzene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 14:49
Chlorobenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 14:49
Ethylbenzene	U		0.00030	0.0010	mg/L	1	03-Aug-2019 14:49
Methylene chloride	U		0.0010	0.0020	mg/L	1	03-Aug-2019 14:49
Toluene	U		0.00020	0.0010	mg/L	1	03-Aug-2019 14:49
Xylenes, Total	U		0.00030	0.0010	mg/L	1	03-Aug-2019 14:49
<i>Surr: 1,2-Dichloroethane-d4</i>		99.6		70-126	%REC	1	03-Aug-2019 14:49
<i>Surr: 4-Bromofluorobenzene</i>		94.2		81-113	%REC	1	03-Aug-2019 14:49
<i>Surr: Dibromofluoromethane</i>		97.0		77-123	%REC	1	03-Aug-2019 14:49
<i>Surr: Toluene-d8</i>		102		82-127	%REC	1	03-Aug-2019 14:49

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW67B-20190731  
 Collection Date: 31-Jul-2019 16:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 19:11
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 19:11
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 19:11
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 19:11
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 19:11
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 19:11
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 19:11
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 19:11
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 19:11
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 19:11
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 19:11
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 19:11
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 19:11
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 19:11
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 19:11
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 19:11
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 19:11
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 19:11
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 19:11
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 19:11
<b>Naphthalene</b>	<b>0.000079</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>06-Aug-2019 19:11</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 19:11
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 19:11
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 19:11
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 19:11
Phenol	U		0.000035	0.00020	mg/L	1	06-Aug-2019 19:11
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 19:11
<i>Surr: 2,4,6-Tribromophenol</i>	<i>66.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:11</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>72.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:11</i>
<i>Surr: 2-Fluorophenol</i>	<i>63.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:11</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:11</i>
<i>Surr: Nitrobenzene-d5</i>	<i>58.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:11</i>
<i>Surr: Phenol-d6</i>	<i>68.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:11</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 07-Aug-2019		Analyst: JC	
<b>Arsenic</b>	<b>0.000494</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>07-Aug-2019 18:34</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49A-20190731  
 Collection Date: 31-Jul-2019 17:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:51
Benzene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:51
Chlorobenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:51
Ethylbenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:51
Methylene chloride	U		0.0010	0.0020	mg/L	1	04-Aug-2019 04:51
Toluene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:51
Vinyl chloride	U		0.00020	0.0010	mg/L	1	04-Aug-2019 04:51
Xylenes, Total	U		0.00030	0.0010	mg/L	1	04-Aug-2019 04:51
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:51</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:51</i>
<i>Surr: Dibromofluoromethane</i>	<i>94.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:51</i>
<i>Surr: Toluene-d8</i>	<i>103</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 04:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49A-20190731  
 Collection Date: 31-Jul-2019 17:45

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 05-Aug-2019		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	06-Aug-2019 19:30
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	06-Aug-2019 19:30
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	06-Aug-2019 19:30
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	06-Aug-2019 19:30
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	06-Aug-2019 19:30
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 19:30
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	06-Aug-2019 19:30
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	06-Aug-2019 19:30
Acenaphthene	U		0.000027	0.00010	mg/L	1	06-Aug-2019 19:30
Acenaphthylene	U		0.000015	0.00010	mg/L	1	06-Aug-2019 19:30
Anthracene	U		0.000014	0.00010	mg/L	1	06-Aug-2019 19:30
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	06-Aug-2019 19:30
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 19:30
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	06-Aug-2019 19:30
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	06-Aug-2019 19:30
Chrysene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 19:30
Dibenzofuran	U		0.000020	0.00010	mg/L	1	06-Aug-2019 19:30
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	06-Aug-2019 19:30
Fluoranthene	U		0.000010	0.00010	mg/L	1	06-Aug-2019 19:30
Fluorene	U		0.000030	0.00010	mg/L	1	06-Aug-2019 19:30
Naphthalene	U		0.000020	0.00010	mg/L	1	06-Aug-2019 19:30
Nitrobenzene	U		0.000024	0.00020	mg/L	1	06-Aug-2019 19:30
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	06-Aug-2019 19:30
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	06-Aug-2019 19:30
Phenanthrene	U		0.000021	0.00010	mg/L	1	06-Aug-2019 19:30
<b>Phenol</b>	<b>0.00064</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	06-Aug-2019 19:30
Pyrene	U		0.000019	0.00010	mg/L	1	06-Aug-2019 19:30
<i>Surr: 2,4,6-Tribromophenol</i>	<i>56.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:30</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>73.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:30</i>
<i>Surr: 2-Fluorophenol</i>	<i>65.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:30</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>87.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:30</i>
<i>Surr: Nitrobenzene-d5</i>	<i>58.7</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:30</i>
<i>Surr: Phenol-d6</i>	<i>67.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>06-Aug-2019 19:30</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 07-Aug-2019		Analyst: JC	
<b>Arsenic</b>	<b>0.000658</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	07-Aug-2019 18:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB06-20190731  
 Collection Date: 31-Jul-2019 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS19080113  
 Lab ID:HS19080113-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	04-Aug-2019 00:40
Benzene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 00:40
Chlorobenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 00:40
Ethylbenzene	U		0.00030	0.0010	mg/L	1	04-Aug-2019 00:40
Methylene chloride	U		0.0010	0.0020	mg/L	1	04-Aug-2019 00:40
Toluene	U		0.00020	0.0010	mg/L	1	04-Aug-2019 00:40
Vinyl chloride	U		0.00020	0.0010	mg/L	1	04-Aug-2019 00:40
Xylenes, Total	U		0.00030	0.0010	mg/L	1	04-Aug-2019 00:40
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 00:40</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 00:40</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 00:40</i>
<i>Surr: Toluene-d8</i>	<i>105</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>04-Aug-2019 00:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## WEIGHT LOG

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**Batch ID:** 143759      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D      **Prep:** 3510\_B\_LOW

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19080113-01	1	1000	1 (mL)	0.001
HS19080113-02	1	1000	1 (mL)	0.001
HS19080113-03	1	1000	1 (mL)	0.001
HS19080113-04	1	1000	1 (mL)	0.001
HS19080113-05	1	1000	1 (mL)	0.001
HS19080113-06	1	1000	1 (mL)	0.001
HS19080113-07	1	1000	1 (mL)	0.001
HS19080113-08	1	1000	1 (mL)	0.001
HS19080113-09	1	1000	1 (mL)	0.001
HS19080113-10	1	1000	1 (mL)	0.001
HS19080113-11	1	1000	1 (mL)	0.001

**Batch ID:** 143819      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19080113-01	1	10	10 (mL)	1
HS19080113-02	1	10	10 (mL)	1
HS19080113-03	1	10	10 (mL)	1
HS19080113-04	1	10	10 (mL)	1
HS19080113-05	1	10	10 (mL)	1

**Batch ID:** 143865      **Method:** ICP-MS METALS BY SW6020A      **Prep:** 3010A

SampID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS19080113-06	1	10	10 (mL)	1
HS19080113-07	1	10	10 (mL)	1
HS19080113-08	1	10	10 (mL)	1
HS19080113-09	1	10	10 (mL)	1
HS19080113-10	1	10	10 (mL)	1
HS19080113-11	1	10	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: 143759 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS19080113-01	WG-1620-MW36D-20190731	31 Jul 2019 07:40		05 Aug 2019 09:58	05 Aug 2019 19:43	1
HS19080113-02	WG-1620-MW38A-20190731	31 Jul 2019 08:35		05 Aug 2019 09:58	05 Aug 2019 20:02	1
HS19080113-03	WG-1620-MW65D-20190731	31 Jul 2019 09:45		05 Aug 2019 09:58	05 Aug 2019 20:21	1
HS19080113-04	WG-1620-MW66D-20190731	31 Jul 2019 10:45		05 Aug 2019 09:58	06 Aug 2019 13:29	1
HS19080113-05	WG-1620-MW59D-20190731	31 Jul 2019 11:50		05 Aug 2019 09:58	06 Aug 2019 13:48	1
HS19080113-06	WG-1620-FD05-20190731	31 Jul 2019 11:50		05 Aug 2019 09:58	06 Aug 2019 14:07	1
HS19080113-07	WG-1620-MW22AR-20190731	31 Jul 2019 13:05		05 Aug 2019 09:58	06 Aug 2019 14:26	1
HS19080113-08	WG-1620-MW22BR-20190731	31 Jul 2019 13:50		05 Aug 2019 09:58	06 Aug 2019 18:33	1
HS19080113-09	WG-1620-MW88C-20190731	31 Jul 2019 15:45		05 Aug 2019 09:58	06 Aug 2019 18:52	1
HS19080113-10	WG-1620-MW67B-20190731	31 Jul 2019 16:45		05 Aug 2019 09:58	06 Aug 2019 19:11	1
HS19080113-11	WG-1620-MW49A-20190731	31 Jul 2019 17:45		05 Aug 2019 09:58	06 Aug 2019 19:30	1
<b>Batch ID: 143819 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19080113-01	WG-1620-MW36D-20190731	31 Jul 2019 07:40		06 Aug 2019 09:00	08 Aug 2019 00:06	1
HS19080113-02	WG-1620-MW38A-20190731	31 Jul 2019 08:35		06 Aug 2019 09:00	08 Aug 2019 00:08	1
HS19080113-03	WG-1620-MW65D-20190731	31 Jul 2019 09:45		06 Aug 2019 09:00	07 Aug 2019 23:50	1
HS19080113-04	WG-1620-MW66D-20190731	31 Jul 2019 10:45		06 Aug 2019 09:00	08 Aug 2019 00:10	1
HS19080113-05	WG-1620-MW59D-20190731	31 Jul 2019 11:50		06 Aug 2019 09:00	08 Aug 2019 00:13	1
<b>Batch ID: 143865 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS19080113-06	WG-1620-FD05-20190731	31 Jul 2019 11:50		07 Aug 2019 09:00	07 Aug 2019 18:26	1
HS19080113-07	WG-1620-MW22AR-20190731	31 Jul 2019 13:05		07 Aug 2019 09:00	07 Aug 2019 18:28	1
HS19080113-08	WG-1620-MW22BR-20190731	31 Jul 2019 13:50		07 Aug 2019 09:00	07 Aug 2019 18:30	1
HS19080113-09	WG-1620-MW88C-20190731	31 Jul 2019 15:45		07 Aug 2019 09:00	07 Aug 2019 18:32	1
HS19080113-10	WG-1620-MW67B-20190731	31 Jul 2019 16:45		07 Aug 2019 09:00	07 Aug 2019 18:34	1
HS19080113-11	WG-1620-MW49A-20190731	31 Jul 2019 17:45		07 Aug 2019 09:00	07 Aug 2019 18:50	1
<b>Batch ID: R343626 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19080113-01	WG-1620-MW36D-20190731	31 Jul 2019 07:40			03 Aug 2019 19:20	1
HS19080113-02	WG-1620-MW38A-20190731	31 Jul 2019 08:35			03 Aug 2019 19:44	1
HS19080113-04	WG-1620-MW66D-20190731	31 Jul 2019 10:45			03 Aug 2019 20:09	1
HS19080113-05	WG-1620-MW59D-20190731	31 Jul 2019 11:50			03 Aug 2019 20:34	1
HS19080113-06	WG-1620-FD05-20190731	31 Jul 2019 11:50			03 Aug 2019 20:58	1
HS19080113-10	WG-1620-MW67B-20190731	31 Jul 2019 16:45			03 Aug 2019 14:49	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	TCLP Date	Prep Date	Analysis Date	DF
<b>Batch ID: R343632 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19080113-07	WG-1620-MW22AR-20190731	31 Jul 2019 13:05			04 Aug 2019 01:54	1
HS19080113-08	WG-1620-MW22BR-20190731	31 Jul 2019 13:50			04 Aug 2019 04:02	1
HS19080113-09	WG-1620-MW88C-20190731	31 Jul 2019 15:45			04 Aug 2019 04:27	1
HS19080113-11	WG-1620-MW49A-20190731	31 Jul 2019 17:45			04 Aug 2019 04:51	1
HS19080113-12	WG-1620-TB06-20190731	31 Jul 2019 00:00			04 Aug 2019 00:40	1
<b>Batch ID: R343643 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS19080113-03	WG-1620-MW65D-20190731	31 Jul 2019 09:45			03 Aug 2019 14:42	1

WorkOrder: HS19080113  
InstrumentID: ICPMS05  
Test Code: ICP\_TW  
Test Number: SW6020  
Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000547	0.000400	0.00200

WorkOrder: HS19080113  
 InstrumentID: ICPMS04  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.00100	0.000579	0.000400	0.00200

WorkOrder: HS19080113  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

Matrix: Aqueous

Units: mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00011	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.00010	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000096	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00012	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00011	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000049	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000013	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.00013	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000051	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000055	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000068	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000053	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.000086	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000066	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000052	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.000099	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000052	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000061	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00013	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000085	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000055	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000094	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000061	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS19080113  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00064	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00060	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00061	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00077	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00058	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00070	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00062	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS19080113  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00067	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00056	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00059	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00051	0.00030	0.0010
A	Methylene chloride	75-09-2	0.00050	0.00054	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00056	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00052	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143819 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-143819</b>	Units: <b>mg/L</b>		Analysis Date: <b>07-Aug-2019 23:45</b>					
Client ID:		Run ID: <b>ICPMS05_343857</b>	SeqNo: <b>5201038</b>	PrepDate: <b>06-Aug-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-143819</b>	Units: <b>mg/L</b>		Analysis Date: <b>07-Aug-2019 23:48</b>					
Client ID:		Run ID: <b>ICPMS05_343857</b>	SeqNo: <b>5201039</b>	PrepDate: <b>06-Aug-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.04979	0.00200	0.05	0	99.6	80 - 120			
<b>MS</b>	Sample ID: <b>HS19080113-03MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>07-Aug-2019 23:55</b>					
Client ID: <b>WG-1620-MW65D-20190731</b>		Run ID: <b>ICPMS05_343857</b>	SeqNo: <b>5201042</b>	PrepDate: <b>06-Aug-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.0508	0.00200	0.05	0.001348	98.9	80 - 120			
<b>MSD</b>	Sample ID: <b>HS19080113-03MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>07-Aug-2019 23:57</b>					
Client ID: <b>WG-1620-MW65D-20190731</b>		Run ID: <b>ICPMS05_343857</b>	SeqNo: <b>5201043</b>	PrepDate: <b>06-Aug-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05054	0.00200	0.05	0.001348	98.4	80 - 120	0.0508	0.507	20
<b>PDS</b>	Sample ID: <b>HS19080113-03PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>07-Aug-2019 23:59</b>					
Client ID: <b>WG-1620-MW65D-20190731</b>		Run ID: <b>ICPMS05_343857</b>	SeqNo: <b>5201044</b>	PrepDate: <b>06-Aug-2019</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.1012	0.00200	0.1	0.001348	99.9	75 - 125			
<b>SD</b>	Sample ID: <b>HS19080113-03SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>07-Aug-2019 23:52</b>					
Client ID: <b>WG-1620-MW65D-20190731</b>		Run ID: <b>ICPMS05_343857</b>	SeqNo: <b>5201041</b>	PrepDate: <b>06-Aug-2019</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	U	0.0100					0.001348	0	10
The following samples were analyzed in this batch:									
	HS19080113-01	HS19080113-02	HS19080113-03	HS19080113-04					
	HS19080113-05								

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

<b>Batch ID:</b> 143865 ( 0 )	<b>Instrument:</b> ICPMS04	<b>Method:</b> ICP-MS METALS BY SW6020A
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<b>MBLK</b>	Sample ID: <b>MBLK-143865</b>	Units: <b>mg/L</b>	Analysis Date: <b>07-Aug-2019 18:22</b>							
Client ID:	Run ID: <b>ICPMS04_343800</b>	SeqNo: <b>5200782</b>	PrepDate: <b>07-Aug-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic U 0.00200

<b>LCS</b>	Sample ID: <b>LCS-143865</b>	Units: <b>mg/L</b>	Analysis Date: <b>07-Aug-2019 18:24</b>							
Client ID:	Run ID: <b>ICPMS04_343800</b>	SeqNo: <b>5200783</b>	PrepDate: <b>07-Aug-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.0481 0.00200 0.05 0 96.2 80 - 120

<b>MS</b>	Sample ID: <b>HS19080113-10MS</b>	Units: <b>mg/L</b>	Analysis Date: <b>07-Aug-2019 18:38</b>							
Client ID: <b>WG-1620-MW67B-20190731</b>	Run ID: <b>ICPMS04_343800</b>	SeqNo: <b>5200920</b>	PrepDate: <b>07-Aug-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.04857 0.00200 0.05 0.000494 96.2 80 - 120

<b>MSD</b>	Sample ID: <b>HS19080113-10MSD</b>	Units: <b>mg/L</b>	Analysis Date: <b>07-Aug-2019 18:46</b>							
Client ID: <b>WG-1620-MW67B-20190731</b>	Run ID: <b>ICPMS04_343800</b>	SeqNo: <b>5200924</b>	PrepDate: <b>07-Aug-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.0492 0.00200 0.05 0.000494 97.4 80 - 120 0.04857 1.28 20

<b>PDS</b>	Sample ID: <b>HS19080113-10PDS</b>	Units: <b>mg/L</b>	Analysis Date: <b>07-Aug-2019 18:48</b>							
Client ID: <b>WG-1620-MW67B-20190731</b>	Run ID: <b>ICPMS04_343800</b>	SeqNo: <b>5200925</b>	PrepDate: <b>07-Aug-2019</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.1001 0.00200 0.1 0.000494 99.7 75 - 125

<b>SD</b>	Sample ID: <b>HS19080113-10SD</b>	Units: <b>mg/L</b>	Analysis Date: <b>07-Aug-2019 18:36</b>							
Client ID: <b>WG-1620-MW67B-20190731</b>	Run ID: <b>ICPMS04_343800</b>	SeqNo: <b>5200919</b>	PrepDate: <b>07-Aug-2019</b> DF: <b>5</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual

Arsenic U 0.0100 0.000494 0 10

<b>The following samples were analyzed in this batch:</b>	HS19080113-06	HS19080113-07	HS19080113-08	HS19080113-09
	HS19080113-10	HS19080113-11		



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-143759	Units: ug/L			Analysis Date: 05-Aug-2019 13:03					
Client ID:	Run ID: SV-7_343719	SeqNo: 5197915		PrepDate: 05-Aug-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>2.803</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>56.1</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.478</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.6</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.943</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.9</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>3.689</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.8</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.625</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.5</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.474</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.5</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-143759	Units: ug/L			Analysis Date: 05-Aug-2019 13:23					
Client ID:	Run ID: SV-7_343719	SeqNo: 5197916		PrepDate: 05-Aug-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.596	0.20	5	0	91.9	39 - 127				
2,4-Dimethylphenol	4.075	0.20	5	0	81.5	35 - 120				
2,4-Dinitrotoluene	4.768	0.20	5	0	95.4	50 - 122				
2,6-Dinitrotoluene	4.413	0.20	5	0	88.3	50 - 120				
2-Chloronaphthalene	4.117	0.20	5	0	82.3	50 - 120				
2-Methylnaphthalene	4.007	0.10	5	0	80.1	50 - 120				
4,6-Dinitro-2-methylphenol	3.484	0.20	5	0	69.7	25 - 121				
4-Nitrophenol	3.601	1.0	5	0	72.0	30 - 130				
Acenaphthene	4.114	0.10	5	0	82.3	45 - 120				
Acenaphthylene	4.278	0.10	5	0	85.6	47 - 120				
Anthracene	3.67	0.10	5	0	73.4	45 - 120				
Benz(a)anthracene	3.909	0.10	5	0	78.2	40 - 120				
Benzo(a)pyrene	3.355	0.10	5	0	67.1	45 - 120				
Bis(2-chloroethoxy)methane	3.992	0.20	5	0	79.8	45 - 120				
Bis(2-ethylhexyl)phthalate	4.151	0.20	5	0	83.0	40 - 139				
Chrysene	3.817	0.10	5	0	76.3	43 - 120				
Dibenzofuran	4.17	0.10	5	0	83.4	50 - 120				
Di-n-butyl phthalate	4.069	0.20	5	0	81.4	45 - 123				
Fluoranthene	3.867	0.10	5	0	77.3	45 - 125				
Fluorene	4.032	0.10	5	0	80.6	49 - 120				
Naphthalene	4.025	0.10	5	0	80.5	45 - 120				
Nitrobenzene	4.643	0.20	5	0	92.9	44 - 120				
N-Nitrosodiphenylamine	4.413	0.20	5	0	88.3	40 - 125				
Pentachlorophenol	2.175	0.20	5	0	43.5	19 - 121				
Phenanthrene	3.605	0.10	5	0	72.1	45 - 121				
Phenol	4.362	0.20	5	0	87.2	20 - 124				
Pyrene	4.031	0.10	5	0	80.6	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.862</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.2</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.737</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.7</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.407</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.1</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>3.975</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>79.5</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.04</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.673</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.5</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS19080113-10MS		Units: ug/L		Analysis Date: 06-Aug-2019 19:49				
Client ID: WG-1620-MW67B-20190731		Run ID: SV-7_343746		SeqNo: 5199678		PrepDate: 05-Aug-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	3.613	0.20	5	0	72.3	39 - 127				
2,4-Dimethylphenol	2.852	0.20	5	0	57.0	35 - 120				
2,4-Dinitrotoluene	1.958	0.20	5	0	39.2	50 - 122			S	
2,6-Dinitrotoluene	2.042	0.20	5	0	40.8	50 - 120			S	
2-Chloronaphthalene	2.937	0.20	5	0	58.7	50 - 120				
2-Methylnaphthalene	4.153	0.10	5	0	83.1	50 - 120				
4,6-Dinitro-2-methylphenol	0.1571	0.20	5	0	3.14	25 - 121			JS	
4-Nitrophenol	1.936	1.0	5	0	38.7	30 - 130				
Acenaphthene	2.841	0.10	5	0	56.8	45 - 120				
Acenaphthylene	2.95	0.10	5	0	59.0	47 - 120				
Anthracene	3.441	0.10	5	0	68.8	45 - 120				
Benz(a)anthracene	4.006	0.10	5	0	80.1	40 - 120				
Benzo(a)pyrene	5.331	0.10	5	0	107	45 - 120				
Bis(2-chloroethoxy)methane	2.695	0.20	5	0	53.9	45 - 120				
Bis(2-ethylhexyl)phthalate	4.215	0.20	5	0	84.3	40 - 139				
Chrysene	4.288	0.10	5	0	85.8	43 - 120				
Dibenzofuran	3.094	0.10	5	0	61.9	50 - 120				
Di-n-butyl phthalate	3.623	0.20	5	0	72.5	45 - 123				
Fluoranthene	3.444	0.10	5	0	68.9	45 - 125				
Fluorene	3.039	0.10	5	0	60.8	49 - 120				
Naphthalene	2.939	0.10	5	0.07887	57.2	45 - 120				
Nitrobenzene	3.422	0.20	5	0	68.4	44 - 120				
N-Nitrosodiphenylamine	4.15	0.20	5	0	83.0	40 - 125				
Pentachlorophenol	U	0.20	5	0	0	19 - 121			S	
Phenanthrene	3.357	0.10	5	0	67.1	45 - 121				
Phenol	3.387	0.20	5	0	67.7	20 - 124				
Pyrene	4.191	0.10	5	0	83.8	40 - 130				
Surr: 2,4,6-Tribromophenol	3.616	0.20	5	0	72.3	34 - 129				
Surr: 2-Fluorobiphenyl	3.153	0.20	5	0	63.1	40 - 125				
Surr: 2-Fluorophenol	3.019	0.20	5	0	60.4	20 - 120				
Surr: 4-Terphenyl-d14	4.415	0.20	5	0	88.3	40 - 135				
Surr: Nitrobenzene-d5	3.077	0.20	5	0	61.5	41 - 120				
Surr: Phenol-d6	3.356	0.20	5	0	67.1	20 - 120				

**Client:** Golder Associates Inc.  
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**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS19080113-03MS		Units: ug/L		Analysis Date: 05-Aug-2019 20:40				
Client ID: WG-1620-MW65D-20190731		Run ID: SV-7_343719		SeqNo: 5197927		PrepDate: 05-Aug-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.676	0.20	5	0	93.5	39 - 127				
2,4-Dimethylphenol	3.928	0.20	5	0	78.6	35 - 120				
2,4-Dinitrotoluene	4.766	0.20	5	0	95.3	50 - 122				
2,6-Dinitrotoluene	4.315	0.20	5	0	86.3	50 - 120				
2-Chloronaphthalene	4.479	0.20	5	0	89.6	50 - 120				
2-Methylnaphthalene	4.225	0.10	5	0	84.5	50 - 120				
4,6-Dinitro-2-methylphenol	1.14	0.20	5	0	22.8	25 - 121			S	
4-Nitrophenol	3.606	1.0	5	0	72.1	30 - 130				
Acenaphthene	4.307	0.10	5	0	86.1	45 - 120				
Acenaphthylene	4.323	0.10	5	0	86.5	47 - 120				
Anthracene	3.817	0.10	5	0	76.3	45 - 120				
Benz(a)anthracene	4.159	0.10	5	0	83.2	40 - 120				
Benzo(a)pyrene	5.611	0.10	5	0	112	45 - 120				
Bis(2-chloroethoxy)methane	4.08	0.20	5	0	81.6	45 - 120				
Bis(2-ethylhexyl)phthalate	4.747	0.20	5	0.1617	91.7	40 - 139				
Chrysene	4.332	0.10	5	0	86.6	43 - 120				
Dibenzofuran	4.302	0.10	5	0	86.0	50 - 120				
Di-n-butyl phthalate	4.32	0.20	5	0	86.4	45 - 123				
Fluoranthene	4.236	0.10	5	0	84.7	45 - 125				
Fluorene	4.113	0.10	5	0	82.3	49 - 120				
Naphthalene	4.197	0.10	5	0	83.9	45 - 120				
Nitrobenzene	4.907	0.20	5	0	98.1	44 - 120				
N-Nitrosodiphenylamine	4.519	0.20	5	0	90.4	40 - 125				
Pentachlorophenol	2.946	0.20	5	0	58.9	19 - 121				
Phenanthrene	3.762	0.10	5	0	75.2	45 - 121				
Phenol	4.887	0.20	5	0	97.7	20 - 124				
Pyrene	4.342	0.10	5	0	86.8	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.707</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.1</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.927</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.5</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.542</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.8</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.579</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.221</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.092</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS	Sample ID: HS19071516-01MS	Units: ug/L			Analysis Date: 05-Aug-2019 15:55					
Client ID:	Run ID: SV-7_343719	SeqNo: 5197918	PrepDate: 05-Aug-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.399	0.20	5	0	88.0	39 - 127				
2,4-Dimethylphenol	3.824	0.20	5	0	76.5	35 - 120				
2,4-Dinitrotoluene	4.519	0.20	5	0	90.4	50 - 122				
2,6-Dinitrotoluene	4.262	0.20	5	0	85.2	50 - 120				
2-Chloronaphthalene	3.69	0.20	5	0	73.8	50 - 120				
2-Methylnaphthalene	3.94	0.10	5	0	78.8	50 - 120				
4,6-Dinitro-2-methylphenol	3.498	0.20	5	0	70.0	25 - 121				
4-Nitrophenol	4.386	1.0	5	0	87.7	30 - 130				
Acenaphthene	3.859	0.10	5	0	77.2	45 - 120				
Acenaphthylene	3.976	0.10	5	0	79.5	47 - 120				
Anthracene	3.561	0.10	5	0	71.2	45 - 120				
Benz(a)anthracene	3.993	0.10	5	0	79.9	40 - 120				
Benzo(a)pyrene	5.251	0.10	5	0	105	45 - 120				
Bis(2-chloroethoxy)methane	3.715	0.20	5	0	74.3	45 - 120				
Bis(2-ethylhexyl)phthalate	4.166	0.20	5	0	83.3	40 - 139				
Chrysene	3.847	0.10	5	0	76.9	43 - 120				
Dibenzofuran	3.979	0.10	5	0	79.6	50 - 120				
Di-n-butyl phthalate	4.067	0.20	5	0	81.3	45 - 123				
Fluoranthene	3.908	0.10	5	0	78.2	45 - 125				
Fluorene	3.81	0.10	5	0	76.2	49 - 120				
Naphthalene	3.784	0.10	5	0.04766	74.7	45 - 120				
Nitrobenzene	4.386	0.20	5	0	87.7	44 - 120				
N-Nitrosodiphenylamine	4.007	0.20	5	0	80.1	40 - 125				
Pentachlorophenol	2.421	0.20	5	0	48.4	19 - 121				
Phenanthrene	3.562	0.10	5	0	71.2	45 - 121				
Phenol	4.081	0.20	5	0	81.6	20 - 124				
Pyrene	4.149	0.10	5	0	83.0	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.919</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.4</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.26</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.798</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>3.946</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.9</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.496</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.9</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.355</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.1</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS19080113-10MSD		Units: ug/L		Analysis Date: 06-Aug-2019 20:08				
Client ID: WG-1620-MW67B-20190731		Run ID: SV-7_343746		SeqNo: 5199679		PrepDate: 05-Aug-2019		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.667	0.20	5	0	73.3	39 - 127	3.613	1.47	20	
2,4-Dimethylphenol	2.94	0.20	5	0	58.8	35 - 120	2.852	3.03	20	
2,4-Dinitrotoluene	1.928	0.20	5	0	38.6	50 - 122	1.958	1.51	20	S
2,6-Dinitrotoluene	2.041	0.20	5	0	40.8	50 - 120	2.042	0.0463	20	S
2-Chloronaphthalene	2.621	0.20	5	0	52.4	50 - 120	2.937	11.4	20	
2-Methylnaphthalene	4.352	0.10	5	0	87.0	50 - 120	4.153	4.67	20	
4,6-Dinitro-2-methylphenol	0.1604	0.20	5	0	3.21	25 - 121	0.1571	0	30	JS
4-Nitrophenol	7.878	1.0	5	0	158	30 - 130	1.936	121	20	SR
Acenaphthene	3.025	0.10	5	0	60.5	45 - 120	2.841	6.28	20	
Acenaphthylene	3.102	0.10	5	0	62.0	47 - 120	2.95	5.03	20	
Anthracene	3.501	0.10	5	0	70.0	45 - 120	3.441	1.73	20	
Benz(a)anthracene	3.825	0.10	5	0	76.5	40 - 120	4.006	4.62	20	
Benzo(a)pyrene	5.442	0.10	5	0	109	45 - 120	5.331	2.07	20	
Bis(2-chloroethoxy)methane	2.818	0.20	5	0	56.4	45 - 120	2.695	4.49	20	
Bis(2-ethylhexyl)phthalate	4.233	0.20	5	0	84.7	40 - 139	4.215	0.428	20	
Chrysene	4.383	0.10	5	0	87.7	43 - 120	4.288	2.19	20	
Dibenzofuran	3.285	0.10	5	0	65.7	50 - 120	3.094	5.97	20	
Di-n-butyl phthalate	3.876	0.20	5	0	77.5	45 - 123	3.623	6.75	20	
Fluoranthene	3.519	0.10	5	0	70.4	45 - 125	3.444	2.15	20	
Fluorene	3.179	0.10	5	0	63.6	49 - 120	3.039	4.5	20	
Naphthalene	3.058	0.10	5	0.07887	59.6	45 - 120	2.939	3.98	20	
Nitrobenzene	3.425	0.20	5	0	68.5	44 - 120	3.422	0.0745	20	
N-Nitrosodiphenylamine	4.252	0.20	5	0	85.0	40 - 125	4.15	2.43	20	
Pentachlorophenol	0.1573	0.20	5	0	3.15	19 - 121	0	0	20	JS
Phenanthrene	3.467	0.10	5	0	69.3	45 - 121	3.357	3.24	20	
Phenol	3.699	0.20	5	0	74.0	20 - 124	3.387	8.79	20	
Pyrene	4.308	0.10	5	0	86.2	40 - 130	4.191	2.74	20	
Surr: 2,4,6-Tribromophenol	3.722	0.20	5	0	74.4	34 - 129	3.616	2.89	20	
Surr: 2-Fluorobiphenyl	3.234	0.20	5	0	64.7	40 - 125	3.153	2.54	20	
Surr: 2-Fluorophenol	3.12	0.20	5	0	62.4	20 - 120	3.019	3.3	20	
Surr: 4-Terphenyl-d14	4.422	0.20	5	0	88.4	40 - 135	4.415	0.152	20	
Surr: Nitrobenzene-d5	3.095	0.20	5	0	61.9	41 - 120	3.077	0.565	20	
Surr: Phenol-d6	3.408	0.20	5	0	68.2	20 - 120	3.356	1.52	20	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS19080113-03MSD	Units: ug/L			Analysis Date: 05-Aug-2019 20:59					
Client ID: WG-1620-MW65D-20190731	Run ID: SV-7_343719	SeqNo: 5197928	PrepDate: 05-Aug-2019	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.528	0.20	5	0	90.6	39 - 127	4.676	3.22	20	
2,4-Dimethylphenol	3.961	0.20	5	0	79.2	35 - 120	3.928	0.822	20	
2,4-Dinitrotoluene	4.569	0.20	5	0	91.4	50 - 122	4.766	4.2	20	
2,6-Dinitrotoluene	4.034	0.20	5	0	80.7	50 - 120	4.315	6.73	20	
2-Chloronaphthalene	3.744	0.20	5	0	74.9	50 - 120	4.479	17.9	20	
2-Methylnaphthalene	4.133	0.10	5	0	82.7	50 - 120	4.225	2.19	20	
4,6-Dinitro-2-methylphenol	1.177	0.20	5	0	23.5	25 - 121	1.14	3.14	30	S
4-Nitrophenol	4.763	1.0	5	0	95.3	30 - 130	3.606	27.6	20	R
Acenaphthene	4.07	0.10	5	0	81.4	45 - 120	4.307	5.64	20	
Acenaphthylene	4.207	0.10	5	0	84.1	47 - 120	4.323	2.72	20	
Anthracene	3.8	0.10	5	0	76.0	45 - 120	3.817	0.46	20	
Benz(a)anthracene	4.432	0.10	5	0	88.6	40 - 120	4.159	6.35	20	
Benzo(a)pyrene	6.166	0.10	5	0	123	45 - 120	5.611	9.42	20	S
Bis(2-chloroethoxy)methane	4.019	0.20	5	0	80.4	45 - 120	4.08	1.5	20	
Bis(2-ethylhexyl)phthalate	4.852	0.20	5	0.1617	93.8	40 - 139	4.747	2.19	20	
Chrysene	4.549	0.10	5	0	91.0	43 - 120	4.332	4.9	20	
Dibenzofuran	4.121	0.10	5	0	82.4	50 - 120	4.302	4.3	20	
Di-n-butyl phthalate	4.609	0.20	5	0	92.2	45 - 123	4.32	6.47	20	
Fluoranthene	4.319	0.10	5	0	86.4	45 - 125	4.236	1.94	20	
Fluorene	3.934	0.10	5	0	78.7	49 - 120	4.113	4.45	20	
Naphthalene	4.063	0.10	5	0	81.3	45 - 120	4.197	3.24	20	
Nitrobenzene	4.783	0.20	5	0	95.7	44 - 120	4.907	2.57	20	
N-Nitrosodiphenylamine	4.317	0.20	5	0	86.3	40 - 125	4.519	4.59	20	
Pentachlorophenol	2.948	0.20	5	0	59.0	19 - 121	2.946	0.0694	20	
Phenanthrene	3.74	0.10	5	0	74.8	45 - 121	3.762	0.598	20	
Phenol	4.613	0.20	5	0	92.3	20 - 124	4.887	5.78	20	
Pyrene	4.298	0.10	5	0	86.0	40 - 130	4.342	1.03	20	
Surr: 2,4,6-Tribromophenol	4.088	0.20	5	0	81.8	34 - 129	4.707	14.1	20	
Surr: 2-Fluorobiphenyl	4.35	0.20	5	0	87.0	40 - 125	4.927	12.4	20	
Surr: 2-Fluorophenol	3.992	0.20	5	0	79.8	20 - 120	4.542	12.9	20	
Surr: 4-Terphenyl-d14	4.195	0.20	5	0	83.9	40 - 135	4.579	8.76	20	
Surr: Nitrobenzene-d5	4.806	0.20	5	0	96.1	41 - 120	5.221	8.28	20	
Surr: Phenol-d6	4.535	0.20	5	0	90.7	20 - 120	5.092	11.6	20	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

Batch ID: 143759 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS19071516-01MSD	Units: ug/L			Analysis Date: 05-Aug-2019 16:14					
Client ID:	Run ID: SV-7_343719	SeqNo: 5197919		PrepDate: 05-Aug-2019		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.873	0.20	5	0	97.5	39 - 127	4.399	10.2	20	
2,4-Dimethylphenol	4.141	0.20	5	0	82.8	35 - 120	3.824	7.97	20	
2,4-Dinitrotoluene	4.836	0.20	5	0	96.7	50 - 122	4.519	6.77	20	
2,6-Dinitrotoluene	4.569	0.20	5	0	91.4	50 - 120	4.262	6.96	20	
2-Chloronaphthalene	4.261	0.20	5	0	85.2	50 - 120	3.69	14.4	20	
2-Methylnaphthalene	4.215	0.10	5	0	84.3	50 - 120	3.94	6.75	20	
4,6-Dinitro-2-methylphenol	3.803	0.20	5	0	76.1	25 - 121	3.498	8.35	30	
4-Nitrophenol	4.882	1.0	5	0	97.6	30 - 130	4.386	10.7	20	
Acenaphthene	4.269	0.10	5	0	85.4	45 - 120	3.859	10.1	20	
Acenaphthylene	4.432	0.10	5	0	88.6	47 - 120	3.976	10.9	20	
Anthracene	3.893	0.10	5	0	77.9	45 - 120	3.561	8.92	20	
Benz(a)anthracene	4.137	0.10	5	0	82.7	40 - 120	3.993	3.54	20	
Benzo(a)pyrene	5.708	0.10	5	0	114	45 - 120	5.251	8.34	20	
Bis(2-chloroethoxy)methane	4.091	0.20	5	0	81.8	45 - 120	3.715	9.62	20	
Bis(2-ethylhexyl)phthalate	4.551	0.20	5	0	91.0	40 - 139	4.166	8.83	20	
Chrysene	4.159	0.10	5	0	83.2	43 - 120	3.847	7.8	20	
Dibenzofuran	4.355	0.10	5	0	87.1	50 - 120	3.979	9.04	20	
Di-n-butyl phthalate	4.434	0.20	5	0	88.7	45 - 123	4.067	8.63	20	
Fluoranthene	4.184	0.10	5	0	83.7	45 - 125	3.908	6.82	20	
Fluorene	4.148	0.10	5	0	83.0	49 - 120	3.81	8.5	20	
Naphthalene	4.222	0.10	5	0.04766	83.5	45 - 120	3.784	11	20	
Nitrobenzene	4.847	0.20	5	0	96.9	44 - 120	4.386	9.97	20	
N-Nitrosodiphenylamine	4.594	0.20	5	0	91.9	40 - 125	4.007	13.6	20	
Pentachlorophenol	2.472	0.20	5	0	49.4	19 - 121	2.421	2.08	20	
Phenanthrene	3.882	0.10	5	0	77.6	45 - 121	3.562	8.6	20	
Phenol	4.803	0.20	5	0	96.1	20 - 124	4.081	16.2	20	
Pyrene	4.443	0.10	5	0	88.9	40 - 130	4.149	6.84	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.157</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.1</i>	<i>34 - 129</i>	<i>3.919</i>	<i>5.88</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.613</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.3</i>	<i>40 - 125</i>	<i>4.26</i>	<i>7.95</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.256</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.1</i>	<i>20 - 120</i>	<i>3.798</i>	<i>11.4</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.073</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>81.5</i>	<i>40 - 135</i>	<i>3.946</i>	<i>3.18</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>4.919</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.4</i>	<i>41 - 120</i>	<i>4.496</i>	<i>8.99</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.413</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.3</i>	<i>20 - 120</i>	<i>4.355</i>	<i>1.3</i>	<i>20</i>	



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

**Batch ID:** 143759 ( 0 )      **Instrument:** SV-7      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

The following samples were analyzed in this batch:

HS19080113-01	HS19080113-02	HS19080113-03	HS19080113-04
HS19080113-05	HS19080113-06	HS19080113-07	HS19080113-08
HS19080113-09	HS19080113-10	HS19080113-11	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

**Batch ID:** R343626 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190803</b>		Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 11:32</b>			
Client ID:		Run ID: <b>VOA4_343626</b>		SeqNo: <b>5195368</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.2</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>47.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.3</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>52.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>81 - 120</i>			

<b>LCS</b>		Sample ID: <b>VLCSW-190803</b>		Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 10:43</b>			
Client ID:		Run ID: <b>VOA4_343626</b>		SeqNo: <b>5195366</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	17.53	1.0	20	0	87.7	70 - 124			
Benzene	17.53	1.0	20	0	87.6	74 - 120			
Chlorobenzene	19.08	1.0	20	0	95.4	76 - 113			
Ethylbenzene	20.7	1.0	20	0	104	77 - 117			
Methylene chloride	17.58	2.0	20	0	87.9	70 - 127			
Toluene	19.46	1.0	20	0	97.3	77 - 118			
Xylenes, Total	62.78	1.0	60	0	105	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.06</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.1</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.13</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>46.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.2</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>52.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

**Batch ID:** R343626 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS19080113-10MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 15:14</b>			
Client ID: <b>WG-1620-MW67B-20190731</b>		Run ID: <b>VOA4_343626</b>			SeqNo: <b>5195377</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.8	1.0	20	0	74.0	70 - 127				
Benzene	15.06	1.0	20	0	75.3	70 - 127				
Chlorobenzene	16.65	1.0	20	0	83.3	70 - 114				
Ethylbenzene	17.5	1.0	20	0	87.5	70 - 124				
Methylene chloride	14.86	2.0	20	0	74.3	70 - 128				
Toluene	16.77	1.0	20	0	83.8	70 - 123				
Xylenes, Total	53.32	1.0	60	0	88.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.56</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.1</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.61</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.2</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>52.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS19080113-10MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 15:38</b>			
Client ID: <b>WG-1620-MW67B-20190731</b>		Run ID: <b>VOA4_343626</b>			SeqNo: <b>5195378</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	14.38	1.0	20	0	71.9	70 - 127	14.8	2.86	20	
Benzene	14.66	1.0	20	0	73.3	70 - 127	15.06	2.74	20	
Chlorobenzene	16.08	1.0	20	0	80.4	70 - 114	16.65	3.48	20	
Ethylbenzene	17.23	1.0	20	0	86.1	70 - 124	17.5	1.53	20	
Methylene chloride	14.35	2.0	20	0	71.7	70 - 128	14.86	3.49	20	
Toluene	16.37	1.0	20	0	81.8	70 - 123	16.77	2.43	20	
Xylenes, Total	51.69	1.0	60	0	86.1	70 - 130	53.32	3.11	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.0</i>	<i>70 - 126</i>	<i>47.56</i>	<i>0.156</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.71</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>81 - 113</i>	<i>48.65</i>	<i>0.105</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>77 - 123</i>	<i>47.61</i>	<i>0.52</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>52.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>82 - 127</i>	<i>52.52</i>	<i>0.972</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19080113-01    HS19080113-02    HS19080113-04    HS19080113-05  
 HS19080113-06    HS19080113-10

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

<b>Batch ID:</b> R343632 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-190803</b>	Units: <b>ug/L</b>			Analysis Date: <b>03-Aug-2019 23:50</b>				
Client ID:	Run ID: <b>VOA4_343632</b>	SeqNo: <b>5195438</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.1</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>52.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-190803</b>	Units: <b>ug/L</b>			Analysis Date: <b>03-Aug-2019 23:01</b>				
Client ID:	Run ID: <b>VOA4_343632</b>	SeqNo: <b>5195436</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	16.76	1.0	20	0	83.8	70 - 124			
Benzene	17.21	1.0	20	0	86.0	74 - 120			
Chlorobenzene	18.92	1.0	20	0	94.6	76 - 113			
Ethylbenzene	19.92	1.0	20	0	99.6	77 - 117			
Methylene chloride	16.69	2.0	20	0	83.4	70 - 127			
Toluene	18.81	1.0	20	0	94.1	77 - 118			
Vinyl chloride	16.62	1.0	20	0	83.1	70 - 130			
Xylenes, Total	59.58	1.0	60	0	99.3	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.68</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.4</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.82</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>46.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.3</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>52.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>106</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

<b>Batch ID:</b> R343632 ( 0 )	<b>Instrument:</b> VOA4	<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C								
<b>MS</b>	Sample ID: <b>HS19080113-07MS</b>	Units: <b>ug/L</b>	Analysis Date: <b>04-Aug-2019 02:45</b>							
Client ID: <b>WG-1620-MW22AR-20190731</b>	Run ID: <b>VOA4_343632</b>	SeqNo: <b>5195445</b>	PrepDate: <b>DF: 1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

1,2-Dichloroethane	15.58	1.0	20	0	77.9	70 - 127				
Benzene	16.65	1.0	20	0	83.2	70 - 127				
Chlorobenzene	17.56	1.0	20	0	87.8	70 - 114				
Ethylbenzene	19.01	1.0	20	0	95.0	70 - 124				
Methylene chloride	16.97	2.0	20	0	84.9	70 - 128				
Toluene	18.42	1.0	20	0	92.1	70 - 123				
Vinyl chloride	17.57	1.0	20	0	87.9	70 - 130				
Xylenes, Total	56.24	1.0	60	0	93.7	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.4</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.44</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.38</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>52.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>82 - 127</i>				

<b>MSD</b>	Sample ID: <b>HS19080113-07MSD</b>	Units: <b>ug/L</b>	Analysis Date: <b>04-Aug-2019 03:10</b>							
Client ID: <b>WG-1620-MW22AR-20190731</b>	Run ID: <b>VOA4_343632</b>	SeqNo: <b>5195446</b>	PrepDate: <b>DF: 1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.43	1.0	20	0	77.1	70 - 127	15.58	0.975	20	
Benzene	16.18	1.0	20	0	80.9	70 - 127	16.65	2.85	20	
Chlorobenzene	17.27	1.0	20	0	86.3	70 - 114	17.56	1.65	20	
Ethylbenzene	19.24	1.0	20	0	96.2	70 - 124	19.01	1.23	20	
Methylene chloride	15.86	2.0	20	0	79.3	70 - 128	16.97	6.8	20	
Toluene	17.3	1.0	20	0	86.5	70 - 123	18.42	6.27	20	
Vinyl chloride	16.33	1.0	20	0	81.7	70 - 130	17.57	7.31	20	
Xylenes, Total	56.89	1.0	60	0	94.8	70 - 130	56.24	1.15	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.3</i>	<i>70 - 126</i>	<i>46.2</i>	<i>0.937</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.18</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 113</i>	<i>49.44</i>	<i>1.47</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>45.82</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>91.6</i>	<i>77 - 123</i>	<i>47.38</i>	<i>3.35</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>52.67</i>	<i>3.91</i>	<i>20</i>	

The following samples were analyzed in this batch: 

HS19080113-07	HS19080113-08	HS19080113-09	HS19080113-11
HS19080113-12			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

**Batch ID:** R343643 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-190803</b>		Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 11:02</b>			
Client ID:		Run ID: <b>VOA2_343643</b>		SeqNo: <b>5195692</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>52.08</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>			

<b>LCS</b>		Sample ID: <b>VLCSW-190803</b>		Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 10:38</b>			
Client ID:		Run ID: <b>VOA2_343643</b>		SeqNo: <b>5195691</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	18.9	1.0	20	0	94.5	70 - 124			
Benzene	19.91	1.0	20	0	99.5	74 - 120			
Chlorobenzene	20.14	1.0	20	0	101	76 - 113			
Ethylbenzene	20.43	1.0	20	0	102	77 - 117			
Methylene chloride	20.9	2.0	20	0	105	70 - 127			
Toluene	20.47	1.0	20	0	102	77 - 118			
Xylenes, Total	61.73	1.0	60	0	103	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>52.93</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>106</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QC BATCH REPORT**

**Batch ID:** R343643 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS19080113-03MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 15:07</b>			
Client ID: <b>WG-1620-MW65D-20190731</b>		Run ID: <b>VOA2_343643</b>			SeqNo: <b>5195702</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.74	1.0	20	0	78.7	70 - 127				
Benzene	17.04	1.0	20	0	85.2	70 - 127				
Chlorobenzene	17.56	1.0	20	0	87.8	70 - 114				
Ethylbenzene	17.79	1.0	20	0	89.0	70 - 124				
Methylene chloride	16.71	2.0	20	0	83.5	70 - 128				
Toluene	17.73	1.0	20	0	88.7	70 - 123				
Xylenes, Total	54.55	1.0	60	0	90.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.62</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.18</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS19080113-03MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>03-Aug-2019 15:31</b>			
Client ID: <b>WG-1620-MW65D-20190731</b>		Run ID: <b>VOA2_343643</b>			SeqNo: <b>5195703</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.52	1.0	20	0	77.6	70 - 127	15.74	1.42	20	
Benzene	16.55	1.0	20	0	82.7	70 - 127	17.04	2.94	20	
Chlorobenzene	17.09	1.0	20	0	85.4	70 - 114	17.56	2.75	20	
Ethylbenzene	17.37	1.0	20	0	86.8	70 - 124	17.79	2.42	20	
Methylene chloride	16.66	2.0	20	0	83.3	70 - 128	16.71	0.3	20	
Toluene	17.45	1.0	20	0	87.3	70 - 123	17.73	1.61	20	
Xylenes, Total	52.42	1.0	60	0	87.4	70 - 130	54.55	3.99	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.41</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 126</i>	<i>50.62</i>	<i>0.421</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>81 - 113</i>	<i>48.95</i>	<i>0.612</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>50.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>77 - 123</i>	<i>50.18</i>	<i>0.568</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>82 - 127</i>	<i>50.5</i>	<i>1.53</i>	<i>20</i>	

The following samples were analyzed in this batch: HS19080113-03

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS19080113

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter



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**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2019	31-Dec-2019
Oklahoma	2018-156	31-Aug-2019
Texas	TX104704231-19-23	30-Apr-2020

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS19080113

**SAMPLE TRACKING**

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Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS19080113-01	WG-1620-MW36D-20190731	Login	02/08/2019 12:41:31	JRM	MET001
HS19080113-01	WG-1620-MW36D-20190731	Login	02/08/2019 12:41:31	JRM	EXT002
HS19080113-01	WG-1620-MW36D-20190731	Login	02/08/2019 12:41:31	JRM	VOA088

Sample Receipt Checklist

Client Name: PBW
Work Order: HS19080113

Date/Time Received: 02-Aug-2019 12:00
Received by: JRM

Checklist completed by: Jared R. Makan
eSignature
Date: 2-Aug-2019

Reviewed by: Dane J. Wacasey
eSignature
Date: 6-Aug-2019

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition?
Custody seals intact on shipping container/cooler?
Custody seals intact on sample bottles?
VOA/TX1005/TX1006 Solids in hermetically sealed vials?
Chain of custody present?
Chain of custody signed when relinquished and received?
Samplers name present on COC?
Chain of custody agrees with sample labels?
Samples in proper container/bottle?
Sample containers intact?
Sufficient sample volume for indicated test?
All samples received within holding time?
Container/Temp Blank temperature in compliance?

- Yes/No checkboxes for each item in the list above.

2 Page(s)
COC IDs:196136, 196124

Temperature(s)/Thermometer(s): 0.3c/0.3c, 0.5c/0.5c, 0.6c/0.6c UC/C IR25

Cooler(s)/Kit(s): 45155, 444481, 7183

Date/Time sample(s) sent to storage: 08/02/2019 14:10

Water - VOA vials have zero headspace? Yes/No checkboxes and No VOA vials submitted checkbox.

Water - pH acceptable upon receipt? Yes/No checkboxes and N/A checkbox.

pH adjusted? Yes/No checkboxes and N/A checkbox.

pH adjusted by: [Empty text box]

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments: [Empty text box]

Corrective Action: [Empty text box]



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# Chain of Custody Form

Page 1 of 2

COC ID: 196136

HS19080113

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive	Address	1400 Douglas Street
	Suite 4004		Stop 0756
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwilc.com	e-Mail Address	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TBO-201801</del>			Water	1	2		X									
2	WG-1620-MW36D-20190731	7-31-19	0740			6	X		X	X							
3	WG-1620-MW38A-20190731		0835			6	X		X	X							
4	WG-1620-MW65D-20190731		0945			12	X		X	X	X						
5	WG-1620-MW66D-20190731		1045			6	X		X	X							
6	WG-1620-MW59D-20190731		1150			6	X		X	X							
7	WG-1620-FDDS-20190731		1150			6	X		X	X							
8	WG-1620-MW22AR-20190731		1305			6	X		X	X							
9	WG-1620-MW22BR-20190731		1350			6	X		X	X							
10	WG-1620-MW88C-20190731		1545			6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24-hour		Results Due Date:
Relinquished by: <b>John B</b>	Date: <b>8-2-19</b>	Time: <b>12:00</b>	Received by:	Notes: UPRR, HOLISTON MWPW	
Relinquished by:	Date: <b>8/2/19</b>	Time: <b>12:00</b>	Received by (Laboratory): <b>J. M. MURPHY</b>	Cooler ID	Cooler Temp.
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	<b>45155</b>	<b>0.3</b>
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				<b>44481</b>	<b>0.5</b>
				<b>7183</b>	<b>0.6</b>
				<b>1025</b>	<b>CF00</b>

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.



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# Chain of Custody Form

Page 2 of 2

COC ID: 196124

## HS19080113

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UFRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TV (5636002 5652646 Metals - As, Pb)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	ms/msD
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TBO-201801			Water	1	2		X									
2	WG-1620-MW67B-20190731	7-31-19	1645			12	X		X	X	X						
3	WG-1620-MW49A-20190731	↓	1745			6		X	X	X							
4	WG-1620-TB06-20190731	↓				2		X									
5																	
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour		Results Due Date:	
Relinquished by: <b>John K</b>	Date: <b>8-2-19</b>	Time: <b>12:00</b>	Received by:	Notes: <b>UPRR Houston MWPW</b>			
Relinquished by: <b>John K</b>	Date: <b>3/2/19</b>	Time: <b>12:00</b>	Received by (Laboratory): <b>J. Murray</b>	Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)	
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	<input type="checkbox"/> Level II Std QC	<input checked="" type="checkbox"/> TRRP Checklist		
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				<input type="checkbox"/> Level III Std QC/Rat Date	<input type="checkbox"/> TRRP Level IV		
				<input type="checkbox"/> Level IV SW/84/CLP			

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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**JANUARY-MARCH 2020**



# Memorandum

March 13, 2020

Revision: April 29, 2020

To: Eric Matzner Ref. No.: 11183954-1620

From: <sup>ck</sup> Chris G. Knight/eew/557-NF Tel: 512-506-8803

CC: Jesse Orth, Jon Lang; Julie Lidstone

**Subject: Data Usability Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

## 1. Scope of Data Usability Study

This document details a Data Usability Summary (DUS) of analytical results for groundwater samples collected in support of the HWPW - Site-Wide Monitoring at the Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works site during January-February 2020. Samples were submitted to ALS Environmental (ALS), located in Houston, Texas and are reported in data packages HS20010407, HS20010442, HS20010596, HS20010713, HS20010767, HS20010930, HS20010958, HS20011116, HS20011249, and HS20020456. The intended use of the data is to support the HWPW - Site-Wide Monitoring at the site by providing current concentration of chemicals of concern.

Data were reviewed and validated by Chris G. Knight of GHD, in accordance with Title 30 of the Texas Administrative Code Section 350.54 (30 TAC 350.54) as described in the Texas Commission on Environmental Quality (TCEQ) Regulatory Guidance document entitled "Review and Reporting of COC Concentration Data under TRRP", (RG-366/TRRP-13), revised May 2010, herein referred to as "TRRP-13 Guidance". Evaluation of the data was based on information obtained from the chain of custody forms, the finished report forms, method blank data, recovery data from surrogate spikes/laboratory control samples (LCS)/matrix spikes (MS), duplicate data, field quality assurance/quality control (QA/QC) samples, the laboratory review checklists (LRC), and the laboratory exception reports (ER).

A sample collection and analysis summary is presented in Table 1. This summary provides a cross-reference of field sample identification numbers and location identification. Each sample is assigned a unique field identification number.

The validated sample results are presented in Table 2. A summary of the analytical methodology is presented in Table 3.



## 2. Laboratory Qualifications

The Laboratory's quality assurance program is consistent with the quality standards outlined in the National Environmental Laboratory Accreditation Program (NELAP). This laboratory was accredited under Texas Certification number # TX104704231 at the time the analysis was performed and the certificate is included in Attachment A.

## 3. Project Objectives

### 3.1 Sampling/Analytical QA/QC Objectives

The QA/QC program was designed to identify contamination resulting from the sampling, sample transport and analytical process through the analysis of trip blank samples, field blank samples, field duplicate sample sets, and method blanks. The QA/QC program was designed to evaluate the quality of the resulting data with respect to bias and precision through analysis of LCS, MS, and duplicate analyses.

## 4. Data Review/Validation Results

### 4.1 Sample Holding Time and Preservation

Samples were shipped with a chain of custody and the paper work was filled out properly with the following exceptions:

- i) HS20010407 – The chain of custody listed nine sample containers submitted for all samples, however only six were received. No sample vials for total petroleum hydrocarbons (TPH) were received for WG1620-MW49A-20200107 and WG-1620-MW-80B-20200107. The laboratory logged in two sample vials for volatile organic compounds (VOCs) and one sample vial TPH analyses. The sample date differed on the chain of custody from the sample date listed on the sample container for sample WG-1620-FB03-20200109. The sample date listed on the sample container was used for login. No further actions were required.
- ii) HS20010442 - The sample time differed on the chain of custody from the sample time listed on the sample container for sample WG-1620 FB 0420200110. The sample time listed on the chain of custody was used for login. No further action was required.
- iii) HS20010713 - The sample time differed on the chain of custody from the sample time listed on the sample container for sample WG-1620-MW25C-20200115. The sample time listed on the chain of custody was used for login. No further action was required.
- iv) HS20010767 – The following samples were received but not listed on the chain of custody: MW48C (01/16/20 10:20), MW47C (01/16/20 11:05), and MW61A MS & MSD (01/16/20 12:00). The laboratory added these samples to the chain of custody. No further actions were required.
- v) HS20010958 – WG-1620-MW38B-20200121 was misidentified on the chain of custody as WG-1620-MW38C-20200121. The sample label was used for login. No further action was required.





All samples were properly preserved, delivered on ice, and stored by the laboratory at the required temperature (0-6°C).

- i) WG-1620-MW18C-20200108 was received in a properly preserved sample container for metals analysis. The included preservative was insufficient to reduce the pH to <2. Additional preservative was added by the laboratory bringing the pH to <2. No further action was required.

The sample chain of custody documents and the analytical report were used to determine sample holding times. All samples were prepared and analyzed within the required holding times.

#### **4.2 Sample Containers**

Sample containers used were certified pre-cleaned glass and plastic containers provided by the laboratory. These containers meet or exceed analyte specifications established in the United States Environmental Protection Agency (USEPA) *Specifications and Guidance for Contaminant-free Sample Containers*.

#### **4.3 Calibrations**

According to the LRC, initial calibration and continuing calibration data met the criteria for the selected method.

#### **4.4 Laboratory Method Blank Analyses**

Method blanks are prepared from a purified matrix and analyzed with investigative samples to determine the existence and magnitude of sample contamination introduced during the analytical procedures. As these were not discrete samples handled in the field, these blanks are not listed on the sample identification cross-reference list found in the data packages.

For this study, laboratory method blanks were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch and results are reported in the laboratory data packages.

The method blank results were non-detect or below the method quantitation limit (MQL), indicating that laboratory contamination was not a factor for this investigation.

#### **4.5 Internal Standard and Surrogate Spike Recoveries**

Recoveries of internal standards are addressed in the LRC of the data packages. All internal standard recoveries associated with the compounds of interest were acceptable per the LRC.

In accordance with the methods employed, all samples, blanks, and QC samples analyzed for VOCs, semi-volatile organic compounds (SVOCs), and TPH are spiked with surrogate compounds prior to sample analysis. Surrogate recoveries provide a means to evaluate the effects of laboratory performance on individual sample matrices. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project. Each individual surrogate compound is expected to meet the laboratory control limits. According to the TRRP-13 Guidelines, one outlying surrogate is acceptable for methods with



multiple surrogate spike compounds as long as the recovery is at least ten percent. Sample analyzed at elevated sample dilutions (five times or greater) were not assessed.

Surrogate recoveries were assessed against laboratory control limits and/or the guidance in TRRP-13. All surrogate recoveries met the above criteria.

#### **4.6 Laboratory Control Sample Analysis**

LCS or LCS/laboratory control sample duplicate (LCSD) are prepared and analyzed as samples to assess the analytical efficiencies of the methods employed, independent of sample matrix effects. The relative percent difference (RPD) of the LCS/LCSD recoveries is used to evaluate analytical precision. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project.

For this study, LCS or LCS/LCSD were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch.

The LCS or LCS/LCSD contained all analytes specified in the methods. All LCS recoveries and/or RPDs were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision (where applicable) with the following exceptions (see Table 4):

- i) One LCS was reported with an elevated recovery for vinyl chloride. The associated sample results was non-detect and not affected by the indicated high bias. No further action was required.
- ii) One LCS was reported with elevated recoveries for 2,4-dinitrotoluene and 2,6-dinitrotoluene. Associated non-detect sample results would not be affected by the indicated high bias. No further actions were required. One associated detected sample results was qualified as estimated; biased high.
- iii) One LCS/LCSD was reported with an elevated RPD for 2-methylnaphthalene. All associated sample results were qualified as estimated.

#### **4.7 Matrix Spike Analysis**

To evaluate the effects of sample matrices on the preparation process, measurement procedures, and accuracy of a particular analysis, samples are spiked with known concentrations of the analytes of interest and analyzed as MS/matrix spike duplicate (MSD) samples. The RPD between the MS and MSD is used to assess analytical precision.

MS/MSD analyses were performed as specified in Table 1. The recovery ranges established by the laboratory is adopted as the acceptance criteria for the project.

The MS/MSD samples were spiked with all analytes specified in the methods. All percent recoveries and the RPD value were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision with the following exceptions:

- i) One MS/MSD was reported with an elevated RPD for 4,6-dinitro-2-methylphenol. The associated sample result was non-detect and not affected by the indicated variability. No further action was required.



- ii) One MS/MSD was reported with elevated MS recoveries for 4,6-dinitro-2-methylphenol and di-n-butyl phthalate. If only the MS or MSD recovery was outside of control limits, no qualification of the data was performed based on the acceptable recovery of the companion spike and the acceptable RPD. No further action was required.
- iii) One MS/MSD was reported with elevated recoveries and/or elevated RPDs for the following SVOCs due to possible matrix interferences and were not assessed: 4-Nitrophenol, bis(2-ethylhexyl)phthalate, and di-n-butyl phthalate

The laboratory also performed additional MS/MSD on non-site samples. These cannot be used to assess accuracy and precision for the site samples.

#### **4.8 Duplicate Sample Analyses**

Analytical precision is evaluated based on the analysis of laboratory duplicate samples. For this study, duplicate samples were prepared and analyzed by the laboratory as specified in Table 1 for arsenic. The RPD established by the laboratory are adopted as the acceptance criteria for the project.

All duplicate analyses performed were acceptable, demonstrating acceptable analytical precision with the following exception (see Table 5):

- i) One duplicate analysis was reported with an elevated RPD for arsenic. All associated sample results were qualified as estimated.

The laboratory also performed additional duplicate analyses on non-site samples. These cannot be used to assess precision for the site samples.

#### **4.9 Field QA/QC Samples**

The field QA/QC consisted of ten trip blank samples, thirteen field blank samples, and five field duplicate sample sets.

##### ***Trip Blank Sample Analysis***

To evaluate contamination from sample collection, transportation, storage, and analytical activities, ten trip blank samples were submitted to the laboratory for VOCs analysis. All results were non-detect for the compounds of interest.

##### ***Field Blank Sample Analysis***

To assess ambient conditions at the site, thirteen field blank samples were submitted for analysis, as identified in Table 1. All results were non-detect for the compounds of interest with the following exceptions (see Table 6):

- i) The following field blanks submitted yielded low level detections for various SVOCs: WG-1620-FB01-20200107, WG-1620-FB02-20200108, WG-1620-FB03-20200109, WG-1620-FB04-20200110, WG-1620-FB05-20200113, WG-1620-FB08-20200116, WG-1620-FB09-20200117, WG-1620-FB10-20200120, WG-1620-FB11-20200121,



WG-1620-FB12-20200123, and WG-1620-FB14-20200210. Associated sample results that were significantly greater than the concentrations found in the field blanks or were non-detect were not impacted. No further action was required. Associated sample results with comparable concentrations to the field blank detections were qualified as non-detect.

### ***Field Duplicate Sample Analysis***

To assess the analytical and sampling protocol precision, five field duplicate sample sets were collected and submitted "blind" to the laboratory, as specified in Table 1. The RPDs associated with these duplicate samples must be less than thirty percent for water samples. The RPDs are only used when sample concentrations are above the estimated regions of detection.

Field duplicate summary data are presented in Table 2. All field duplicate results were within acceptable agreement, demonstrating acceptable sampling and analytical precision with the following exceptions (see Table 7):

- i) WG-1620-MW35B-20200110 and WG-1620-FD01-20200110 did show some variability in VOCs and SVOCs results and were qualified as estimated.
- ii) WG-1620-MW21C-20200114 and WG-1620-FD02-20200114 did show some variability in 2,6-dinitrotoluene results and were qualified as estimated.
- iii) WG-1620-MW33A-20200120 and WG-1620-FD05-20200120 did show some variability in acenaphthene, fluoranthene, and pyrene results and were qualified as estimated.

### **4.10 Field Procedures**

Golder Associates, Inc. collected groundwater samples in accordance with their Standard Operating Procedures (SOP) for sample collection.

### **4.11 Analyte Reporting**

The laboratory reported detected results for each analyte down to the sample detection limit (SDL), which is defined as the method detection limit (MDL) with sample-specific adjustments for dilutions, aliquot size, volumes, etc. Positive analyte detections less than the MQL but greater than the SDL were qualified as estimated (J) in Table 2 unless qualified otherwise in this memorandum.

- i) The following samples were analyzed at the lowest practical dilution for VOCs determinations due to sample matrix: WG-1620-MW18C-20200108, WG-1620-MW18A-20200108. No further actions were required.
- ii) The following samples were analyzed at the lowest practical dilution for VOCs and/or SVOCs determinations due to elevated concentration of non-target analytes: WG-1620-MW32B-20200120, WG-1620-MW23C-20200109, WG-1620-MW49A-20200107, WG-1620-MW57B-20200108, WG-1620-MW70B-20200120, and WG-1620-MW72B-20200109. No further actions were required.



- iii) The following samples were analyzed at the lowest practical dilution for VOCs determinations due to matrix interferences: WG-1620-MW57B-20200108, WG-1620-MW72B-20200109, WG-1620-MW23C-20200109. No further actions were required.

The detectability check standard (DCS) results supported the laboratory MDLs.

## 5. Conclusion

Based on the assessment detailed in the foregoing, the data summarized in Table 2 are usable for the purpose of supporting the HWPW - Site-Wide Monitoring at the site by providing current concentration of chemicals of concern with the specific qualifications noted herein.

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW49A-20200107	MW-49A	Water	01/07/2020	13:10	X	X	X	X	
WG-1620-MW80B-20200107	MW-80B	Water	01/07/2020	16:00	X	X	X	X	
WG-1620-FB01-20200107	-	Water	01/07/2020	16:30	X	X		X	Field Blank
WG-1620-FB02-20200108	-	Water	01/08/2020	09:00	X	X		X	Field Blank
WG-1620-MW18C-20200108	MW-18C	Water	01/08/2020	10:10	X	X		X	
WG-1620-MW18A-20200108	MW-18A	Water	01/08/2020	11:10	X	X		X	
WG-1620-MW58A-20200108	MW-58A	Water	01/08/2020	12:15	X	X		X	
WG-1620-MW57B-20200108	MW-57B	Water	01/08/2020	13:30	X	X		X	
WG-1620-MW57A-20200108	MW-57A	Water	01/08/2020	14:55	X	X		X	
WG-1620-FB03-20200109	-	Water	01/09/2020	07:30	X	X		X	Field Blank
WG-1620-MW72B-20200109	MW-72B	Water	01/09/2020	08:35	X	X		X	
WG-1620-MW23C-20200109	MW-23C	Water	01/09/2020	09:40	X	X		X	
WG-1620-MW19C-20200109	MW-19C	Water	01/09/2020	10:45	X	X		X	MS/MSD-P; DUP
WG-1620-MW53C-20200109	MW-53C	Water	01/09/2020	11:15	X	X		X	MS/MSD-P; DUP
WG-1620-MW51A-20200109	MW-51A	Water	01/09/2020	11:50	X	X	X	X	MS/MSD-P

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**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW54C-20200109	MW-54C	Water	01/09/2020	12:35	X	X		X	
WG-1620-MW51C-20200109	MW-51C	Water	01/09/2020	12:40	X	X	X	X	
WG-1620-MW44A-20200109	MW-44A	Water	01/09/2020	13:30	X	X		X	MS/MSD-P
WG-1620-MW85C-20200109	MW-85C	Water	01/09/2020	13:35	X	X	X	X	
WG-1620-MW77A-20200109	MW-77A	Water	01/09/2020	14:25	X	X	X	X	
WG-1620-MW36B-20200109	MW-36B	Water	01/09/2020	14:40	X	X		X	
WG-1620-MW76C-20200109	MW-76C	Water	01/09/2020	15:20	X	X	X	X	
WG-1620-MW36A-20200109	MW-36A	Water	01/09/2020	16:00	X	X		X	
WG-1620-MW81B-20200109	MW-81B	Water	01/09/2020	16:20	X	X	X	X	
WG-1620-MW50A-20200109	MW-50A	Water	01/09/2020	17:15	X	X	X	X	
WG-1620-TB01-20200109	-	Water	01/09/2020	-	X				Trip Blank
WG-1620-MW83C-20200110	MW-83C	Water	01/10/2020	09:00	X	X		X	
WG-1620-MW83B-20200110	MW-83B	Water	01/10/2020	10:00	X	X		X	
WG-1620-FB04-20200110	-	Water	01/10/2020	10:30	X	X		X	Field Blank
WG-1620-MW35A-20200110	MW-35A	Water	01/10/2020	11:20	X	X		X	

Table 1

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**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW35B-20200110	MW-35B	Water	01/10/2020	12:45	X	X		X	
WG-1620-FD01-20200110	MW-35B	Water	01/10/2020	12:45	X	X		X	Field duplicate of MW35B
WQ-1620-TB01-20200110	-	Water	01/10/2020	-	X				Trip Blank
WG-1620-MW40B-20200113	MW-40B	Water	01/13/2020	12:10	X	X		X	MS/MSD-P; DUP
WG-1620-MW42B-20200113	MW-42B	Water	01/13/2020	13:10	X	X		X	
WG-1620-TW41B-20200113	TW-41B	Water	01/13/2020	14:00	X	X		X	
WG-1620-MW05-20200113	MW-05	Water	01/13/2020	14:55	X	X		X	
WG-1620-MW12C-20200113	MW-12C	Water	01/13/2020	15:55	X	X		X	
WG-1620-MW12A-20200113	MW-12A	Water	01/13/2020	16:45	X	X		X	
WG-1620-MW39B-20200113	MW-39B	Water	01/13/2020	17:35	X	X		X	
WG-1620-FB05-20200113	-	Water	01/13/2020	17:50	X	X		X	Field Blank
WG-1620-MW13-20200114	MW-13	Water	01/14/2020	08:10	X	X		X	MS/MSD-P
WG-1620-MW09-20200114	MW-09	Water	01/14/2020	09:00	X	X		X	
WG-1620-MW21C-20200114	MW-21C	Water	01/14/2020	10:00	X	X		X	
WG-1620-FD02-20200114	MW-21C	Water	01/14/2020	10:00	X	X		X	Field duplicate of MW21C



Table 1

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**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-P11-20200114	P-11	Water	01/14/2020	10:55	X	X		X	
WG-1620-MW88C-20200114	MW-88C	Water	01/14/2020	11:50	X	X		X	
WG-1620-MW20A-20200114	MW-20A	Water	01/14/2020	12:55	X	X		X	
WG-1620-MW15A-20200114	MW-15A	Water	01/14/2020	13:45	X	X		X	
WG-1620-MW15B-20200114	MW-15B	Water	01/14/2020	14:30	X	X		X	
WG-1620-MW15C-20200114	MW-15C	Water	01/14/2020	15:20	X	X		X	
WG-1620-MW14-20200114	MW-14	Water	01/14/2020	16:15	X	X		X	
WG-1620-MW17-20200114	MW-17	Water	01/14/2020	17:10	X	X		X	
WG-1620-FB06-20200114	-	Water	01/14/2020	17:25	X	X		X	Field Blank
WG-1620-TB03-20200114	-	Water	01/14/2020	-	X				Trip Blank
WG-1620-MW17C-20200115	MW-17C	Water	01/15/2020	08:15	X	X		X	
WG-1620-MW67B-20200115	MW-67B	Water	01/15/2020	09:30	X	X		X	MS/MSD; DUP
WG-1620-MW46C-20200115	MW-46C	Water	01/15/2020	10:25	X	X		X	
WG-1620-MW25A-20200115	MW-25A	Water	01/15/2020	11:20	X	X		X	
WG-1620-MW44C-20200115	MW-44C	Water	01/15/2020	13:15	X	X		X	

Table 1

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**HWPW - Site-Wide Monitoring**  
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**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW71B-20200115	MW-71B	Water	01/15/2020	14:00	X	X		X	
WG-1620-MW33BR-20200115	MW-33BR	Water	01/15/2020	15:00	X	X		X	
WG-1620-MW25C-20200115	MW-25C	Water	01/15/2020	15:50	X	X		X	
WG-1620-MW27A-20200115	MW-27A	Water	01/15/2020	16:40	X	X		X	
WG-1620-MW27C-20200115	MW-27C	Water	01/15/2020	17:20	X	X		X	
WQ-1620-TB04-20200115	-	Water	01/15/2020	-	X				Trip Blank
WG-1620-MW59A-20200116	MW-59A	Water	01/16/2020	08:05	X	X		X	
WG-1620-MW59B-20200116	MW-59B	Water	01/16/2020	09:00	X	X		X	
WG-1620-MW28A-20200116	MW-28A	Water	01/16/2020	09:10	X	X		X	
WG-1620-MW28C-20200116	MW-28C	Water	01/16/2020	10:15	X	X		X	MS/MSD; DUP
WG-1620-MW48C-20200116	MW-48C	Water	01/16/2020	10:20	X	X		X	
WG-1620-MW47C-20200116	MW-47C	Water	01/16/2020	11:05	X	X		X	
WG-1620-MW61A-20200116	MW-61A	Water	01/16/2020	12:00	X	X		X	MS/MSD; DUP
WG-1620-FB08-20200116	-	Water	01/16/2020	12:00	X	X		X	Field Blank
WG-1620-MW89B-20200116	MW-89B	Water	01/16/2020	12:45	X	X		X	

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW84B-20200116	MW-84B	Water	01/16/2020	13:45	X	X		X	MS/MSD-P
WG-1620-MW59D-20200116	MW-59D	Water	01/16/2020	13:50	X	X		X	
WG-1620-FD03-20200116	MW-59D	Water	01/16/2020	13:50	X	X		X	Field duplicate of MW59D
WG-1620-MW63B-20200116	MW-63B	Water	01/16/2020	14:45	X	X		X	
WG-1620-MW66D-20200116	MW-66D	Water	01/16/2020	15:00	X	X		X	
WG-1620-MW26A-20200116	MW-26A	Water	01/16/2020	16:00	X	X		X	
WG-1620-MW65D-20200116	MW-65D	Water	01/16/2020	16:20	X	X		X	MS/MSD; DUP
WG-1620-MW36D-20200116	MW-36D	Water	01/16/2020	17:25	X	X		X	
WG-1620-MW86C-20200117	MW-86C	Water	01/17/2020	07:25	X	X	X	X	
WG-1620-FD04-20200117	MW-86C	Water	01/17/2020	07:25	X	X	X	X	Field duplicate of MW86C
WG-1620-MW75B-20200117	MW-75B	Water	01/17/2020	08:15	X	X	X	X	
WG-1620-MW68C-20200117	MW-68C	Water	01/17/2020	09:10	X	X		X	
WG-1620-MW79A-20200117	MW-79A	Water	01/17/2020	09:20	X	X	X	X	
WG-1620-FB09-20200117	-	Water	01/17/2020	10:00	X	X		X	Field Blank
WG-1620-MW68A-20200117	MW-68A	Water	01/17/2020	10:15	X	X		X	

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW74B-20200117	MW-74B	Water	01/17/2020	10:20	X	X	X	X	
WG-1620-MW78A-20200117	MW-78A	Water	01/17/2020	11:25	X	X	X	X	
WQ-1620-TB05-2020117	-	Water	01/17/2020	-	X				Trip Blank
WG-1620-MW32AR-20200120	MW-32AR	Water	01/20/2020	09:10	X	X		X	MS/MSD-P; DUP
WG-1620-FB10-20200120	-	Water	01/20/2020	09:30	X	X		X	Field Blank
WG-1620-MW32B-20200120	MW-32B	Water	01/20/2020	10:15	X	X		X	
WG-1620-MW70B-20200120	MW-70B	Water	01/20/2020	11:30	X	X		X	
WG-1620-MW33A-20200120	MW-33A	Water	01/20/2020	12:45	X	X		X	
WG-1620-FD05-20200120	MW-33A	Water	01/20/2020	12:45	X	X		X	Field duplicate of MW-33A
WG-1620-MW87C-20200120	MW-87C	Water	01/20/2020	14:00	X	X		X	
WG-1620-MW90B-20200120	MW-90B	Water	01/20/2020	15:15	X	X		X	
WG-1620-MW45C-20200120	MW-45C	Water	01/20/2020	16:30	X	X		X	
WQ-1620-TB06-20200120	-	Water	01/20/2020	-	X				Trip Blank
WG-1620-MW82B-20200121	MW-82B	Water	01/21/2020	09:10	X	X		X	MS/MSD; DUP
WG-1620-MW38A-20200121	MW-38A	Water	01/21/2020	11:10	X	X		X	

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW38B-20200121	MW-38B	Water	01/21/2020	13:45	X	X		X	
WG-1620-MW03-20200121	MW-03	Water	01/21/2020	15:00	X	X		X	
WG-1620-FB11-20200121	-	Water	01/21/2020	15:30	X	X		X	Field Blank
WG-1620-MW04-20200121	MW-04	Water	01/21/2020	16:00	X	X		X	
WQ-1620-TB07-20200121	-	Water	01/21/2020	-	X				Trip Blank
WG-1620-MW41B-20200123	MW-41B	Water	01/23/2020	12:40	X	X		X	MS/MSD-P; DUP
WG-1620-FB12-20200123	-	Water	01/23/2020	13:00	X	X		X	Field Blank
WG-1620-MW12B-20200123	MW-12B	Water	01/23/2020	14:00	X	X		X	
WG-1620-MW68B-20200123	MW-68B	Water	01/23/2020	15:25	X	X		X	
WQ-1620-TB08-20200123	-	Water	01/23/2020	-	X				Trip Blank
WG-1620-MW62B-20200127	MW-62B	Water	01/27/2020	09:25	X	X		X	
WG-1620-MW64A-20200127	MW-64A	Water	01/27/2020	11:15	X	X		X	
WG-1620-FB13-20200127	-	Water	01/27/2020	12:00	X	X		X	Field Blank
WG-1620-MW49B-20200127	MW-49B	Water	01/27/2020	13:30	X	X		X	
WQ-1620-TB09-20200127	-	Water	01/27/2020	-	X				Trip Blank

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW22AR-20200210	MW-22AR	Water	02/10/2020	10:00	X	X		X	
WG-1620-MW22BR-20200210	MW-22BR	Water	02/10/2020	11:00	X	X		X	
WG-1620-FB14-20200210	-	Water	02/10/2020	11:30	X	X		X	Field Blank
WQ-1620-TB11-20200211	-	Water	02/11/2020	-	X	X		X	Trip Blank

## Notes:

- VOCs - Volatile Organic Compounds  
SVOCs - Semi-volatile Organic Compounds  
TPH - Total Petroleum Hydrocarbons  
MS/MSD - Matrix Spike/ Matrix Spike Duplicate  
MS/MSD-P - Matrix Spike/ Matrix Spike Duplicate (partial parameters)  
DUP - Laboratory Duplicate  
"- " - Not Applicable

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-03	MW-04	MW-05	MW-09	MW-12A
Sample Name:	WG-1620-MW03-20200121	WG-1620-MW04-20200121	WG-1620-MW05-20200113	WG-1620-MW09-20200114	WG-1620-MW12A-20200113
Sample Date:	01/21/2020	01/21/2020	01/13/2020	01/14/2020	01/13/2020

Parameters	Unit	MW-03	MW-04	MW-05	MW-09	MW-12A
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.000045 J	0.000026	<0.000019	<0.000019	0.0023
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.000061 J	0.000015	<0.000027	<0.000027	0.28
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015	0.0012
Anthracene	mg/L	0.000069 J	0.000010	<0.000014	<0.000014	0.0097
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050	0.00015
Benzo(a)pyrene	mg/L	0.000024 J	0.000033 J	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00011 J	0.0018	<0.000037	<0.000037	0.00011 J
Chrysene	mg/L	<0.000021	0.000034 J	<0.000021	<0.000021	0.00015
Di-n-butylphthalate (DBP)	mg/L	0.000029 J	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.000097 J	<0.000020	<0.000020	0.23
Fluoranthene	mg/L	0.000055 J	0.000029	<0.000010	<0.000010	0.0067
Fluorene	mg/L	<0.000030	0.000067 J	<0.000030	<0.000030	0.28
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-03	MW-04	MW-05	MW-09	MW-12A
Sample Name:	WG-1620-MW03-20200121	WG-1620-MW04-20200121	WG-1620-MW05-20200113	WG-1620-MW09-20200114	WG-1620-MW12A-20200113
Sample Date:	01/21/2020	01/21/2020	01/13/2020	01/14/2020	01/13/2020

Parameters	Unit	MW-03	MW-04	MW-05	MW-09	MW-12A
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.00024	0.0013	<0.00012	0.00011	<0.00051
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	0.00017	<0.000021	<0.000021	0.082
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.000036 J	0.00020	<0.000019	<0.000019	0.0031
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00207	0.00223	0.00146 J	0.00430	0.00134 J



Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-12B	MW-12C	MW-13	MW-14	MW-15A
Sample Name:	WG-1620-MW12B-20200123	WG-1620-MW12C-20200113	WG-1620-MW13-20200114	WG-1620-MW14-20200114	WG-1620-MW15A-20200114
Sample Date:	01/23/2020	01/13/2020	01/14/2020	01/14/2020	01/14/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.0050	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.030	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.0075	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.014 J	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.025	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.0050	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.0050	--	--	--
Xylenes (total)	mg/L	<0.0075	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.0021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.0040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.0058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.0042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.0021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	320 J	0.00018	0.000054 J	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.0020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.0047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	350	0.00015	0.00022	0.000091 J
Acenaphthylene	mg/L	5.7	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	160	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	79	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	22	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.0030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.0037	<0.000037	<0.000037	<0.000037
Chrysene	mg/L	64	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.0020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	320	0.00012	0.00016	<0.000020
Fluoranthene	mg/L	520	0.000091 J	<0.000010	0.000064 J
Fluorene	mg/L	420	0.00019	0.00016	0.000057 J
N-Nitrosodiphenylamine	mg/L	<0.0025	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-12B	MW-12C	MW-13	MW-14	MW-15A
	Sample Name:	WG-1620-MW12B-20200123	WG-1620-MW12C-20200113	WG-1620-MW13-20200114	WG-1620-MW14-20200114	WG-1620-MW15A-20200114
	Sample Date:	01/23/2020	01/13/2020	01/14/2020	01/14/2020	01/14/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	760	0.0017	0.00026	<0.000020	0.00022
Nitrobenzene	mg/L	<0.0024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.0079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	1200	0.00015	<0.000021	<0.000021	0.025
Phenol	mg/L	<0.0035	0.00012 J	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	330	0.000068 J	<0.000019	0.000060 J	0.0013
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.0491 J	0.00195 J	0.0642	0.00185 J	0.0441

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-15B	MW-15C	MW-17	MW-17C	MW-18A	
Sample Name:	WG-1620-MW15B-20200114	WG-1620-MW15C-20200114	WG-1620-MW17-20200114	WG-1620-MW17C-20200115	WG-1620-MW18A-20200108	
Sample Date:	01/14/2020	01/14/2020	01/14/2020	01/15/2020	01/08/2020	
Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.0050
Benzene	mg/L	<0.00020	0.00059 J	<0.00020	<0.00020	0.41
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.0075
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	0.32
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010	<0.025
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	0.11
Vinyl chloride	mg/L	--	--	--	--	<0.0050
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	0.50
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040	1.8
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	<0.000019	0.000073 J	0.36
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047	0.0042
Acenaphthene	mg/L	0.000073 J	0.016	<0.000027	0.00023	0.33
Acenaphthylene	mg/L	<0.000015	0.00089	<0.000015	<0.000015	0.0081
Anthracene	mg/L	0.00011	0.00057	<0.000014	<0.000014	0.0082
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000055 J	<0.000037	<0.000037	<0.000037	<0.00043
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.0045	<0.000020	0.000065 J	0.20
Fluoranthene	mg/L	0.00011	0.00066	0.00011	0.00034	0.0023
Fluorene	mg/L	0.000062 J	0.00051	<0.000030	0.000061 J	0.16
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-15B	MW-15C	MW-17	MW-17C	MW-18A
Sample Name:	WG-1620-MW15B-20200114	WG-1620-MW15C-20200114	WG-1620-MW17-20200114	WG-1620-MW17C-20200115	WG-1620-MW18A-20200108
Sample Date:	01/14/2020	01/14/2020	01/14/2020	01/15/2020	01/08/2020

Parameters	Unit	MW-15B	MW-15C	MW-17	MW-17C	MW-18A
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.000020	0.00024	<0.000020	0.00016	4.6
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.000050 J	0.000080 J	<0.000021	<0.000021	0.10
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	0.0014
Pyrene	mg/L	0.000060 J	0.00038	0.000066 J	0.00021	0.0013
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00343	0.000773 J	0.00154 J	0.00184 J	0.0236

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-18C	MW-19C	MW-20A	MW-21C	MW-21C
	Sample Name:	WG-1620-MW18C-20200108	WG-1620-MW19C-20200109	WG-1620-MW20A-20200114	WG-1620-MW21C-20200114	WG-1620-FD02-20200114
	Sample Date:	01/08/2020	01/09/2020	01/14/2020	01/14/2020	01/14/2020 Duplicate
Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.0020	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	1.1	0.00032 J	0.020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.0030	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.31	0.00064 J	0.025	<0.00030	<0.00030
Methylene chloride	mg/L	<0.010	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.97	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.0020	<0.00020	--	--	--
Xylenes (total)	mg/L	0.92	0.0012	0.024	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	0.000078 J	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.023	<0.000040	0.0018	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042 J	0.0030 J
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.31	<0.000019	0.075	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.093	0.00099	0.11	<0.000027	0.00020
Acenaphthylene	mg/L	0.0026	0.000050 J	0.00092	<0.000015	<0.000015
Anthracene	mg/L	0.020	<0.00017	0.0075	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	0.0025	0.00026	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	0.00053	0.00010	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000071	<0.000037	0.00035	0.000065 J
Chrysene	mg/L	0.0027	0.00024	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	0.000021 J	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.085	0.00069	0.079	<0.000020	0.00017
Fluoranthene	mg/L	0.013	0.0011	0.00067	<0.000010	<0.000010
Fluorene	mg/L	0.043	0.00047	0.080	<0.000030	0.00018
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-18C	MW-19C	MW-20A	MW-21C	MW-21C
	Sample Name:	WG-1620-MW18C-20200108	WG-1620-MW19C-20200109	WG-1620-MW20A-20200114	WG-1620-MW21C-20200114	WG-1620-FD02-20200114
	Sample Date:	01/08/2020	01/09/2020	01/14/2020	01/14/2020	01/14/2020 Duplicate
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	14	<0.00011	2.0	0.00011	0.00013
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	0.023	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.065	<0.00045	0.028	<0.000021	0.000061 J
Phenol	mg/L	0.0062	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.0066	0.0014	0.00030	<0.000019	<0.000019
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00283	0.00140 J	0.00808	0.00109 J	0.00109 J

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-22AR	MW-22BR	MW-23C	MW-25A	MW-25C
Sample Name:	WG-1620-MW22AR-20200210	WG-1620-MW22BR-20200210	WG-1620-MW23C-20200109	WG-1620-MW25A-20200115	WG-1620-MW25C-20200115
Sample Date:	02/10/2020	02/10/2020	01/09/2020	01/15/2020	01/15/2020

Parameters	Unit	MW-22AR	MW-22BR	MW-23C	MW-25A	MW-25C
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.0010	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	0.0027 J	<0.00020	0.0017
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.0015	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	0.028	<0.00030	0.033
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0050	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.0010	<0.00020	0.010
Vinyl chloride	mg/L	--	--	--	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	<0.00030	0.025	<0.00030	0.22
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.00021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.047	<0.000040	0.0062
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.00058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.00042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	0.00011 J	<0.00021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.00012	<0.000043	0.12	<0.000019	0.96
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.00020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.00047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.00015	0.018	0.21	<0.000027	0.36
Acenaphthylene	mg/L	<0.000015	0.00018	0.0029	<0.000015	0.0025
Anthracene	mg/L	0.00020	0.00042	0.035	0.000089 J	0.027
Benzo(a)anthracene	mg/L	0.00029	<0.000050	0.0080	0.00012	0.00057
Benzo(a)pyrene	mg/L	0.00010	<0.000020	0.0022	0.000038 J	0.00019
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.00030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00011 J	0.000075 J	<0.00037	<0.000037	0.00066
Chrysene	mg/L	0.00032	<0.000021	0.0072	0.000097 J	0.00056
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.00020	<0.000020	0.00068
Dibenzofuran	mg/L	<0.00015	<0.00029	0.17	<0.000020	0.34
Fluoranthene	mg/L	0.0015	0.0017	0.068	0.00045	0.0058
Fluorene	mg/L	0.00019	0.0026	0.12	<0.000030	0.15
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.00025	<0.000025	<0.000025

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-22AR	MW-22BR	MW-23C	MW-25A	MW-25C
Sample Name:	WG-1620-MW22AR-20200210	WG-1620-MW22BR-20200210	WG-1620-MW23C-20200109	WG-1620-MW25A-20200115	WG-1620-MW25C-20200115
Sample Date:	02/10/2020	02/10/2020	01/09/2020	01/15/2020	01/15/2020

Parameters	Unit	MW-22AR	MW-22BR	MW-23C	MW-25A	MW-25C
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.00044	<0.00013	0.66	0.00010	7.5
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.00024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.00079	<0.000079	0.17
Phenanthrene	mg/L	0.0015	<0.00013	0.25	0.00017	0.18
Phenol	mg/L	<0.000041	<0.000035	0.0079	<0.000035	<0.000035
Pyrene	mg/L	0.0010	0.00082	0.043	0.00035	0.0049
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00352	0.0160	0.00333	0.00216	0.00391



Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-26A	MW-27A	MW-27C	MW-28A	MW-28C
Sample Name:	WG-1620-MW26A-20200116	WG-1620-MW27A-20200115	WG-1620-MW27C-20200115	WG-1620-MW28A-20200116	WG-1620-MW28C-20200116
Sample Date:	01/16/2020	01/15/2020	01/15/2020	01/16/2020	01/16/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	0.000034 J
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000042	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.00017	<0.000027	<0.000027	<0.00014
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000087	<0.000014	<0.000014	<0.00013
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	0.000073 J
Benzo(a)pyrene	mg/L	<0.000020	0.000052 J	<0.000020	0.000057 J
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.000062 J	<0.000037	0.000090 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	0.000065 J
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	0.000032 J
Dibenzofuran	mg/L	<0.000044	<0.000020	<0.000020	<0.00010
Fluoranthene	mg/L	0.00073	<0.000010	<0.000010	0.00030
Fluorene	mg/L	<0.00029	<0.000030	<0.000030	<0.00014
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-26A	MW-27A	MW-27C	MW-28A	MW-28C
	Sample Name:	WG-1620-MW26A-20200116	WG-1620-MW27A-20200115	WG-1620-MW27C-20200115	WG-1620-MW28A-20200116	WG-1620-MW28C-20200116
	Sample Date:	01/16/2020	01/15/2020	01/15/2020	01/16/2020	01/16/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.00017	0.000053 J	0.000055 J	<0.00060	<0.00016
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000030	<0.000021	<0.000021	0.00048	<0.000021
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	0.000086 J
Pyrene	mg/L	0.00035	<0.000019	<0.000019	0.00021	<0.000019
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.0217	0.000859 J	0.000623 J	0.00664	0.000937 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-32AR	MW-32B	MW-33A	MW-33A	MW-33BR
Sample Name:	WG-1620-MW32AR-20200120	WG-1620-MW32B-20200120	WG-1620-MW33A-20200120	WG-1620-FD05-20200120	WG-1620-MW33BR-20200115
Sample Date:	01/20/2020	01/20/2020	01/20/2020	01/20/2020 Duplicate	01/15/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.0050	<0.00020	<0.00020
Benzene	mg/L	<0.00020	1.8	<0.00020	0.12
Chlorobenzene	mg/L	<0.00030	<0.0075	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	0.73	<0.00030	0.043
Methylene chloride	mg/L	<0.0010	<0.025	<0.0010	<0.0010
Toluene	mg/L	<0.00020	2.3	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	<0.0050	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	2.1	<0.00030	0.011
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.00042	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	26	<0.000040	0.00018 J
2,4-Dinitrotoluene	mg/L	<0.000058	<0.0012	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.00084	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.00042	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000028	50	<0.000063	0.00025
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.00040	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.00094	<0.000047	<0.000047
Acenaphthene	mg/L	0.0014	26	0.00073 J	0.0051
Acenaphthylene	mg/L	<0.000015	<0.00030	0.000049 J	0.00015
Anthracene	mg/L	<0.000014	23	0.000057 J	0.00046
Benzo(a)anthracene	mg/L	<0.000050	2.2	0.00012	0.00010
Benzo(a)pyrene	mg/L	0.000081 J	0.74	0.000066 J	0.000039 J
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.00060	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000043 J	<0.00074	<0.000037	0.00067
Chrysene	mg/L	0.000042 J	2.4	0.000078 J	0.00011
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.00040	<0.000020	0.00099
Dibenzofuran	mg/L	<0.000020	28	0.000096 J	0.0039
Fluoranthene	mg/L	<0.00029	21	0.00035 J	0.00082
Fluorene	mg/L	0.000038 J	23	0.000059 J	0.0018
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.00050	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-32AR	MW-32B	MW-33A	MW-33A	MW-33BR
	Sample Name:	WG-1620-MW32AR-20200120	WG-1620-MW32B-20200120	WG-1620-MW33A-20200120	WG-1620-FD05-20200120	WG-1620-MW33BR-20200115
	Sample Date:	01/20/2020	01/20/2020	01/20/2020	01/20/2020 Duplicate	01/15/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.00014	300	<0.00050	<0.00066	0.0051
Nitrobenzene	mg/L	<0.000024	<0.00048	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.0016	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000032	69	<0.000056	<0.000046	0.0013
Phenol	mg/L	<0.000035	17	<0.000035	<0.000035	0.00018 J
Pyrene	mg/L	0.00042	14	<0.00019 J	0.00052 J	0.00057
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00770	0.00193 J	0.00810	0.00755	0.000877 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-35A	MW-35B	MW-35B	MW-36A	MW-36B
Sample Name:	WG-1620-MW35A-20200110	WG-1620-MW35B-20200110	WG-1620-FD01-20200110	WG-1620-MW36A-20200109	WG-1620-MW36B-20200109
Sample Date:	01/10/2020	01/10/2020	01/10/2020 Duplicate	01/09/2020	01/09/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	0.025 J	0.016 J	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	0.096	0.083	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	0.058 J	0.041 J	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	0.00021	<0.000052
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.14 J	0.062 J	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.059 J	0.033 J	<0.000027
Acenaphthylene	mg/L	<0.000015	0.00038	0.00027	<0.000015
Anthracene	mg/L	<0.000014	0.0029 J	0.0018 J	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000061 J	<0.000037	<0.000037	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.054 J	0.031 J	<0.000020
Fluoranthene	mg/L	<0.000010	0.0015 J	0.00093 J	<0.000010
Fluorene	mg/L	0.000065 J	0.028 J	0.015 J	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-35A	MW-35B	MW-35B	MW-36A	MW-36B
	Sample Name:	WG-1620-MW35A-20200110	WG-1620-MW35B-20200110	WG-1620-FD01-20200110	WG-1620-MW36A-20200109	WG-1620-MW36B-20200109
	Sample Date:	01/10/2020	01/10/2020	01/10/2020 Duplicate	01/09/2020	01/09/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.00020	2.0 J	0.13 J	<0.00034	<0.00093
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	0.024 J	0.013 J	<0.000021	<0.00010
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	<0.000075
Pyrene	mg/L	<0.000019	0.00070 J	0.00045 J	<0.000019	<0.000019
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.0219	0.0120	0.0159	0.00354	0.00100 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

	Location ID:	MW-36D	MW-38A	MW-38B	MW-39B	MW-40B
	Sample Name:	WG-1620-MW36D-20200116	WG-1620-MW38A-20200121	WG-1620-MW38B-20200121	WG-1620-MW39B-20200113	WG-1620-MW40B-20200113
	Sample Date:	01/16/2020	01/21/2020	01/21/2020	01/13/2020	01/13/2020
Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	0.011
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	0.089
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020	0.022
Vinyl chloride	mg/L	--	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030	0.16
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	0.00025	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040	0.0013
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.00017	0.00017	0.00021	0.34
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.00068	0.054	0.0012	0.39
Acenaphthylene	mg/L	<0.000015	<0.000015	0.00045	<0.000015	0.0021
Anthracene	mg/L	<0.000033	0.00015	0.0012	0.00016	0.040
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	0.000055 J	<0.000050	0.000094 J
Benzo(a)pyrene	mg/L	<0.000020	0.000032 J	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.000063 J	<0.000037	<0.000037	<0.000037
Chrysene	mg/L	<0.000021	0.000060 J	0.000057 J	<0.000021	0.000051 J
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	0.000036 J	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.00036	0.0054	0.00019	0.29
Fluoranthene	mg/L	<0.000010	0.00018	0.0025	0.00023	0.017
Fluorene	mg/L	<0.000030	0.00029	0.015	0.00032	0.30
N-Nitrosodiphenylamine	mg/L	<0.000025	0.00011 J	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-36D	MW-38A	MW-38B	MW-39B	MW-40B
	Sample Name:	WG-1620-MW36D-20200116	WG-1620-MW38A-20200121	WG-1620-MW38B-20200121	WG-1620-MW39B-20200113	WG-1620-MW40B-20200113
	Sample Date:	01/16/2020	01/21/2020	01/21/2020	01/13/2020	01/13/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.000033	0.0011	0.00098	0.0023	7.9
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000047	0.00039	0.00095	0.00019	0.22
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	0.00014	0.0015	0.00022	0.0037
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.000533 J	0.0177	0.0173	0.00248	0.0523



Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-41B	MW-42B	MW-44A	MW-44C	MW-45C
Sample Name:	WG-1620-MW41B-20200123	WG-1620-MW42B-20200113	WG-1620-MW44A-20200109	WG-1620-MW44C-20200115	WG-1620-MW45C-20200120
Sample Date:	01/23/2020	01/13/2020	01/09/2020	01/15/2020	01/20/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.012	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.066	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.087	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	--	<0.00020
Xylenes (total)	mg/L	0.16	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.043	<0.000040	<0.000042	0.00011 J
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.67 J	0.000069 J	<0.000035	0.00012
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.49	0.00012	0.064	0.00045
Acenaphthylene	mg/L	0.0089	<0.000015	0.00097	0.00065
Anthracene	mg/L	0.045	0.000049 J	0.00027	0.00086
Benzo(a)anthracene	mg/L	0.0040	<0.000050	0.000054 J	0.0018
Benzo(a)pyrene	mg/L	0.00090	<0.000020	<0.000020	0.00065
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00032	0.00013 J	<0.000037	0.0012
Chrysene	mg/L	0.0033	<0.000021	<0.000021	0.0022
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	0.0016
Dibenzofuran	mg/L	0.39	0.000065 J	<0.000033	0.00034
Fluoranthene	mg/L	0.052	0.00018	0.0059	0.0095
Fluorene	mg/L	0.35	0.000067 J	0.0078	0.00067
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-41B	MW-42B	MW-44A	MW-44C	MW-45C
Sample Name:	WG-1620-MW41B-20200123	WG-1620-MW42B-20200113	WG-1620-MW44A-20200109	WG-1620-MW44C-20200115	WG-1620-MW45C-20200120
Sample Date:	01/23/2020	01/13/2020	01/09/2020	01/15/2020	01/20/2020

Parameters	Unit	MW-41B	MW-42B	MW-44A	MW-44C	MW-45C
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	10	<0.00076	<0.0030	0.000097 J	<0.00061
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.41	0.00012	<0.00022	0.0077	<0.00027
Phenol	mg/L	<0.000035	<0.000035	<0.00013	<0.000035	<0.000035
Pyrene	mg/L	0.029	0.00013	0.0045	0.0060	0.00068
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.0842 J	0.00133 J	0.00966	0.00314	0.000730 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-46C	MW-47C	MW-48C	MW-49A	MW-49B
Sample Name:	WG-1620-MW46C-20200115	WG-1620-MW47C-20200116	WG-1620-MW48C-20200116	WG-1620-MW49A-20200107	WG-1620-MW49B-20200127
Sample Date:	01/15/2020	01/16/2020	01/16/2020	01/07/2020	01/27/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.010
Benzene	mg/L	0.028	<0.00020	<0.00020	0.60
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	0.0063 J
Ethylbenzene	mg/L	0.051	<0.00030	<0.00030	0.10
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.010
Toluene	mg/L	0.0047	<0.00020	<0.00020	0.19
Vinyl chloride	mg/L	--	--	--	<0.0020
Xylenes (total)	mg/L	0.11	<0.00030	<0.00030	0.24
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.0013
2,4-Dimethylphenol	mg/L	0.00016 J	0.00020	<0.000040	5.7
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.0035
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.0025
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.0013
2-Methylnaphthalene	mg/L	0.20	<0.00010	<0.000076	0.13
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.0012
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.0028
Acenaphthene	mg/L	0.16	<0.000057	<0.000073	0.057
Acenaphthylene	mg/L	0.0016	<0.000015	<0.000015	0.00097
Anthracene	mg/L	0.073	<0.000014	<0.000014	0.0028
Benzo(a)anthracene	mg/L	0.0071	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	0.0021	0.000040 J	0.000021 J	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.0011	0.00079	0.000039 J	<0.00011
Chrysene	mg/L	0.0068	0.000060 J	0.000029 J	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	0.0014	<0.000020	<0.000020
Dibenzofuran	mg/L	0.16	<0.000049	<0.000054	0.037
Fluoranthene	mg/L	0.084	<0.00011	<0.000061	0.00056
Fluorene	mg/L	0.12	<0.000042	<0.000050	0.027
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.0015

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-46C	MW-47C	MW-48C	MW-49A	MW-49B
Sample Name:	WG-1620-MW46C-20200115	WG-1620-MW47C-20200116	WG-1620-MW48C-20200116	WG-1620-MW49A-20200107	WG-1620-MW49B-20200127
Sample Date:	01/15/2020	01/16/2020	01/16/2020	01/07/2020	01/27/2020

Parameters	Unit	MW-46C	MW-47C	MW-48C	MW-49A	MW-49B
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	1.8	<0.00056	<0.00052	4.5	1200
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.00024	<0.0014
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.0047
Phenanthrene	mg/L	0.24	<0.000088	<0.000052	0.016	500
Phenol	mg/L	<0.000035	<0.000035	<0.000035	0.17	<0.0021
Pyrene	mg/L	0.053	<0.00010	<0.000065	0.00052	96
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	32.0	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	11	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	<0.19	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	21	--
<b>Metals</b>						
Arsenic	mg/L	0.00272	0.00234	0.00126 J	0.00463	0.0107

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-50A	MW-51A	MW-51C	MW-53C	MW-54C
Sample Name:	WG-1620-MW50A-20200109	WG-1620-MW51A-20200109	WG-1620-MW51C-20200109	WG-1620-MW53C-20200109	WG-1620-MW54C-20200109
Sample Date:	01/09/2020	01/09/2020	01/09/2020	01/09/2020	01/09/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.00071	<0.000040	<0.000040	<0.0011
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.00019	<0.000044	<0.000055	<0.00025
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.00037	<0.000040	<0.000094	<0.00032
Acenaphthylene	mg/L	0.000022 J	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000063	<0.00010	<0.000028
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	0.000027 J	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000085	<0.00014	<0.000085	<0.000037
Chrysene	mg/L	<0.000021	0.000050 J	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	0.000034 J	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.00013	<0.000049	<0.00011	<0.00027
Fluoranthene	mg/L	0.000052 J	0.00013	0.00018	0.000071 J
Fluorene	mg/L	<0.00011	<0.000043	<0.00011	<0.00020
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	0.000056 J

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-50A	MW-51A	MW-51C	MW-53C	MW-54C
	Sample Name:	WG-1620-MW50A-20200109	WG-1620-MW51A-20200109	WG-1620-MW51C-20200109	WG-1620-MW53C-20200109	WG-1620-MW54C-20200109
	Sample Date:	01/09/2020	01/09/2020	01/09/2020	01/09/2020	01/09/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.0020	<0.00032	<0.00030	<0.0012	<0.035
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000070	<0.00021	<0.00056	<0.00025	0.0056
Phenol	mg/L	<0.000066	<0.000035	<0.000035	0.00050	<0.000064
Pyrene	mg/L	<0.000034	<0.000083	<0.00011	<0.000061	0.0014
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	<0.19	<0.19	<0.19	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	<0.19	<0.19	<0.19	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	<0.19	<0.19	<0.19	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	<0.19	<0.19	<0.19	--	--
<b>Metals</b>						
Arsenic	mg/L	0.000718 J	0.00188 J	0.000452 J	0.000728 J	0.00122 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-57A	MW-57B	MW-58A	MW-59A	MW-59B
Sample Name:	WG-1620-MW57A-20200108	WG-1620-MW57B-20200108	WG-1620-MW58A-20200108	WG-1620-MW59A-20200116	WG-1620-MW59B-20200116
Sample Date:	01/08/2020	01/08/2020	01/08/2020	01/16/2020	01/16/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.0020	<0.00020	<0.00020
Benzene	mg/L	0.038	0.52	0.0053	<0.00020
Chlorobenzene	mg/L	0.00046 J	<0.0030	0.00072 J	<0.00030
Ethylbenzene	mg/L	0.015	0.25	0.0098	<0.00030
Methylene chloride	mg/L	<0.0010	<0.010	<0.0010	<0.0010
Toluene	mg/L	0.010	0.64	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	<0.00020	<0.00020
Xylenes (total)	mg/L	0.026	0.70	0.028	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.00021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.41	2.8	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.00058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.00042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.00021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.18	1.1	0.080	<0.00016
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.00020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.00047	<0.000047	<0.000047
Acenaphthene	mg/L	0.19	0.55	0.27	<0.00029
Acenaphthylene	mg/L	0.0025	0.0086	0.0015	<0.000015
Anthracene	mg/L	0.048	0.19	0.018	<0.000027
Benzo(a)anthracene	mg/L	0.0036	0.040	0.000091 J	<0.000050
Benzo(a)pyrene	mg/L	0.0011	0.012	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.00030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.00037	<0.00037	<0.00011	<0.000037
Chrysene	mg/L	0.0030	0.042	0.000071 J	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.00020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.15	0.49	0.13	<0.00016
Fluoranthene	mg/L	0.038	0.34	0.0099	<0.000032
Fluorene	mg/L	0.13	0.46	0.19	<0.00011
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.00025	<0.000025	<0.000025

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-57A	MW-57B	MW-58A	MW-59A	MW-59B
Sample Name:	WG-1620-MW57A-20200108	WG-1620-MW57B-20200108	WG-1620-MW58A-20200108	WG-1620-MW59A-20200116	WG-1620-MW59B-20200116
Sample Date:	01/08/2020	01/08/2020	01/08/2020	01/16/2020	01/16/2020

Parameters	Unit	MW-57A	MW-57B	MW-58A	MW-59A	MW-59B
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	1.0	13	1.2	<0.0012	<0.00024
Nitrobenzene	mg/L	<0.000024	<0.00024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.00079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.17	1.1	0.062	<0.00012	<0.000099
Phenol	mg/L	0.0049	0.32	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.025	0.21	0.0050	<0.000023	<0.000021
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.0545	0.00404	0.000906 J	0.00368	0.000486 J



Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-59D	MW-59D	MW-61A	MW-62B	MW-63B
Sample Name:	WG-1620-MW59D-20200116	WG-1620-FD03-20200116	WG-1620-MW61A-20200116	WG-1620-MW62B-20200127	WG-1620-MW63B-20200116
Sample Date:	01/16/2020	01/16/2020 Duplicate	01/16/2020	01/27/2020	01/16/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	0.018
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	0.051
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	<0.00020	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	0.013
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	0.00016 J
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	<0.000019	0.00011
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000036	<0.000027	<0.000052	0.040
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	0.00029
Anthracene	mg/L	<0.000014	<0.000014	<0.000014	0.00096
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	0.000081 J	0.000095 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	0.000027 J
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	0.00014 J
Dibenzofuran	mg/L	<0.000028	<0.000020	<0.000020	0.0042
Fluoranthene	mg/L	<0.000010	<0.000010	<0.000060	0.0020
Fluorene	mg/L	<0.000031	<0.000030	<0.000030	0.011
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-59D	MW-59D	MW-61A	MW-62B	MW-63B
	Sample Name:	WG-1620-MW59D-20200116	WG-1620-FD03-20200116	WG-1620-MW61A-20200116	WG-1620-MW62B-20200127	WG-1620-MW63B-20200116
	Sample Date:	01/16/2020	01/16/2020 Duplicate	01/16/2020	01/27/2020	01/16/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	<0.00013	<0.000020	<0.00021	0.0010	<0.00035
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000043	<0.000021	<0.00010	0.00087	<0.000072
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000019	<0.000059	0.0013	<0.000021
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.000508 J	0.000478 J	0.00107 J	0.0161	0.00204

Table 2

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-64A	MW-65D	MW-66D	MW-67B	MW-68A
Sample Name:	WG-1620-MW64A-20200127	WG-1620-MW65D-20200116	WG-1620-MW66D-20200116	WG-1620-MW67B-20200115	WG-1620-MW68A-20200117
Sample Date:	01/27/2020	01/16/2020	01/16/2020	01/15/2020	01/17/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.00037	<0.000019	<0.000095	0.000048 J
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.00016	<0.000027	<0.000027	0.00030
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	0.000054 J
Anthracene	mg/L	0.000096 J	<0.000014	<0.000014	0.000059 J
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00031	0.00026	0.000046 J	0.000090 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.000079 J	<0.000020	<0.000020	0.000034 J
Dibenzofuran	mg/L	0.000100 J	<0.000020	<0.000020	<0.000020
Fluoranthene	mg/L	0.00028	<0.000010	<0.000010	0.00013
Fluorene	mg/L	0.000082 J	<0.000030	<0.000030	0.000069 J
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

	Location ID:	MW-64A	MW-65D	MW-66D	MW-67B	MW-68A
	Sample Name:	WG-1620-MW64A-20200127	WG-1620-MW65D-20200116	WG-1620-MW66D-20200116	WG-1620-MW67B-20200115	WG-1620-MW68A-20200117
	Sample Date:	01/27/2020	01/16/2020	01/16/2020	01/15/2020	01/17/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	0.0020	<0.000020	<0.00013	<0.000020	<0.00022
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021	0.00016
Phenol	mg/L	<0.000035	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.00015	<0.000019	<0.000019	<0.000019	0.000068 J
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00126 J	0.000507 J	0.00138 J	0.000467 J	0.0423

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-68B	MW-68C	MW-70B	MW-71B	MW-72B	
Sample Name:	WG-1620-MW68B-20200123	WG-1620-MW68C-20200117	WG-1620-MW70B-20200120	WG-1620-MW71B-20200115	WG-1620-MW72B-20200109	
Sample Date:	01/23/2020	01/17/2020	01/20/2020	01/15/2020	01/09/2020	
Parameters	Unit					
<b>Volatile Organic Compounds</b>						
1,2-Dichloroethane	mg/L	<0.0050	<0.00020	<0.010	<0.00020	<0.0020
Benzene	mg/L	1.4	0.00056 J	1.9	0.0021	0.70
Chlorobenzene	mg/L	<0.0075	<0.00030	<0.015	<0.00030	<0.0030
Ethylbenzene	mg/L	0.39	0.00047 J	0.57	0.0013	0.13
Methylene chloride	mg/L	<0.025	<0.0010	<0.050	<0.0010	<0.010
Toluene	mg/L	0.23	<0.00020	2.1	<0.00020	0.58
Vinyl chloride	mg/L	<0.0050	--	<0.010	--	--
Xylenes (total)	mg/L	1.1	0.00062 J	1.5	0.0030	0.36
<b>Semi-volatile Organic Compounds</b>						
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.00042	<0.000021	<0.00021
2,4-Dimethylphenol	mg/L	0.075	<0.000040	49	<0.000040	13
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.0012	<0.000058	<0.00058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.00084	<0.000042	<0.00042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.00042	<0.000021	<0.00021
2-Methylnaphthalene	mg/L	1.1 J	0.00033	10	0.0060	0.22
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.00040	<0.000020	<0.00020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.00094	<0.000047	<0.00047
Acenaphthene	mg/L	0.44	0.000036 J	7.1	0.0053	0.079
Acenaphthylene	mg/L	0.0037	0.000023 J	0.079	0.00015	0.0023
Anthracene	mg/L	0.15	0.000036 J	5.8	0.0039	0.020
Benzo(a)anthracene	mg/L	0.047	0.000055 J	0.52	0.0011	0.0031
Benzo(a)pyrene	mg/L	0.0089	0.000063 J	0.13	0.00052	0.0011
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.00060	<0.000030	<0.00030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.0011	0.00029	0.0039 J	0.00063	<0.00037
Chrysene	mg/L	0.040	0.000054 J	0.63	0.0011	0.0033
Di-n-butylphthalate (DBP)	mg/L	0.00018 J	<0.000020	<0.00040	0.00098	<0.00020
Dibenzofuran	mg/L	0.50	0.000030 J	6.2	0.0049	0.061
Fluoranthene	mg/L	0.31	0.000074 J	4.7	0.0057	0.024
Fluorene	mg/L	0.32	0.000036 J	6.7	0.0038	0.049
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.00050	<0.000025	<0.00025

**Analytical Results Summary  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Location ID:	MW-68B	MW-68C	MW-70B	MW-71B	MW-72B
Sample Name:	WG-1620-MW68B-20200123	WG-1620-MW68C-20200117	WG-1620-MW70B-20200120	WG-1620-MW71B-20200115	WG-1620-MW72B-20200109
Sample Date:	01/23/2020	01/17/2020	01/20/2020	01/15/2020	01/09/2020

Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	12	<0.00052	71	0.0064	2.0
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.00048	<0.000024	<0.00024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.0016	<0.000079	<0.00079
Phenanthrene	mg/L	0.95	0.000042 J	14	0.017	0.075
Phenol	mg/L	<0.000035	<0.000035	3.4	<0.000035	5.2
Pyrene	mg/L	0.19	0.000092 J	3.2	0.0042	0.013
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00944 J	<0.000400	0.00120 J	0.00279	0.00105 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-74B	MW-75B	MW-76C	MW-77A	MW-78A
Sample Name:	WG-1620-MW74B-20200117	WG-1620-MW75B-20200117	WG-1620-MW76C-20200109	WG-1620-MW77A-20200109	WG-1620-MW78A-20200117
Sample Date:	01/17/2020	01/17/2020	01/09/2020	01/09/2020	01/17/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.12	0.12	<0.00020	0.058
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.076	0.037	<0.00030	0.063
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.29	0.13	<0.00020	0.0069
Vinyl chloride	mg/L	--	--	--	--
Xylenes (total)	mg/L	0.20	0.11	<0.00030	0.065
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.21	0.86	<0.00038	2.5
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	1.3	0.81	<0.00045	0.22
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.98	0.67	<0.00041	0.13
Acenaphthylene	mg/L	0.018	0.0086	0.000035 J	0.0013
Anthracene	mg/L	0.42	0.39	<0.00017	0.0036
Benzo(a)anthracene	mg/L	0.13	0.12	0.00014	<0.000050
Benzo(a)pyrene	mg/L	0.034	0.035	0.000037 J	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.0014	0.0021	<0.00016	<0.000077
Chrysene	mg/L	0.082	0.095	0.00013	0.000041 J
Di-n-butylphthalate (DBP)	mg/L	0.0022	0.0016	<0.000020	<0.000020
Dibenzofuran	mg/L	0.72	0.54	<0.00029	0.079
Fluoranthene	mg/L	0.67	0.75	0.0010	0.00071
Fluorene	mg/L	0.88	0.61	<0.00032	0.067
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-74B	MW-75B	MW-76C	MW-77A	MW-78A	
Sample Name:	WG-1620-MW74B-20200117	WG-1620-MW75B-20200117	WG-1620-MW76C-20200109	WG-1620-MW77A-20200109	WG-1620-MW78A-20200117	
Sample Date:	01/17/2020	01/17/2020	01/09/2020	01/09/2020	01/17/2020	
<b>Parameters</b>	<b>Unit</b>					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	8.5	7.5	<0.0038	4.7	3.1
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	2.3	1.9	<0.00069	0.031	0.031
Phenol	mg/L	0.048	0.0029	<0.000035	<0.000035	0.13
Pyrene	mg/L	0.44	0.51	0.00070	0.00046	0.0018
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	25.5	34.9	<0.19	12.1	4.30
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	13	21	<0.19	3.8	1.6
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	0.47 J	0.91	<0.19	<0.19	<0.19
Total Petroleum Hydrocarbons (C6-C12)	mg/L	12	13	<0.19	8.3	2.7
<b>Metals</b>						
Arsenic	mg/L	0.000443 J	0.00167 J	0.00299	0.0237	0.00991



Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-79A	MW-80B	MW-81B	MW-82B	MW-83B
Sample Name:	WG-1620-MW79A-20200117	WG-1620-MW80B-20200107	WG-1620-MW81B-20200109	WG-1620-MW82B-20200121	WG-1620-MW83B-20200110
Sample Date:	01/17/2020	01/07/2020	01/09/2020	01/21/2020	01/10/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.12	<0.00020	<0.00020	0.021
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.075	<0.00030	<0.00030	0.078
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.26	<0.00020	<0.00020	0.0052
Vinyl chloride	mg/L	--	--	--	<0.00020
Xylenes (total)	mg/L	0.20	<0.00030	<0.00030	0.098
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	6.7	0.00028	<0.000040	0.000072 J
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.43	<0.000036	<0.000019	0.00010
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.19	0.000046 J	<0.000077	0.000043 J
Acenaphthylene	mg/L	0.0036	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	0.0094	0.000026 J	<0.000014	0.000031 J
Benzo(a)anthracene	mg/L	0.00081	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	0.00029	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000069	<0.000066	<0.000037
Chrysene	mg/L	0.00072	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.00098	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.14	0.000048 J	<0.000020	0.000040 J
Fluoranthene	mg/L	0.0076	<0.000010	<0.000010	<0.000010
Fluorene	mg/L	0.094	0.000032 J	<0.000030	0.000037 J
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

	Location ID:	MW-79A	MW-80B	MW-81B	MW-82B	MW-83B
	Sample Name:	WG-1620-MW79A-20200117	WG-1620-MW80B-20200107	WG-1620-MW81B-20200109	WG-1620-MW82B-20200121	WG-1620-MW83B-20200110
	Sample Date:	01/17/2020	01/07/2020	01/09/2020	01/21/2020	01/10/2020
Parameters	Unit					
<b>Semi-volatile Organic Compounds (Continued)</b>						
Naphthalene	mg/L	9.0	<0.00034	<0.00023	<0.00041	1.2
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.081	0.000027 J	<0.000021	0.00012	0.021
Phenol	mg/L	0.35	<0.000035	<0.000035	<0.000035	<0.000035
Pyrene	mg/L	0.0035	<0.000019	<0.000019	<0.000019	0.0022
<b>Total Petroleum Hydrocarbons</b>						
Total Petroleum Hydrocarbons	mg/L	37.0	<0.19	<0.19	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	16	<0.19	<0.19	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	<0.19	<0.19	<0.19	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	21	<0.19	<0.19	--	--
<b>Metals</b>						
Arsenic	mg/L	0.00893	0.00212	0.00130 J	0.00484	0.0709

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-83C	MW-84B	MW-85C	MW-86C
Sample Name:	WG-1620-MW83C-20200110	WG-1620-MW84B-20200116	WG-1620-MW85C-20200109	WG-1620-MW86C-20200117
Sample Date:	01/10/2020	01/16/2020	01/09/2020	01/17/2020
Parameters	Unit			
<b>Volatile Organic Compounds</b>				
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	0.0064	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	0.0094	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	0.0039
Vinyl chloride	mg/L	<0.00020	--	--
Xylenes (total)	mg/L	<0.00030	0.0032	<0.00030
<b>Semi-volatile Organic Compounds</b>				
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	0.00048	0.00018 J
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.00021	0.0065	<0.000041
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	0.00035	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	0.00077 J
Acenaphthene	mg/L	0.00018	0.025	<0.000077
Acenaphthylene	mg/L	<0.000015	0.00031	<0.000015
Anthracene	mg/L	<0.000014	0.00095	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	0.000031 J
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	<0.00016
Chrysene	mg/L	<0.000021	<0.000021	0.000053 J
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.00015	0.015	<0.000059
Fluoranthene	mg/L	<0.000010	0.00037	0.00025
Fluorene	mg/L	0.00012	0.0063	<0.000055
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

	Location ID:	MW-83C	MW-84B	MW-85C	MW-86C
	Sample Name:	WG-1620-MW83C-20200110	WG-1620-MW84B-20200116	WG-1620-MW85C-20200109	WG-1620-MW86C-20200117
	Sample Date:	01/10/2020	01/16/2020	01/09/2020	01/17/2020
Parameters	Unit				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Naphthalene	mg/L	0.0018	0.13	<0.00046	<0.00075
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.00017	0.0065	<0.000067	<0.000021
Phenol	mg/L	<0.000035	<0.000035	<0.000042	<0.000035
Pyrene	mg/L	<0.000019	0.00025	0.00021	<0.000019
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	--	--	<0.20	<0.19
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	<0.20	<0.19
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	<0.20	<0.19
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	<0.20	<0.19
<b>Metals</b>					
Arsenic	mg/L	0.00564	0.00363	0.00272	0.000645 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Location ID:	MW-86C	MW-87C	MW-88C	MW-89B
Sample Name:	WG-1620-FD04-20200117	WG-1620-MW87C-20200120	WG-1620-MW88C-20200114	WG-1620-MW89B-20200116
Sample Date:	01/17/2020 Duplicate	01/20/2020	01/14/2020	01/16/2020
Parameters	Unit			
<b>Volatile Organic Compounds</b>				
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.0045	<0.00020	<0.00020
Vinyl chloride	mg/L	--	<0.00020	--
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>				
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.000066 J	0.00023	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.00011	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	0.00012	0.000088 J	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	0.000066 J	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000053	<0.000037	0.000064 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	0.00011	<0.000020
Fluoranthene	mg/L	<0.000010	<0.000086	<0.000010
Fluorene	mg/L	<0.000030	0.000093 J	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

	Location ID:	MW-86C	MW-87C	MW-88C	MW-89B
	Sample Name:	WG-1620-FD04-20200117	WG-1620-MW87C-20200120	WG-1620-MW88C-20200114	WG-1620-MW89B-20200116
	Sample Date:	01/17/2020	01/20/2020	01/14/2020	01/16/2020
		Duplicate			
Parameters	Unit				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Naphthalene	mg/L	<0.00022	0.0011	<0.000020	<0.00029
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.00036	<0.00028	<0.000021	0.00083
Phenol	mg/L	<0.000035	0.00048	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000057	<0.000019	0.00081
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	<0.19	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	<0.19	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	<0.19	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	<0.19	--	--	--
<b>Metals</b>					
Arsenic	mg/L	0.000699 J	0.00135 J	0.000862 J	0.000463 J

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

	Location ID:	MW-90B	P-11	TW-41B
	Sample Name:	WG-1620-MW90B-20200120	WG-1620-P11-20200114	WG-1620-TW41B-20200113
	Sample Date:	01/20/2020	01/14/2020	01/13/2020
Parameters	Unit			
<b>Volatile Organic Compounds</b>				
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	0.00039 J
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020
Vinyl chloride	mg/L	<0.00020	--	--
Xylenes (total)	mg/L	<0.00030	<0.00030	0.0021
<b>Semi-volatile Organic Compounds</b>				
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	0.000075 J	<0.000040	0.00017 J
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	0.00012 J
2-Methylnaphthalene	mg/L	0.00034	0.000076 J	0.0071
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000047
Acenaphthene	mg/L	0.00015	0.0046	0.12
Acenaphthylene	mg/L	<0.000015	<0.000015	0.0011
Anthracene	mg/L	0.00014	0.000081 J	0.0031
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	0.000052 J	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	<0.000037	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	0.00011	0.00016	0.055
Fluoranthene	mg/L	<0.00012	<0.00010	0.0021
Fluorene	mg/L	0.00013	0.0013	0.078
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000025

Table 2

**Analytical Results Summary**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

	Location ID:	MW-90B	P-11	TW-41B
	Sample Name:	WG-1620-MW90B-20200120	WG-1620-P11-20200114	WG-1620-TW41B-20200113
	Sample Date:	01/20/2020	01/14/2020	01/13/2020
Parameters	Unit			
<b>Semi-volatile Organic Compounds (Continued)</b>				
Naphthalene	mg/L	0.0030	0.00048	0.058
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000079
Phenanthrene	mg/L	0.00040	0.00035	0.0084
Phenol	mg/L	0.00019 J	<0.000035	<0.000035
Pyrene	mg/L	<0.000077	<0.000019	0.00082
<b>Total Petroleum Hydrocarbons</b>				
Total Petroleum Hydrocarbons	mg/L	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--
<b>Metals</b>				
Arsenic	mg/L	0.00290	0.0360	0.0557

## Notes:

- < - Not detected at the associated reporting limit
- J - Estimated concentration
- JH - Estimated concentration; biased high
- J - Estimated concentration
- "--" - Not analyzed



Table 3

**Analytical Methods**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Method	Matrix	Holding Time	
			Collection to Extraction (Days)	Extraction to Analysis (Days)
VOCs	SW-846 8260C	Water	-	14
SVOCs	SW-846 8270D	Water	7	40
TPH	TX1005	Water	14	40
Arsenic	SW-846 6020A	Water	-	180

## Notes:

- VOCs - Volatile Organic Compounds  
SVOCs - Semi-volatile Organic Compounds  
TPH - Total Petroleum Hydrocarbons  
"-" - Not Applicable

## Method References:

- SW-846 - "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition, 1986, with subsequent revisions

Table 4

**Qualified Sample Results Due to Outlying LCS/LCSD Results**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Analyte	LCS Date (mm/dd/yyyy)	LCS % Recovery	LCSD % Recovery	RPD (percent)	Control Limits		Associated Sample ID	Qualified Result	Units
						% Recovery	RPD			
SVOCs	2-Methylnaphthalene	01/31/2020	95.4	117	20.1	50-120	20	WG-1620-MW12B-20200123	320 J	mg/L
								WG-1620-MW41B-20200123	0.67 J	mg/L
								WG-1620-MW68B-20200123	1.1 J	mg/L

## Notes:

- LCS - Laboratory Control Sample
- LCSD - Laboratory Control Sample Duplicate
- RPD - Relative Percent Difference
- SVOCs - Semi-volatile Organic Compounds
- J - Estimated concentration

**Table 5**

**Qualified Sample Data Due to Outlying Laboratory Duplicate Results**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Sample ID	Analyte	RPD		Associated Sample IDs	Qualified Result	Units
			RPD (percent)	Control Limit (percent)			
Metals	WG-1620-MW41B-20200123	Arsenic	11.1	10	WG-1620-MW12B-20200123	0.0491 J	mg/L
					WG-1620-MW41B-20200123	0.0842 J	mg/L
					WG-1620-MW68B-20200123	0.00944 J	mg/L

Notes:

- RPD - Relative Percent Difference
- J - Estimated concentration

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB01-20200107	01/07/2020	2-Methylnaphthalene	0.000031 J	WG-1620-MW80B-20200107	0.000036 J	<0.000036	mg/L
			bis(2-Ethylhexyl)phthalate (DEHP)	0.00017 J	WG-1620-MW49A-20200107	0.00011 J	<0.00011	mg/L
					WG-1620-MW80B-20200107	0.000069 J	<0.000069	mg/L
			Naphthalene	0.00019	WG-1620-MW80B-20200107	0.00034	<0.00034	mg/L
SVOCs	WG-1620-FB02-20200108	01/08/2020	bis(2-Ethylhexyl)phthalate (DEHP)	0.000099 J	WG-1620-MW18A-20200108	0.00043	<0.00043	mg/L
					WG-1620-MW57A-20200108	0.00037	<0.00037	mg/L
					WG-1620-MW58A-20200108	0.00011 J	<0.00011	mg/L
SVOCs	WG-1620-FB03-20200109	01/09/2020	2,4-Dimethylphenol	0.001	WG-1620-MW23C-20200109	0.047	<0.047	mg/L
					WG-1620-MW36A-20200109	0.000052 J	<0.000052	mg/L
					WG-1620-MW36B-20200109	0.000054 J	<0.000054	mg/L
					WG-1620-MW44A-20200109	0.00042	<0.00042	mg/L
					WG-1620-MW50A-20200109	0.00071	<0.00071	mg/L
					WG-1620-MW53C-20200109	0.0011	<0.0011	mg/L
			2-Methylnaphthalene		WG-1620-MW76C-20200109	0.00038	<0.00038	mg/L
					WG-1620-MW36B-20200109	0.00013	<0.00013	mg/L
					WG-1620-MW44A-20200109	0.00035	<0.00035	mg/L
	WG-1620-MW50A-20200109	0.00019	<0.00019	mg/L				

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB03-20200109	01/09/2020	2-Methylnaphthalene	0.00038	WG-1620-MW51A-20200109	0.000044 J	<0.000044	mg/L
					WG-1620-MW51C-20200109	0.000055 J	<0.000055	mg/L
					WG-1620-MW53C-20200109	0.00025	<0.00025	mg/L
					WG-1620-MW76C-20200109	0.00045	<0.00045	mg/L
					WG-1620-MW85C-20200109	0.000041 J	<0.000041	mg/L
			Acenaphthene	0.00015	WG-1620-MW36B-20200109	0.00017	<0.00017	mg/L
					WG-1620-MW50A-20200109	0.00037	<0.00037	mg/L
					WG-1620-MW51A-20200109	0.000040 J	<0.000040	mg/L
					WG-1620-MW51C-20200109	0.000094 J	<0.000094	mg/L
					WG-1620-MW53C-20200109	0.00032	<0.00032	mg/L
			Anthracene	0.000047 J	WG-1620-MW76C-20200109	0.00041	<0.00041	mg/L
					WG-1620-MW81B-20200109	0.000077 J	<0.000077	mg/L
					WG-1620-MW85C-20200109	0.000077 J	<0.000077	mg/L
					WG-1620-MW19C-20200109	0.00017	<0.00017	mg/L
					WG-1620-MW51A-20200109	0.000063 J	<0.000063	mg/L
					WG-1620-MW51C-20200109	0.00010	<0.00010	mg/L
					WG-1620-MW53C-20200109	0.000028 J	<0.000028	mg/L
					WG-1620-MW76C-20200109	0.00017	<0.00017	mg/L

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units		
SVOCs	WG-1620-FB03-20200109	01/09/2020	bis(2-Ethylhexyl)phthalate (DEHP)	0.00011 J	WG-1620-MW19C-20200109	0.000071 J	<0.000071	mg/L		
					WG-1620-MW50A-20200109	0.000085 J	<0.000085	mg/L		
					WG-1620-MW51A-20200109	0.00014 J	<0.00014	mg/L		
					WG-1620-MW51C-20200109	0.000085 J	<0.000085	mg/L		
					WG-1620-MW54C-20200109	0.00018 J	<0.00018	mg/L		
					WG-1620-MW76C-20200109	0.00016 J	<0.00016	mg/L		
					WG-1620-MW77A-20200109	0.000077 J	<0.000077	mg/L		
					WG-1620-MW81B-20200109	0.000066 J	<0.000066	mg/L		
					WG-1620-MW85C-20200109	0.00016 J	<0.00016	mg/L		
					Dibenzofuran	0.00012	WG-1620-MW36B-20200109	0.00012	<0.00012	mg/L
							WG-1620-MW44A-20200109	0.00033	<0.00033	mg/L
							WG-1620-MW50A-20200109	0.00013	<0.00013	mg/L
							WG-1620-MW51A-20200109	0.000049 J	<0.000049	mg/L
							WG-1620-MW51C-20200109	0.00011	<0.00011	mg/L
							WG-1620-MW53C-20200109	0.00027	<0.00027	mg/L
SVOCs	WG-1620-FB03-20200109	01/09/2020	Dibenzofuran	0.00012	WG-1620-MW85C-20200109	0.000059 J	<0.000059	mg/L		
			Fluorene		0.000092 J	WG-1620-MW36B-20200109	0.00012	<0.00012	mg/L	
						WG-1620-MW50A-20200109	0.00011	<0.00011	mg/L	

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units			
SVOCs	WG-1620-FB03-20200109	01/09/2020	Fluorene	0.000092 J	WG-1620-MW51A-20200109	0.000043 J	<0.000043	mg/L			
					WG-1620-MW51C-20200109	0.00011	<0.00011	mg/L			
					WG-1620-MW53C-20200109	0.00020	<0.00020	mg/L			
					WG-1620-MW76C-20200109	0.00032	<0.00032	mg/L			
					WG-1620-MW85C-20200109	0.000055 J	<0.000055	mg/L			
					Naphthalene	0.0078	WG-1620-MW19C-20200109	0.00011	<0.00011	mg/L	
							WG-1620-MW36A-20200109	0.00034	<0.00034	mg/L	
							WG-1620-MW36B-20200109	0.00093	<0.00093	mg/L	
							WG-1620-MW44A-20200109	0.0030	<0.0030	mg/L	
							WG-1620-MW50A-20200109	0.0020	<0.0020	mg/L	
			WG-1620-MW51A-20200109	0.00032			<0.00032	mg/L			
			WG-1620-MW51C-20200109	0.00030			<0.00030	mg/L			
			WG-1620-MW53C-20200109	0.0012			<0.0012	mg/L			
			WG-1620-MW54C-20200109	0.035			<0.035	mg/L			
			WG-1620-MW76C-20200109	0.0038			<0.0038	mg/L			
			Phenanthrene	0.00019	WG-1620-MW81B-20200109	0.00023	<0.00023	mg/L			
					WG-1620-MW85C-20200109	0.00046	<0.00046	mg/L			
								WG-1620-MW19C-20200109	0.00045	<0.00045	mg/L
								WG-1620-MW36B-20200109	0.00010	<0.00010	mg/L

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB03-20200109	01/09/2020	Phenanthrene	0.00019	WG-1620-MW44A-20200109	0.00022	<0.00022	mg/L
					WG-1620-MW50A-20200109	0.000070 J	<0.000070	mg/L
					WG-1620-MW51A-20200109	0.00021	<0.00021	mg/L
					WG-1620-MW51C-20200109	0.00056	<0.00056	mg/L
					WG-1620-MW53C-20200109	0.00025	<0.00025	mg/L
					WG-1620-MW76C-20200109	0.00069	<0.00069	mg/L
					WG-1620-MW85C-20200109	0.000067 J	<0.000067	mg/L
					Phenol	0.000053 J	WG-1620-MW36B-20200109	0.000075 J
			WG-1620-MW44A-20200109	0.00013 J	<0.00013		mg/L	
			WG-1620-MW50A-20200109	0.000066 J	<0.000066		mg/L	
			WG-1620-MW50A-20200109	0.000034 J	<0.000034		mg/L	
			WG-1620-MW51A-20200109	0.000083 J	<0.000083		mg/L	
			WG-1620-MW51C-20200109	0.00011	<0.00011		mg/L	
			WG-1620-MW53C-20200109	0.000061 J	<0.000061		mg/L	
			WG-1620-MW54C-20200109	0.000064 J	<0.000064		mg/L	
			WG-1620-MW85C-20200109	0.000042 J	<0.000042	mg/L		
SVOCs	WG-1620-FB04-20200110	01/10/2020	Naphthalene	0.00020	WG-1620-MW35A-20200110	0.00020	<0.00020	mg/L



Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units	
SVOCs	WG-1620-FB05-20200113	01/13/2020	Naphthalene	0.00033	WG-1620-MW05-20200113	0.00012	<0.00012	mg/L	
					WG-1620-MW12A-20200113	0.00051	<0.00051	mg/L	
					WG-1620-MW42B-20200113	0.00076	<0.00076	mg/L	
SVOCs	WG-1620-FB08-20200116	01/16/2020	2-Methylnaphthalene	0.000052 J	WG-1620-MW26A-20200116	0.000042 J	<0.000042	mg/L	
					WG-1620-MW47C-20200116	0.00010	<0.00010	mg/L	
					WG-1620-MW48C-20200116	0.000076 J	<0.000076	mg/L	
					WG-1620-MW59A-20200116	0.00016	<0.00016	mg/L	
					WG-1620-MW59B-20200116	0.000038 J	<0.000038	mg/L	
					WG-1620-MW63B-20200116	0.000056 J	<0.000056	mg/L	
					WG-1620-MW66D-20200116	0.000095 J	<0.000095	mg/L	
SVOCs	WG-1620-FB08-20200116	01/16/2020	2-Methylnaphthalene	0.000052 J	WG-1620-MW89B-20200116	0.000032 J	<0.000032	mg/L	
					Acenaphthene	WG-1620-MW26A-20200116	0.00017	<0.00017	mg/L
						WG-1620-MW28A-20200116	0.00014	<0.00014	mg/L
						WG-1620-MW47C-20200116	0.000057 J	<0.000057	mg/L
						WG-1620-MW48C-20200116	0.000073 J	<0.000073	mg/L
						WG-1620-MW59A-20200116	0.00029	<0.00029	mg/L
						WG-1620-MW59B-20200116	0.000051 J	<0.000051	mg/L
WG-1620-MW59D-20200116	0.000036 J	<0.000036	mg/L						

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB08-20200116	01/16/2020	Acenaphthene	0.000077 J	WG-1620-MW61A-20200116	0.000052 J	<0.000052	mg/L
					WG-1620-MW63B-20200116	0.000059 J	<0.000059	mg/L
					WG-1620-MW89B-20200116	0.000083 J	<0.000083	mg/L
			Anthracene	0.000027 J	WG-1620-MW26A-20200116	0.000087 J	<0.000087	mg/L
					WG-1620-MW28A-20200116	0.00013	<0.00013	mg/L
					WG-1620-MW36D-20200116	0.000033 J	<0.000033	mg/L
					WG-1620-MW59A-20200116	0.000027 J	<0.000027	mg/L
					WG-1620-MW59B-20200116	0.000031 J	<0.000031	mg/L
					WG-1620-MW63B-20200116	0.000081 J	<0.000081	mg/L
			Dibenzofuran	0.000066 J	WG-1620-MW26A-20200116	0.000044 J	<0.000044	mg/L
					WG-1620-MW28A-20200116	0.00010	<0.00010	mg/L
					WG-1620-MW47C-20200116	0.000049 J	<0.000049	mg/L
					WG-1620-MW48C-20200116	0.000054 J	<0.000054	mg/L
					WG-1620-MW59A-20200116	0.00016	<0.00016	mg/L
					WG-1620-MW59B-20200116	0.000050 J	<0.000050	mg/L
					WG-1620-MW59D-20200116	0.000028 J	<0.000028	mg/L
					WG-1620-MW63B-20200116	0.000063 J	<0.000063	mg/L
					WG-1620-MW89B-20200116	0.00011	<0.00011	mg/L

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB08-20200116	01/16/2020	Fluoranthene	0.000033 J	WG-1620-MW47C-20200116	0.00011	<0.00011	mg/L
					WG-1620-MW48C-20200116	0.000061 J	<0.000061	mg/L
					WG-1620-MW59A-20200116	0.000032 J	<0.000032	mg/L
					WG-1620-MW59B-20200116	0.000025 J	<0.000025	mg/L
					WG-1620-MW61A-20200116	0.000060 J	<0.000060	mg/L
					WG-1620-MW63B-20200116	0.000023 J	<0.000023	mg/L
			Fluorene	0.000058 J	WG-1620-MW26A-20200116	0.00029	<0.00029	mg/L
					WG-1620-MW28A-20200116	0.00014	<0.00014	mg/L
					WG-1620-MW47C-20200116	0.000042 J	<0.000042	mg/L
					WG-1620-MW48C-20200116	0.000050 J	<0.000050	mg/L
					WG-1620-MW59A-20200116	0.00011	<0.00011	mg/L
					WG-1620-MW59B-20200116	0.000042 J	<0.000042	mg/L
			Naphthalene	0.00054	WG-1620-MW59D-20200116	0.000031 J	<0.000031	mg/L
					WG-1620-MW63B-20200116	0.00012	<0.00012	mg/L
					WG-1620-MW89B-20200116	0.00018	<0.00018	mg/L
					WG-1620-MW36D-20200116	0.000033 J	<0.000033	mg/L
					WG-1620-MW66D-20200116	0.00013	<0.00013	mg/L
					WG-1620-MW28C-20200116	0.00016	<0.00016	mg/L
				WG-1620-MW26A-20200116	0.00017	<0.00017	mg/L	

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB08-20200116	01/16/2020	Naphthalene	0.00054	WG-1620-MW61A-20200116	0.00021	<0.00021	mg/L
					WG-1620-MW59B-20200116	0.00024	<0.00024	mg/L
					WG-1620-MW89B-20200116	0.00029	<0.00029	mg/L
					WG-1620-MW63B-20200116	0.00035	<0.00035	mg/L
					WG-1620-MW59D-20200116	0.00013	<0.00013	mg/L
					WG-1620-MW48C-20200116	0.00052	<0.00052	mg/L
					WG-1620-MW47C-20200116	0.00056	<0.00056	mg/L
					WG-1620-MW28A-20200116	0.00060	<0.00060	mg/L
					WG-1620-MW59A-20200116	0.0012	<0.0012	mg/L
					Phenanthrene	0.000067 J	WG-1620-MW26A-20200116	0.000030 J
			WG-1620-MW36D-20200116	0.000047 J			<0.000047	mg/L
			WG-1620-MW47C-20200116	0.000088 J			<0.000088	mg/L
			WG-1620-MW48C-20200116	0.000052 J			<0.000052	mg/L
			WG-1620-MW59A-20200116	0.00012			<0.00012	mg/L
			WG-1620-MW59B-20200116	0.000099 J			<0.000099	mg/L
			WG-1620-MW59D-20200116	0.000043 J			<0.000043	mg/L
			WG-1620-MW61A-20200116	0.00010			<0.00010	mg/L
			WG-1620-MW63B-20200116	0.000072 J			<0.000072	mg/L

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB08-20200116	01/16/2020	Pyrene	0.000022 J	WG-1620-MW47C-20200116	0.00010	<0.00010	mg/L
					WG-1620-MW48C-20200116	0.000065 J	<0.000065	mg/L
					WG-1620-MW59A-20200116	0.000023 J	<0.000023	mg/L
					WG-1620-MW59B-20200116	0.000021 J	<0.000021	mg/L
					WG-1620-MW61A-20200116	0.000059 J	<0.000059	mg/L
					WG-1620-MW63B-20200116	0.000021 J	<0.000021	mg/L
SVOCs	WG-1620-FB09-20200117	01/17/2020	bis(2-Ethylhexyl)phthalate (DEHP)	0.000050 J	WG-1620-FD04-20200117	0.000053 J	<0.000053	mg/L
					WG-1620-MW68A-20200117	0.000061 J	<0.000061	mg/L
			Naphthalene	0.00017	WG-1620-FD04-20200117	0.00022	<0.00022	mg/L
					WG-1620-MW68A-20200117	0.00022	<0.00022	mg/L
					WG-1620-MW68C-20200117	0.00052	<0.00052	mg/L
					WG-1620-MW86C-20200117	0.00075	<0.00075	mg/L
SVOCs	WG-1620-FB10-20200120	01/20/2020	1,2-Diphenylhydrazine	0.000047 J	WG-1620-MW45C-20200120	0.000085 J	<0.000085	mg/L
			2-Methylnaphthalene	0.000035 J	WG-1620-FD05-20200120	0.000095 J	<0.000095	mg/L
SVOCs	WG-1620-FB10-20200120	01/20/2020	1,2-Diphenylhydrazine	0.000047 J	WG-1620-MW32AR-20200120	0.000028 J	<0.000028	mg/L
			2-Methylnaphthalene	0.000035 J	WG-1620-MW33A-20200120	0.000063 J	<0.000063	mg/L
					WG-1620-MW45C-20200120	0.00013	<0.00013	mg/L

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks  
HWPW - Site-Wide Monitoring  
Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works  
Houston, Texas  
January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units		
SVOCs	WG-1620-FB10-20200120	01/20/2020	Fluoranthene	0.000060 J	WG-1620-MW32AR-20200120	0.00029	<0.00029	mg/L		
					WG-1620-MW87C-20200120	0.000086 J	<0.000086	mg/L		
					WG-1620-MW90B-20200120	0.00012	<0.00012	mg/L		
			Naphthalene	0.00019	WG-1620-FD05-20200120	0.00066	<0.00066	mg/L		
					WG-1620-MW32AR-20200120	0.00014	<0.00014	mg/L		
					WG-1620-MW33A-20200120	0.00050	<0.00050	mg/L		
			Naphthalene	0.00019	WG-1620-MW45C-20200120	0.00061	<0.00061	mg/L		
					Phenanthrene	0.000066 J	WG-1620-FD05-20200120	0.00046 J	<0.00046	mg/L
							WG-1620-MW32AR-20200120	0.00032 J	<0.00032	mg/L
			WG-1620-MW33A-20200120	0.00056 J			<0.00056	mg/L		
			Phenanthrene	0.000066 J	WG-1620-MW45C-20200120	0.00027	<0.00027	mg/L		
					WG-1620-MW87C-20200120	0.00028	<0.00028	mg/L		
					Pyrene	0.000040 J	WG-1620-MW33A-20200120	0.00019	<0.00019 J	mg/L
			WG-1620-MW87C-20200120	0.000057 J			<0.000057	mg/L		
			WG-1620-MW90B-20200120	0.000077 J			<0.000077	mg/L		
SVOCs	WG-1620-FB11-20200121	01/21/2020	Naphthalene	0.00018	WG-1620-MW03-20200121	0.00024	<0.00024	mg/L		
					WG-1620-MW82B-20200121	0.00041	<0.00041	mg/L		

Table 6

**Qualified Sample Data Due to Analyte Concentrations in the Field Blanks**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Field Blank ID	Blank Date (mm/dd/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	WG-1620-FB14-20200210	02/10/2020	2-Methylnaphthalene	0.00038	WG-1620-MW22AR-20200210	0.00012	<0.00012	mg/L
					WG-1620-MW22BR-20200210	0.000043 J	<0.000043	mg/L
SVOCs	WG-1620-FB14-20200210	02/10/2020	Acenaphthene	0.00020	WG-1620-MW22AR-20200210	0.00015	<0.00015	mg/L
			Dibenzofuran	0.00019	WG-1620-MW22AR-20200210	0.00015	<0.00015	mg/L
SVOCs	WG-1620-FB14-20200210	02/10/2020	Dibenzofuran	0.00019	WG-1620-MW22BR-20200210	0.00029	<0.00029	mg/L
			Naphthalene	0.0034	WG-1620-MW22AR-20200210	0.00044	<0.00044	mg/L
				0.0034	WG-1620-MW22BR-20200210	0.00013	<0.00013	mg/L
			Phenanthrene	0.000029 J	WG-1620-MW22BR-20200210	0.00013	<0.00013	mg/L
			Phenol	0.00053	WG-1620-MW22AR-20200210	0.000041 J	<0.000041	mg/L

## Notes:

SVOCs - Semi-volatile Organic Compounds

VOCs - Volatile Organic Compounds

J - Estimated concentration

&lt; - Not detected at the associated reporting limit

Table 7

**Qualified Sample Data Due to Variability in Field Duplicate Results**  
**HWPW - Site-Wide Monitoring**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January-February 2020**

Parameter	Analyte	RPD	Diff	Sample ID	Qualified Result	Field Duplicate Sample ID	Qualified Result	Units
VOCs	Benzene	43.9	-	WG-1620-MW35B-20200110	0.025 J	WG-1620-FD01-20200110	0.016 J	mg/L
	Xylenes (total)	34.3	-		0.058 J		0.041 J	mg/L
SVOCs	2-Methylnaphthalene	77.2	-	WG-1620-MW35B-20200110	0.14 J	WG-1620-FD01-20200110	0.062 J	mg/L
	Acenaphthene	56.5	-		0.059 J		0.033 J	mg/L
	Anthracene	46.8	-		0.0029 J		0.0018 J	mg/L
	Dibenzofuran	54.1	-		0.054 J		0.031 J	mg/L
	Fluoranthene	46.9	-		0.0015 J		0.00093 J	mg/L
	Fluorene	60.5	-		0.028 J		0.015 J	mg/L
	Naphthalene	175	-		2.0 J		0.13 J	mg/L
	Phenanthrene	59.5	-		0.024 J		0.013 J	mg/L
	Pyrene	43.5	-		0.00070 J		0.00045 J	mg/L
SVOCs	2,6-Dinitrotoluene	-	0.0028	WG-1620-MW21C-20200114	<0.00020 J	WG-1620-FD02-20200114	0.0030 J	mg/L
SVOCs	Acenaphthene	40.4	-	WG-1620-MW33A-20200120	0.00073 J	WG-1620-FD05-20200120	0.0011 J	mg/L
	Fluoranthene	47.8	-		0.00035 J		0.00057 J	mg/L
	Pyrene	-	0.00042		<0.00010 J		0.00052 J	mg/L

## Notes:

RPD - Relative Percent Difference

Diff - Difference

VOCs - Volatile Organic Compounds

SVOCs - Semi-volatile Organic Compounds

J - Estimated concentration

&lt; - Not detected at the associated reporting limit

"- " - Not Applicable



# Attachment A

## Laboratory NELAP Certificate



# Texas Commission on Environmental Quality

## NELAP - Recognized Laboratory Fields of Accreditation



ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: *Drinking Water***

**Method** EPA 1613

Analyte	AB	Analyte ID	Method ID
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408

**Method** EPA 200.8

Analyte	AB	Analyte ID	Method ID
Copper	TX	1055	10014605
Lead	TX	1075	10014605



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**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 120.1			
Analyte Conductivity	TX	1610	10006403
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 160.4			
Analyte Residue-volatile	TX	1970	10010409
Method EPA 1613			
Analyte 1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10120408
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10120408
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10120408
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10120408
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10120408
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10120408
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10120408
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10120408
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,6,7,8-HxCDD)	TX	9456	10120408
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10120408
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10120408
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10120408
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10120408
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10120408



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**Matrix: Non-Potable Water**

2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10120408
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10120408
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10120408
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10120408
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10120408
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10120408
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10120408
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10120408
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10120408
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10120408
<b>Method EPA 1664</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10127807
<b>Method EPA 180.1</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Turbidity	TX	2055	10011606
<b>Method EPA 200.8</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Aluminum	TX	1000	10014605
Antimony	TX	1005	10014605
Arsenic	TX	1010	10014605
Barium	TX	1015	10014605
Beryllium	TX	1020	10014605
Boron	TX	1025	10014605
Cadmium	TX	1030	10014605
Calcium	TX	1035	10014605
Chromium	TX	1040	10014605
Cobalt	TX	1050	10014605
Copper	TX	1055	10014605
Iron	TX	1070	10014605



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**Matrix: Non-Potable Water**

Lead	TX	1075	10014605
Magnesium	TX	1085	10014605
Manganese	TX	1090	10014605
Molybdenum	TX	1100	10014605
Nickel	TX	1105	10014605
Potassium	TX	1125	10014605
Selenium	TX	1140	10014605
Silver	TX	1150	10014605
Sodium	TX	1155	10014605
Strontium	TX	1160	10014605
Thallium	TX	1165	10014605
Tin	TX	1175	10014605
Titanium	TX	1180	10014605
Uranium	TX	3035	10014605
Vanadium	TX	1185	10014605
Zinc	TX	1190	10014605

**Method EPA 245.1**

Analyte	AB	Analyte ID	Method ID
Mercury	TX	1095	10036609

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 325.1**

Analyte	AB	Analyte ID	Method ID
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**Matrix: Non-Potable Water**

Chloride	TX	1575	10056801
<b>Method</b> EPA 335.1			
<b>Analyte</b> Amenable cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1510	<b>Method ID</b> 10060001
<b>Method</b> EPA 335.2			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10278203
<b>Method</b> EPA 335.4			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10061402
<b>Method</b> EPA 350.3			
<b>Analyte</b> Ammonia as N	<b>AB</b> TX	<b>Analyte ID</b> 1515	<b>Method ID</b> 10064401
<b>Method</b> EPA 365.3			
<b>Analyte</b> Orthophosphate as P Phosphorus	<b>AB</b> TX TX	<b>Analyte ID</b> 1870 1910	<b>Method ID</b> 10070801 10070801
<b>Method</b> EPA 375.4			
<b>Analyte</b> Sulfate	<b>AB</b> TX	<b>Analyte ID</b> 2000	<b>Method ID</b> 10073800
<b>Method</b> EPA 376.1			
<b>Analyte</b> Sulfide	<b>AB</b> TX	<b>Analyte ID</b> 2005	<b>Method ID</b> 10074201
<b>Method</b> EPA 410.4			
<b>Analyte</b> Chemical oxygen demand (COD)	<b>AB</b> TX	<b>Analyte ID</b> 1565	<b>Method ID</b> 10077404
<b>Method</b> EPA 415.1			
<b>Analyte</b> Total Organic Carbon (TOC)	<b>AB</b> TX	<b>Analyte ID</b> 2040	<b>Method ID</b> 10078407
<b>Method</b> EPA 420.1			
<b>Analyte</b> Total phenolics	<b>AB</b> TX	<b>Analyte ID</b> 1905	<b>Method ID</b> 10079400



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**Matrix: Non-Potable Water**

**Method EPA 420.4**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10080203

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156419
Antimony	TX	1005	10156419
Arsenic	TX	1010	10156419
Barium	TX	1015	10156419
Beryllium	TX	1020	10156419
Boron	TX	1025	10156419
Cadmium	TX	1030	10156419
Calcium	TX	1035	10156419
Chromium	TX	1040	10156419
Cobalt	TX	1050	10156419
Copper	TX	1055	10156419
Iron	TX	1070	10156419
Lead	TX	1075	10156419
Lithium	TX	1080	10156419
Magnesium	TX	1085	10156419
Manganese	TX	1090	10156419
Molybdenum	TX	1100	10156419
Nickel	TX	1105	10156419
Potassium	TX	1125	10156419
Selenium	TX	1140	10156419
Silver	TX	1150	10156419
Sodium	TX	1155	10156419
Strontium	TX	1160	10156419
Thallium	TX	1165	10156419
Tin	TX	1175	10156419
Titanium	TX	1180	10156419



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
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Issue Date: 5/1/2019

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**Matrix: Non-Potable Water**

Vanadium	TX	1185	10156419
Zinc	TX	1190	10156419
<b>Method EPA 608</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
4,4'-DDD	TX	7355	10103603
4,4'-DDE	TX	7360	10103603
4,4'-DDT	TX	7365	10103603
Aldrin	TX	7025	10103603
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10103603
alpha-Chlordane	TX	7240	10103603
Aroclor-1016 (PCB-1016)	TX	8880	10103603
Aroclor-1221 (PCB-1221)	TX	8885	10103603
Aroclor-1232 (PCB-1232)	TX	8890	10103603
Aroclor-1242 (PCB-1242)	TX	8895	10103603
Aroclor-1248 (PCB-1248)	TX	8900	10103603
Aroclor-1254 (PCB-1254)	TX	8905	10103603
Aroclor-1260 (PCB-1260)	TX	8910	10103603
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10103603
Chlordane (tech.)	TX	7250	10103603
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10103603
Dieldrin	TX	7470	10103603
Endosulfan I	TX	7510	10103603
Endosulfan II	TX	7515	10103603
Endosulfan sulfate	TX	7520	10103603
Endrin	TX	7540	10103603
Endrin aldehyde	TX	7530	10103603
Endrin ketone	TX	7535	10103603
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10103603
gamma-Chlordane	TX	7245	10103603
Heptachlor	TX	7685	10103603





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**Matrix: Non-Potable Water**

Heptachlor epoxide	TX	7690	10103603
Methoxychlor	TX	7810	10103603
Toxaphene (Chlorinated camphene)	TX	8250	10103603

**Method EPA 624**

Analyte	AB	Analyte ID	Method ID
1,1,1-Trichloroethane	TX	5160	10107207
1,1,2,2-Tetrachloroethane	TX	5110	10107207
1,1,2-Trichloroethane	TX	5165	10107207
1,1-Dichloroethane	TX	4630	10107207
1,1-Dichloroethylene	TX	4640	10107207
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10107207
1,2-Dichlorobenzene	TX	4610	10107207
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10107207
1,2-Dichloropropane	TX	4655	10107207
1,3-Dichlorobenzene	TX	4615	10107207
1,4-Dichlorobenzene	TX	4620	10107207
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10107207
2-Chloroethyl vinyl ether	TX	4500	10107207
Acetone (2-Propanone)	TX	4315	10107207
Acrolein (Propenal)	TX	4325	10107207
Acrylonitrile	TX	4340	10107207
Benzene	TX	4375	10107207
Bromodichloromethane	TX	4395	10107207
Bromoform	TX	4400	10107207
Carbon tetrachloride	TX	4455	10107207
Chlorobenzene	TX	4475	10107207
Chlorodibromomethane	TX	4575	10107207
Chloroethane (Ethyl chloride)	TX	4485	10107207
Chloroform	TX	4505	10107207
cis-1,2-Dichloroethylene	TX	4645	10107207



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**Matrix: Non-Potable Water**

cis-1,3-Dichloropropene	TX	4680	10107207
Ethylbenzene	TX	4765	10107207
m+p-xylene	TX	5240	10107207
Methyl bromide (Bromomethane)	TX	4950	10107207
Methyl chloride (Chloromethane)	TX	4960	10107207
Methyl tert-butyl ether (MTBE)	TX	5000	10107207
Methylene chloride (Dichloromethane)	TX	4975	10107207
Naphthalene	TX	5005	10107207
o-Xylene	TX	5250	10107207
Tetrachloroethylene (Perchloroethylene)	TX	5115	10107207
Toluene	TX	5140	10107207
trans-1,2-Dichloroethylene	TX	4700	10107207
trans-1,3-Dichloropropylene	TX	4685	10107207
Trichloroethene (Trichloroethylene)	TX	5170	10107207
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10107207
Vinyl chloride	TX	5235	10107207
Xylene (total)	TX	5260	10107207

**Method EPA 625**

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10107401
1,2,4-Trichlorobenzene	TX	5155	10107401
1,2-Dichlorobenzene	TX	4610	10107401
1,2-Diphenylhydrazine	TX	6220	10107401
1,3-Dichlorobenzene	TX	4615	10107401
1,4-Dichlorobenzene	TX	4620	10107401
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10107401
2,4,5-Trichlorophenol	TX	6835	10107401
2,4,6-Trichlorophenol	TX	6840	10107401
2,4-Dichlorophenol	TX	6000	10107401
2,4-Dimethylphenol	TX	6130	10107401



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**Matrix: Non-Potable Water**

2,4-Dinitrophenol	TX	6175	10107401
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10107401
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10107401
2-Chloronaphthalene	TX	5795	10107401
2-Chlorophenol	TX	5800	10107401
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10107401
2-Methylphenol (o-Cresol)	TX	6400	10107401
2-Nitrophenol	TX	6490	10107401
3,3'-Dichlorobenzidine	TX	5945	10107401
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10107401
4-Chloro-3-methylphenol	TX	5700	10107401
4-Chlorophenyl phenylether	TX	5825	10107401
4-Methylphenol (p-Cresol)	TX	6410	10107401
4-Nitrophenol	TX	6500	10107401
Acenaphthene	TX	5500	10107401
Acenaphthylene	TX	5505	10107401
Anthracene	TX	5555	10107401
Benzidine	TX	5595	10107401
Benzo(a)anthracene	TX	5575	10107401
Benzo(a)pyrene	TX	5580	10107401
Benzo(b)fluoranthene	TX	5585	10107401
Benzo(g,h,i)perylene	TX	5590	10107401
Benzo(k)fluoranthene	TX	5600	10107401
bis(2-Chloroethoxy)methane	TX	5760	10107401
bis(2-Chloroethyl) ether	TX	5765	10107401
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10107401
Butyl benzyl phthalate	TX	5670	10107401
Chrysene	TX	5855	10107401
Dibenz(a,h) anthracene	TX	5895	10107401
Diethyl phthalate	TX	6070	10107401



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**Matrix: Non-Potable Water**

Dimethyl phthalate	TX	6135	10107401
Di-n-butyl phthalate	TX	5925	10107401
Di-n-octyl phthalate	TX	6200	10107401
Fluoranthene	TX	6265	10107401
Fluorene	TX	6270	10107401
Hexachlorobenzene	TX	6275	10107401
Hexachlorobutadiene	TX	4835	10107401
Hexachlorocyclopentadiene	TX	6285	10107401
Hexachloroethane	TX	4840	10107401
Indeno(1,2,3-cd) pyrene	TX	6315	10107401
Isophorone	TX	6320	10107401
Naphthalene	TX	5005	10107401
Nitrobenzene	TX	5015	10107401
n-Nitrosodiethylamine	TX	6525	10107401
n-Nitrosodimethylamine	TX	6530	10107401
n-Nitrosodi-n-butylamine	TX	5025	10107401
n-Nitrosodi-n-propylamine	TX	6545	10107401
n-Nitrosodiphenylamine	TX	6535	10107401
Pentachlorobenzene	TX	6590	10107401
Pentachlorophenol	TX	6605	10107401
Phenanthrene	TX	6615	10107401
Phenol	TX	6625	10107401
Pyrene	TX	6665	10107401
Pyridine	TX	5095	10107401
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603



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**Matrix: Non-Potable Water**

**Method EPA 8011**

Analyte	AB	Analyte ID	Method ID
1,2,3-Trichloropropane	TX	5180	10173009
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10173009
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10173009

**Method EPA 8015**

Analyte	AB	Analyte ID	Method ID
Diesel range organics (DRO)	TX	9369	10173203
Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402



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**Matrix: Non-Potable Water**

4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402
Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Hexachlorobenzene	TX	6275	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201





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**Matrix: Non-Potable Water**

Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8151**

Analyte	AB	Analyte ID	Method ID
2,4,5-T	TX	8655	10183003
2,4-D	TX	8545	10183003
2,4-DB	TX	8560	10183003
Dalapon	TX	8555	10183003
Dicamba	TX	8595	10183003
Dichloroprop (Dichloroprop, Weedone)	TX	8605	10183003
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10183003
MCPA	TX	7775	10183003
MCPP	TX	7780	10183003
Silvex (2,4,5-TP)	TX	8650	10183003

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404
1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404



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**Matrix: Non-Potable Water**

1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
2-Pentanone	TX	5045	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404
Allyl alcohol	TX	4350	10184404
Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404





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**Matrix: Non-Potable Water**

Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Di-isopropylether (DIPE)	TX	9375	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Ethyl-t-butylether (ETBE) (2-Ethoxy-2-methylpropane)	TX	4770	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404
Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404



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**Matrix: Non-Potable Water**

Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
T-amylmethylether (TAME)	TX	4370	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404
Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

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Issue Date: 5/1/2019

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**Matrix: Non-Potable Water**

Method EPA 8270

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203
2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203



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**Matrix: Non-Potable Water**

2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Dimethyl aminoazobenzene	TX	6105	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrobiphenyl	TX	6480	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Chloro-2-methylaniline	TX	5695	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203



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**Matrix: Non-Potable Water**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Captan	TX	7190	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203



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**Matrix: Non-Potable Water**

Chrysene	TX	5855	10185203
Coumaphos	TX	7315	10185203
Demeton	TX	7390	10185203
Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Dioxathion	TX	7495	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethion	TX	7565	10185203
Ethyl methanesulfonate	TX	6260	10185203
Famphur	TX	7580	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203





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**Matrix: Non-Potable Water**

Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203
Kepone	TX	7740	10185203
Maleic anhydride	TX	6335	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naled	TX	7905	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203



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**Matrix: Non-Potable Water**

Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Phosmet (Imidan)	TX	8000	10185203
Phthalic anhydride	TX	6640	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203
Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Resorcinol	TX	6680	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203
Trifluralin (Treflan)	TX	8295	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209





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**Matrix: Non-Potable Water**

1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209

**Method EPA 8316**

Analyte	AB	Analyte ID	Method ID
Acrylamide	TX	4330	10188202

**Method EPA 8330**

Analyte	AB	Analyte ID	Method ID
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807



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**Matrix: Non-Potable Water**

Methyl-2,4,6-trinitrophenylamine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807
<b>Method EPA 9014</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803
<b>Method EPA 9038</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Sulfate	TX	2000	10196608
<b>Method EPA 9040</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
pH	TX	1900	10196802
<b>Method EPA 9050</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	10198604
<b>Method EPA 9056</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209
<b>Method EPA 9060</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	10200201



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**Matrix: Non-Potable Water**

Method	Analyte	AB	Analyte ID	Method ID
EPA 9065	Total phenolics	TX	1905	10200405
EPA 9066	Total phenolics	TX	1905	10200609
EPA 9250	Chloride	TX	1575	10207202
EPA RSK 175	2-methylpropane (Isobutane)	TX	4942	10212905
	Ethane	TX	4747	10212905
	Ethene	TX	4752	10212905
	Methane	TX	4926	10212905
	n-Butane	TX	5007	10212905
	n-Propane	TX	5029	10212905
HACH 8000	Chemical oxygen demand (COD)	TX	1565	60003001
SM 2120 B	Color	TX	1605	20223807
SM 2310 B (4a)	Acidity, as CaCO <sub>3</sub>	TX	1500	20002806
SM 2320 B	Alkalinity as CaCO <sub>3</sub>	TX	1505	20045005
SM 2340 B		AB	Analyte ID	Method ID



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**Matrix: Non-Potable Water**

Total hardness as CaCO <sub>3</sub>	TX	1755	20046008
<b>Method</b> SM 2510 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	20048004
<b>Method</b> SM 2540 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-total (total solids)	TX	1950	20004608
<b>Method</b> SM 2540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-filterable (TDS)	TX	1955	20049803
<b>Method</b> SM 2540 D			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-nonfilterable (TSS)	TX	1960	20004802
<b>Method</b> SM 3500-Cr B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	20065809
<b>Method</b> SM 4500-Cl F			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total residual chlorine	TX	1940	20080482
<b>Method</b> SM 4500-Cl <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chloride	TX	1575	20019209
<b>Method</b> SM 4500-CN <sup>-</sup> C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20020808
<b>Method</b> SM 4500-CN <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20021209
<b>Method</b> SM 4500-CN <sup>-</sup> G			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	20021607



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**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method SM 4500-H+ B			
Analyte	AB	Analyte ID	Method ID
pH	TX	1900	20104603
Method SM 4500-NH3 D			
Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	20108809
Kjeldahl Nitrogen (Total Kjeldahl Nitrogen-TKN)	TX	1790	20108809
Method SM 4500-NH3 F			
Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	20023001
Method SM 4500-O G			
Analyte	AB	Analyte ID	Method ID
Oxygen, dissolved	TX	1880	20025405
Method SM 4500-P E			
Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	20025803
Phosphorus	TX	1910	20025803
Method SM 4500-S2 <sup>-</sup> F			
Analyte	AB	Analyte ID	Method ID
Sulfide	TX	2005	20126209
Method SM 4500-SiO2 D			
Analyte	AB	Analyte ID	Method ID
Silica as SiO2	TX	1990	20127202
Method SM 4500-SO3 <sup>-</sup> B			
Analyte	AB	Analyte ID	Method ID
Sulfite	TX	2015	20026806
Method SM 5210 B			
Analyte	AB	Analyte ID	Method ID
Biochemical oxygen demand (BOD)	TX	1530	20027401
Carbonaceous BOD, CBOD	TX	1555	20027401
Method SM 5310 B			
Analyte	AB	Analyte ID	Method ID



# Texas Commission on Environmental Quality



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10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

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Issue Date: 5/1/2019

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**Matrix: Non-Potable Water**

Total Organic Carbon (TOC)	TX	2040	20137206
<b>Method</b> SM 5310 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	20138209
<b>Method</b> SM 5540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Surfactants - MBAS	TX	2025	20144405
<b>Method</b> TCEQ 1005			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208





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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
Method ASTM D2216			
Analyte Moisture	TX	10337	ASTM D2216-05
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 1030			
Analyte Ignitability	TX	1780	10117201
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 1668			
Analyte Decachlorobiphenyls	TX	10332	10262007
Dichlorobiphenyls	TX	464	10262007
Heptachlorobiphenyls	TX	486	10262007
Hexachlorobiphenyls	TX	487	10262007
Monochlorobiphenyls	TX	501	10262007
Nonachlorobiphenyls	TX	507	10262007
Octachlorobiphenyls	TX	508	10262007
Pentachlorobiphenyls	TX	515	10262007
Tetrachlorobiphenyls	TX	528	10262007
Trichlorobiphenyls	TX	541	10262007
Method EPA 200.8			
Analyte Uranium	TX	3035	10014605



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**Matrix: Solid & Chemical Materials**

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 310.1**

Analyte	AB	Analyte ID	Method ID
Alkalinity as CaCO3	TX	1505	10054805

**Method EPA 350.3**

Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	10064401

**Method EPA 365.3**

Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	10070801
Phosphorus	TX	1910	10070801

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156204
Antimony	TX	1005	10156204
Arsenic	TX	1010	10156204
Barium	TX	1015	10156204
Beryllium	TX	1020	10156204
Boron	TX	1025	10156204
Cadmium	TX	1030	10156204
Calcium	TX	1035	10156204
Chromium	TX	1040	10156204





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**Matrix: Solid & Chemical Materials**

Cobalt	TX	1050	10156204
Copper	TX	1055	10156204
Iron	TX	1070	10156204
Lead	TX	1075	10156204
Lithium	TX	1080	10156204
Magnesium	TX	1085	10156204
Manganese	TX	1090	10156204
Molybdenum	TX	1100	10156204
Nickel	TX	1105	10156204
Potassium	TX	1125	10156204
Selenium	TX	1140	10156204
Silver	TX	1150	10156204
Sodium	TX	1155	10156204
Strontium	TX	1160	10156204
Thallium	TX	1165	10156204
Tin	TX	1175	10156204
Titanium	TX	1180	10156204
Vanadium	TX	1185	10156204
Zinc	TX	1190	10156204
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603
<b>Method EPA 7471</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10166004
<b>Method EPA 8015</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Diesel range organics (DRO)	TX	9369	10173203



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**Matrix: Solid & Chemical Materials**

Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402
4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402



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**Matrix: Solid & Chemical Materials**

Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201
Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404



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**Matrix: Solid & Chemical Materials**

1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404



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**Matrix: Solid & Chemical Materials**

Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404
Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404



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**Matrix: Solid & Chemical Materials**

Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404
Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404





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**Matrix: Solid & Chemical Materials**

Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404

**Method EPA 8270**

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203



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**Matrix: Solid & Chemical Materials**

2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203
2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203





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**Matrix: Solid & Chemical Materials**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203
Chrysene	TX	5855	10185203



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

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**Matrix: Solid & Chemical Materials**

Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzo(a,e) pyrene	TX	5890	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethyl methanesulfonate	TX	6260	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203
Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203



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**Matrix: Solid & Chemical Materials**

Kepona	TX	7740	10185203
Malathion	TX	7770	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203
Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203



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**Matrix: Solid & Chemical Materials**

Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209



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**Matrix: Solid & Chemical Materials**

Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209
<b>Method EPA 8316</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Acrylamide	TX	4330	10188202
<b>Method EPA 8330</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807
Methyl-2,4,6-trinitrophenylnitramine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807
<b>Method EPA 9014</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803
<b>Method EPA 9038</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Sulfate	TX	2000	10196608



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**Matrix: Solid & Chemical Materials**

**Method EPA 9040**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197203
pH	TX	1900	10196802

**Method EPA 9045**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197805
pH	TX	1900	10197805

**Method EPA 9050**

Analyte	AB	Analyte ID	Method ID
Conductivity	TX	1610	10198604

**Method EPA 9056**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209

**Method EPA 9060**

Analyte	AB	Analyte ID	Method ID
Total Organic Carbon (TOC)	TX	2040	10200201

**Method EPA 9065**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10200405

**Method EPA 9071**

Analyte	AB	Analyte ID	Method ID
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10201204





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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
EPA 9095			
<b>Analyte</b> Paint Filter Liquids Test	TX	10312	10204009
EPA 9250			
<b>Analyte</b> Chloride	TX	1575	10207202
SM 2320 B			
<b>Analyte</b> Alkalinity as CaCO3	TX	1505	20045005
SM 2510 B			
<b>Analyte</b> Conductivity	TX	1610	20048004
SM 2540 G			
<b>Analyte</b> Residue-total (total solids)	TX	1950	20005203
SSA/ASA Part 3:34			
<b>Analyte</b> Carbon, organic (Walkley-Black)	TX	10340	SSA/ASA Pt 3:34
TCEQ 1005			
<b>Analyte</b> Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
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January 20, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010407**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 27 sample(s) on Jan 10, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "Dane J. Wacasey".

Generated By: DAYNA.FISHER  
Dane J. Wacasey



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

**Laboratory Review Checklist: Reportable Data**

Laboratory Name: ALS Laboratory Group			LRC Date: 01/20/2020				
Project Name: Houston TX-Wood Preserving Works			Laboratory Job Number: HS20010407				
Reviewer Name: Dane Wacasey			Prep Batch Number(s): 149585, 149592, 149618, 149624, 149670, R354349, R354396, R354399, R354465				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				3
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data								
Laboratory Name: ALS Laboratory Group					LRC Date: 01/20/2020			
Project Name: Houston TX-Wood Preserving Works					Laboratory Job Number: HS20010407			
Reviewer Name: Dane Wacasey					Prep Batch Number(s): 149585, 149592, 149618, 149624, 149670, R354349, R354396, R354399, R354465			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER <sup>5</sup>	
S1	OI	<b>Initial calibration (ICAL)</b>						
		Were response factors and/or relative response factors for each analyte within QC limits?	X					
		Were percent RSDs or correlation coefficient criteria met?	X					
		Was the number of standards recommended in the method used for all analytes?	X					
		Were all points generated between the lowest and highest standard used to calculate the curve?	X					
		Are ICAL data available for all instruments used?	X					
		Has the initial calibration curve been verified using an appropriate second source standard?	X					
S2	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>						
		Was the CCV analyzed at the method-required frequency?	X					
		Were percent differences for each analyte within the method-required QC limits?	X					
		Was the ICAL curve verified for each analyte?	X					
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X					
S3	O	<b>Mass spectral tuning:</b>						
		Was the appropriate compound for the method used for tuning?	X					
		Were ion abundance data within the method-required QC limits?	X					
S4	O	<b>Internal standards (IS):</b>						
		Were IS area counts and retention times within the method-required QC limits?	X					
S5	OI	<b>Raw data (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section</b>						
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X					
		Were data associated with manual integrations flagged on the raw data?	X					
S6	O	<b>Dual column confirmation</b>						
		Did dual column confirmation results meet the method-required QC?			X			
S7	O	<b>Tentatively identified compounds (TICs):</b>						
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X			
S8	I	<b>Interference Check Sample (ICS) results:</b>						
		Were percent recoveries within method QC limits?	X					
S9	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>						
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X					
S10	OI	<b>Method detection limit (MDL) studies</b>						
		Was a MDL study performed for each reported analyte?	X					
		Is the MDL either adjusted or supported by the analysis of DCSs?	X					
S11	OI	<b>Proficiency test reports:</b>						
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X					
S12	OI	<b>Standards documentation</b>						
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X					
S13	OI	<b>Compound/analyte identification procedures</b>						
		Are the procedures for compound/analyte identification documented?	X					
S14	OI	<b>Demonstration of analyst competency (DOC)</b>						
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X					
		Is documentation of the analyst's competency up-to-date and on file?	X					
S15	OI	<b>Verification/validation documentation for methods (NELAC Chap 5 or ISO/IEC 17025 Section 5)</b>						
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X					
S16	OI	<b>Laboratory standard operating procedures (SOPs):</b>						
		Are laboratory SOPs current and on file for each method performed?	X					

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 01/20/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20010407
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 149585, 149592, 149618, 149624, 149670, R354349, R354396, R354399, R354465

ER# <sup>5</sup>	Description
1	<p>Batch 149624 and 149670, Semivolatiles by Method SW8270, Samples WG-1620-MW77A-20200109, WG-1620-MW49A-20200107, WG-1620-MW18C-20200108, WG-1620-MW18A-20200108, WG-1620-MW58A-20200108, WG-1620-MW57A-20200108, WG-1620-MW57B-20200108, WG-1620-MW72B-20200109: Surrogate recoveries could not be determined due to dilution below the calibration range.</p> <p>Batch 149670, Semivolatiles by Method SW8270, Sample WG-1620-MW23C-20200109, Surrogate Nitrobenzene-d5 recovered above upper control limits. This was due to a dilution required for sample analysis.</p>
2	<p>Batch R354396, Volatiles by Method SW8260, Sample HS20010411-10, MS and MSD were performed on an unrelated sample.</p>
3	<p>Batch R354396, Volatiles by Method SW8260, Samples WG-1620-MW18C-20200108, WG-1620-MW18A-20200108: Lowest practical dilution due to sample matrix.</p> <p>Batch R354465, Volatiles by Method SW8260, Samples WG-1620-MW23C-20200109, WG-1620-MW49A-20200107, WG-1620-MW57B-20200108, WG-1620-MW72B-20200109: Lowest practical dilution due to high concentration of non-target analyte(s).</p> <p>Batch 149670, Semivolatiles by Method SW8270, Sample WG-1620-MW57B-20200108, WG-1620-MW72B-20200109, WG-1620-MW23C-20200109: The GCMS semi-volatile extract of this sample was run at a dilution due to a high level of matrix interference.</p>

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010407

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010407-01	WG-1620-MW19C-20200109	Water		09-Jan-2020 10:45	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-02	WG-1620-MW51A-20200109	Water		09-Jan-2020 11:50	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-03	WG-1620-MW51C-20200109	Water		09-Jan-2020 12:40	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-04	WG-1620-MW85C-20200109	Water		09-Jan-2020 13:35	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-05	WG-1620-MW77A-20200109	Water		09-Jan-2020 14:25	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-06	WG-1620-MW76C-20200109	Water		09-Jan-2020 15:20	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-07	WG-1620-MW81B-20200109	Water		09-Jan-2020 16:20	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-08	WG-1620-MW50A-20200109	Water		09-Jan-2020 17:15	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-09	WG-1620-MW49A-20200107	Water		07-Jan-2020 13:10	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-10	WG-1620-MW80B-20200107	Water		07-Jan-2020 16:00	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-11	WG-1620-FB01-20200107	Water		07-Jan-2020 16:30	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-12	WG-1620-FB02-20200108	Water		08-Jan-2020 09:00	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-13	WG-1620-MW18C-20200108	Water		08-Jan-2020 10:10	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-14	WG-1620-MW18A-20200108	Water		08-Jan-2020 11:10	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-15	WG-1620-MW58A-20200108	Water		08-Jan-2020 12:15	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-16	WG-1620-MW57A-20200108	Water		08-Jan-2020 14:55	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-17	WG-1620-MW57B-20200108	Water		08-Jan-2020 13:30	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-18	WG-1620-FB03-20200109	Water		09-Jan-2020 07:30	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-19	WG-1620-MW49B-20200107	Water		07-Jan-2020 14:30	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-20	WG-1620-MW72B-20200109	Water		09-Jan-2020 08:35	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-21	WG-1620-MW23C-20200109	Water		09-Jan-2020 09:40	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-22	WG-1620-MW53C-20200109	Water		09-Jan-2020 11:15	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-23	WG-1620-MW54C-20200109	Water		09-Jan-2020 12:35	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-24	WG-1620-MW44A-20200109	Water		09-Jan-2020 13:30	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-25	WG-1620-MW36B-20200109	Water		09-Jan-2020 14:40	10-Jan-2020 09:35	<input type="checkbox"/>
HS20010407-26	WG-1620-MW36A-20200109	Water		09-Jan-2020 16:00	10-Jan-2020 09:35	<input type="checkbox"/>

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010407

**SAMPLE SUMMARY**

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Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010407-27	WG-1620-TB01-20200109	Water		09-Jan-2020 00:00	10-Jan-2020 09:35	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW19C-20200109  
 Collection Date: 09-Jan-2020 10:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:06
<b>Benzene</b>	<b>0.00032</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jan-2020 23:06
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:06
<b>Ethylbenzene</b>	<b>0.00064</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jan-2020 23:06
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 23:06
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:06
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:06
<b>Xylenes, Total</b>	<b>0.0012</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Jan-2020 23:06
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.3</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:06</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:06</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:06</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:06</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW19C-20200109  
 Collection Date: 09-Jan-2020 10:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	0.000078	J	0.000021	0.00020	mg/L	1	15-Jan-2020 08:50
2,4-Dimethylphenol		U	0.000040	0.00020	mg/L	1	15-Jan-2020 08:50
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jan-2020 08:50
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jan-2020 08:50
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jan-2020 08:50
2-Methylnaphthalene		U	0.000019	0.00010	mg/L	1	15-Jan-2020 08:50
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jan-2020 08:50
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jan-2020 08:50
Acenaphthene	0.000099		0.000027	0.00010	mg/L	1	15-Jan-2020 08:50
Acenaphthylene	0.000050	J	0.000015	0.00010	mg/L	1	15-Jan-2020 08:50
Anthracene	0.00017		0.000014	0.00010	mg/L	1	15-Jan-2020 08:50
Benz(a)anthracene	0.00026		0.000050	0.00010	mg/L	1	15-Jan-2020 08:50
Benzo(a)pyrene	0.00010		0.000020	0.00010	mg/L	1	15-Jan-2020 08:50
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jan-2020 08:50
Bis(2-ethylhexyl)phthalate	0.000071	J	0.000037	0.00020	mg/L	1	15-Jan-2020 08:50
Chrysene	0.00024		0.000021	0.00010	mg/L	1	15-Jan-2020 08:50
Dibenzofuran	0.00069		0.000020	0.00010	mg/L	1	15-Jan-2020 08:50
Di-n-butyl phthalate	0.000021	J	0.000020	0.00020	mg/L	1	15-Jan-2020 08:50
Fluoranthene	0.0011		0.000010	0.00010	mg/L	1	15-Jan-2020 08:50
Fluorene	0.00047		0.000030	0.00010	mg/L	1	15-Jan-2020 08:50
Naphthalene	0.00011		0.000020	0.00010	mg/L	1	15-Jan-2020 08:50
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jan-2020 08:50
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jan-2020 08:50
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jan-2020 08:50
Phenanthrene	0.00045		0.000021	0.00010	mg/L	1	15-Jan-2020 08:50
Phenol		U	0.000035	0.00020	mg/L	1	15-Jan-2020 08:50
Pyrene	0.0014		0.000019	0.00010	mg/L	1	15-Jan-2020 08:50
Surr: 2,4,6-Tribromophenol	65.3			34-129	%REC	1	15-Jan-2020 08:50
Surr: 2-Fluorobiphenyl	61.5			40-125	%REC	1	15-Jan-2020 08:50
Surr: 2-Fluorophenol	54.7			20-120	%REC	1	15-Jan-2020 08:50
Surr: 4-Terphenyl-d14	82.1			40-135	%REC	1	15-Jan-2020 08:50
Surr: Nitrobenzene-d5	50.1			41-120	%REC	1	15-Jan-2020 08:50
Surr: Phenol-d6	52.1			20-120	%REC	1	15-Jan-2020 08:50
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00140	J	0.000400	0.00200	mg/L	1	13-Jan-2020 22:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51A-20200109  
 Collection Date: 09-Jan-2020 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 01:27
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 01:27
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 01:27
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 01:27
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 01:27
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 01:27
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 01:27
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 01:27</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 01:27</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 01:27</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 01:27</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51A-20200109  
 Collection Date: 09-Jan-2020 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 09:09
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 09:09
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 09:09
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 09:09
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 09:09
<b>2-Methylnaphthalene</b>	<b>0.000044</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 09:09
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 09:09
<b>Acenaphthene</b>	<b>0.000040</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jan-2020 09:09
<b>Anthracene</b>	<b>0.000063</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 09:09
<b>Benzo(a)pyrene</b>	<b>0.000027</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 09:09
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00014</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<b>Chrysene</b>	<b>0.000050</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<b>Dibenzofuran</b>	<b>0.000049</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<b>Di-n-butyl phthalate</b>	<b>0.000034</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<b>Fluoranthene</b>	<b>0.00013</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<b>Fluorene</b>	<b>0.000043</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<b>Naphthalene</b>	<b>0.00032</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 09:09
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 09:09
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 09:09
<b>Phenanthrene</b>	<b>0.00021</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 09:09
<b>Pyrene</b>	<b>0.000083</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:09
<i>Surr: 2,4,6-Tribromophenol</i>	74.3			34-129	%REC	1	15-Jan-2020 09:09
<i>Surr: 2-Fluorobiphenyl</i>	79.5			40-125	%REC	1	15-Jan-2020 09:09
<i>Surr: 2-Fluorophenol</i>	53.3			20-120	%REC	1	15-Jan-2020 09:09
<i>Surr: 4-Terphenyl-d14</i>	91.9			40-135	%REC	1	15-Jan-2020 09:09
<i>Surr: Nitrobenzene-d5</i>	63.9			41-120	%REC	1	15-Jan-2020 09:09
<i>Surr: Phenol-d6</i>	66.7			20-120	%REC	1	15-Jan-2020 09:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51A-20200109  
 Collection Date: 09-Jan-2020 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 13-Jan-2020		Analyst: MBG	
nC6 to nC12	U		0.19	0.48	mg/L	1	13-Jan-2020 21:36
>nC12 to nC28	U		0.19	0.48	mg/L	1	13-Jan-2020 21:36
>nC28 to nC35	U		0.19	0.48	mg/L	1	13-Jan-2020 21:36
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	13-Jan-2020 21:36
Surr: 2-Fluorobiphenyl	104			70-130	%REC	1	13-Jan-2020 21:36
Surr: Trifluoromethyl benzene	106			70-130	%REC	1	13-Jan-2020 21:36
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00188	J	0.000400	0.00200	mg/L	1	13-Jan-2020 22:25

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51C-20200109  
 Collection Date: 09-Jan-2020 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 03:29
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 03:29
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 03:29
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 03:29
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 03:29
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 03:29
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 03:29
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:29</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:29</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:29</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:29</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51C-20200109  
 Collection Date: 09-Jan-2020 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 09:28
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 09:28
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 09:28
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 09:28
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 09:28
<b>2-Methylnaphthalene</b>	<b>0.000055</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 09:28
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 09:28
<b>Acenaphthene</b>	<b>0.000094</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jan-2020 09:28
<b>Anthracene</b>	<b>0.00010</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 09:28
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 09:28
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 09:28
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000085</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 09:28
<b>Dibenzofuran</b>	<b>0.00011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 09:28
<b>Fluoranthene</b>	<b>0.00018</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
<b>Fluorene</b>	<b>0.00011</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
<b>Naphthalene</b>	<b>0.00030</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 09:28
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 09:28
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 09:28
<b>Phenanthrene</b>	<b>0.00056</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 09:28
<b>Pyrene</b>	<b>0.00011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:28
<i>Surr: 2,4,6-Tribromophenol</i>	<i>64.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:28</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>73.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:28</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:28</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>80.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:28</i>
<i>Surr: Nitrobenzene-d5</i>	<i>55.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:28</i>
<i>Surr: Phenol-d6</i>	<i>61.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:28</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW51C-20200109  
 Collection Date: 09-Jan-2020 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>			Prep:TX1005PR / 13-Jan-2020		Analyst: MBG
nC6 to nC12	U		0.19	0.48	mg/L	1	13-Jan-2020 22:05
>nC12 to nC28	U		0.19	0.48	mg/L	1	13-Jan-2020 22:05
>nC28 to nC35	U		0.19	0.48	mg/L	1	13-Jan-2020 22:05
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	13-Jan-2020 22:05
Surr: 2-Fluorobiphenyl	108			70-130	%REC	1	13-Jan-2020 22:05
Surr: Trifluoromethyl benzene	110			70-130	%REC	1	13-Jan-2020 22:05
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 13-Jan-2020		Analyst: JHD
Arsenic	0.000452	J	0.000400	0.00200	mg/L	1	13-Jan-2020 22:27

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW85C-20200109  
 Collection Date: 09-Jan-2020 13:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:31
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:31
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:31
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:31
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 23:31
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:31
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:31
<i>Surr: 1,2-Dichloroethane-d4</i>		96.4		70-126	%REC	1	14-Jan-2020 23:31
<i>Surr: 4-Bromofluorobenzene</i>		95.7		81-113	%REC	1	14-Jan-2020 23:31
<i>Surr: Dibromofluoromethane</i>		102		77-123	%REC	1	14-Jan-2020 23:31
<i>Surr: Toluene-d8</i>		101		82-127	%REC	1	14-Jan-2020 23:31

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW85C-20200109  
 Collection Date: 09-Jan-2020 13:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 09:46
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 09:46
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 09:46
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 09:46
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 09:46
<b>2-Methylnaphthalene</b>	<b>0.000041</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 09:46
<b>4-Nitrophenol</b>	<b>0.00077</b>	J	<b>0.000047</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Acenaphthene</b>	<b>0.000077</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jan-2020 09:46
Anthracene	U		0.000014	0.00010	mg/L	1	15-Jan-2020 09:46
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 09:46
<b>Benzo(a)pyrene</b>	<b>0.000031</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 09:46
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00016</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Chrysene</b>	<b>0.000053</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Dibenzofuran</b>	<b>0.000059</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 09:46
<b>Fluoranthene</b>	<b>0.00025</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Fluorene</b>	<b>0.000055</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Naphthalene</b>	<b>0.00046</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 09:46
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 09:46
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 09:46
<b>Phenanthrene</b>	<b>0.000067</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Phenol</b>	<b>0.000042</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<b>Pyrene</b>	<b>0.00021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 09:46
<i>Surr: 2,4,6-Tribromophenol</i>	<i>84.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:46</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>78.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:46</i>
<i>Surr: 2-Fluorophenol</i>	<i>74.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:46</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>98.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:46</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:46</i>
<i>Surr: Phenol-d6</i>	<i>69.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 09:46</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW85C-20200109  
 Collection Date: 09-Jan-2020 13:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>			Prep:TX1005PR / 13-Jan-2020		Analyst: MBG
nC6 to nC12	U		0.20	0.49	mg/L	1	13-Jan-2020 22:34
>nC12 to nC28	U		0.20	0.49	mg/L	1	13-Jan-2020 22:34
>nC28 to nC35	U		0.20	0.49	mg/L	1	13-Jan-2020 22:34
Total Petroleum Hydrocarbon	U		0.20	0.49	mg/L	1	13-Jan-2020 22:34
Surr: 2-Fluorobiphenyl	108			70-130	%REC	1	13-Jan-2020 22:34
Surr: Trifluoromethyl benzene	108			70-130	%REC	1	13-Jan-2020 22:34
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 13-Jan-2020		Analyst: JHD
Arsenic	0.00272		0.000400	0.00200	mg/L	1	13-Jan-2020 22:29

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW77A-20200109  
 Collection Date: 09-Jan-2020 14:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:58
<b>Benzene</b>	<b>0.058</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 12:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:58
<b>Ethylbenzene</b>	<b>0.063</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 12:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 12:58
<b>Toluene</b>	<b>0.0069</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 12:58
<b>Xylenes, Total</b>	<b>0.065</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 12:58
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:58</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>103</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:58</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:58</i>
<i>Surr: Toluene-d8</i>	<i>99.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW77A-20200109  
 Collection Date: 09-Jan-2020 14:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 10:05
<b>2,4-Dimethylphenol</b>	<b>0.055</b>		<b>0.00080</b>	<b>0.0040</b>	<b>mg/L</b>	20	15-Jan-2020 14:11
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 10:05
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 10:05
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 10:05
<b>2-Methylnaphthalene</b>	<b>0.22</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jan-2020 16:04
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 10:05
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 10:05
<b>Acenaphthene</b>	<b>0.13</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 14:11
<b>Acenaphthylene</b>	<b>0.0013</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:05
<b>Anthracene</b>	<b>0.0036</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:05
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 10:05
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 10:05
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 10:05
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000077</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 10:05
<b>Chrysene</b>	<b>0.000041</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:05
<b>Dibenzofuran</b>	<b>0.079</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 14:11
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 10:05
<b>Fluoranthene</b>	<b>0.00071</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:05
<b>Fluorene</b>	<b>0.067</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 14:11
<b>Naphthalene</b>	<b>4.7</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	15-Jan-2020 17:57
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 10:05
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 10:05
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 10:05
<b>Phenanthrene</b>	<b>0.031</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 14:11
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 10:05
<b>Pyrene</b>	<b>0.00046</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:05
<i>Surr: 2,4,6-Tribromophenol</i>	<i>90.3</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>15-Jan-2020 16:04</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>15-Jan-2020 17:57</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:05</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>96.4</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 14:11</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>80.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:05</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>96.1</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 14:11</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>101</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>15-Jan-2020 16:04</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>15-Jan-2020 17:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>79.1</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 14:11</i>
<i>Surr: 2-Fluorophenol</i>	<i>110</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>15-Jan-2020 16:04</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>15-Jan-2020 17:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:05</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW77A-20200109  
 Collection Date: 09-Jan-2020 14:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	102			40-135	%REC	1	15-Jan-2020 10:05
Surr: 4-Terphenyl-d14	112			40-135	%REC	20	15-Jan-2020 14:11
Surr: 4-Terphenyl-d14	115	J		40-135	%REC	100	15-Jan-2020 16:04
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	15-Jan-2020 17:57
Surr: Nitrobenzene-d5	93.9			41-120	%REC	20	15-Jan-2020 14:11
Surr: Nitrobenzene-d5	66.8	J		41-120	%REC	100	15-Jan-2020 16:04
Surr: Nitrobenzene-d5	95.0			41-120	%REC	1	15-Jan-2020 10:05
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	15-Jan-2020 17:57
Surr: Phenol-d6	72.0			20-120	%REC	1	15-Jan-2020 10:05
Surr: Phenol-d6	0	JS		20-120	%REC	1000	15-Jan-2020 17:57
Surr: Phenol-d6	77.0	J		20-120	%REC	20	15-Jan-2020 14:11
Surr: Phenol-d6	60.7	J		20-120	%REC	100	15-Jan-2020 16:04
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 13-Jan-2020		Analyst: MBG	
nC6 to nC12	8.3		0.19	0.48	mg/L	1	13-Jan-2020 23:03
>nC12 to nC28	3.8		0.19	0.48	mg/L	1	13-Jan-2020 23:03
>nC28 to nC35	U		0.19	0.48	mg/L	1	13-Jan-2020 23:03
<b>Total Petroleum Hydrocarbon</b>	<b>12.1</b>		<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	<b>1</b>	<b>13-Jan-2020 23:03</b>
Surr: 2-Fluorobiphenyl	108			70-130	%REC	1	13-Jan-2020 23:03
Surr: Trifluoromethyl benzene	108			70-130	%REC	1	13-Jan-2020 23:03
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.0237		0.000400	0.00200	mg/L	1	13-Jan-2020 22:31

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76C-20200109  
 Collection Date: 09-Jan-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:55
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:55
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:55
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:55
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 23:55
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 23:55
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 23:55
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:55</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:55</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:55</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 23:55</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76C-20200109  
 Collection Date: 09-Jan-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jan-2020 10:24
<b>2,4-Dimethylphenol</b>	<b>0.00038</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jan-2020 10:24
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jan-2020 10:24
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jan-2020 10:24
<b>2-Methylnaphthalene</b>	<b>0.00045</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jan-2020 10:24
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jan-2020 10:24
<b>Acenaphthene</b>	<b>0.00041</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Acenaphthylene</b>	<b>0.000035</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Anthracene</b>	<b>0.00017</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Benz(a)anthracene</b>	<b>0.00014</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Benzo(a)pyrene</b>	<b>0.000037</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jan-2020 10:24
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00016</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Chrysene</b>	<b>0.00013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Dibenzofuran</b>	<b>0.00029</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jan-2020 10:24
<b>Fluoranthene</b>	<b>0.0010</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Fluorene</b>	<b>0.00032</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<b>Naphthalene</b>	<b>0.0038</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jan-2020 10:24
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jan-2020 10:24
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jan-2020 10:24
<b>Phenanthrene</b>	<b>0.00069</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
Phenol		U	0.000035	0.00020	mg/L	1	15-Jan-2020 10:24
<b>Pyrene</b>	<b>0.00070</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:24
<i>Surr: 2,4,6-Tribromophenol</i>	<i>78.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>72.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>52.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:24</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:24</i>
<i>Surr: Nitrobenzene-d5</i>	<i>56.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:24</i>
<i>Surr: Phenol-d6</i>	<i>58.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 10:24</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76C-20200109  
 Collection Date: 09-Jan-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 13-Jan-2020		Analyst: MBG	
nC6 to nC12	U		0.19	0.48	mg/L	1	13-Jan-2020 23:33
>nC12 to nC28	U		0.19	0.48	mg/L	1	13-Jan-2020 23:33
>nC28 to nC35	U		0.19	0.48	mg/L	1	13-Jan-2020 23:33
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	13-Jan-2020 23:33
Surr: 2-Fluorobiphenyl	116			70-130	%REC	1	13-Jan-2020 23:33
Surr: Trifluoromethyl benzene	116			70-130	%REC	1	13-Jan-2020 23:33
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00299		0.000400	0.00200	mg/L	1	13-Jan-2020 22:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW81B-20200109  
 Collection Date: 09-Jan-2020 16:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 03:54
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 03:54
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 03:54
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 03:54
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 03:54
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 03:54
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 03:54
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:54</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:54</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:54</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 03:54</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW81B-20200109  
 Collection Date: 09-Jan-2020 16:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 10:43
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 10:43
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 10:43
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 10:43
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 10:43
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	15-Jan-2020 10:43
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 10:43
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 10:43
<b>Acenaphthene</b>	<b>0.000077</b>	<b>J</b>	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:43
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jan-2020 10:43
Anthracene	U		0.000014	0.00010	mg/L	1	15-Jan-2020 10:43
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 10:43
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 10:43
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 10:43
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000066</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 10:43
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 10:43
Dibenzofuran	U		0.000020	0.00010	mg/L	1	15-Jan-2020 10:43
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 10:43
Fluoranthene	U		0.000010	0.00010	mg/L	1	15-Jan-2020 10:43
Fluorene	U		0.000030	0.00010	mg/L	1	15-Jan-2020 10:43
<b>Naphthalene</b>	<b>0.00023</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 10:43
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 10:43
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 10:43
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 10:43
Phenanthrene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 10:43
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 10:43
Pyrene	U		0.000019	0.00010	mg/L	1	15-Jan-2020 10:43
<i>Surr: 2,4,6-Tribromophenol</i>	65.6			34-129	%REC	1	15-Jan-2020 10:43
<i>Surr: 2-Fluorobiphenyl</i>	65.2			40-125	%REC	1	15-Jan-2020 10:43
<i>Surr: 2-Fluorophenol</i>	63.5			20-120	%REC	1	15-Jan-2020 10:43
<i>Surr: 4-Terphenyl-d14</i>	85.7			40-135	%REC	1	15-Jan-2020 10:43
<i>Surr: Nitrobenzene-d5</i>	53.1			41-120	%REC	1	15-Jan-2020 10:43
<i>Surr: Phenol-d6</i>	56.4			20-120	%REC	1	15-Jan-2020 10:43

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW81B-20200109  
 Collection Date: 09-Jan-2020 16:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>			Prep:TX1005PR / 13-Jan-2020		Analyst: MBG
nC6 to nC12	U		0.19	0.48	mg/L	1	14-Jan-2020 00:02
>nC12 to nC28	U		0.19	0.48	mg/L	1	14-Jan-2020 00:02
>nC28 to nC35	U		0.19	0.48	mg/L	1	14-Jan-2020 00:02
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	14-Jan-2020 00:02
Surr: 2-Fluorobiphenyl	105			70-130	%REC	1	14-Jan-2020 00:02
Surr: Trifluoromethyl benzene	106			70-130	%REC	1	14-Jan-2020 00:02
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 13-Jan-2020		Analyst: JHD
Arsenic	0.00130	J	0.000400	0.00200	mg/L	1	13-Jan-2020 22:36

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50A-20200109  
 Collection Date: 09-Jan-2020 17:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 04:18
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 04:18
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 04:18
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 04:18
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 04:18
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 04:18
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 04:18
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 04:18</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 04:18</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 04:18</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 04:18</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50A-20200109  
 Collection Date: 09-Jan-2020 17:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jan-2020 11:02
<b>2,4-Dimethylphenol</b>	<b>0.00071</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jan-2020 11:02
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jan-2020 11:02
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jan-2020 11:02
<b>2-Methylnaphthalene</b>	<b>0.00019</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jan-2020 11:02
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jan-2020 11:02
<b>Acenaphthene</b>	<b>0.00037</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
<b>Acenaphthylene</b>	<b>0.000022</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
Anthracene		U	0.000014	0.00010	mg/L	1	15-Jan-2020 11:02
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	15-Jan-2020 11:02
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	15-Jan-2020 11:02
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jan-2020 11:02
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000085</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
Chrysene		U	0.000021	0.00010	mg/L	1	15-Jan-2020 11:02
<b>Dibenzofuran</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jan-2020 11:02
<b>Fluoranthene</b>	<b>0.000052</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
<b>Fluorene</b>	<b>0.00011</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
<b>Naphthalene</b>	<b>0.0020</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jan-2020 11:02
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jan-2020 11:02
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jan-2020 11:02
<b>Phenanthrene</b>	<b>0.000070</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
<b>Phenol</b>	<b>0.000066</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
<b>Pyrene</b>	<b>0.000034</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:02
<i>Surr: 2,4,6-Tribromophenol</i>	83.6			34-129	%REC	1	15-Jan-2020 11:02
<i>Surr: 2-Fluorobiphenyl</i>	81.2			40-125	%REC	1	15-Jan-2020 11:02
<i>Surr: 2-Fluorophenol</i>	57.4			20-120	%REC	1	15-Jan-2020 11:02
<i>Surr: 4-Terphenyl-d14</i>	93.2			40-135	%REC	1	15-Jan-2020 11:02
<i>Surr: Nitrobenzene-d5</i>	64.4			41-120	%REC	1	15-Jan-2020 11:02
<i>Surr: Phenol-d6</i>	63.4			20-120	%REC	1	15-Jan-2020 11:02

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50A-20200109  
 Collection Date: 09-Jan-2020 17:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>			Prep:TX1005PR / 13-Jan-2020		Analyst: MBG
nC6 to nC12	U		0.19	0.48	mg/L	1	14-Jan-2020 00:31
>nC12 to nC28	U		0.19	0.48	mg/L	1	14-Jan-2020 00:31
>nC28 to nC35	U		0.19	0.48	mg/L	1	14-Jan-2020 00:31
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	14-Jan-2020 00:31
Surr: 2-Fluorobiphenyl	103			70-130	%REC	1	14-Jan-2020 00:31
Surr: Trifluoromethyl benzene	103			70-130	%REC	1	14-Jan-2020 00:31
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 13-Jan-2020		Analyst: JHD
Arsenic	0.000718	J	0.000400	0.00200	mg/L	1	13-Jan-2020 22:38

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49A-20200107  
 Collection Date: 07-Jan-2020 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.0020	0.010	mg/L	10	15-Jan-2020 14:41
<b>Benzene</b>	<b>0.30</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 14:41
<b>Chlorobenzene</b>	<b>0.0063</b>	J	<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 14:41
<b>Ethylbenzene</b>	<b>0.10</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 14:41
Methylene chloride		U	0.010	0.020	mg/L	10	15-Jan-2020 14:41
<b>Toluene</b>	<b>0.19</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 14:41
Vinyl chloride		U	0.0020	0.010	mg/L	10	15-Jan-2020 14:41
<b>Xylenes, Total</b>	<b>0.24</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 14:41
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.4</i>			<i>70-126</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 14:41</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.4</i>			<i>81-113</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 14:41</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.4</i>			<i>77-123</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 14:41</i>
<i>Surr: Toluene-d8</i>	<i>98.8</i>			<i>82-127</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 14:41</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49A-20200107  
 Collection Date: 07-Jan-2020 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 11:21
<b>2,4-Dimethylphenol</b>	<b>5.7</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	1000	15-Jan-2020 18:16
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 11:21
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 11:21
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 11:21
<b>2-Methylnaphthalene</b>	<b>0.13</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jan-2020 16:23
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 11:21
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 11:21
<b>Acenaphthene</b>	<b>0.057</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:30
<b>Acenaphthylene</b>	<b>0.00097</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:21
<b>Anthracene</b>	<b>0.0028</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:21
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 11:21
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 11:21
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 11:21
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:21
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 11:21
<b>Dibenzofuran</b>	<b>0.037</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:30
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 11:21
<b>Fluoranthene</b>	<b>0.00056</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:21
<b>Fluorene</b>	<b>0.027</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:30
<b>Naphthalene</b>	<b>4.5</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	15-Jan-2020 18:16
Nitrobenzene	U		0.00024	0.0020	mg/L	10	15-Jan-2020 14:30
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 11:21
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 11:21
<b>Phenanthrene</b>	<b>0.016</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:30
<b>Phenol</b>	<b>0.17</b>		<b>0.0035</b>	<b>0.020</b>	<b>mg/L</b>	100	15-Jan-2020 16:23
<b>Pyrene</b>	<b>0.00052</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:21
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	1000	15-Jan-2020 18:16
Surr: 2,4,6-Tribromophenol	69.4			34-129	%REC	1	15-Jan-2020 11:21
Surr: 2,4,6-Tribromophenol	101			34-129	%REC	10	15-Jan-2020 14:30
Surr: 2,4,6-Tribromophenol	86.8	J		34-129	%REC	100	15-Jan-2020 16:23
Surr: 2-Fluorobiphenyl	59.3			40-125	%REC	1	15-Jan-2020 11:21
Surr: 2-Fluorobiphenyl	107			40-125	%REC	10	15-Jan-2020 14:30
Surr: 2-Fluorobiphenyl	107	J		40-125	%REC	100	15-Jan-2020 16:23
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	1000	15-Jan-2020 18:16
Surr: 2-Fluorophenol	75.3			20-120	%REC	1	15-Jan-2020 11:21
Surr: 2-Fluorophenol	83.0			20-120	%REC	10	15-Jan-2020 14:30
Surr: 2-Fluorophenol	86.4	J		20-120	%REC	100	15-Jan-2020 16:23
Surr: 2-Fluorophenol	0	JS		20-120	%REC	1000	15-Jan-2020 18:16

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49A-20200107  
 Collection Date: 07-Jan-2020 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	84.6			40-135	%REC	1	15-Jan-2020 11:21
Surr: 4-Terphenyl-d14	112			40-135	%REC	10	15-Jan-2020 14:30
Surr: 4-Terphenyl-d14	120	J		40-135	%REC	100	15-Jan-2020 16:23
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	15-Jan-2020 18:16
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	15-Jan-2020 18:16
Surr: Nitrobenzene-d5	55.3			41-120	%REC	1	15-Jan-2020 11:21
Surr: Nitrobenzene-d5	65.6			41-120	%REC	10	15-Jan-2020 14:30
Surr: Nitrobenzene-d5	65.6	J		41-120	%REC	100	15-Jan-2020 16:23
Surr: Phenol-d6	66.3			20-120	%REC	1	15-Jan-2020 11:21
Surr: Phenol-d6	87.2			20-120	%REC	10	15-Jan-2020 14:30
Surr: Phenol-d6	82.2	J		20-120	%REC	100	15-Jan-2020 16:23
Surr: Phenol-d6	0	JS		20-120	%REC	1000	15-Jan-2020 18:16
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 13-Jan-2020		Analyst: MBG	
nC6 to nC12	21		0.19	0.48	mg/L	1	14-Jan-2020 01:00
>nC12 to nC28	11		0.19	0.48	mg/L	1	14-Jan-2020 01:00
>nC28 to nC35	U		0.19	0.48	mg/L	1	14-Jan-2020 01:00
<b>Total Petroleum Hydrocarbon</b>	<b>32.0</b>		<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	14-Jan-2020 01:00
Surr: 2-Fluorobiphenyl	112			70-130	%REC	1	14-Jan-2020 01:00
Surr: Trifluoromethyl benzene	113			70-130	%REC	1	14-Jan-2020 01:00
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00463		0.000400	0.00200	mg/L	1	13-Jan-2020 22:40

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW80B-20200107  
 Collection Date: 07-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 04:42
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 04:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 04:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 04:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 04:42
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 04:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 04:42
<i>Surr: 1,2-Dichloroethane-d4</i>	98.3			70-126	%REC	1	14-Jan-2020 04:42
<i>Surr: 4-Bromofluorobenzene</i>	95.5			81-113	%REC	1	14-Jan-2020 04:42
<i>Surr: Dibromofluoromethane</i>	102			77-123	%REC	1	14-Jan-2020 04:42
<i>Surr: Toluene-d8</i>	99.7			82-127	%REC	1	14-Jan-2020 04:42

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW80B-20200107  
 Collection Date: 07-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	15-Jan-2020 11:40
<b>2,4-Dimethylphenol</b>	<b>0.00028</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	15-Jan-2020 11:40
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	15-Jan-2020 11:40
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	15-Jan-2020 11:40
<b>2-Methylnaphthalene</b>	<b>0.000036</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	15-Jan-2020 11:40
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	15-Jan-2020 11:40
<b>Acenaphthene</b>	<b>0.000046</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
Acenaphthylene		U	0.000015	0.00010	mg/L	1	15-Jan-2020 11:40
<b>Anthracene</b>	<b>0.000026</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	15-Jan-2020 11:40
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	15-Jan-2020 11:40
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	15-Jan-2020 11:40
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000069</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
Chrysene		U	0.000021	0.00010	mg/L	1	15-Jan-2020 11:40
<b>Dibenzofuran</b>	<b>0.000048</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	15-Jan-2020 11:40
Fluoranthene		U	0.000010	0.00010	mg/L	1	15-Jan-2020 11:40
<b>Fluorene</b>	<b>0.000032</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
<b>Naphthalene</b>	<b>0.00034</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
Nitrobenzene		U	0.000024	0.00020	mg/L	1	15-Jan-2020 11:40
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	15-Jan-2020 11:40
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	15-Jan-2020 11:40
<b>Phenanthrene</b>	<b>0.000027</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:40
Phenol		U	0.000035	0.00020	mg/L	1	15-Jan-2020 11:40
Pyrene		U	0.000019	0.00010	mg/L	1	15-Jan-2020 11:40
<i>Surr: 2,4,6-Tribromophenol</i>	<i>85.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>74.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>62.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:40</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>93.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:40</i>
<i>Surr: Nitrobenzene-d5</i>	<i>56.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:40</i>
<i>Surr: Phenol-d6</i>	<i>63.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW80B-20200107  
 Collection Date: 07-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 13-Jan-2020		Analyst: MBG	
nC6 to nC12	U		0.19	0.48	mg/L	1	14-Jan-2020 01:29
>nC12 to nC28	U		0.19	0.48	mg/L	1	14-Jan-2020 01:29
>nC28 to nC35	U		0.19	0.48	mg/L	1	14-Jan-2020 01:29
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	14-Jan-2020 01:29
Surr: 2-Fluorobiphenyl	105			70-130	%REC	1	14-Jan-2020 01:29
Surr: Trifluoromethyl benzene	104			70-130	%REC	1	14-Jan-2020 01:29
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00212		0.000400	0.00200	mg/L	1	13-Jan-2020 22:42

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB01-20200107  
 Collection Date: 07-Jan-2020 16:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:30
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:30
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 17:30
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 17:30
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 17:30
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:30
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:30
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 17:30
<i>Surr: 1,2-Dichloroethane-d4</i>		116		70-126	%REC	1	14-Jan-2020 17:30
<i>Surr: 4-Bromofluorobenzene</i>		103		81-113	%REC	1	14-Jan-2020 17:30
<i>Surr: Dibromofluoromethane</i>		102		77-123	%REC	1	14-Jan-2020 17:30
<i>Surr: Toluene-d8</i>		102		82-127	%REC	1	14-Jan-2020 17:30

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB01-20200107  
 Collection Date: 07-Jan-2020 16:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
<b>1,2-Diphenylhydrazine</b>	<b>0.000041</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:59
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 11:59
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 11:59
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 11:59
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 11:59
<b>2-Methylnaphthalene</b>	<b>0.000031</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:59
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 11:59
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 11:59
Acenaphthene	U		0.000027	0.00010	mg/L	1	15-Jan-2020 11:59
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jan-2020 11:59
Anthracene	U		0.000014	0.00010	mg/L	1	15-Jan-2020 11:59
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 11:59
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 11:59
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 11:59
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00017</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 11:59
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 11:59
Dibenzofuran	U		0.000020	0.00010	mg/L	1	15-Jan-2020 11:59
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 11:59
Fluoranthene	U		0.000010	0.00010	mg/L	1	15-Jan-2020 11:59
Fluorene	U		0.000030	0.00010	mg/L	1	15-Jan-2020 11:59
<b>Naphthalene</b>	<b>0.00019</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 11:59
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 11:59
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 11:59
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 11:59
Phenanthrene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 11:59
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 11:59
Pyrene	U		0.000019	0.00010	mg/L	1	15-Jan-2020 11:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>77.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>86.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>77.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>68.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:59</i>
<i>Surr: Phenol-d6</i>	<i>73.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 11:59</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	13-Jan-2020 22:45

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB02-20200108  
 Collection Date: 08-Jan-2020 09:00

**ANALYTICAL REPORT**

WorkOrder:HS20010407  
 Lab ID:HS20010407-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:55
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:55
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 17:55
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 17:55
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 17:55
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:55
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 17:55
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 17:55
<i>Surr: 1,2-Dichloroethane-d4</i>		118		70-126	%REC	1	14-Jan-2020 17:55
<i>Surr: 4-Bromofluorobenzene</i>		101		81-113	%REC	1	14-Jan-2020 17:55
<i>Surr: Dibromofluoromethane</i>		102		77-123	%REC	1	14-Jan-2020 17:55
<i>Surr: Toluene-d8</i>		99.3		82-127	%REC	1	14-Jan-2020 17:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB02-20200108  
 Collection Date: 08-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG
<b>1,2-Diphenylhydrazine</b>	<b>0.000044</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 12:17
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 12:17
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 12:17
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 12:17
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 12:17
<b>2-Methylnaphthalene</b>	<b>0.000038</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:17
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 12:17
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 12:17
<b>Acenaphthene</b>	<b>0.000053</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:17
Acenaphthylene	U		0.000015	0.00010	mg/L	1	15-Jan-2020 12:17
Anthracene	U		0.000014	0.00010	mg/L	1	15-Jan-2020 12:17
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 12:17
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 12:17
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 12:17
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000099</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 12:17
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 12:17
Dibenzofuran	U		0.000020	0.00010	mg/L	1	15-Jan-2020 12:17
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 12:17
Fluoranthene	U		0.000010	0.00010	mg/L	1	15-Jan-2020 12:17
Fluorene	U		0.000030	0.00010	mg/L	1	15-Jan-2020 12:17
<b>Naphthalene</b>	<b>0.00030</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:17
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 12:17
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 12:17
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 12:17
Phenanthrene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 12:17
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 12:17
Pyrene	U		0.000019	0.00010	mg/L	1	15-Jan-2020 12:17
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>84.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:17</i>
<i>Surr: 2-Fluorophenol</i>	<i>75.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:17</i>
<i>Surr: Nitrobenzene-d5</i>	<i>71.7</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:17</i>
<i>Surr: Phenol-d6</i>	<i>70.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:17</i>
<b>ICP-MS METALS BY SW6020A</b>			<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD
Arsenic	U		0.000400	0.00200	mg/L	1	13-Jan-2020 22:58

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18C-20200108  
 Collection Date: 08-Jan-2020 10:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0020	0.010	mg/L	10	14-Jan-2020 17:22
<b>Benzene</b>	<b>1.1</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jan-2020 17:22
Chlorobenzene	U		0.0030	0.010	mg/L	10	14-Jan-2020 17:22
<b>Ethylbenzene</b>	<b>0.31</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jan-2020 17:22
Methylene chloride	U		0.010	0.020	mg/L	10	14-Jan-2020 17:22
<b>Toluene</b>	<b>0.97</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jan-2020 17:22
Vinyl chloride	U		0.0020	0.010	mg/L	10	14-Jan-2020 17:22
<b>Xylenes, Total</b>	<b>0.92</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	14-Jan-2020 17:22
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.1</i>			<i>70-126</i>	<i>%REC</i>	<i>10</i>	<i>14-Jan-2020 17:22</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.9</i>			<i>81-113</i>	<i>%REC</i>	<i>10</i>	<i>14-Jan-2020 17:22</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>10</i>	<i>14-Jan-2020 17:22</i>
<i>Surr: Toluene-d8</i>	<i>98.9</i>			<i>82-127</i>	<i>%REC</i>	<i>10</i>	<i>14-Jan-2020 17:22</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18C-20200108  
 Collection Date: 08-Jan-2020 10:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 12:36
<b>2,4-Dimethylphenol</b>	<b>0.023</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 12:36
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 12:36
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 12:36
<b>2-Methylnaphthalene</b>	<b>0.31</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	15-Jan-2020 16:41
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 12:36
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 12:36
<b>Acenaphthene</b>	<b>0.093</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
<b>Acenaphthylene</b>	<b>0.0026</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:36
<b>Anthracene</b>	<b>0.020</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
<b>Benz(a)anthracene</b>	<b>0.0025</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:36
<b>Benzo(a)pyrene</b>	<b>0.00053</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:36
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 12:36
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	15-Jan-2020 12:36
<b>Chrysene</b>	<b>0.0027</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:36
<b>Dibenzofuran</b>	<b>0.085</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 12:36
<b>Fluoranthene</b>	<b>0.013</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
<b>Fluorene</b>	<b>0.043</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
<b>Naphthalene</b>	<b>14</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	15-Jan-2020 18:35
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 12:36
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 12:36
<b>Pentachlorophenol</b>	<b>0.023</b>		<b>0.00079</b>	<b>0.0020</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
<b>Phenanthrene</b>	<b>0.065</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	15-Jan-2020 14:48
<b>Phenol</b>	<b>0.0062</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 12:36
<b>Pyrene</b>	<b>0.0066</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:36
Surr: 2,4,6-Tribromophenol	66.4			34-129	%REC	1	15-Jan-2020 12:36
Surr: 2,4,6-Tribromophenol	86.5			34-129	%REC	10	15-Jan-2020 14:48
Surr: 2,4,6-Tribromophenol	112	J		34-129	%REC	100	15-Jan-2020 16:41
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	2000	15-Jan-2020 18:35
Surr: 2-Fluorobiphenyl	64.7			40-125	%REC	1	15-Jan-2020 12:36
Surr: 2-Fluorobiphenyl	92.7			40-125	%REC	10	15-Jan-2020 14:48
Surr: 2-Fluorobiphenyl	108	J		40-125	%REC	100	15-Jan-2020 16:41
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	2000	15-Jan-2020 18:35
Surr: 2-Fluorophenol	97.6			20-120	%REC	1	15-Jan-2020 12:36
Surr: 2-Fluorophenol	56.1			20-120	%REC	10	15-Jan-2020 14:48
Surr: 2-Fluorophenol	71.0	J		20-120	%REC	100	15-Jan-2020 16:41
Surr: 2-Fluorophenol	0	JS		20-120	%REC	2000	15-Jan-2020 18:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18C-20200108  
 Collection Date: 08-Jan-2020 10:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	70.5			40-135	%REC	1	15-Jan-2020 12:36
Surr: 4-Terphenyl-d14	105			40-135	%REC	10	15-Jan-2020 14:48
Surr: 4-Terphenyl-d14	104	J		40-135	%REC	100	15-Jan-2020 16:41
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	15-Jan-2020 18:35
Surr: Nitrobenzene-d5	108			41-120	%REC	1	15-Jan-2020 12:36
Surr: Nitrobenzene-d5	78.6			41-120	%REC	10	15-Jan-2020 14:48
Surr: Nitrobenzene-d5	97.5	J		41-120	%REC	100	15-Jan-2020 16:41
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	15-Jan-2020 18:35
Surr: Phenol-d6	59.5	J		20-120	%REC	100	15-Jan-2020 16:41
Surr: Phenol-d6	0	JS		20-120	%REC	2000	15-Jan-2020 18:35
Surr: Phenol-d6	63.0			20-120	%REC	1	15-Jan-2020 12:36
Surr: Phenol-d6	73.0			20-120	%REC	10	15-Jan-2020 14:48
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00283		0.000400	0.00200	mg/L	1	13-Jan-2020 23:00

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18A-20200108  
 Collection Date: 08-Jan-2020 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.0050	0.025	mg/L	25	14-Jan-2020 17:49
<b>Benzene</b>	<b>0.41</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	14-Jan-2020 17:49
Chlorobenzene		U	0.0075	0.025	mg/L	25	14-Jan-2020 17:49
<b>Ethylbenzene</b>	<b>0.32</b>		<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	14-Jan-2020 17:49
Methylene chloride		U	0.025	0.050	mg/L	25	14-Jan-2020 17:49
<b>Toluene</b>	<b>0.11</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	14-Jan-2020 17:49
Vinyl chloride		U	0.0050	0.025	mg/L	25	14-Jan-2020 17:49
<b>Xylenes, Total</b>	<b>0.50</b>		<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	14-Jan-2020 17:49
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.3</i>			<i>70-126</i>	<i>%REC</i>	25	<i>14-Jan-2020 17:49</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.1</i>			<i>81-113</i>	<i>%REC</i>	25	<i>14-Jan-2020 17:49</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.1</i>			<i>77-123</i>	<i>%REC</i>	25	<i>14-Jan-2020 17:49</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	25	<i>14-Jan-2020 17:49</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18A-20200108  
 Collection Date: 08-Jan-2020 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 12:55
<b>2,4-Dimethylphenol</b>	<b>1.8</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	500	15-Jan-2020 17:00
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 12:55
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 12:55
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 12:55
<b>2-Methylnaphthalene</b>	<b>0.36</b>		<b>0.00095</b>	<b>0.0050</b>	<b>mg/L</b>	50	15-Jan-2020 15:07
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 12:55
<b>4-Nitrophenol</b>	<b>0.0042</b>		<b>0.000047</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
<b>Acenaphthene</b>	<b>0.33</b>		<b>0.0014</b>	<b>0.0050</b>	<b>mg/L</b>	50	15-Jan-2020 15:07
<b>Acenaphthylene</b>	<b>0.0081</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
<b>Anthracene</b>	<b>0.0082</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	15-Jan-2020 12:55
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 12:55
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 12:55
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00043</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
Chrysene	U		0.000021	0.00010	mg/L	1	15-Jan-2020 12:55
<b>Dibenzofuran</b>	<b>0.20</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	50	15-Jan-2020 15:07
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 12:55
<b>Fluoranthene</b>	<b>0.0023</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
<b>Fluorene</b>	<b>0.16</b>		<b>0.0015</b>	<b>0.0050</b>	<b>mg/L</b>	50	15-Jan-2020 15:07
<b>Naphthalene</b>	<b>4.6</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	500	15-Jan-2020 17:00
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 12:55
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 12:55
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 12:55
<b>Phenanthrene</b>	<b>0.10</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	50	15-Jan-2020 15:07
<b>Phenol</b>	<b>0.0014</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
<b>Pyrene</b>	<b>0.0013</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 12:55
Surr: 2,4,6-Tribromophenol	70.4			34-129	%REC	1	15-Jan-2020 12:55
Surr: 2,4,6-Tribromophenol	97.3	J		34-129	%REC	50	15-Jan-2020 15:07
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	500	15-Jan-2020 17:00
Surr: 2-Fluorobiphenyl	97.7	J		40-125	%REC	50	15-Jan-2020 15:07
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	500	15-Jan-2020 17:00
Surr: 2-Fluorobiphenyl	62.0			40-125	%REC	1	15-Jan-2020 12:55
Surr: 2-Fluorophenol	81.6			20-120	%REC	1	15-Jan-2020 12:55
Surr: 2-Fluorophenol	97.2	J		20-120	%REC	50	15-Jan-2020 15:07
Surr: 2-Fluorophenol	0	JS		20-120	%REC	500	15-Jan-2020 17:00
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	500	15-Jan-2020 17:00
Surr: 4-Terphenyl-d14	86.6			40-135	%REC	1	15-Jan-2020 12:55
Surr: 4-Terphenyl-d14	108	J		40-135	%REC	50	15-Jan-2020 15:07

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW18A-20200108  
 Collection Date: 08-Jan-2020 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	88.0			41-120	%REC	1	15-Jan-2020 12:55
Surr: Nitrobenzene-d5	77.8	J		41-120	%REC	50	15-Jan-2020 15:07
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	500	15-Jan-2020 17:00
Surr: Phenol-d6	0	JS		20-120	%REC	500	15-Jan-2020 17:00
Surr: Phenol-d6	66.3			20-120	%REC	1	15-Jan-2020 12:55
Surr: Phenol-d6	79.6	J		20-120	%REC	50	15-Jan-2020 15:07
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.0236		0.000400	0.00200	mg/L	1	13-Jan-2020 23:03

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW58A-20200108  
 Collection Date: 08-Jan-2020 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	15-Jan-2020 13:22
<b>Benzene</b>	<b>0.0053</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:22
<b>Chlorobenzene</b>	<b>0.00072</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:22
<b>Ethylbenzene</b>	<b>0.0098</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:22
Methylene chloride		U	0.0010	0.0020	mg/L	1	15-Jan-2020 13:22
Toluene		U	0.00020	0.0010	mg/L	1	15-Jan-2020 13:22
Vinyl chloride		U	0.00020	0.0010	mg/L	1	15-Jan-2020 13:22
<b>Xylenes, Total</b>	<b>0.028</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:22
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:22</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:22</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:22</i>
<i>Surr: Toluene-d8</i>	<i>98.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:22</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW58A-20200108  
 Collection Date: 08-Jan-2020 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 13:14
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	15-Jan-2020 13:14
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 13:14
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 13:14
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 13:14
<b>2-Methylnaphthalene</b>	<b>0.080</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:26
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 13:14
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 13:14
<b>Acenaphthene</b>	<b>0.27</b>		<b>0.0054</b>	<b>0.020</b>	<b>mg/L</b>	200	15-Jan-2020 17:19
<b>Acenaphthylene</b>	<b>0.0015</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:14
<b>Anthracene</b>	<b>0.018</b>		<b>0.00028</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:26
<b>Benz(a)anthracene</b>	<b>0.000091</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:14
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	15-Jan-2020 13:14
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 13:14
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 13:14
<b>Chrysene</b>	<b>0.000071</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:14
<b>Dibenzofuran</b>	<b>0.13</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:26
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 13:14
<b>Fluoranthene</b>	<b>0.0099</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:14
<b>Fluorene</b>	<b>0.19</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:26
<b>Naphthalene</b>	<b>1.2</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	15-Jan-2020 17:19
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 13:14
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 13:14
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 13:14
<b>Phenanthrene</b>	<b>0.062</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:26
Phenol	U		0.000035	0.00020	mg/L	1	15-Jan-2020 13:14
<b>Pyrene</b>	<b>0.0050</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:14
Surr: 2,4,6-Tribromophenol	82.3			34-129	%REC	20	15-Jan-2020 15:26
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	200	15-Jan-2020 17:19
Surr: 2,4,6-Tribromophenol	73.5			34-129	%REC	1	15-Jan-2020 13:14
Surr: 2-Fluorobiphenyl	59.5			40-125	%REC	1	15-Jan-2020 13:14
Surr: 2-Fluorobiphenyl	87.0			40-125	%REC	20	15-Jan-2020 15:26
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	200	15-Jan-2020 17:19
Surr: 2-Fluorophenol	52.6	J		20-120	%REC	20	15-Jan-2020 15:26
Surr: 2-Fluorophenol	0	JS		20-120	%REC	200	15-Jan-2020 17:19
Surr: 2-Fluorophenol	50.4			20-120	%REC	1	15-Jan-2020 13:14
Surr: 4-Terphenyl-d14	87.2			40-135	%REC	1	15-Jan-2020 13:14
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	15-Jan-2020 17:19
Surr: 4-Terphenyl-d14	104			40-135	%REC	20	15-Jan-2020 15:26

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW58A-20200108  
 Collection Date: 08-Jan-2020 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	15-Jan-2020 17:19
Surr: Nitrobenzene-d5	54.3			41-120	%REC	1	15-Jan-2020 13:14
Surr: Nitrobenzene-d5	66.9	J		41-120	%REC	20	15-Jan-2020 15:26
Surr: Phenol-d6	52.1			20-120	%REC	1	15-Jan-2020 13:14
Surr: Phenol-d6	56.2	J		20-120	%REC	20	15-Jan-2020 15:26
Surr: Phenol-d6	0	JS		20-120	%REC	200	15-Jan-2020 17:19
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.000906	J	0.000400	0.00200	mg/L	1	13-Jan-2020 23:05

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57A-20200108  
 Collection Date: 08-Jan-2020 14:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	15-Jan-2020 13:47
<b>Benzene</b>	<b>0.038</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:47
<b>Chlorobenzene</b>	<b>0.00046</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:47
<b>Ethylbenzene</b>	<b>0.015</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:47
Methylene chloride		U	0.0010	0.0020	mg/L	1	15-Jan-2020 13:47
<b>Toluene</b>	<b>0.010</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:47
Vinyl chloride		U	0.00020	0.0010	mg/L	1	15-Jan-2020 13:47
<b>Xylenes, Total</b>	<b>0.026</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 13:47
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>92.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:47</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:47</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:47</i>
<i>Surr: Toluene-d8</i>	<i>97.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:47</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57A-20200108  
 Collection Date: 08-Jan-2020 14:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	15-Jan-2020 13:33
<b>2,4-Dimethylphenol</b>	<b>0.41</b>		<b>0.0080</b>	<b>0.040</b>	<b>mg/L</b>	200	15-Jan-2020 17:38
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	15-Jan-2020 13:33
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	15-Jan-2020 13:33
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	15-Jan-2020 13:33
<b>2-Methylnaphthalene</b>	<b>0.18</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	15-Jan-2020 13:33
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	15-Jan-2020 13:33
<b>Acenaphthene</b>	<b>0.19</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
<b>Acenaphthylene</b>	<b>0.0025</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:33
<b>Anthracene</b>	<b>0.048</b>		<b>0.00028</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
<b>Benz(a)anthracene</b>	<b>0.0036</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:33
<b>Benzo(a)pyrene</b>	<b>0.0011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:33
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	15-Jan-2020 13:33
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00037</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 13:33
<b>Chrysene</b>	<b>0.0030</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	15-Jan-2020 13:33
<b>Dibenzofuran</b>	<b>0.15</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	15-Jan-2020 13:33
<b>Fluoranthene</b>	<b>0.038</b>		<b>0.00020</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
<b>Fluorene</b>	<b>0.13</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
<b>Naphthalene</b>	<b>1.0</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	15-Jan-2020 17:38
Nitrobenzene	U		0.000024	0.00020	mg/L	1	15-Jan-2020 13:33
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	15-Jan-2020 13:33
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	15-Jan-2020 13:33
<b>Phenanthrene</b>	<b>0.17</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
<b>Phenol</b>	<b>0.0049</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	15-Jan-2020 13:33
<b>Pyrene</b>	<b>0.025</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	15-Jan-2020 15:45
<i>Surr: 2,4,6-Tribromophenol</i>	<i>66.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:33</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>102</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>15-Jan-2020 17:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>93.8</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>15-Jan-2020 17:38</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>56.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:33</i>
<i>Surr: 2-Fluorophenol</i>	<i>65.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:33</i>
<i>Surr: 2-Fluorophenol</i>	<i>85.2</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>15-Jan-2020 17:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>115</i>			<i>40-135</i>	<i>%REC</i>	<i>20</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>200</i>	<i>15-Jan-2020 17:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>87.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 13:33</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57A-20200108  
 Collection Date: 08-Jan-2020 14:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 14-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	67.2			41-120	%REC	1	15-Jan-2020 13:33
Surr: Nitrobenzene-d5	75.6	J		41-120	%REC	20	15-Jan-2020 15:45
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	15-Jan-2020 17:38
Surr: Phenol-d6	59.7			20-120	%REC	1	15-Jan-2020 13:33
Surr: Phenol-d6	80.7			20-120	%REC	20	15-Jan-2020 15:45
Surr: Phenol-d6	0	JS		20-120	%REC	200	15-Jan-2020 17:38
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.0545		0.000400	0.00200	mg/L	1	13-Jan-2020 23:07

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57B-20200108  
 Collection Date: 08-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0020	0.010	mg/L	10	15-Jan-2020 15:07
<b>Benzene</b>	<b>0.52</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:07
Chlorobenzene	U		0.0030	0.010	mg/L	10	15-Jan-2020 15:07
<b>Ethylbenzene</b>	<b>0.25</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:07
Methylene chloride	U		0.010	0.020	mg/L	10	15-Jan-2020 15:07
<b>Toluene</b>	<b>0.64</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:07
<b>Xylenes, Total</b>	<b>0.70</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:07
<i>Surr: 1,2-Dichloroethane-d4</i>	95.6			70-126	%REC	10	15-Jan-2020 15:07
<i>Surr: 4-Bromofluorobenzene</i>	99.3			81-113	%REC	10	15-Jan-2020 15:07
<i>Surr: Dibromofluoromethane</i>	99.6			77-123	%REC	10	15-Jan-2020 15:07
<i>Surr: Toluene-d8</i>	98.9			82-127	%REC	10	15-Jan-2020 15:07

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57B-20200108  
 Collection Date: 08-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	17-Jan-2020 04:40
<b>2,4-Dimethylphenol</b>	<b>2.8</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	1000	17-Jan-2020 12:49
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	17-Jan-2020 04:40
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	17-Jan-2020 04:40
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	17-Jan-2020 04:40
<b>2-Methylnaphthalene</b>	<b>1.1</b>		<b>0.019</b>	<b>0.10</b>	<b>mg/L</b>	1000	17-Jan-2020 12:49
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	17-Jan-2020 04:40
4-Nitrophenol	U		0.00047	0.010	mg/L	10	17-Jan-2020 04:40
<b>Acenaphthene</b>	<b>0.55</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
<b>Acenaphthylene</b>	<b>0.0086</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 04:40
<b>Anthracene</b>	<b>0.19</b>		<b>0.0014</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
<b>Benz(a)anthracene</b>	<b>0.040</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 04:40
<b>Benzo(a)pyrene</b>	<b>0.012</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 04:40
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	17-Jan-2020 04:40
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	17-Jan-2020 04:40
<b>Chrysene</b>	<b>0.042</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 04:40
<b>Dibenzofuran</b>	<b>0.49</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	17-Jan-2020 04:40
<b>Fluoranthene</b>	<b>0.34</b>		<b>0.0010</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
<b>Fluorene</b>	<b>0.46</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
<b>Naphthalene</b>	<b>13</b>		<b>0.20</b>	<b>1.0</b>	<b>mg/L</b>	10000	17-Jan-2020 16:07
Nitrobenzene	U		0.00024	0.0020	mg/L	10	17-Jan-2020 04:40
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	17-Jan-2020 04:40
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	17-Jan-2020 04:40
<b>Phenanthrene</b>	<b>1.1</b>		<b>0.021</b>	<b>0.10</b>	<b>mg/L</b>	1000	17-Jan-2020 12:49
<b>Phenol</b>	<b>0.32</b>		<b>0.0035</b>	<b>0.020</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
<b>Pyrene</b>	<b>0.21</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 12:29
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>10000</i>	<i>17-Jan-2020 16:07</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>17-Jan-2020 12:29</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>17-Jan-2020 12:49</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>91.9</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>17-Jan-2020 04:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>80.9</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>17-Jan-2020 04:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>17-Jan-2020 12:29</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>17-Jan-2020 12:49</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>10000</i>	<i>17-Jan-2020 16:07</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>10000</i>	<i>17-Jan-2020 16:07</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>17-Jan-2020 12:29</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>17-Jan-2020 12:49</i>
<i>Surr: 2-Fluorophenol</i>	<i>117</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>17-Jan-2020 04:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW57B-20200108  
 Collection Date: 08-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	17-Jan-2020 12:29
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	17-Jan-2020 12:49
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	10000	17-Jan-2020 16:07
Surr: 4-Terphenyl-d14	112			40-135	%REC	10	17-Jan-2020 04:40
Surr: Nitrobenzene-d5	119			41-120	%REC	10	17-Jan-2020 04:40
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	17-Jan-2020 12:49
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	10000	17-Jan-2020 16:07
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	17-Jan-2020 12:29
Surr: Phenol-d6	0	JS		20-120	%REC	10000	17-Jan-2020 16:07
Surr: Phenol-d6	0	JS		20-120	%REC	100	17-Jan-2020 12:29
Surr: Phenol-d6	0	JS		20-120	%REC	1000	17-Jan-2020 12:49
Surr: Phenol-d6	99.2			20-120	%REC	10	17-Jan-2020 04:40
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00404		0.000400	0.00200	mg/L	1	13-Jan-2020 23:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB03-20200109  
 Collection Date: 09-Jan-2020 07:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 20:40
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 20:40
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 20:40
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 20:40
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 20:40
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 20:40
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 20:40
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 20:40
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>97.6</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 20:40</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.5</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 20:40</i>
<i>Surr: Dibromofluoromethane</i>		<i>101</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 20:40</i>
<i>Surr: Toluene-d8</i>		<i>103</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 20:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB03-20200109  
 Collection Date: 09-Jan-2020 07:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	17-Jan-2020 04:59
<b>2,4-Dimethylphenol</b>	<b>0.0010</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	17-Jan-2020 04:59
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	17-Jan-2020 04:59
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	17-Jan-2020 04:59
<b>2-Methylnaphthalene</b>	<b>0.00038</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	17-Jan-2020 04:59
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	17-Jan-2020 04:59
<b>Acenaphthene</b>	<b>0.00015</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
Acenaphthylene		U	0.000015	0.00010	mg/L	1	17-Jan-2020 04:59
<b>Anthracene</b>	<b>0.000047</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	17-Jan-2020 04:59
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	17-Jan-2020 04:59
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	17-Jan-2020 04:59
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
Chrysene		U	0.000021	0.00010	mg/L	1	17-Jan-2020 04:59
<b>Dibenzofuran</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	17-Jan-2020 04:59
<b>Fluoranthene</b>	<b>0.000069</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
<b>Fluorene</b>	<b>0.000092</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
<b>Naphthalene</b>	<b>0.0078</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
Nitrobenzene		U	0.000024	0.00020	mg/L	1	17-Jan-2020 04:59
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	17-Jan-2020 04:59
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	17-Jan-2020 04:59
<b>Phenanthrene</b>	<b>0.00019</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
<b>Phenol</b>	<b>0.000053</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
<b>Pyrene</b>	<b>0.000039</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 04:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 04:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>77.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 04:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 04:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>81.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 04:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>78.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 04:59</i>
<i>Surr: Phenol-d6</i>	<i>78.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 04:59</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic		U	0.000400	0.00200	mg/L	1	13-Jan-2020 23:11

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW72B-20200109  
 Collection Date: 09-Jan-2020 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0020	0.010	mg/L	10	15-Jan-2020 15:34
<b>Benzene</b>	<b>0.70</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:34
Chlorobenzene	U		0.0030	0.010	mg/L	10	15-Jan-2020 15:34
<b>Ethylbenzene</b>	<b>0.13</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:34
Methylene chloride	U		0.010	0.020	mg/L	10	15-Jan-2020 15:34
<b>Toluene</b>	<b>0.58</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:34
<b>Xylenes, Total</b>	<b>0.36</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	10	15-Jan-2020 15:34
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.1</i>			<i>70-126</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 15:34</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.7</i>			<i>81-113</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 15:34</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.8</i>			<i>77-123</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 15:34</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>10</i>	<i>15-Jan-2020 15:34</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW72B-20200109  
 Collection Date: 09-Jan-2020 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	17-Jan-2020 13:08
<b>2,4-Dimethylphenol</b>	<b>13</b>		<b>0.40</b>	<b>2.0</b>	<b>mg/L</b>	10000	17-Jan-2020 16:26
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	17-Jan-2020 13:08
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	17-Jan-2020 13:08
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	17-Jan-2020 13:08
<b>2-Methylnaphthalene</b>	<b>0.22</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 13:28
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	17-Jan-2020 13:08
4-Nitrophenol	U		0.00047	0.010	mg/L	10	17-Jan-2020 13:08
<b>Acenaphthene</b>	<b>0.079</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Acenaphthylene</b>	<b>0.0023</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Anthracene</b>	<b>0.020</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Benz(a)anthracene</b>	<b>0.0031</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Benzo(a)pyrene</b>	<b>0.0011</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	17-Jan-2020 13:08
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	17-Jan-2020 13:08
<b>Chrysene</b>	<b>0.0033</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Dibenzofuran</b>	<b>0.061</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	17-Jan-2020 13:08
<b>Fluoranthene</b>	<b>0.024</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Fluorene</b>	<b>0.049</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Naphthalene</b>	<b>2.0</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	17-Jan-2020 13:48
Nitrobenzene	U		0.00024	0.0020	mg/L	10	17-Jan-2020 13:08
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	17-Jan-2020 13:08
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	17-Jan-2020 13:08
<b>Phenanthrene</b>	<b>0.075</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<b>Phenol</b>	<b>5.2</b>		<b>0.035</b>	<b>0.20</b>	<b>mg/L</b>	1000	17-Jan-2020 13:48
<b>Pyrene</b>	<b>0.013</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 13:08
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>10000</i>	<i>17-Jan-2020 16:26</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.1</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>17-Jan-2020 13:08</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>17-Jan-2020 13:28</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>17-Jan-2020 13:48</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>88.2</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>17-Jan-2020 13:08</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>17-Jan-2020 13:28</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>17-Jan-2020 13:48</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>10000</i>	<i>17-Jan-2020 16:26</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>10000</i>	<i>17-Jan-2020 16:26</i>
<i>Surr: 2-Fluorophenol</i>	<i>108</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>17-Jan-2020 13:08</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>17-Jan-2020 13:28</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>17-Jan-2020 13:48</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW72B-20200109  
 Collection Date: 09-Jan-2020 08:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
Surr: 4-Terphenyl-d14	87.4			40-135	%REC	10	17-Jan-2020 13:08
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	17-Jan-2020 13:28
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	17-Jan-2020 13:48
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	10000	17-Jan-2020 16:26
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	10000	17-Jan-2020 16:26
Surr: Nitrobenzene-d5	82.9			41-120	%REC	10	17-Jan-2020 13:08
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	17-Jan-2020 13:28
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	17-Jan-2020 13:48
Surr: Phenol-d6	81.9			20-120	%REC	10	17-Jan-2020 13:08
Surr: Phenol-d6	0	JS		20-120	%REC	100	17-Jan-2020 13:28
Surr: Phenol-d6	0	JS		20-120	%REC	1000	17-Jan-2020 13:48
Surr: Phenol-d6	0	JS		20-120	%REC	10000	17-Jan-2020 16:26
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Jan-2020		Analyst: JHD	
Arsenic	0.00105	J	0.000400	0.00200	mg/L	1	13-Jan-2020 23:14

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW23C-20200109  
 Collection Date: 09-Jan-2020 09:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0010	0.0050	mg/L	5	15-Jan-2020 14:14
<b>Benzene</b>	<b>0.0027</b>	J	<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	5	15-Jan-2020 14:14
Chlorobenzene	U		0.0015	0.0050	mg/L	5	15-Jan-2020 14:14
<b>Ethylbenzene</b>	<b>0.028</b>		<b>0.0015</b>	<b>0.0050</b>	<b>mg/L</b>	5	15-Jan-2020 14:14
Methylene chloride	U		0.0050	0.010	mg/L	5	15-Jan-2020 14:14
Toluene	U		0.0010	0.0050	mg/L	5	15-Jan-2020 14:14
<b>Xylenes, Total</b>	<b>0.025</b>		<b>0.0015</b>	<b>0.0050</b>	<b>mg/L</b>	5	15-Jan-2020 14:14
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.0</i>			<i>70-126</i>	<i>%REC</i>	5	<i>15-Jan-2020 14:14</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	5	<i>15-Jan-2020 14:14</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	5	<i>15-Jan-2020 14:14</i>
<i>Surr: Toluene-d8</i>	<i>98.9</i>			<i>82-127</i>	<i>%REC</i>	5	<i>15-Jan-2020 14:14</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW23C-20200109  
 Collection Date: 09-Jan-2020 09:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.00021	0.0020	mg/L	10	17-Jan-2020 14:08
<b>2,4-Dimethylphenol</b>	<b>0.047</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
2,4-Dinitrotoluene	U		0.00058	0.0020	mg/L	10	17-Jan-2020 14:08
2,6-Dinitrotoluene	U		0.00042	0.0020	mg/L	10	17-Jan-2020 14:08
2-Chloronaphthalene	U		0.00021	0.0020	mg/L	10	17-Jan-2020 14:08
<b>2-Methylnaphthalene</b>	<b>0.12</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 16:46
4,6-Dinitro-2-methylphenol	U		0.00020	0.0020	mg/L	10	17-Jan-2020 14:08
4-Nitrophenol	U		0.00047	0.010	mg/L	10	17-Jan-2020 14:08
<b>Acenaphthene</b>	<b>0.21</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 16:46
<b>Acenaphthylene</b>	<b>0.0029</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
<b>Anthracene</b>	<b>0.035</b>		<b>0.00014</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
<b>Benz(a)anthracene</b>	<b>0.0080</b>		<b>0.00050</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
<b>Benzo(a)pyrene</b>	<b>0.0022</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
Bis(2-chloroethoxy)methane	U		0.00030	0.0020	mg/L	10	17-Jan-2020 14:08
Bis(2-ethylhexyl)phthalate	U		0.00037	0.0020	mg/L	10	17-Jan-2020 14:08
<b>Chrysene</b>	<b>0.0072</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
<b>Dibenzofuran</b>	<b>0.17</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 16:46
Di-n-butyl phthalate	U		0.00020	0.0020	mg/L	10	17-Jan-2020 14:08
<b>Fluoranthene</b>	<b>0.068</b>		<b>0.00010</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
<b>Fluorene</b>	<b>0.12</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 16:46
<b>Naphthalene</b>	<b>0.66</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 16:46
Nitrobenzene	U		0.00024	0.0020	mg/L	10	17-Jan-2020 14:08
N-Nitrosodiphenylamine	U		0.00025	0.0020	mg/L	10	17-Jan-2020 14:08
Pentachlorophenol	U		0.00079	0.0020	mg/L	10	17-Jan-2020 14:08
<b>Phenanthrene</b>	<b>0.25</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	17-Jan-2020 16:46
<b>Phenol</b>	<b>0.0079</b>		<b>0.00035</b>	<b>0.0020</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
<b>Pyrene</b>	<b>0.043</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 14:08
Surr: 2,4,6-Tribromophenol	87.3			34-129	%REC	10	17-Jan-2020 14:08
Surr: 2,4,6-Tribromophenol	103	J		34-129	%REC	100	17-Jan-2020 16:46
Surr: 2-Fluorobiphenyl	67.3	J		40-125	%REC	100	17-Jan-2020 16:46
Surr: 2-Fluorobiphenyl	72.8			40-125	%REC	10	17-Jan-2020 14:08
Surr: 2-Fluorophenol	81.0			20-120	%REC	10	17-Jan-2020 14:08
Surr: 2-Fluorophenol	80.8	J		20-120	%REC	100	17-Jan-2020 16:46
Surr: 4-Terphenyl-d14	97.6	J		40-135	%REC	100	17-Jan-2020 16:46
Surr: 4-Terphenyl-d14	89.4			40-135	%REC	10	17-Jan-2020 14:08
Surr: Nitrobenzene-d5	131	JS		41-120	%REC	100	17-Jan-2020 16:46
Surr: Nitrobenzene-d5	116			41-120	%REC	10	17-Jan-2020 14:08
Surr: Phenol-d6	71.5			20-120	%REC	10	17-Jan-2020 14:08
Surr: Phenol-d6	84.0	J		20-120	%REC	100	17-Jan-2020 16:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW23C-20200109  
 Collection Date: 09-Jan-2020 09:40

**ANALYTICAL REPORT**

WorkOrder:HS20010407  
 Lab ID:HS20010407-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 13-Jan-2020		Analyst: JHD
Arsenic	0.00333		0.000400	0.00200	mg/L	1	13-Jan-2020 23:16

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW53C-20200109  
 Collection Date: 09-Jan-2020 11:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 05:56
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 05:56
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 05:56
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 05:56
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 05:56
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 05:56
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 05:56
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 05:56</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 05:56</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 05:56</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 05:56</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW53C-20200109  
 Collection Date: 09-Jan-2020 11:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	17-Jan-2020 14:27
<b>2,4-Dimethylphenol</b>	<b>0.0011</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	17-Jan-2020 14:27
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	17-Jan-2020 14:27
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	17-Jan-2020 14:27
<b>2-Methylnaphthalene</b>	<b>0.00025</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	17-Jan-2020 14:27
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	17-Jan-2020 14:27
<b>Acenaphthene</b>	<b>0.00032</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
Acenaphthylene		U	0.000015	0.00010	mg/L	1	17-Jan-2020 14:27
<b>Anthracene</b>	<b>0.000028</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	17-Jan-2020 14:27
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	17-Jan-2020 14:27
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	17-Jan-2020 14:27
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	17-Jan-2020 14:27
Chrysene		U	0.000021	0.00010	mg/L	1	17-Jan-2020 14:27
<b>Dibenzofuran</b>	<b>0.00027</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	17-Jan-2020 14:27
<b>Fluoranthene</b>	<b>0.000071</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
<b>Fluorene</b>	<b>0.00020</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
<b>Naphthalene</b>	<b>0.0012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
Nitrobenzene		U	0.000024	0.00020	mg/L	1	17-Jan-2020 14:27
<b>N-Nitrosodiphenylamine</b>	<b>0.000056</b>	J	<b>0.000025</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	17-Jan-2020 14:27
<b>Phenanthrene</b>	<b>0.00025</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
<b>Phenol</b>	<b>0.00050</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
<b>Pyrene</b>	<b>0.000061</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:27
Surr: 2,4,6-Tribromophenol	65.4			34-129	%REC	1	17-Jan-2020 14:27
Surr: 2-Fluorobiphenyl	64.0			40-125	%REC	1	17-Jan-2020 14:27
Surr: 2-Fluorophenol	66.8			20-120	%REC	1	17-Jan-2020 14:27
Surr: 4-Terphenyl-d14	91.4			40-135	%REC	1	17-Jan-2020 14:27
Surr: Nitrobenzene-d5	64.9			41-120	%REC	1	17-Jan-2020 14:27
Surr: Phenol-d6	68.0			20-120	%REC	1	17-Jan-2020 14:27
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000728</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	14-Jan-2020 21:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54C-20200109  
 Collection Date: 09-Jan-2020 12:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 06:20
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 06:20
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 06:20
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 06:20
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 06:20
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 06:20
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 06:20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 06:20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 06:20</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 06:20</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 06:20</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54C-20200109  
 Collection Date: 09-Jan-2020 12:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 14:47
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 14:47
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 14:47
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 14:47
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 14:47
<b>2-Methylnaphthalene</b>	<b>0.0020</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 14:47
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 14:47
<b>Acenaphthene</b>	<b>0.021</b>		<b>0.00014</b>	<b>0.00050</b>	<b>mg/L</b>	5	17-Jan-2020 17:06
<b>Acenaphthylene</b>	<b>0.00023</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Anthracene</b>	<b>0.0015</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 14:47
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 14:47
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 14:47
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00018</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Chrysene</b>	<b>0.000054</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Dibenzofuran</b>	<b>0.020</b>		<b>0.00010</b>	<b>0.00050</b>	<b>mg/L</b>	5	17-Jan-2020 17:06
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 14:47
<b>Fluoranthene</b>	<b>0.0026</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Fluorene</b>	<b>0.0095</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Naphthalene</b>	<b>0.035</b>		<b>0.00010</b>	<b>0.00050</b>	<b>mg/L</b>	5	17-Jan-2020 17:06
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 14:47
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 14:47
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 14:47
<b>Phenanthrene</b>	<b>0.0056</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Phenol</b>	<b>0.000064</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<b>Pyrene</b>	<b>0.0014</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 14:47
<i>Surr: 2,4,6-Tribromophenol</i>	58.6			34-129	%REC	1	17-Jan-2020 14:47
<i>Surr: 2,4,6-Tribromophenol</i>	71.5			34-129	%REC	5	17-Jan-2020 17:06
<i>Surr: 2-Fluorobiphenyl</i>	65.7			40-125	%REC	5	17-Jan-2020 17:06
<i>Surr: 2-Fluorobiphenyl</i>	52.3			40-125	%REC	1	17-Jan-2020 14:47
<i>Surr: 2-Fluorophenol</i>	66.8			20-120	%REC	5	17-Jan-2020 17:06
<i>Surr: 2-Fluorophenol</i>	52.0			20-120	%REC	1	17-Jan-2020 14:47
<i>Surr: 4-Terphenyl-d14</i>	81.0			40-135	%REC	1	17-Jan-2020 14:47
<i>Surr: 4-Terphenyl-d14</i>	102			40-135	%REC	5	17-Jan-2020 17:06
<i>Surr: Nitrobenzene-d5</i>	58.3			41-120	%REC	1	17-Jan-2020 14:47
<i>Surr: Nitrobenzene-d5</i>	71.5			41-120	%REC	5	17-Jan-2020 17:06
<i>Surr: Phenol-d6</i>	54.6			20-120	%REC	1	17-Jan-2020 14:47
<i>Surr: Phenol-d6</i>	69.4			20-120	%REC	5	17-Jan-2020 17:06

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54C-20200109  
 Collection Date: 09-Jan-2020 12:35

**ANALYTICAL REPORT**

WorkOrder:HS20010407  
 Lab ID:HS20010407-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
Arsenic	0.00122	J	0.000400	0.00200	mg/L	1	14-Jan-2020 21:52

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44A-20200109  
 Collection Date: 09-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:33
Benzene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:33
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:33
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:33
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 12:33
Toluene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:33
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:33
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>100</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:33</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:33</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:33</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:33</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44A-20200109  
 Collection Date: 09-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	17-Jan-2020 15:07
<b>2,4-Dimethylphenol</b>	<b>0.00042</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	17-Jan-2020 15:07
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	17-Jan-2020 15:07
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	17-Jan-2020 15:07
<b>2-Methylnaphthalene</b>	<b>0.00035</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	17-Jan-2020 15:07
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	17-Jan-2020 15:07
<b>Acenaphthene</b>	<b>0.064</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	17-Jan-2020 17:26
<b>Acenaphthylene</b>	<b>0.00097</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
<b>Anthracene</b>	<b>0.00027</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
<b>Benz(a)anthracene</b>	<b>0.000054</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	17-Jan-2020 15:07
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	17-Jan-2020 15:07
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	17-Jan-2020 15:07
Chrysene		U	0.000021	0.00010	mg/L	1	17-Jan-2020 15:07
<b>Dibenzofuran</b>	<b>0.00033</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	17-Jan-2020 15:07
<b>Fluoranthene</b>	<b>0.0059</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
<b>Fluorene</b>	<b>0.0078</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
<b>Naphthalene</b>	<b>0.0030</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
Nitrobenzene		U	0.000024	0.00020	mg/L	1	17-Jan-2020 15:07
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	17-Jan-2020 15:07
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	17-Jan-2020 15:07
<b>Phenanthrene</b>	<b>0.00022</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
<b>Phenol</b>	<b>0.00013</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
<b>Pyrene</b>	<b>0.0045</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:07
Surr: 2,4,6-Tribromophenol	75.5			34-129	%REC	1	17-Jan-2020 15:07
Surr: 2,4,6-Tribromophenol	95.3			34-129	%REC	10	17-Jan-2020 17:26
Surr: 2-Fluorobiphenyl	64.7			40-125	%REC	1	17-Jan-2020 15:07
Surr: 2-Fluorobiphenyl	79.8			40-125	%REC	10	17-Jan-2020 17:26
Surr: 2-Fluorophenol	59.8			20-120	%REC	1	17-Jan-2020 15:07
Surr: 2-Fluorophenol	83.8			20-120	%REC	10	17-Jan-2020 17:26
Surr: 4-Terphenyl-d14	82.7			40-135	%REC	1	17-Jan-2020 15:07
Surr: 4-Terphenyl-d14	114			40-135	%REC	10	17-Jan-2020 17:26
Surr: Nitrobenzene-d5	87.6			41-120	%REC	10	17-Jan-2020 17:26
Surr: Nitrobenzene-d5	61.3			41-120	%REC	1	17-Jan-2020 15:07
Surr: Phenol-d6	64.2			20-120	%REC	1	17-Jan-2020 15:07
Surr: Phenol-d6	82.5			20-120	%REC	10	17-Jan-2020 17:26

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44A-20200109  
 Collection Date: 09-Jan-2020 13:30

**ANALYTICAL REPORT**

WorkOrder:HS20010407  
 Lab ID:HS20010407-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 14-Jan-2020		Analyst: ALR
Arsenic	0.00966		0.000400	0.00200	mg/L	1	14-Jan-2020 21:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36B-20200109  
 Collection Date: 09-Jan-2020 14:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:20
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:20
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 18:20
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 18:20
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 18:20
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:20
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:20
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 18:20
<i>Surr: 1,2-Dichloroethane-d4</i>		112		70-126	%REC	1	14-Jan-2020 18:20
<i>Surr: 4-Bromofluorobenzene</i>		98.0		81-113	%REC	1	14-Jan-2020 18:20
<i>Surr: Dibromofluoromethane</i>		105		77-123	%REC	1	14-Jan-2020 18:20
<i>Surr: Toluene-d8</i>		101		82-127	%REC	1	14-Jan-2020 18:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36B-20200109  
 Collection Date: 09-Jan-2020 14:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 15:27
<b>2,4-Dimethylphenol</b>	<b>0.000054</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 15:27
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 15:27
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 15:27
<b>2-Methylnaphthalene</b>	<b>0.00013</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 15:27
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 15:27
<b>Acenaphthene</b>	<b>0.00017</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 15:27
Anthracene	U		0.000014	0.00010	mg/L	1	17-Jan-2020 15:27
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 15:27
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 15:27
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 15:27
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 15:27
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 15:27
<b>Dibenzofuran</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 15:27
<b>Fluoranthene</b>	<b>0.000051</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
<b>Fluorene</b>	<b>0.00012</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
<b>Naphthalene</b>	<b>0.00093</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 15:27
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 15:27
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 15:27
<b>Phenanthrene</b>	<b>0.00010</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
<b>Phenol</b>	<b>0.000075</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 15:27
Pyrene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 15:27
<i>Surr: 2,4,6-Tribromophenol</i>	70.4			34-129	%REC	1	17-Jan-2020 15:27
<i>Surr: 2-Fluorobiphenyl</i>	56.0			40-125	%REC	1	17-Jan-2020 15:27
<i>Surr: 2-Fluorophenol</i>	50.6			20-120	%REC	1	17-Jan-2020 15:27
<i>Surr: 4-Terphenyl-d14</i>	79.3			40-135	%REC	1	17-Jan-2020 15:27
<i>Surr: Nitrobenzene-d5</i>	56.4			41-120	%REC	1	17-Jan-2020 15:27
<i>Surr: Phenol-d6</i>	50.2			20-120	%REC	1	17-Jan-2020 15:27
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00100</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	14-Jan-2020 21:57

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36A-20200109  
 Collection Date: 09-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-26  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:44
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:44
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 18:44
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 18:44
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 18:44
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:44
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 18:44
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 18:44
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>118</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 18:44</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.4</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 18:44</i>
<i>Surr: Dibromofluoromethane</i>		<i>109</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 18:44</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 18:44</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36A-20200109  
 Collection Date: 09-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-26  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 15-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 15:47
<b>2,4-Dimethylphenol</b>	<b>0.000052</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 15:47
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 15:47
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 15:47
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 15:47
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 15:47
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 15:47
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 15:47
Acenaphthene	U		0.000027	0.00010	mg/L	1	17-Jan-2020 15:47
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 15:47
Anthracene	U		0.000014	0.00010	mg/L	1	17-Jan-2020 15:47
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 15:47
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 15:47
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 15:47
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 15:47
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 15:47
Dibenzofuran	U		0.000020	0.00010	mg/L	1	17-Jan-2020 15:47
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 15:47
Fluoranthene	U		0.000010	0.00010	mg/L	1	17-Jan-2020 15:47
Fluorene	U		0.000030	0.00010	mg/L	1	17-Jan-2020 15:47
<b>Naphthalene</b>	<b>0.00034</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 15:47
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 15:47
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 15:47
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 15:47
Phenanthrene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 15:47
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 15:47
Pyrene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 15:47
<i>Surr: 2,4,6-Tribromophenol</i>	<i>54.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 15:47</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>61.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 15:47</i>
<i>Surr: 2-Fluorophenol</i>	<i>61.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 15:47</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>79.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 15:47</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 15:47</i>
<i>Surr: Phenol-d6</i>	<i>60.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 15:47</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00354</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	14-Jan-2020 21:59

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB01-20200109  
 Collection Date: 09-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010407  
 Lab ID:HS20010407-27  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:09
Benzene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:09
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:09
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:09
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 12:09
Toluene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:09
Vinyl chloride	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:09
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:09
<i>Surr: 1,2-Dichloroethane-d4</i>		99.6		70-126	%REC	1	15-Jan-2020 12:09
<i>Surr: 4-Bromofluorobenzene</i>		95.3		81-113	%REC	1	15-Jan-2020 12:09
<i>Surr: Dibromofluoromethane</i>		101		77-123	%REC	1	15-Jan-2020 12:09
<i>Surr: Toluene-d8</i>		102		82-127	%REC	1	15-Jan-2020 12:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**Batch ID:** 149585      **Start Date:** 13 Jan 2020 10:30      **End Date:** 13 Jan 2020 13:30  
**Method:** TX 1005 PREP      **Prep Code:** TX 1005\_W PR

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010407-02	1	31.02 (g)	3 (mL)	0.09671
HS20010407-03	1	31.1 (g)	3 (mL)	0.09646
HS20010407-04	1	30.67 (g)	3 (mL)	0.09782
HS20010407-05	1	31.05 (g)	3 (mL)	0.09662
HS20010407-06	1	31.3 (g)	3 (mL)	0.09585
HS20010407-07	1	31.21 (g)	3 (mL)	0.09612
HS20010407-08	1	31.12 (g)	3 (mL)	0.0964
HS20010407-09	1	31.54 (g)	3 (mL)	0.09512
HS20010407-10	1	31.32 (g)	3 (mL)	0.09579

**Batch ID:** 149592      **Start Date:** 13 Jan 2020 11:00      **End Date:** 13 Jan 2020 15:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010407-01		10 (mL)	10 (mL)	1
HS20010407-02		10 (mL)	10 (mL)	1
HS20010407-03		10 (mL)	10 (mL)	1
HS20010407-04		10 (mL)	10 (mL)	1
HS20010407-05		10 (mL)	10 (mL)	1
HS20010407-06		10 (mL)	10 (mL)	1
HS20010407-07		10 (mL)	10 (mL)	1
HS20010407-08		10 (mL)	10 (mL)	1
HS20010407-09		10 (mL)	10 (mL)	1
HS20010407-10		10 (mL)	10 (mL)	1
HS20010407-11		10 (mL)	10 (mL)	1
HS20010407-12		10 (mL)	10 (mL)	1
HS20010407-13		10 (mL)	10 (mL)	1
HS20010407-14		10 (mL)	10 (mL)	1
HS20010407-15		10 (mL)	10 (mL)	1
HS20010407-16		10 (mL)	10 (mL)	1
HS20010407-17		10 (mL)	10 (mL)	1
HS20010407-18		10 (mL)	10 (mL)	1
HS20010407-20		10 (mL)	10 (mL)	1
HS20010407-21		10 (mL)	10 (mL)	1

**Batch ID:** 149618      **Start Date:** 14 Jan 2020 08:30      **End Date:** 14 Jan 2020 16:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010407-22		10 (mL)	10 (mL)	1
HS20010407-23		10 (mL)	10 (mL)	1
HS20010407-24		10 (mL)	10 (mL)	1
HS20010407-25		10 (mL)	10 (mL)	1
HS20010407-26		10 (mL)	10 (mL)	1

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**Batch ID:** 149624      **Start Date:** 14 Jan 2020 07:00      **End Date:** 14 Jan 2020 12:00  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010407-01	1	1000 (mL)	1 (mL)	0.001
HS20010407-02	1	1000 (mL)	1 (mL)	0.001
HS20010407-03	1	1000 (mL)	1 (mL)	0.001
HS20010407-04	1	1000 (mL)	1 (mL)	0.001
HS20010407-05	1	1000 (mL)	1 (mL)	0.001
HS20010407-06	1	1000 (mL)	1 (mL)	0.001
HS20010407-07	1	1000 (mL)	1 (mL)	0.001
HS20010407-08	1	1000 (mL)	1 (mL)	0.001
HS20010407-09	1	1000 (mL)	1 (mL)	0.001
HS20010407-10	1	1000 (mL)	1 (mL)	0.001
HS20010407-11	1	1000 (mL)	1 (mL)	0.001
HS20010407-12	1	1000 (mL)	1 (mL)	0.001
HS20010407-13	1	1000 (mL)	1 (mL)	0.001
HS20010407-14	1	1000 (mL)	1 (mL)	0.001
HS20010407-15	1	1000 (mL)	1 (mL)	0.001
HS20010407-16	1	1000 (mL)	1 (mL)	0.001

**Batch ID:** 149670      **Start Date:** 15 Jan 2020 08:00      **End Date:** 15 Jan 2020 14:00  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010407-17		1000 (mL)	1 (mL)	0.001
HS20010407-18		1000 (mL)	1 (mL)	0.001
HS20010407-20		1000 (mL)	1 (mL)	0.001
HS20010407-21		1000 (mL)	1 (mL)	0.001
HS20010407-22		1000 (mL)	1 (mL)	0.001
HS20010407-23		1000 (mL)	1 (mL)	0.001
HS20010407-24		1000 (mL)	1 (mL)	0.001
HS20010407-25		1000 (mL)	1 (mL)	0.001
HS20010407-26		1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149585 ( 0 )</b>		<b>Test Name : LOW-LEVEL TEXAS TPH BY TX1005</b>			<b>Matrix: Water</b>	
HS20010407-02	WG-1620-MW51A-20200109	09 Jan 2020 11:50		13 Jan 2020 10:30	13 Jan 2020 21:36	1
HS20010407-03	WG-1620-MW51C-20200109	09 Jan 2020 12:40		13 Jan 2020 10:30	13 Jan 2020 22:05	1
HS20010407-04	WG-1620-MW85C-20200109	09 Jan 2020 13:35		13 Jan 2020 10:30	13 Jan 2020 22:34	1
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25		13 Jan 2020 10:30	13 Jan 2020 23:03	1
HS20010407-06	WG-1620-MW76C-20200109	09 Jan 2020 15:20		13 Jan 2020 10:30	13 Jan 2020 23:33	1
HS20010407-07	WG-1620-MW81B-20200109	09 Jan 2020 16:20		13 Jan 2020 10:30	14 Jan 2020 00:02	1
HS20010407-08	WG-1620-MW50A-20200109	09 Jan 2020 17:15		13 Jan 2020 10:30	14 Jan 2020 00:31	1
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10		13 Jan 2020 10:30	14 Jan 2020 01:00	1
HS20010407-10	WG-1620-MW80B-20200107	07 Jan 2020 16:00		13 Jan 2020 10:30	14 Jan 2020 01:29	1
<b>Batch ID: 149592 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010407-01	WG-1620-MW19C-20200109	09 Jan 2020 10:45		13 Jan 2020 15:00	13 Jan 2020 22:09	1
HS20010407-02	WG-1620-MW51A-20200109	09 Jan 2020 11:50		13 Jan 2020 15:00	13 Jan 2020 22:25	1
HS20010407-03	WG-1620-MW51C-20200109	09 Jan 2020 12:40		13 Jan 2020 15:00	13 Jan 2020 22:27	1
HS20010407-04	WG-1620-MW85C-20200109	09 Jan 2020 13:35		13 Jan 2020 15:00	13 Jan 2020 22:29	1
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25		13 Jan 2020 15:00	13 Jan 2020 22:31	1
HS20010407-06	WG-1620-MW76C-20200109	09 Jan 2020 15:20		13 Jan 2020 15:00	13 Jan 2020 22:34	1
HS20010407-07	WG-1620-MW81B-20200109	09 Jan 2020 16:20		13 Jan 2020 15:00	13 Jan 2020 22:36	1
HS20010407-08	WG-1620-MW50A-20200109	09 Jan 2020 17:15		13 Jan 2020 15:00	13 Jan 2020 22:38	1
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10		13 Jan 2020 15:00	13 Jan 2020 22:40	1
HS20010407-10	WG-1620-MW80B-20200107	07 Jan 2020 16:00		13 Jan 2020 15:00	13 Jan 2020 22:42	1
HS20010407-11	WG-1620-FB01-20200107	07 Jan 2020 16:30		13 Jan 2020 15:00	13 Jan 2020 22:45	1
HS20010407-12	WG-1620-FB02-20200108	08 Jan 2020 09:00		13 Jan 2020 15:00	13 Jan 2020 22:58	1
HS20010407-13	WG-1620-MW18C-20200108	08 Jan 2020 10:10		13 Jan 2020 15:00	13 Jan 2020 23:00	1
HS20010407-14	WG-1620-MW18A-20200108	08 Jan 2020 11:10		13 Jan 2020 15:00	13 Jan 2020 23:03	1
HS20010407-15	WG-1620-MW58A-20200108	08 Jan 2020 12:15		13 Jan 2020 15:00	13 Jan 2020 23:05	1
HS20010407-16	WG-1620-MW57A-20200108	08 Jan 2020 14:55		13 Jan 2020 15:00	13 Jan 2020 23:07	1
HS20010407-17	WG-1620-MW57B-20200108	08 Jan 2020 13:30		13 Jan 2020 15:00	13 Jan 2020 23:09	1
HS20010407-18	WG-1620-FB03-20200109	09 Jan 2020 07:30		13 Jan 2020 15:00	13 Jan 2020 23:11	1
HS20010407-20	WG-1620-MW72B-20200109	09 Jan 2020 08:35		13 Jan 2020 15:00	13 Jan 2020 23:14	1
HS20010407-21	WG-1620-MW23C-20200109	09 Jan 2020 09:40		13 Jan 2020 15:00	13 Jan 2020 23:16	1
<b>Batch ID: 149618 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010407-22	WG-1620-MW53C-20200109	09 Jan 2020 11:15		14 Jan 2020 16:00	14 Jan 2020 21:37	1
HS20010407-23	WG-1620-MW54C-20200109	09 Jan 2020 12:35		14 Jan 2020 16:00	14 Jan 2020 21:52	1
HS20010407-24	WG-1620-MW44A-20200109	09 Jan 2020 13:30		14 Jan 2020 16:00	14 Jan 2020 21:55	1
HS20010407-25	WG-1620-MW36B-20200109	09 Jan 2020 14:40		14 Jan 2020 16:00	14 Jan 2020 21:57	1
HS20010407-26	WG-1620-MW36A-20200109	09 Jan 2020 16:00		14 Jan 2020 16:00	14 Jan 2020 21:59	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149624 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010407-01	WG-1620-MW19C-20200109	09 Jan 2020 10:45		14 Jan 2020 11:05	15 Jan 2020 08:50	1
HS20010407-02	WG-1620-MW51A-20200109	09 Jan 2020 11:50		14 Jan 2020 11:05	15 Jan 2020 09:09	1
HS20010407-03	WG-1620-MW51C-20200109	09 Jan 2020 12:40		14 Jan 2020 11:05	15 Jan 2020 09:28	1
HS20010407-04	WG-1620-MW85C-20200109	09 Jan 2020 13:35		14 Jan 2020 11:05	15 Jan 2020 09:46	1
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25		14 Jan 2020 11:05	15 Jan 2020 17:57	1000
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25		14 Jan 2020 11:05	15 Jan 2020 16:04	100
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25		14 Jan 2020 11:05	15 Jan 2020 14:11	20
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25		14 Jan 2020 11:05	15 Jan 2020 10:05	1
HS20010407-06	WG-1620-MW76C-20200109	09 Jan 2020 15:20		14 Jan 2020 11:05	15 Jan 2020 10:24	1
HS20010407-07	WG-1620-MW81B-20200109	09 Jan 2020 16:20		14 Jan 2020 11:05	15 Jan 2020 10:43	1
HS20010407-08	WG-1620-MW50A-20200109	09 Jan 2020 17:15		14 Jan 2020 11:05	15 Jan 2020 11:02	1
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10		14 Jan 2020 11:05	15 Jan 2020 16:23	100
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10		14 Jan 2020 11:05	15 Jan 2020 18:16	1000
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10		14 Jan 2020 11:05	15 Jan 2020 14:30	10
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10		14 Jan 2020 11:05	15 Jan 2020 11:21	1
HS20010407-10	WG-1620-MW80B-20200107	07 Jan 2020 16:00		14 Jan 2020 11:05	15 Jan 2020 11:40	1
HS20010407-11	WG-1620-FB01-20200107	07 Jan 2020 16:30		14 Jan 2020 11:05	15 Jan 2020 11:59	1
HS20010407-12	WG-1620-FB02-20200108	08 Jan 2020 09:00		14 Jan 2020 11:05	15 Jan 2020 12:17	1
HS20010407-13	WG-1620-MW18C-20200108	08 Jan 2020 10:10		14 Jan 2020 11:05	15 Jan 2020 16:41	100
HS20010407-13	WG-1620-MW18C-20200108	08 Jan 2020 10:10		14 Jan 2020 11:05	15 Jan 2020 18:35	2000
HS20010407-13	WG-1620-MW18C-20200108	08 Jan 2020 10:10		14 Jan 2020 11:05	15 Jan 2020 14:48	10
HS20010407-13	WG-1620-MW18C-20200108	08 Jan 2020 10:10		14 Jan 2020 11:05	15 Jan 2020 12:36	1
HS20010407-14	WG-1620-MW18A-20200108	08 Jan 2020 11:10		14 Jan 2020 11:05	15 Jan 2020 17:00	500
HS20010407-14	WG-1620-MW18A-20200108	08 Jan 2020 11:10		14 Jan 2020 11:05	15 Jan 2020 15:07	50
HS20010407-14	WG-1620-MW18A-20200108	08 Jan 2020 11:10		14 Jan 2020 11:05	15 Jan 2020 12:55	1
HS20010407-15	WG-1620-MW58A-20200108	08 Jan 2020 12:15		14 Jan 2020 11:05	15 Jan 2020 17:19	200
HS20010407-15	WG-1620-MW58A-20200108	08 Jan 2020 12:15		14 Jan 2020 11:05	15 Jan 2020 15:26	20
HS20010407-15	WG-1620-MW58A-20200108	08 Jan 2020 12:15		14 Jan 2020 11:05	15 Jan 2020 13:14	1
HS20010407-16	WG-1620-MW57A-20200108	08 Jan 2020 14:55		14 Jan 2020 11:05	15 Jan 2020 17:38	200
HS20010407-16	WG-1620-MW57A-20200108	08 Jan 2020 14:55		14 Jan 2020 11:05	15 Jan 2020 15:45	20
HS20010407-16	WG-1620-MW57A-20200108	08 Jan 2020 14:55		14 Jan 2020 11:05	15 Jan 2020 13:33	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149670 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010407-17	WG-1620-MW57B-20200108	08 Jan 2020 13:30		15 Jan 2020 10:11	17 Jan 2020 16:07	1000 0
HS20010407-17	WG-1620-MW57B-20200108	08 Jan 2020 13:30		15 Jan 2020 10:11	17 Jan 2020 12:49	1000
HS20010407-17	WG-1620-MW57B-20200108	08 Jan 2020 13:30		15 Jan 2020 10:11	17 Jan 2020 12:29	100
HS20010407-17	WG-1620-MW57B-20200108	08 Jan 2020 13:30		15 Jan 2020 10:11	17 Jan 2020 04:40	10
HS20010407-18	WG-1620-FB03-20200109	09 Jan 2020 07:30		15 Jan 2020 10:11	17 Jan 2020 04:59	1
HS20010407-20	WG-1620-MW72B-20200109	09 Jan 2020 08:35		15 Jan 2020 10:11	17 Jan 2020 16:26	1000 0
HS20010407-20	WG-1620-MW72B-20200109	09 Jan 2020 08:35		15 Jan 2020 10:11	17 Jan 2020 13:48	1000
HS20010407-20	WG-1620-MW72B-20200109	09 Jan 2020 08:35		15 Jan 2020 10:11	17 Jan 2020 13:28	100
HS20010407-20	WG-1620-MW72B-20200109	09 Jan 2020 08:35		15 Jan 2020 10:11	17 Jan 2020 13:08	10
HS20010407-21	WG-1620-MW23C-20200109	09 Jan 2020 09:40		15 Jan 2020 10:11	17 Jan 2020 16:46	100
HS20010407-21	WG-1620-MW23C-20200109	09 Jan 2020 09:40		15 Jan 2020 10:11	17 Jan 2020 14:08	10
HS20010407-22	WG-1620-MW53C-20200109	09 Jan 2020 11:15		15 Jan 2020 10:11	17 Jan 2020 14:27	1
HS20010407-23	WG-1620-MW54C-20200109	09 Jan 2020 12:35		15 Jan 2020 10:11	17 Jan 2020 17:06	5
HS20010407-23	WG-1620-MW54C-20200109	09 Jan 2020 12:35		15 Jan 2020 10:11	17 Jan 2020 14:47	1
HS20010407-24	WG-1620-MW44A-20200109	09 Jan 2020 13:30		15 Jan 2020 10:11	17 Jan 2020 17:26	10
HS20010407-24	WG-1620-MW44A-20200109	09 Jan 2020 13:30		15 Jan 2020 10:11	17 Jan 2020 15:07	1
HS20010407-25	WG-1620-MW36B-20200109	09 Jan 2020 14:40		15 Jan 2020 10:11	17 Jan 2020 15:27	1
HS20010407-26	WG-1620-MW36A-20200109	09 Jan 2020 16:00		15 Jan 2020 10:11	17 Jan 2020 15:47	1
<b>Batch ID: R354349 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010407-02	WG-1620-MW51A-20200109	09 Jan 2020 11:50			14 Jan 2020 01:27	1
HS20010407-03	WG-1620-MW51C-20200109	09 Jan 2020 12:40			14 Jan 2020 03:29	1
HS20010407-07	WG-1620-MW81B-20200109	09 Jan 2020 16:20			14 Jan 2020 03:54	1
HS20010407-08	WG-1620-MW50A-20200109	09 Jan 2020 17:15			14 Jan 2020 04:18	1
HS20010407-10	WG-1620-MW80B-20200107	07 Jan 2020 16:00			14 Jan 2020 04:42	1
HS20010407-22	WG-1620-MW53C-20200109	09 Jan 2020 11:15			14 Jan 2020 05:56	1
HS20010407-23	WG-1620-MW54C-20200109	09 Jan 2020 12:35			14 Jan 2020 06:20	1
<b>Batch ID: R354396 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010407-01	WG-1620-MW19C-20200109	09 Jan 2020 10:45			14 Jan 2020 23:06	1
HS20010407-04	WG-1620-MW85C-20200109	09 Jan 2020 13:35			14 Jan 2020 23:31	1
HS20010407-06	WG-1620-MW76C-20200109	09 Jan 2020 15:20			14 Jan 2020 23:55	1
HS20010407-13	WG-1620-MW18C-20200108	08 Jan 2020 10:10			14 Jan 2020 17:22	10
HS20010407-14	WG-1620-MW18A-20200108	08 Jan 2020 11:10			14 Jan 2020 17:49	25
HS20010407-18	WG-1620-FB03-20200109	09 Jan 2020 07:30			14 Jan 2020 20:40	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: R354399 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010407-11	WG-1620-FB01-20200107	07 Jan 2020 16:30			14 Jan 2020 17:30	1
HS20010407-12	WG-1620-FB02-20200108	08 Jan 2020 09:00			14 Jan 2020 17:55	1
HS20010407-25	WG-1620-MW36B-20200109	09 Jan 2020 14:40			14 Jan 2020 18:20	1
HS20010407-26	WG-1620-MW36A-20200109	09 Jan 2020 16:00			14 Jan 2020 18:44	1
<b>Batch ID: R354465 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010407-05	WG-1620-MW77A-20200109	09 Jan 2020 14:25			15 Jan 2020 12:58	1
HS20010407-09	WG-1620-MW49A-20200107	07 Jan 2020 13:10			15 Jan 2020 14:41	10
HS20010407-15	WG-1620-MW58A-20200108	08 Jan 2020 12:15			15 Jan 2020 13:22	1
HS20010407-16	WG-1620-MW57A-20200108	08 Jan 2020 14:55			15 Jan 2020 13:47	1
HS20010407-17	WG-1620-MW57B-20200108	08 Jan 2020 13:30			15 Jan 2020 15:07	10
HS20010407-20	WG-1620-MW72B-20200109	09 Jan 2020 08:35			15 Jan 2020 15:34	10
HS20010407-21	WG-1620-MW23C-20200109	09 Jan 2020 09:40			15 Jan 2020 14:14	5
HS20010407-24	WG-1620-MW44A-20200109	09 Jan 2020 13:30			15 Jan 2020 12:33	1
HS20010407-27	WG-1620-TB01-20200109	09 Jan 2020 00:00			15 Jan 2020 12:09	1

WorkOrder: HS20010407  
 InstrumentID: FID-11  
 Test Code: TX1005\_W\_Low  
 Test Number: TX1005  
 Test Name: Low-level Texas TPH by TX1005

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	nC6 to nC12	TPH-1005-1	0.25	0.25	0.20	0.50
A	>nC12 to nC28	TPH-1005-2	0.25	0.29	0.20	0.50
A	>nC28 to nC35	TPH-1005-4	0.25	0.25	0.20	0.50
A	Total Petroleum Hydrocarbon	TPH	0.25	0.25	0.20	0.50
S	2-Fluorobiphenyl	321-60-8	0	0	0	0
S	Trifluoromethyl benzene	98-08-8	0	0	0	0

WorkOrder: HS20010407  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20010407  
 InstrumentID: ICPMS04  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000456	0.000400	0.00200

WorkOrder: HS20010407  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010407  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010407  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010



WorkOrder: HS20010407  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00045	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

**Batch ID:** 149585 ( 0 )      **Instrument:** FID-11      **Method:** LOW-LEVEL TEXAS TPH BY TX1005

<b>MBLK</b>		Sample ID: <b>MBLK-149585</b>		Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 18:11</b>				
Client ID:		Run ID: <b>FID-11_354328</b>		SeqNo: <b>5433689</b>		PrepDate: <b>13-Jan-2020</b>		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	U	0.50								
>nC12 to nC28	U	0.50								
>nC28 to nC35	U	0.50								
Total Petroleum Hydrocarbon	U	0.50								
<i>Surr: 2-Fluorobiphenyl</i>	2.667	0	2.5	0	107	70 - 130				
<i>Surr: Trifluoromethyl benzene</i>	2.689	0	2.5	0	108	70 - 130				

<b>LCS</b>		Sample ID: <b>LCS-149585</b>		Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 18:40</b>				
Client ID:		Run ID: <b>FID-11_354328</b>		SeqNo: <b>5433690</b>		PrepDate: <b>13-Jan-2020</b>		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	25.62	0.50	25	0	102	75 - 125				
>nC12 to nC28	23.57	0.50	25	0	94.3	75 - 125				
<i>Surr: 2-Fluorobiphenyl</i>	2.945	0	2.5	0	118	70 - 130				
<i>Surr: Trifluoromethyl benzene</i>	2.74	0	2.5	0	110	70 - 130				

<b>LCSD</b>		Sample ID: <b>LCSD-149585</b>		Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 19:09</b>				
Client ID:		Run ID: <b>FID-11_354328</b>		SeqNo: <b>5433691</b>		PrepDate: <b>13-Jan-2020</b>		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	27.82	0.50	25	0	111	75 - 125	25.62	8.24	20	
>nC12 to nC28	22.91	0.50	25	0	91.6	75 - 125	23.57	2.87	20	
<i>Surr: 2-Fluorobiphenyl</i>	3.003	0	2.5	0	120	70 - 130	2.945	1.96	20	
<i>Surr: Trifluoromethyl benzene</i>	2.941	0	2.5	0	118	70 - 130	2.74	7.09	20	

<b>MS</b>		Sample ID: <b>HS20010449-05MS</b>		Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 20:08</b>				
Client ID:		Run ID: <b>FID-11_354328</b>		SeqNo: <b>5433693</b>		PrepDate: <b>13-Jan-2020</b>		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	25.84	0.48	23.79	0	109	75 - 125				
>nC12 to nC28	21.53	0.48	23.79	0	90.5	75 - 125				
<i>Surr: 2-Fluorobiphenyl</i>	2.988	0	2.379	0	126	70 - 130				
<i>Surr: Trifluoromethyl benzene</i>	2.864	0	2.379	0	120	70 - 130				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

**Batch ID:** 149585 ( 0 )      **Instrument:** FID-11      **Method:** LOW-LEVEL TEXAS TPH BY TX1005

**MSD**      Sample ID: **HS20010449-05MSD**      Units: **mg/L**      Analysis Date: **13-Jan-2020 20:37**  
 Client ID:      Run ID: **FID-11\_354328**      SeqNo: **5433694**      PrepDate: **13-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

nC6 to nC12	26.05	0.48	24.06	0	108	75 - 125	25.84	0.8	20
>nC12 to nC28	22.27	0.48	24.06	0	92.6	75 - 125	21.53	3.38	20
Surr: 2-Fluorobiphenyl	3.117	0	2.406	0	130	70 - 130	2.988	4.23	20
Surr: Trifluoromethyl benzene	3.06	0	2.406	0	127	70 - 130	2.864	6.62	20

The following samples were analyzed in this batch:

HS20010407-02	HS20010407-03	HS20010407-04	HS20010407-05
HS20010407-06	HS20010407-07	HS20010407-08	HS20010407-09
HS20010407-10			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

<b>Batch ID:</b> 149592 ( 0 )		<b>Instrument:</b> ICPMS05		<b>Method:</b> ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-149592</b>	Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 22:05</b>					
Client ID:		Run ID: <b>ICPMS05_354278</b>	SeqNo: <b>5433457</b>	PrepDate: <b>13-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual

Arsenic U 0.00200

<b>LCS</b>	Sample ID: <b>LCS-149592</b>	Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 22:07</b>					
Client ID:		Run ID: <b>ICPMS05_354278</b>	SeqNo: <b>5433458</b>	PrepDate: <b>13-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual

Arsenic 0.04766 0.00200 0.05 0 95.3 80 - 120

<b>MS</b>	Sample ID: <b>HS20010407-01MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 22:14</b>					
Client ID: <b>WG-1620-MW19C-20200109</b>		Run ID: <b>ICPMS05_354278</b>	SeqNo: <b>5433463</b>	PrepDate: <b>13-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual

Arsenic 0.05662 0.00200 0.05 0.001405 110 80 - 120

<b>MSD</b>	Sample ID: <b>HS20010407-01MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 22:16</b>					
Client ID: <b>WG-1620-MW19C-20200109</b>		Run ID: <b>ICPMS05_354278</b>	SeqNo: <b>5433464</b>	PrepDate: <b>13-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual

Arsenic 0.0493 0.00200 0.05 0.001405 95.8 80 - 120 0.05662 13.8 20

<b>PDS</b>	Sample ID: <b>HS20010407-01PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 22:18</b>					
Client ID: <b>WG-1620-MW19C-20200109</b>		Run ID: <b>ICPMS05_354278</b>	SeqNo: <b>5433465</b>	PrepDate: <b>13-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual

Arsenic 0.1219 0.00200 0.1 0.001405 121 75 - 125

<b>SD</b>	Sample ID: <b>HS20010407-01SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>13-Jan-2020 22:11</b>					
Client ID: <b>WG-1620-MW19C-20200109</b>		Run ID: <b>ICPMS05_354278</b>	SeqNo: <b>5433460</b>	PrepDate: <b>13-Jan-2020</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual

Arsenic U 0.0100 0.001405 0 10

<b>The following samples were analyzed in this batch:</b>	HS20010407-01	HS20010407-02	HS20010407-03	HS20010407-04
	HS20010407-05	HS20010407-06	HS20010407-07	HS20010407-08
	HS20010407-09	HS20010407-10	HS20010407-11	HS20010407-12
	HS20010407-13	HS20010407-14	HS20010407-15	HS20010407-16
	HS20010407-17	HS20010407-18	HS20010407-20	HS20010407-21

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149618 ( 0 )		Instrument: ICPMS04		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-149618</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Jan-2020 21:28</b>					
Client ID:		Run ID: <b>ICPMS04_354392</b>	SeqNo: <b>5434990</b>	PrepDate: <b>14-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-149618</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Jan-2020 21:30</b>					
Client ID:		Run ID: <b>ICPMS04_354392</b>	SeqNo: <b>5434991</b>	PrepDate: <b>14-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05218	0.00200	0.05	0	104	80 - 120			
<b>MS</b>	Sample ID: <b>HS20010407-22MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Jan-2020 21:41</b>					
Client ID: <b>WG-1620-MW53C-20200109</b>		Run ID: <b>ICPMS04_354392</b>	SeqNo: <b>5434996</b>	PrepDate: <b>14-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05347	0.00200	0.05	0.000728	105	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20010407-22MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Jan-2020 21:44</b>					
Client ID: <b>WG-1620-MW53C-20200109</b>		Run ID: <b>ICPMS04_354392</b>	SeqNo: <b>5434997</b>	PrepDate: <b>14-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05102	0.00200	0.05	0.000728	101	80 - 120	0.05347	4.68	20
<b>PDS</b>	Sample ID: <b>HS20010407-22PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Jan-2020 21:46</b>					
Client ID: <b>WG-1620-MW53C-20200109</b>		Run ID: <b>ICPMS04_354392</b>	SeqNo: <b>5434998</b>	PrepDate: <b>14-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.1134	0.00200	0.1	0.000728	113	75 - 125			
<b>SD</b>	Sample ID: <b>HS20010407-22SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Jan-2020 21:39</b>					
Client ID: <b>WG-1620-MW53C-20200109</b>		Run ID: <b>ICPMS04_354392</b>	SeqNo: <b>5434995</b>	PrepDate: <b>14-Jan-2020</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	U	0.0100					0.000728	0	10
The following samples were analyzed in this batch:									
	HS20010407-22	HS20010407-23	HS20010407-24	HS20010407-25					
	HS20010407-26								

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149624 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149624	Units: ug/L			Analysis Date: 14-Jan-2020 11:54					
Client ID:	Run ID: SV-6_354345	SeqNo: 5435594	PrepDate: 14-Jan-2020	DF: 1						
Analyte	Result	SQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	3.342	0.20	5	0	66.8	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.576	0.20	5	0	91.5	40 - 125				
<i>Surr: 2-Fluorophenol</i>	3.643	0.20	5	0	72.9	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.605	0.20	5	0	92.1	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	3.631	0.20	5	0	72.6	41 - 120				
<i>Surr: Phenol-d6</i>	3.457	0.20	5	0	69.1	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149624 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149624	Units: ug/L			Analysis Date: 14-Jan-2020 12:13					
Client ID:	Run ID: SV-6_354345	SeqNo: 5435595	PrepDate: 14-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.509	0.20	5	0	90.2	39 - 127				
2,4-Dimethylphenol	4.222	0.20	5	0	84.4	35 - 120				
2,4-Dinitrotoluene	5.427	0.20	5	0	109	50 - 122				
2,6-Dinitrotoluene	5.379	0.20	5	0	108	50 - 120				
2-Chloronaphthalene	4.621	0.20	5	0	92.4	50 - 120				
2-Methylnaphthalene	4.461	0.10	5	0	89.2	50 - 120				
4,6-Dinitro-2-methylphenol	5.218	0.20	5	0	104	25 - 121				
4-Nitrophenol	4.957	1.0	5	0	99.1	30 - 130				
Acenaphthene	4.67	0.10	5	0	93.4	45 - 120				
Acenaphthylene	4.934	0.10	5	0	98.7	47 - 120				
Anthracene	5.352	0.10	5	0	107	45 - 120				
Benz(a)anthracene	4.963	0.10	5	0	99.3	40 - 120				
Benzo(a)pyrene	4.889	0.10	5	0	97.8	45 - 120				
Bis(2-chloroethoxy)methane	4.092	0.20	5	0	81.8	45 - 120				
Bis(2-ethylhexyl)phthalate	5.661	0.20	5	0	113	40 - 139				
Chrysene	5.074	0.10	5	0	101	43 - 120				
Dibenzofuran	4.846	0.10	5	0	96.9	50 - 120				
Di-n-butyl phthalate	5.647	0.20	5	0	113	45 - 123				
Fluoranthene	4.974	0.10	5	0	99.5	45 - 125				
Fluorene	5.112	0.10	5	0	102	49 - 120				
Naphthalene	4.743	0.10	5	0	94.9	45 - 120				
Nitrobenzene	4	0.20	5	0	80.0	44 - 120				
N-Nitrosodiphenylamine	5.745	0.20	5	0	115	40 - 125				
Pentachlorophenol	3.506	0.20	5	0	70.1	19 - 121				
Phenanthrene	4.852	0.10	5	0	97.0	45 - 121				
Phenol	3.759	0.20	5	0	75.2	20 - 124				
Pyrene	5.037	0.10	5	0	101	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.25</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.482</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.6</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.707</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.1</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.481</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.639</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.8</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.704</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.1</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149624 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-149624		Units: ug/L		Analysis Date: 14-Jan-2020 12:31				
Client ID:		Run ID: SV-6_354345		SeqNo: 5435596		PrepDate: 14-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.829	0.20	5	0	96.6	39 - 127	4.509	6.83	20	
2,4-Dimethylphenol	4.34	0.20	5	0	86.8	35 - 120	4.222	2.76	20	
2,4-Dinitrotoluene	5.633	0.20	5	0	113	50 - 122	5.427	3.73	20	
2,6-Dinitrotoluene	5.545	0.20	5	0	111	50 - 120	5.379	3.02	20	
2-Chloronaphthalene	5.172	0.20	5	0	103	50 - 120	4.621	11.2	20	
2-Methylnaphthalene	4.729	0.10	5	0	94.6	50 - 120	4.461	5.83	20	
4,6-Dinitro-2-methylphenol	5.955	0.20	5	0	119	25 - 121	5.218	13.2	30	
4-Nitrophenol	5.539	1.0	5	0	111	30 - 130	4.957	11.1	20	
Acenaphthene	4.912	0.10	5	0	98.2	45 - 120	4.67	5.04	20	
Acenaphthylene	5.119	0.10	5	0	102	47 - 120	4.934	3.68	20	
Anthracene	5.346	0.10	5	0	107	45 - 120	5.352	0.112	20	
Benz(a)anthracene	5.559	0.10	5	0	111	40 - 120	4.963	11.3	20	
Benzo(a)pyrene	5.228	0.10	5	0	105	45 - 120	4.889	6.72	20	
Bis(2-chloroethoxy)methane	4.243	0.20	5	0	84.9	45 - 120	4.092	3.63	20	
Bis(2-ethylhexyl)phthalate	6.111	0.20	5	0	122	40 - 139	5.661	7.65	20	
Chrysene	5.131	0.10	5	0	103	43 - 120	5.074	1.12	20	
Dibenzofuran	5.09	0.10	5	0	102	50 - 120	4.846	4.91	20	
Di-n-butyl phthalate	6.115	0.20	5	0	122	45 - 123	5.647	7.97	20	
Fluoranthene	5.298	0.10	5	0	106	45 - 125	4.974	6.3	20	
Fluorene	5.396	0.10	5	0	108	49 - 120	5.112	5.4	20	
Naphthalene	4.921	0.10	5	0	98.4	45 - 120	4.743	3.68	20	
Nitrobenzene	4.07	0.20	5	0	81.4	44 - 120	4	1.75	20	
N-Nitrosodiphenylamine	5.768	0.20	5	0	115	40 - 125	5.745	0.406	20	
Pentachlorophenol	3.936	0.20	5	0	78.7	19 - 121	3.506	11.6	20	
Phenanthrene	5.216	0.10	5	0	104	45 - 121	4.852	7.23	20	
Phenol	4.067	0.20	5	0	81.3	20 - 124	3.759	7.87	20	
Pyrene	5.47	0.10	5	0	109	40 - 130	5.037	8.23	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.446</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.9</i>	<i>34 - 129</i>	<i>4.25</i>	<i>4.5</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.64</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.8</i>	<i>40 - 125</i>	<i>4.482</i>	<i>3.46</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.254</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.1</i>	<i>20 - 120</i>	<i>3.707</i>	<i>13.8</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.914</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.3</i>	<i>40 - 135</i>	<i>4.481</i>	<i>9.21</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.552</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>71.0</i>	<i>41 - 120</i>	<i>3.639</i>	<i>2.42</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.871</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.4</i>	<i>20 - 120</i>	<i>3.704</i>	<i>4.4</i>	<i>20</i>	



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

**Batch ID:** 149624 ( 0 )      **Instrument:** SV-6      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

The following samples were analyzed in this batch:

HS20010407-01	HS20010407-02	HS20010407-03	HS20010407-04
HS20010407-05	HS20010407-06	HS20010407-07	HS20010407-08
HS20010407-09	HS20010407-10	HS20010407-11	HS20010407-12
HS20010407-13	HS20010407-14	HS20010407-15	HS20010407-16

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149670 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149670	Units: ug/L			Analysis Date: 15-Jan-2020 10:42					
Client ID:	Run ID: SV-7_354430	SeqNo: 5436172		PrepDate: 15-Jan-2020		DF: 1				
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.283</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>106</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.701</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.0</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.816</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.3</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.833</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.7</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.164</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.547</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>70.9</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149670 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149670	Units: ug/L			Analysis Date: 15-Jan-2020 11:21					
Client ID:	Run ID: SV-7_354430	SeqNo: 5436173		PrepDate: 15-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.467	0.20	5	0	69.3	39 - 127				
2,4-Dimethylphenol	3.994	0.20	5	0	79.9	35 - 120				
2,4-Dinitrotoluene	4.458	0.20	5	0	89.2	50 - 122				
2,6-Dinitrotoluene	4.495	0.20	5	0	89.9	50 - 120				
2-Chloronaphthalene	4.578	0.20	5	0	91.6	50 - 120				
2-Methylnaphthalene	4.353	0.10	5	0	87.1	50 - 120				
4,6-Dinitro-2-methylphenol	4.277	0.20	5	0	85.5	25 - 121				
4-Nitrophenol	4.553	1.0	5	0	91.1	30 - 130				
Acenaphthene	4.129	0.10	5	0	82.6	45 - 120				
Acenaphthylene	4.347	0.10	5	0	86.9	47 - 120				
Anthracene	4.583	0.10	5	0	91.7	45 - 120				
Benz(a)anthracene	4.026	0.10	5	0	80.5	40 - 120				
Benzo(a)pyrene	4.379	0.10	5	0	87.6	45 - 120				
Bis(2-chloroethoxy)methane	3.263	0.20	5	0	65.3	45 - 120				
Bis(2-ethylhexyl)phthalate	4.686	0.20	5	0	93.7	40 - 139				
Chrysene	4.461	0.10	5	0	89.2	43 - 120				
Dibenzofuran	4.41	0.10	5	0	88.2	50 - 120				
Di-n-butyl phthalate	4.842	0.20	5	0	96.8	45 - 123				
Fluoranthene	4.447	0.10	5	0	88.9	45 - 125				
Fluorene	4.442	0.10	5	0	88.8	49 - 120				
Naphthalene	4.37	0.10	5	0	87.4	45 - 120				
Nitrobenzene	2.972	0.20	5	0	59.4	44 - 120				
N-Nitrosodiphenylamine	4.559	0.20	5	0	91.2	40 - 125				
Pentachlorophenol	2.864	0.20	5	0	57.3	19 - 121				
Phenanthrene	4.457	0.10	5	0	89.1	45 - 121				
Phenol	3.476	0.20	5	0	69.5	20 - 124				
Pyrene	4.412	0.10	5	0	88.2	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.275</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.513</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.534</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>70.7</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.504</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.1</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.162</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.2</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.471</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.4</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: 149670 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-149670		Units: ug/L		Analysis Date: 15-Jan-2020 11:40				
Client ID:		Run ID: SV-7_354430		SeqNo: 5436174		PrepDate: 15-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.453	0.20	5	0	69.1	39 - 127	3.467	0.41	20	
2,4-Dimethylphenol	3.884	0.20	5	0	77.7	35 - 120	3.994	2.81	20	
2,4-Dinitrotoluene	4.441	0.20	5	0	88.8	50 - 122	4.458	0.366	20	
2,6-Dinitrotoluene	4.362	0.20	5	0	87.2	50 - 120	4.495	3.01	20	
2-Chloronaphthalene	4.546	0.20	5	0	90.9	50 - 120	4.578	0.706	20	
2-Methylnaphthalene	4.339	0.10	5	0	86.8	50 - 120	4.353	0.309	20	
4,6-Dinitro-2-methylphenol	4.522	0.20	5	0	90.4	25 - 121	4.277	5.58	30	
4-Nitrophenol	4.659	1.0	5	0	93.2	30 - 130	4.553	2.31	20	
Acenaphthene	4.083	0.10	5	0	81.7	45 - 120	4.129	1.1	20	
Acenaphthylene	4.304	0.10	5	0	86.1	47 - 120	4.347	1	20	
Anthracene	4.551	0.10	5	0	91.0	45 - 120	4.583	0.688	20	
Benz(a)anthracene	4.096	0.10	5	0	81.9	40 - 120	4.026	1.71	20	
Benzo(a)pyrene	4.392	0.10	5	0	87.8	45 - 120	4.379	0.304	20	
Bis(2-chloroethoxy)methane	3.282	0.20	5	0	65.6	45 - 120	3.263	0.583	20	
Bis(2-ethylhexyl)phthalate	4.83	0.20	5	0	96.6	40 - 139	4.686	3.03	20	
Chrysene	4.427	0.10	5	0	88.5	43 - 120	4.461	0.775	20	
Dibenzofuran	4.305	0.10	5	0	86.1	50 - 120	4.41	2.43	20	
Di-n-butyl phthalate	4.945	0.20	5	0	98.9	45 - 123	4.842	2.11	20	
Fluoranthene	4.517	0.10	5	0	90.3	45 - 125	4.447	1.55	20	
Fluorene	4.431	0.10	5	0	88.6	49 - 120	4.442	0.249	20	
Naphthalene	4.337	0.10	5	0	86.7	45 - 120	4.37	0.757	20	
Nitrobenzene	2.885	0.20	5	0	57.7	44 - 120	2.972	2.99	20	
N-Nitrosodiphenylamine	4.481	0.20	5	0	89.6	40 - 125	4.559	1.74	20	
Pentachlorophenol	2.789	0.20	5	0	55.8	19 - 121	2.864	2.66	20	
Phenanthrene	4.556	0.10	5	0	91.1	45 - 121	4.457	2.19	20	
Phenol	3.665	0.20	5	0	73.3	20 - 124	3.476	5.3	20	
Pyrene	4.528	0.10	5	0	90.6	40 - 130	4.412	2.6	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.124</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>34 - 129</i>	<i>5.275</i>	<i>2.9</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.48</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.6</i>	<i>40 - 125</i>	<i>4.513</i>	<i>0.72</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.744</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.9</i>	<i>20 - 120</i>	<i>3.534</i>	<i>5.78</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.601</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.0</i>	<i>40 - 135</i>	<i>4.504</i>	<i>2.14</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.042</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>60.8</i>	<i>41 - 120</i>	<i>3.162</i>	<i>3.86</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.574</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>71.5</i>	<i>20 - 120</i>	<i>3.471</i>	<i>2.92</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

<b>Batch ID:</b> 149670 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS20010407-17	HS20010407-18	HS20010407-20	HS20010407-21
HS20010407-22	HS20010407-23	HS20010407-24	HS20010407-25
HS20010407-26			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

<b>Batch ID:</b> R354349 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200113</b>	Units: <b>ug/L</b>			Analysis Date: <b>13-Jan-2020 23:49</b>				
Client ID:	Run ID: <b>VOA2_354349</b>	SeqNo: <b>5434111</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200113</b>	Units: <b>ug/L</b>			Analysis Date: <b>13-Jan-2020 23:00</b>				
Client ID:	Run ID: <b>VOA2_354349</b>	SeqNo: <b>5434110</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	18.96	1.0	20	0	94.8	70 - 124			
Benzene	18.2	1.0	20	0	91.0	74 - 120			
Chlorobenzene	19.65	1.0	20	0	98.2	76 - 113			
Ethylbenzene	19.26	1.0	20	0	96.3	77 - 117			
Methylene chloride	18.51	2.0	20	0	92.5	70 - 127			
Toluene	20.36	1.0	20	0	102	77 - 118			
Xylenes, Total	57.72	1.0	60	0	96.2	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.8</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: R354349 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>		Sample ID: <b>HS20010407-02MS</b>		Units: <b>ug/L</b>		Analysis Date: <b>14-Jan-2020 01:51</b>				
Client ID: <b>WG-1620-MW51A-20200109</b>		Run ID: <b>VOA2_354349</b>		SeqNo: <b>5434116</b>		PrepDate:		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	17.57	1.0	20	0	87.8	70 - 127				
Benzene	17.81	1.0	20	0	89.0	70 - 127				
Chlorobenzene	18.54	1.0	20	0	92.7	70 - 114				
Ethylbenzene	18.26	1.0	20	0	91.3	70 - 124				
Methylene chloride	16.96	2.0	20	0	84.8	70 - 128				
Toluene	19.27	1.0	20	0	96.3	70 - 123				
Xylenes, Total	55.25	1.0	60	0	92.1	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>48.69</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010407-02MSD</b>		Units: <b>ug/L</b>		Analysis Date: <b>14-Jan-2020 02:16</b>			
Client ID: <b>WG-1620-MW51A-20200109</b>		Run ID: <b>VOA2_354349</b>		SeqNo: <b>5434117</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	17.38	1.0	20	0	86.9	70 - 127	17.57	1.05	20
Benzene	17.31	1.0	20	0	86.5	70 - 127	17.81	2.84	20
Chlorobenzene	18.55	1.0	20	0	92.8	70 - 114	18.54	0.0788	20
Ethylbenzene	18.19	1.0	20	0	91.0	70 - 124	18.26	0.356	20
Methylene chloride	15.63	2.0	20	0	78.2	70 - 128	16.96	8.15	20
Toluene	19.36	1.0	20	0	96.8	70 - 123	19.27	0.472	20
Xylenes, Total	55.39	1.0	60	0	92.3	70 - 130	55.25	0.244	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>70 - 126</i>	<i>48.95</i>	<i>2.43</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.6</i>	<i>81 - 113</i>	<i>49.03</i>	<i>1.55</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>49.18</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.4</i>	<i>77 - 123</i>	<i>50.01</i>	<i>1.67</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>49.56</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>82 - 127</i>	<i>48.69</i>	<i>1.77</i>	<i>20</i>

The following samples were analyzed in this batch: 

HS20010407-02	HS20010407-03	HS20010407-07	HS20010407-08
HS20010407-10	HS20010407-22	HS20010407-23	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

<b>Batch ID:</b> R354396 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200114</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 14:53</b>				
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435103</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.0</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200214</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 14:04</b>				
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435102</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18	1.0	20	0	90.0	70 - 124			
Benzene	17.81	1.0	20	0	89.1	74 - 120			
Chlorobenzene	18.74	1.0	20	0	93.7	76 - 113			
Ethylbenzene	18.45	1.0	20	0	92.3	77 - 117			
Methylene chloride	18.59	2.0	20	0	93.0	70 - 127			
Toluene	19.5	1.0	20	0	97.5	77 - 118			
Vinyl chloride	17.32	1.0	20	0	86.6	70 - 130			
Xylenes, Total	56.46	1.0	60	0	94.1	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.68</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.44</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>48.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.5</i>	<i>81 - 120</i>			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

Batch ID: R354396 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>	Sample ID: <b>HS20010411-10MS</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 18:13</b>					
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435111</b>		PrepDate:		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	16.57	1.0	20	0	82.9	70 - 127				
Benzene	18.27	1.0	20	0.9903	86.4	70 - 127				
Chlorobenzene	17.64	1.0	20	0	88.2	70 - 114				
Ethylbenzene	22.73	1.0	20	5.13	88.0	70 - 124				
Methylene chloride	16.6	2.0	20	0	83.0	70 - 128				
Toluene	19.16	1.0	20	0	95.8	70 - 123				
Vinyl chloride	16.93	1.0	20	0	84.6	70 - 130				
Xylenes, Total	344.6	1.0	60	319.7	41.5	70 - 130			SO	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.2</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>52.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.53</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>47.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>82 - 127</i>				

<b>MSD</b>	Sample ID: <b>HS20010411-10MSD</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 18:37</b>				
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435112</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	16.84	1.0	20	0	84.2	70 - 127	16.57	1.58	20
Benzene	17.96	1.0	20	0.9903	84.9	70 - 127	18.27	1.68	20
Chlorobenzene	17.63	1.0	20	0	88.1	70 - 114	17.64	0.0493	20
Ethylbenzene	22.1	1.0	20	5.13	84.8	70 - 124	22.73	2.82	20
Methylene chloride	16.24	2.0	20	0	81.2	70 - 128	16.6	2.19	20
Toluene	18.83	1.0	20	0	94.2	70 - 123	19.16	1.72	20
Vinyl chloride	16.95	1.0	20	0	84.8	70 - 130	16.93	0.151	20
Xylenes, Total	335.7	1.0	60	319.7	26.5	70 - 130	344.6	2.64	20 SO
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.9</i>	<i>70 - 126</i>	<i>48.09</i>	<i>1.33</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>52.3</i>	<i>1.73</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.0</i>	<i>77 - 123</i>	<i>48.53</i>	<i>1.09</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>48.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.6</i>	<i>82 - 127</i>	<i>47.83</i>	<i>0.978</i>	<i>20</i>

The following samples were analyzed in this batch: 

HS20010407-01	HS20010407-04	HS20010407-06	HS20010407-13
HS20010407-14	HS20010407-18		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

<b>Batch ID:</b> R354399 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200114</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 12:50</b>				
Client ID:	Run ID: <b>VOA4_354399</b>	SeqNo: <b>5435154</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>57.34</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>115</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.44</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>52.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.53</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200114</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 12:01</b>				
Client ID:	Run ID: <b>VOA4_354399</b>	SeqNo: <b>5435153</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	23.08	1.0	20	0	115	70 - 124			
Benzene	20.28	1.0	20	0	101	74 - 120			
Chlorobenzene	18.89	1.0	20	0	94.5	76 - 113			
Ethylbenzene	19.37	1.0	20	0	96.8	77 - 117			
Methylene chloride	21.61	2.0	20	0	108	70 - 127			
Toluene	21.14	1.0	20	0	106	77 - 118			
Vinyl chloride	21.93	1.0	20	0	110	70 - 130			
Xylenes, Total	61.14	1.0	60	0	102	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>56.63</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>113</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.69</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>53.13</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>106</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

**Batch ID:** R354399 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010411-05MS			Units: ug/L		Analysis Date: 14-Jan-2020 15:47			
Client ID:		Run ID: VOA4_354399			SeqNo: 5435161		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	20.56	1.0	20	0	103	70 - 127				
Benzene	19.01	1.0	20	0	95.0	70 - 127				
Chlorobenzene	17.18	1.0	20	0	85.9	70 - 114				
Ethylbenzene	20.13	1.0	20	0	101	70 - 124				
Methylene chloride	17.46	2.0	20	0	87.3	70 - 128				
Toluene	22.65	1.0	20	0	113	70 - 123				
Vinyl chloride	24.4	1.0	20	0	122	70 - 130				
Xylenes, Total	69.5	1.0	60	0	116	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>57.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>116</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>54.66</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>109</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010411-05MSD			Units: ug/L		Analysis Date: 14-Jan-2020 16:11			
Client ID:		Run ID: VOA4_354399			SeqNo: 5435162		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	22.07	1.0	20	0	110	70 - 127	20.56	7.1	20	
Benzene	19.34	1.0	20	0	96.7	70 - 127	19.01	1.72	20	
Chlorobenzene	17.77	1.0	20	0	88.8	70 - 114	17.18	3.33	20	
Ethylbenzene	18.83	1.0	20	0	94.2	70 - 124	20.13	6.64	20	
Methylene chloride	17.52	2.0	20	0	87.6	70 - 128	17.46	0.372	20	
Toluene	21.23	1.0	20	0	106	70 - 123	22.65	6.47	20	
Vinyl chloride	23.62	1.0	20	0	118	70 - 130	24.4	3.26	20	
Xylenes, Total	63.98	1.0	60	0	107	70 - 130	69.5	8.28	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>57.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>115</i>	<i>70 - 126</i>	<i>57.78</i>	<i>0.68</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>51.46</i>	<i>0.515</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>54.13</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>108</i>	<i>77 - 123</i>	<i>54.66</i>	<i>0.963</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>	<i>50.14</i>	<i>0.198</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010407-11    HS20010407-12    HS20010407-25    HS20010407-26

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

<b>Batch ID:</b> R354465 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200115</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jan-2020 11:45</b>				
Client ID:	Run ID: <b>VOA2_354465</b>	SeqNo: <b>5436426</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.42</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.8</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.35</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.91</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.78</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200215</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jan-2020 10:56</b>				
Client ID:	Run ID: <b>VOA2_354465</b>	SeqNo: <b>5436425</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18.44	1.0	20	0	92.2	70 - 124			
Benzene	17.88	1.0	20	0	89.4	74 - 120			
Chlorobenzene	18.53	1.0	20	0	92.6	76 - 113			
Ethylbenzene	18.15	1.0	20	0	90.8	77 - 117			
Methylene chloride	19.11	2.0	20	0	95.6	70 - 127			
Toluene	19.34	1.0	20	0	96.7	77 - 118			
Vinyl chloride	17.21	1.0	20	0	86.0	70 - 130			
Xylenes, Total	54.69	1.0	60	0	91.2	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.66</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.2</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>48.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QC BATCH REPORT**

**Batch ID:** R354465 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010407-24MS			Units: ug/L		Analysis Date: 15-Jan-2020 16:50			
Client ID: WG-1620-MW44A-20200109		Run ID: VOA2_354465			SeqNo: 5436438		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.97	1.0	20	0	84.8	70 - 127				
Benzene	17.78	1.0	20	0	88.9	70 - 127				
Chlorobenzene	17.83	1.0	20	0	89.2	70 - 114				
Ethylbenzene	17.86	1.0	20	0	89.3	70 - 124				
Methylene chloride	16.86	2.0	20	0	84.3	70 - 128				
Toluene	19.06	1.0	20	0	95.3	70 - 123				
Vinyl chloride	17.26	1.0	20	0	86.3	70 - 130				
Xylenes, Total	54.2	1.0	60	0	90.3	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>47.85</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010407-24MSD			Units: ug/L		Analysis Date: 15-Jan-2020 17:15			
Client ID: WG-1620-MW44A-20200109		Run ID: VOA2_354465			SeqNo: 5436439		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.39	1.0	20	0	87.0	70 - 127	16.97	2.48	20	
Benzene	17.36	1.0	20	0	86.8	70 - 127	17.78	2.36	20	
Chlorobenzene	18.08	1.0	20	0	90.4	70 - 114	17.83	1.36	20	
Ethylbenzene	18.24	1.0	20	0	91.2	70 - 124	17.86	2.1	20	
Methylene chloride	16.44	2.0	20	0	82.2	70 - 128	16.86	2.52	20	
Toluene	19.31	1.0	20	0	96.5	70 - 123	19.06	1.3	20	
Vinyl chloride	16.59	1.0	20	0	82.9	70 - 130	17.26	4	20	
Xylenes, Total	54.24	1.0	60	0	90.4	70 - 130	54.2	0.0713	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.0</i>	<i>70 - 126</i>	<i>48.67</i>	<i>2.44</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>81 - 113</i>	<i>48.96</i>	<i>0.246</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>77 - 123</i>	<i>49.55</i>	<i>2.43</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.82</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>82 - 127</i>	<i>47.85</i>	<i>2.01</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS20010407-05	HS20010407-09	HS20010407-15	HS20010407-16
HS20010407-17	HS20010407-20	HS20010407-21	HS20010407-24
HS20010407-27			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010407

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

---

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010407

Date/Time Received: 10-Jan-2020 09:35
Received by: NDR

Checklist completed by: Nilesh D. Ranchod
eSignature
Date: 10-Jan-2020

Reviewed by: Corey Grandits
eSignature
Date: 13-Jan-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [ ] No [checked]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

Temperature(s)/Thermometer(s): 0.8C/0.8C,1.2C/1.2C,1.0C/1.0C,1.0C/1.0C,0.8C/0.8C IR # 25
C,0.8C/0.8C UC/C

Cooler(s)/Kit(s): 45678/45710/45492/45533/44308/45063

Date/Time sample(s) sent to storage: 01/10/2020 18:00

Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]

Water - pH acceptable upon receipt? Yes [ ] No [checked] N/A [ ]

pH adjusted? Yes [checked] No [ ] N/A [ ]

pH adjusted by: Si Ma

Login Notes: Sx WG1620-MW49A-20200107, WG-1620-MW-80B-20200107 No TPH Vials Rec'd COC= 9 Rec'd =6 Logged in 2 vials for VOC 1 vial for TPH. Sx Date differ WG-1620-FB03-20200109 COC= 1/8/20 Label =1/9/20.

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010407

Date/Time Received: 10-Jan-2020 09:35
Received by: NDR

Checklist completed by: Nilesh D. Ranchod
eSignature
Date: 10-Jan-2020

Reviewed by: Corey Grandits
eSignature
Date: 13-Jan-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [ ] No [checked]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

Temperature(s)/Thermometer(s): C
Cooler(s)/Kit(s):
Date/Time sample(s) sent to storage: 01/10/2020 18:00

- Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [ ] No [checked] N/A [ ]
pH adjusted? Yes [checked] No [ ] N/A [ ]

pH adjusted by: Si Ma

Login Notes: SX # WG-1620-MW18C-20200108
Metal pH>(7) Preserved with 0.5mil HNO3
on 01/10/2020 @ 01:30pm
Lot #312090209 By Sima
After preserved pH (1)

Client Contacted: Date Contacted: Person Contacted:
Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page 1 of 3

COC ID: 206389

HS20010407

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address	

8260\_LL\_W (5632528 Volatile Organics Site Specific)  
 8260\_LL\_W (5632528 VOC Site Specific + V.C.)  
 8270\_LOW\_W (5632532 SemiVolatiles Site specific)  
 ICP\_TW (5636002 5652546 Metals - As)  
 TPH TX 1005

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-MW19C-20200109</del>																
2	WG-1620-MW19C-20200109	1-9-20	1045	W		2		X									
3	WG-1620-MW51A-20200109	↓	1150	W		6		X	X	X							
4	WG-1620-MW51C-20200109		1240	W		9	X		X	X	X						
5	WG-1620-MW85C-20200109		1335	W		9	X		X	X	X						
6	WG-1620-MW77A-20200109		1425	W		9	X		X	X	X						
7	WG-1620-MW76C-20200109		1520	W		9	X		X	X	X						
8	WG-1620-MW81B-20200109		1620	W		9	X		X	X	X						
9	WG-1620-MW50A-20200109		1715	W		9	X		X	X	X						
10	WG-1620-MW49A-20200107		1-7-20	1310	W		9	X		X	X	X					

Sampler(s) Please Print & Sign  
 JOHN BRAYDON  
 Date: 1-10-20 Time: 0935  
 Date: 1-10-20 Time: 0935

Shipment Method: HAND DELIVERED  
 Required Turnaround Time: (Check Box)  
 STD 10 Wk Days  5 Wk Days  2 Wk Days  24 Hour

Results Due Date: \_\_\_\_\_

Notes: UPRR Houston MWPW

Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)
45678	0.3	<input type="checkbox"/> Level II Std QC
45710	1.2	<input type="checkbox"/> Level III Std QC/Raw Data
45492	1.0	<input checked="" type="checkbox"/> TRRP Check st
45535	1.0	<input type="checkbox"/> TRRP Level IV
44303	0.8	<input type="checkbox"/> Other

Reservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

e: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions of the contract.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.



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# Chain of Custody Form

Page 2 of     

COC ID: 206388

HS20010407

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92888	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5632646 Metals - As)
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E	TPH TX 1005
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	F	
Phone	(512) 671-3434	Phone		G	
Fax	(512) 671-3446	Fax		H	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		I	
				J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TB0-202001</del>																
2	WG-1620-MW80B-20200107	1-7-20	1600	Water	+	2		*									
3	WG-1620-FB01-20200107	1-7-20	1630	W		9	X		X	X	X						
4	WG-1620-FB02-20200108	1-8-20	0900	W		6	X		X	X							
5	WG-1620-MW18C-20200108		1010	W		6	X		X	X							
6	WG-1620-MW18A-20200108		1110	W		6		X	X	X							
7	WG-1620-MW58A-20200108		1215	W		6		X	X	X							
8	WG-1620-MW57A-20200108		1455	W		6		X	X	X							
9	WG-1620-MW57B-20200108		1330	W		6		X	X	X							
10	WG-1620-FB03-20200109		0730	W		6	X		X	X							

Sampler(s) Please Print & Sign: JOHN BRAYTON

Relinquished by: [Signature] Date: 1-10-20 Time: 0935

Relinquished by: [Signature] Date: 1-10-20 Time: 0935

Shipment Method: HAND DELIVERY Required Turnaround Time: (Check Box)  STD 10 Wk Days  5 Wk Days  2 Wk Days  24 Hour

Received by: [Signature] Received by (Laboratory): N/A Checked by (Laboratory): [Signature]

Notes: UPRR Houston MWPW

QC Package: (Check One Box Below)

Level II Std QC  TRRP Check st

Level III Std QC/Raw Data  TRRP Level IV

Level IV SV/B4/CLP  Other

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
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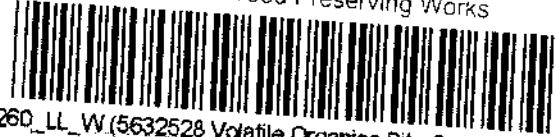
# Chain of Custody Form

Page 3 of 3

COC ID: 206386

HS20010407

Golder Associates Inc.  
Houston TX-Wood Preserving Works



8260\_LL\_W (5632528 Volatile Organics Site Specific)  
8260\_LL\_W (5632528 VOC Site Specific + V.C.)  
8270\_LOW\_W (5632532 SemiVolatiles Site specific)  
ICP\_TW (5636002 5652646 Metals - As)  
TPH TR 1965

m, W  
8

0

ALS Project Manager:

Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR.92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750
City/State/Zip	Round Rock, TX 78684	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TB0-202001																
2	WG-1620-MW49B-20200107	1-7-20	1430	Water	1	2		X									
3	WG-1620-MW72B-20200109	1-9-20	0835	W		9	X	X	X	X							
4	WG-1620-MW23C-20200109		0940	W		6	X	X	X								
5	WG-1620-MW53C-20200109		1115	W		6	X	X	X								
6	WG-1620-MW54C-20200109		1235	W		6	X	X	X								
7	WG-1620-MW44A-20200109		1330	W		6	X	X	X								
8	WG-1620-MW36B-20200109		1440	W		6	X	X	X								
9	WG-1620-MW36A-20200109		1600	W		6	X	X	X								
10	WG-1620-TB01-20200109			W		2	X	X	X								

WILL RESAMPLE WELL  
DUAL IN SAMPLE

Sampler(s) Please Print & Sign: JOHN BEATON

Relinquished by: [Signature] Date: 1-10-20 Time: 09:35

Relinquished by: [Signature] Date: 1-10-20 Time: 07:35

Shipment Method: HAND DELIVER

Received by: [Signature]

Received by (Laboratory): [Signature]

Checked by (Laboratory): [Signature]

Required Turnaround Time: (Check Box)  
 STD 10 Wk Days  
 5 Wk Days  
 2 Wk Days  
 24 Hour  
 Other

Results Due Date: \_\_\_\_\_

Notes: UPRR Houston MWPW

QC Package: (Check One Box Below)  
 Level II Std QC  
 Level III St: QC/Raw Data  
 Level IV SY/643/CLP  
 TRRP Check st  
 TRRP Level IV  
 Other

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

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- Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the...
- The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
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F: +1 281 530 5887

January 20, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010442**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 7 sample(s) on Jan 10, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: DAYNA.FISHER  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

---

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/20/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010442			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149618, 149706, R354396, R354484			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				



<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/20/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010442			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149618, 149706, R354396, R354484			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 01/20/2020
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS20010442
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 149618, 149706, R354396, R354484
ER# <sup>5</sup>	Description	
1	Semivolatiles by Method SW8270, Sample WG-1620-MW35B-20200110, WG-1620-MW83B-20200110: surrogate recoveries could not be determined due to dilution below the calibration range.	
2	Batch R354396, Volatiles by Method SW8260, Sample HS20010411-10, MS and MSD were performed on an unrelated sample	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.</p> <p>O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);          NA = Not Applicable;          NR = Not Reviewed;          R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010442

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010442-02	WQ-1620-TB01-20200110	Water	CG-121719 -74	10-Jan-2020 14:00	10-Jan-2020 15:11	<input type="checkbox"/>
HS20010442-03	WG-1620-FD01-20200110	Water		10-Jan-2020 00:00	10-Jan-2020 15:11	<input type="checkbox"/>
HS20010442-04	WG-1620-MW35B-20200110	Water		10-Jan-2020 12:45	10-Jan-2020 15:11	<input type="checkbox"/>
HS20010442-05	WG-1620-MW35A-20200110	Water		10-Jan-2020 11:20	10-Jan-2020 15:11	<input type="checkbox"/>
HS20010442-06	WG-1620-FB04-20200110	Water		10-Jan-2020 10:30	10-Jan-2020 15:11	<input type="checkbox"/>
HS20010442-07	WG-1620-MW83B-20200110	Water		10-Jan-2020 10:00	10-Jan-2020 15:11	<input type="checkbox"/>
HS20010442-08	WG-1620-MW83C-20200110	Water		10-Jan-2020 09:00	10-Jan-2020 15:11	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB01-20200110  
 Collection Date: 10-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	14-Jan-2020 22:42
Benzene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 22:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 22:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	14-Jan-2020 22:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	14-Jan-2020 22:42
Toluene	U		0.00020	0.0010	mg/L	1	14-Jan-2020 22:42
Vinyl chloride	U		0.00020	0.0010	mg/L	1	14-Jan-2020 22:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	14-Jan-2020 22:42
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>97.4</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 22:42</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>93.4</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 22:42</i>
<i>Surr: Dibromofluoromethane</i>		<i>100</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 22:42</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Jan-2020 22:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200110  
 Collection Date: 10-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	15-Jan-2020 15:45
<b>Benzene</b>	<b>0.016</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 15:45
Chlorobenzene		U	0.00030	0.0010	mg/L	1	15-Jan-2020 15:45
<b>Ethylbenzene</b>	<b>0.083</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 15:45
Methylene chloride		U	0.0010	0.0020	mg/L	1	15-Jan-2020 15:45
Toluene		U	0.00020	0.0010	mg/L	1	15-Jan-2020 15:45
Vinyl chloride		U	0.00020	0.0010	mg/L	1	15-Jan-2020 15:45
<b>Xylenes, Total</b>	<b>0.041</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 15:45
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>121</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: Dibromofluoromethane</i>	<i>110</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 15:45</i>
<i>Surr: Toluene-d8</i>	<i>99.1</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 15:45</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200110  
 Collection Date: 10-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	20-Jan-2020 09:44
<b>2,4-Dimethylphenol</b>	<b>0.00021</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jan-2020 09:44
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	20-Jan-2020 09:44
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	20-Jan-2020 09:44
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	20-Jan-2020 09:44
<b>2-Methylnaphthalene</b>	<b>0.062</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:03
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	20-Jan-2020 09:44
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	20-Jan-2020 09:44
<b>Acenaphthene</b>	<b>0.033</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:03
<b>Acenaphthylene</b>	<b>0.00027</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 09:44
<b>Anthracene</b>	<b>0.0018</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 09:44
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	20-Jan-2020 09:44
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	20-Jan-2020 09:44
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	20-Jan-2020 09:44
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	20-Jan-2020 09:44
Chrysene		U	0.000021	0.00010	mg/L	1	20-Jan-2020 09:44
<b>Dibenzofuran</b>	<b>0.031</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:03
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	20-Jan-2020 09:44
<b>Fluoranthene</b>	<b>0.00093</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 09:44
<b>Fluorene</b>	<b>0.015</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:03
<b>Naphthalene</b>	<b>0.13</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	20-Jan-2020 11:20
Nitrobenzene		U	0.000024	0.00020	mg/L	1	20-Jan-2020 09:44
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	20-Jan-2020 09:44
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	20-Jan-2020 09:44
<b>Phenanthrene</b>	<b>0.013</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:03
Phenol		U	0.000035	0.00020	mg/L	1	20-Jan-2020 09:44
<b>Pyrene</b>	<b>0.00045</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 09:44
Surr: 2,4,6-Tribromophenol	49.5	J		34-129	%REC	20	20-Jan-2020 11:20
Surr: 2,4,6-Tribromophenol	81.1			34-129	%REC	1	20-Jan-2020 09:44
Surr: 2,4,6-Tribromophenol	62.7			34-129	%REC	10	20-Jan-2020 10:03
Surr: 2-Fluorobiphenyl	80.1			40-125	%REC	1	20-Jan-2020 09:44
Surr: 2-Fluorobiphenyl	72.0			40-125	%REC	10	20-Jan-2020 10:03
Surr: 2-Fluorobiphenyl	72.9	J		40-125	%REC	20	20-Jan-2020 11:20
Surr: 2-Fluorophenol	78.7	J		20-120	%REC	20	20-Jan-2020 11:20
Surr: 2-Fluorophenol	85.0			20-120	%REC	1	20-Jan-2020 09:44
Surr: 2-Fluorophenol	71.8			20-120	%REC	10	20-Jan-2020 10:03
Surr: 4-Terphenyl-d14	94.6			40-135	%REC	1	20-Jan-2020 09:44
Surr: 4-Terphenyl-d14	78.2			40-135	%REC	10	20-Jan-2020 10:03
Surr: 4-Terphenyl-d14	79.1	J		40-135	%REC	20	20-Jan-2020 11:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200110  
 Collection Date: 10-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	107			41-120	%REC	1	20-Jan-2020 09:44
Surr: Nitrobenzene-d5	83.7			41-120	%REC	10	20-Jan-2020 10:03
Surr: Nitrobenzene-d5	96.3			41-120	%REC	20	20-Jan-2020 11:20
Surr: Phenol-d6	72.7	J		20-120	%REC	20	20-Jan-2020 11:20
Surr: Phenol-d6	74.0			20-120	%REC	1	20-Jan-2020 09:44
Surr: Phenol-d6	73.6			20-120	%REC	10	20-Jan-2020 10:03
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
Arsenic	0.0159		0.000400	0.00200	mg/L	1	14-Jan-2020 22:01

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35B-20200110  
 Collection Date: 10-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	15-Jan-2020 16:10
<b>Benzene</b>	<b>0.025</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:10
Chlorobenzene		U	0.00030	0.0010	mg/L	1	15-Jan-2020 16:10
<b>Ethylbenzene</b>	<b>0.096</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:10
Methylene chloride		U	0.0010	0.0020	mg/L	1	15-Jan-2020 16:10
Toluene		U	0.00020	0.0010	mg/L	1	15-Jan-2020 16:10
Vinyl chloride		U	0.00020	0.0010	mg/L	1	15-Jan-2020 16:10
<b>Xylenes, Total</b>	<b>0.058</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:10
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>121</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:10</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>103</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:10</i>
<i>Surr: Dibromofluoromethane</i>	<i>110</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:10</i>
<i>Surr: Toluene-d8</i>	<i>98.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:10</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35B-20200110  
 Collection Date: 10-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 18:05
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 18:05
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 18:05
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 18:05
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 18:05
<b>2-Methylnaphthalene</b>	<b>0.14</b>		<b>0.00095</b>	<b>0.0050</b>	<b>mg/L</b>	50	20-Jan-2020 11:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 18:05
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 18:05
<b>Acenaphthene</b>	<b>0.059</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:23
<b>Acenaphthylene</b>	<b>0.00038</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 18:05
<b>Anthracene</b>	<b>0.0029</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 18:05
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 18:05
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 18:05
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 18:05
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 18:05
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 18:05
<b>Dibenzofuran</b>	<b>0.054</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:23
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 18:05
<b>Fluoranthene</b>	<b>0.0015</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 18:05
<b>Fluorene</b>	<b>0.028</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:23
<b>Naphthalene</b>	<b>2.0</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	250	20-Jan-2020 12:18
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 18:05
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 18:05
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 18:05
<b>Phenanthrene</b>	<b>0.024</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 10:23
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 18:05
<b>Pyrene</b>	<b>0.00070</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 18:05
<i>Surr: 2,4,6-Tribromophenol</i>	78.3			34-129	%REC	1	17-Jan-2020 18:05
<i>Surr: 2,4,6-Tribromophenol</i>	86.2			34-129	%REC	10	20-Jan-2020 10:23
<i>Surr: 2,4,6-Tribromophenol</i>	71.2	J		34-129	%REC	50	20-Jan-2020 11:40
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	250	20-Jan-2020 12:18
<i>Surr: 2-Fluorobiphenyl</i>	87.2	J		40-125	%REC	50	20-Jan-2020 11:40
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	250	20-Jan-2020 12:18
<i>Surr: 2-Fluorobiphenyl</i>	73.4			40-125	%REC	10	20-Jan-2020 10:23
<i>Surr: 2-Fluorobiphenyl</i>	74.3			40-125	%REC	1	17-Jan-2020 18:05
<i>Surr: 2-Fluorophenol</i>	70.9			20-120	%REC	1	17-Jan-2020 18:05
<i>Surr: 2-Fluorophenol</i>	76.9			20-120	%REC	10	20-Jan-2020 10:23
<i>Surr: 2-Fluorophenol</i>	83.1	J		20-120	%REC	50	20-Jan-2020 11:40
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	250	20-Jan-2020 12:18

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35B-20200110  
 Collection Date: 10-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
Surr: 4-Terphenyl-d14	88.9	J		40-135	%REC	50	20-Jan-2020 11:40
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	250	20-Jan-2020 12:18
Surr: 4-Terphenyl-d14	88.6			40-135	%REC	10	20-Jan-2020 10:23
Surr: 4-Terphenyl-d14	88.9			40-135	%REC	1	17-Jan-2020 18:05
Surr: Nitrobenzene-d5	95.7			41-120	%REC	10	20-Jan-2020 10:23
Surr: Nitrobenzene-d5	102	J		41-120	%REC	50	20-Jan-2020 11:40
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	250	20-Jan-2020 12:18
Surr: Nitrobenzene-d5	68.2			41-120	%REC	1	17-Jan-2020 18:05
Surr: Phenol-d6	69.0			20-120	%REC	1	17-Jan-2020 18:05
Surr: Phenol-d6	67.6			20-120	%REC	10	20-Jan-2020 10:23
Surr: Phenol-d6	80.1	J		20-120	%REC	50	20-Jan-2020 11:40
Surr: Phenol-d6	0	JS		20-120	%REC	250	20-Jan-2020 12:18
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
Arsenic	0.0120		0.000400	0.00200	mg/L	1	14-Jan-2020 22:03

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35A-20200110  
 Collection Date: 10-Jan-2020 11:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 16:35
Benzene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 16:35
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 16:35
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 16:35
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 16:35
Toluene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 16:35
Vinyl chloride	U		0.00020	0.0010	mg/L	1	15-Jan-2020 16:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jan-2020 16:35
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>117</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:35</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.2</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:35</i>
<i>Surr: Dibromofluoromethane</i>		<i>109</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:35</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:35</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW35A-20200110  
 Collection Date: 10-Jan-2020 11:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 18:25
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 18:25
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 18:25
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 18:25
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 18:25
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 18:25
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 18:25
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 18:25
Acenaphthene	U		0.000027	0.00010	mg/L	1	17-Jan-2020 18:25
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 18:25
Anthracene	U		0.000014	0.00010	mg/L	1	17-Jan-2020 18:25
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 18:25
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 18:25
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 18:25
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000061</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 18:25
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 18:25
Dibenzofuran	U		0.000020	0.00010	mg/L	1	17-Jan-2020 18:25
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 18:25
Fluoranthene	U		0.000010	0.00010	mg/L	1	17-Jan-2020 18:25
<b>Fluorene</b>	<b>0.000065</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 18:25
<b>Naphthalene</b>	<b>0.00020</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 18:25
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 18:25
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 18:25
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 18:25
Phenanthrene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 18:25
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 18:25
Pyrene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 18:25
<i>Surr: 2,4,6-Tribromophenol</i>	<i>81.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 18:25</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>76.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 18:25</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 18:25</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>90.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 18:25</i>
<i>Surr: Nitrobenzene-d5</i>	<i>72.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 18:25</i>
<i>Surr: Phenol-d6</i>	<i>71.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 18:25</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.0219</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	14-Jan-2020 22:06

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB04-20200110  
 Collection Date: 10-Jan-2020 10:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:51
Benzene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:51
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:51
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:51
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 12:51
Toluene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:51
Vinyl chloride	U		0.00020	0.0010	mg/L	1	15-Jan-2020 12:51
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jan-2020 12:51
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>119</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:51</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.3</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:51</i>
<i>Surr: Dibromofluoromethane</i>		<i>109</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:51</i>
<i>Surr: Toluene-d8</i>		<i>99.9</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 12:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB04-20200110  
 Collection Date: 10-Jan-2020 10:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jan-2020 10:42
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	20-Jan-2020 10:42
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jan-2020 10:42
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jan-2020 10:42
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jan-2020 10:42
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	20-Jan-2020 10:42
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jan-2020 10:42
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jan-2020 10:42
Acenaphthene	U		0.000027	0.00010	mg/L	1	20-Jan-2020 10:42
Acenaphthylene	U		0.000015	0.00010	mg/L	1	20-Jan-2020 10:42
Anthracene	U		0.000014	0.00010	mg/L	1	20-Jan-2020 10:42
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jan-2020 10:42
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jan-2020 10:42
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jan-2020 10:42
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jan-2020 10:42
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jan-2020 10:42
Dibenzofuran	U		0.000020	0.00010	mg/L	1	20-Jan-2020 10:42
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jan-2020 10:42
Fluoranthene	U		0.000010	0.00010	mg/L	1	20-Jan-2020 10:42
Fluorene	U		0.000030	0.00010	mg/L	1	20-Jan-2020 10:42
<b>Naphthalene</b>		<b>0.00020</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 10:42
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jan-2020 10:42
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jan-2020 10:42
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jan-2020 10:42
Phenanthrene	U		0.000021	0.00010	mg/L	1	20-Jan-2020 10:42
Phenol	U		0.000035	0.00020	mg/L	1	20-Jan-2020 10:42
Pyrene	U		0.000019	0.00010	mg/L	1	20-Jan-2020 10:42
<i>Surr: 2,4,6-Tribromophenol</i>		<i>67.6</i>		<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 10:42</i>
<i>Surr: 2-Fluorobiphenyl</i>		<i>93.0</i>		<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 10:42</i>
<i>Surr: 2-Fluorophenol</i>		<i>73.7</i>		<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 10:42</i>
<i>Surr: 4-Terphenyl-d14</i>		<i>93.4</i>		<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 10:42</i>
<i>Surr: Nitrobenzene-d5</i>		<i>89.7</i>		<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 10:42</i>
<i>Surr: Phenol-d6</i>		<i>79.9</i>		<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 10:42</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
Arsenic	U		0.000400	0.00200	mg/L	1	14-Jan-2020 22:08

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83B-20200110  
 Collection Date: 10-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 16:59
<b>Benzene</b>	<b>0.021</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:59
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 16:59
<b>Ethylbenzene</b>	<b>0.078</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:59
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 16:59
<b>Toluene</b>	<b>0.0052</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:59
Vinyl chloride	U		0.00020	0.0010	mg/L	1	15-Jan-2020 16:59
<b>Xylenes, Total</b>	<b>0.098</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Jan-2020 16:59
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>124</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:59</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>105</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>109</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:59</i>
<i>Surr: Toluene-d8</i>	<i>99.1</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Jan-2020 16:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83B-20200110  
 Collection Date: 10-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	17-Jan-2020 19:04
<b>2,4-Dimethylphenol</b>	<b>0.000072</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	17-Jan-2020 19:04
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	17-Jan-2020 19:04
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	17-Jan-2020 19:04
<b>2-Methylnaphthalene</b>	<b>0.078</b>		<b>0.00019</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 11:01
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	17-Jan-2020 19:04
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	17-Jan-2020 19:04
<b>Acenaphthene</b>	<b>0.066</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 11:01
<b>Acenaphthylene</b>	<b>0.00047</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
<b>Anthracene</b>	<b>0.0039</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
<b>Benz(a)anthracene</b>	<b>0.000070</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	17-Jan-2020 19:04
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	17-Jan-2020 19:04
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	17-Jan-2020 19:04
<b>Chrysene</b>	<b>0.000065</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
<b>Dibenzofuran</b>	<b>0.023</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 11:01
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	17-Jan-2020 19:04
<b>Fluoranthene</b>	<b>0.0033</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
<b>Fluorene</b>	<b>0.024</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 11:01
<b>Naphthalene</b>	<b>1.2</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	250	20-Jan-2020 11:59
Nitrobenzene		U	0.000024	0.00020	mg/L	1	17-Jan-2020 19:04
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	17-Jan-2020 19:04
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	17-Jan-2020 19:04
<b>Phenanthrene</b>	<b>0.021</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	20-Jan-2020 11:01
Phenol		U	0.000035	0.00020	mg/L	1	17-Jan-2020 19:04
<b>Pyrene</b>	<b>0.0022</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:04
Surr: 2,4,6-Tribromophenol	72.8			34-129	%REC	1	17-Jan-2020 19:04
Surr: 2,4,6-Tribromophenol	59.3			34-129	%REC	10	20-Jan-2020 11:01
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	250	20-Jan-2020 11:59
Surr: 2-Fluorobiphenyl	60.1			40-125	%REC	10	20-Jan-2020 11:01
Surr: 2-Fluorobiphenyl	64.9			40-125	%REC	1	17-Jan-2020 19:04
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	250	20-Jan-2020 11:59
Surr: 2-Fluorophenol	63.7			20-120	%REC	10	20-Jan-2020 11:01
Surr: 2-Fluorophenol	0	JS		20-120	%REC	250	20-Jan-2020 11:59
Surr: 2-Fluorophenol	70.7			20-120	%REC	1	17-Jan-2020 19:04
Surr: 4-Terphenyl-d14	88.4			40-135	%REC	1	17-Jan-2020 19:04
Surr: 4-Terphenyl-d14	79.1			40-135	%REC	10	20-Jan-2020 11:01
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	250	20-Jan-2020 11:59

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83B-20200110  
 Collection Date: 10-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	82.9			41-120	%REC	10	20-Jan-2020 11:01
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	250	20-Jan-2020 11:59
Surr: Nitrobenzene-d5	116			41-120	%REC	1	17-Jan-2020 19:04
Surr: Phenol-d6	62.3			20-120	%REC	1	17-Jan-2020 19:04
Surr: Phenol-d6	54.0			20-120	%REC	10	20-Jan-2020 11:01
Surr: Phenol-d6	0	JS		20-120	%REC	250	20-Jan-2020 11:59
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
Arsenic	0.0709		0.000400	0.00200	mg/L	1	14-Jan-2020 22:10

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83C-20200110  
 Collection Date: 10-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	15-Jan-2020 17:24
Benzene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 17:24
Chlorobenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 17:24
Ethylbenzene	U		0.00030	0.0010	mg/L	1	15-Jan-2020 17:24
Methylene chloride	U		0.0010	0.0020	mg/L	1	15-Jan-2020 17:24
Toluene	U		0.00020	0.0010	mg/L	1	15-Jan-2020 17:24
Vinyl chloride	U		0.00020	0.0010	mg/L	1	15-Jan-2020 17:24
Xylenes, Total	U		0.00030	0.0010	mg/L	1	15-Jan-2020 17:24
<i>Surr: 1,2-Dichloroethane-d4</i>		122		70-126	%REC	1	15-Jan-2020 17:24
<i>Surr: 4-Bromofluorobenzene</i>		98.8		81-113	%REC	1	15-Jan-2020 17:24
<i>Surr: Dibromofluoromethane</i>		110		77-123	%REC	1	15-Jan-2020 17:24
<i>Surr: Toluene-d8</i>		102		82-127	%REC	1	15-Jan-2020 17:24

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW83C-20200110  
 Collection Date: 10-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010442  
 Lab ID:HS20010442-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 19:24
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 19:24
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 19:24
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 19:24
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 19:24
<b>2-Methylnaphthalene</b>	<b>0.00021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:24
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 19:24
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 19:24
<b>Acenaphthene</b>	<b>0.00018</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:24
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 19:24
Anthracene	U		0.000014	0.00010	mg/L	1	17-Jan-2020 19:24
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 19:24
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 19:24
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 19:24
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 19:24
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 19:24
<b>Dibenzofuran</b>	<b>0.00015</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:24
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 19:24
Fluoranthene	U		0.000010	0.00010	mg/L	1	17-Jan-2020 19:24
<b>Fluorene</b>	<b>0.00012</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:24
<b>Naphthalene</b>	<b>0.0018</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:24
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 19:24
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 19:24
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 19:24
<b>Phenanthrene</b>	<b>0.00017</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:24
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 19:24
Pyrene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 19:24
<i>Surr: 2,4,6-Tribromophenol</i>	<i>66.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>59.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>62.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:24</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>80.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:24</i>
<i>Surr: Nitrobenzene-d5</i>	<i>63.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:24</i>
<i>Surr: Phenol-d6</i>	<i>61.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:24</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 14-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00564</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	14-Jan-2020 22:12

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**Batch ID:** 149618      **Start Date:** 14 Jan 2020 08:30      **End Date:** 14 Jan 2020 16:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010442-03		10 (mL)	10 (mL)	1
HS20010442-04		10 (mL)	10 (mL)	1
HS20010442-05		10 (mL)	10 (mL)	1
HS20010442-06		10 (mL)	10 (mL)	1
HS20010442-07		10 (mL)	10 (mL)	1
HS20010442-08		10 (mL)	10 (mL)	1

**Batch ID:** 149706      **Start Date:** 16 Jan 2020 07:00      **End Date:** 16 Jan 2020 11:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010442-03	1	1000 (mL)	1 (mL)	0.001
HS20010442-04	1	1000 (mL)	1 (mL)	0.001
HS20010442-05	1	1000 (mL)	1 (mL)	0.001
HS20010442-06	1	1000 (mL)	1 (mL)	0.001
HS20010442-07	1	1000 (mL)	1 (mL)	0.001
HS20010442-08	1	1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149618 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010442-03	WG-1620-FD01-20200110	10 Jan 2020 00:00		14 Jan 2020 16:00	14 Jan 2020 22:01	1
HS20010442-04	WG-1620-MW35B-20200110	10 Jan 2020 12:45		14 Jan 2020 16:00	14 Jan 2020 22:03	1
HS20010442-05	WG-1620-MW35A-20200110	10 Jan 2020 11:20		14 Jan 2020 16:00	14 Jan 2020 22:06	1
HS20010442-06	WG-1620-FB04-20200110	10 Jan 2020 10:30		14 Jan 2020 16:00	14 Jan 2020 22:08	1
HS20010442-07	WG-1620-MW83B-20200110	10 Jan 2020 10:00		14 Jan 2020 16:00	14 Jan 2020 22:10	1
HS20010442-08	WG-1620-MW83C-20200110	10 Jan 2020 09:00		14 Jan 2020 16:00	14 Jan 2020 22:12	1
<b>Batch ID: 149706 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010442-03	WG-1620-FD01-20200110	10 Jan 2020 00:00		16 Jan 2020 10:59	20 Jan 2020 11:20	20
HS20010442-03	WG-1620-FD01-20200110	10 Jan 2020 00:00		16 Jan 2020 10:59	20 Jan 2020 10:03	10
HS20010442-03	WG-1620-FD01-20200110	10 Jan 2020 00:00		16 Jan 2020 10:59	20 Jan 2020 09:44	1
HS20010442-04	WG-1620-MW35B-20200110	10 Jan 2020 12:45		16 Jan 2020 10:59	20 Jan 2020 11:40	50
HS20010442-04	WG-1620-MW35B-20200110	10 Jan 2020 12:45		16 Jan 2020 10:59	20 Jan 2020 12:18	250
HS20010442-04	WG-1620-MW35B-20200110	10 Jan 2020 12:45		16 Jan 2020 10:59	20 Jan 2020 10:23	10
HS20010442-04	WG-1620-MW35B-20200110	10 Jan 2020 12:45		16 Jan 2020 10:59	17 Jan 2020 18:05	1
HS20010442-05	WG-1620-MW35A-20200110	10 Jan 2020 11:20		16 Jan 2020 10:59	17 Jan 2020 18:25	1
HS20010442-06	WG-1620-FB04-20200110	10 Jan 2020 10:30		16 Jan 2020 10:59	20 Jan 2020 10:42	1
HS20010442-07	WG-1620-MW83B-20200110	10 Jan 2020 10:00		16 Jan 2020 10:59	20 Jan 2020 11:59	250
HS20010442-07	WG-1620-MW83B-20200110	10 Jan 2020 10:00		16 Jan 2020 10:59	20 Jan 2020 11:01	10
HS20010442-07	WG-1620-MW83B-20200110	10 Jan 2020 10:00		16 Jan 2020 10:59	17 Jan 2020 19:04	1
HS20010442-08	WG-1620-MW83C-20200110	10 Jan 2020 09:00		16 Jan 2020 10:59	17 Jan 2020 19:24	1
<b>Batch ID: R354396 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010442-02	WQ-1620-TB01-20200110	10 Jan 2020 14:00			14 Jan 2020 22:42	1
<b>Batch ID: R354484 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010442-03	WG-1620-FD01-20200110	10 Jan 2020 00:00			15 Jan 2020 15:45	1
HS20010442-04	WG-1620-MW35B-20200110	10 Jan 2020 12:45			15 Jan 2020 16:10	1
HS20010442-05	WG-1620-MW35A-20200110	10 Jan 2020 11:20			15 Jan 2020 16:35	1
HS20010442-06	WG-1620-FB04-20200110	10 Jan 2020 10:30			15 Jan 2020 12:51	1
HS20010442-07	WG-1620-MW83B-20200110	10 Jan 2020 10:00			15 Jan 2020 16:59	1
HS20010442-08	WG-1620-MW83C-20200110	10 Jan 2020 09:00			15 Jan 2020 17:24	1

WorkOrder: HS20010442  
 InstrumentID: ICPMS04  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000456	0.000400	0.00200

WorkOrder: HS20010442  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010442  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010



WorkOrder: HS20010442  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00045	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

**Batch ID:** 149618 ( 0)      **Instrument:** ICPMS04      **Method:** ICP-MS METALS BY SW6020A

**MBLK**      Sample ID: **MBLK-149618**      Units: **mg/L**      Analysis Date: **14-Jan-2020 21:28**  
 Client ID:      Run ID: **ICPMS04\_354392**      SeqNo: **5434990**      PrepDate: **14-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      U      0.00200

**LCS**      Sample ID: **LCS-149618**      Units: **mg/L**      Analysis Date: **14-Jan-2020 21:30**  
 Client ID:      Run ID: **ICPMS04\_354392**      SeqNo: **5434991**      PrepDate: **14-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.05218      0.00200      0.05      0      104      80 - 120

**MS**      Sample ID: **HS20010407-22MS**      Units: **mg/L**      Analysis Date: **14-Jan-2020 21:41**  
 Client ID:      Run ID: **ICPMS04\_354392**      SeqNo: **5434996**      PrepDate: **14-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.05347      0.00200      0.05      0.000728      105      80 - 120

**MSD**      Sample ID: **HS20010407-22MSD**      Units: **mg/L**      Analysis Date: **14-Jan-2020 21:44**  
 Client ID:      Run ID: **ICPMS04\_354392**      SeqNo: **5434997**      PrepDate: **14-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.05102      0.00200      0.05      0.000728      101      80 - 120      0.05347      4.68      20

**PDS**      Sample ID: **HS20010407-22PDS**      Units: **mg/L**      Analysis Date: **14-Jan-2020 21:46**  
 Client ID:      Run ID: **ICPMS04\_354392**      SeqNo: **5434998**      PrepDate: **14-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.1134      0.00200      0.1      0.000728      113      75 - 125

**SD**      Sample ID: **HS20010407-22SD**      Units: **mg/L**      Analysis Date: **14-Jan-2020 21:39**  
 Client ID:      Run ID: **ICPMS04\_354392**      SeqNo: **5434995**      PrepDate: **14-Jan-2020**      DF: **5**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %D      %D Limit Qual

Arsenic      U      0.0100      0.000728      0      10

**The following samples were analyzed in this batch:** HS20010442-03      HS20010442-04      HS20010442-05      HS20010442-06  
 HS20010442-07      HS20010442-08

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

Batch ID: 149706 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149706	Units: ug/L			Analysis Date: 17-Jan-2020 11:10					
Client ID:	Run ID: SV-7_354580	SeqNo: 5438870	PrepDate: 16-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	3.801	0.20	5	0	76.0	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.834	0.20	5	0	96.7	40 - 125				
<i>Surr: 2-Fluorophenol</i>	5.028	0.20	5	0	101	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	5.043	0.20	5	0	101	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	4.556	0.20	5	0	91.1	41 - 120				
<i>Surr: Phenol-d6</i>	4.987	0.20	5	0	99.7	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

Batch ID: 149706 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149706	Units: ug/L			Analysis Date: 17-Jan-2020 11:29					
Client ID:	Run ID: SV-7_354580	SeqNo: 5438871		PrepDate: 16-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.272	0.20	5	0	105	39 - 127				
2,4-Dimethylphenol	4.201	0.20	5	0	84.0	35 - 120				
2,4-Dinitrotoluene	4.666	0.20	5	0	93.3	50 - 122				
2,6-Dinitrotoluene	4.42	0.20	5	0	88.4	50 - 120				
2-Chloronaphthalene	4.636	0.20	5	0	92.7	50 - 120				
2-Methylnaphthalene	4.065	0.10	5	0	81.3	50 - 120				
4,6-Dinitro-2-methylphenol	5.471	0.20	5	0	109	25 - 121				
4-Nitrophenol	4.842	1.0	5	0	96.8	30 - 130				
Acenaphthene	4.398	0.10	5	0	88.0	45 - 120				
Acenaphthylene	4.139	0.10	5	0	82.8	47 - 120				
Anthracene	4.609	0.10	5	0	92.2	45 - 120				
Benz(a)anthracene	4.284	0.10	5	0	85.7	40 - 120				
Benzo(a)pyrene	4.099	0.10	5	0	82.0	45 - 120				
Bis(2-chloroethoxy)methane	4.466	0.20	5	0	89.3	45 - 120				
Bis(2-ethylhexyl)phthalate	4.202	0.20	5	0	84.0	40 - 139				
Chrysene	4.745	0.10	5	0	94.9	43 - 120				
Dibenzofuran	4.287	0.10	5	0	85.7	50 - 120				
Di-n-butyl phthalate	4.89	0.20	5	0	97.8	45 - 123				
Fluoranthene	4.866	0.10	5	0	97.3	45 - 125				
Fluorene	4.384	0.10	5	0	87.7	49 - 120				
Naphthalene	4.293	0.10	5	0	85.9	45 - 120				
Nitrobenzene	4.568	0.20	5	0	91.4	44 - 120				
N-Nitrosodiphenylamine	4.806	0.20	5	0	96.1	40 - 125				
Pentachlorophenol	3.923	0.20	5	0	78.5	19 - 121				
Phenanthrene	4.546	0.10	5	0	90.9	45 - 121				
Phenol	4.568	0.20	5	0	91.4	20 - 124				
Pyrene	4.445	0.10	5	0	88.9	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.171</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.4</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.609</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.848</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.705</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.1</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.603</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.1</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.591</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

**Batch ID:** 149706 ( 0 )      **Instrument:** SV-7      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

LCSD		Sample ID: LCSD-149706			Units: ug/L		Analysis Date: 17-Jan-2020 11:49			
Client ID:		Run ID: SV-7_354580			SeqNo: 5438872		PrepDate: 16-Jan-2020		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.285	0.20	5	0	106	39 - 127	5.272	0.234	20	
2,4-Dimethylphenol	4.2	0.20	5	0	84.0	35 - 120	4.201	0.0176	20	
2,4-Dinitrotoluene	4.635	0.20	5	0	92.7	50 - 122	4.666	0.65	20	
2,6-Dinitrotoluene	4.325	0.20	5	0	86.5	50 - 120	4.42	2.17	20	
2-Chloronaphthalene	4.49	0.20	5	0	89.8	50 - 120	4.636	3.19	20	
2-Methylnaphthalene	4.163	0.10	5	0	83.3	50 - 120	4.065	2.37	20	
4,6-Dinitro-2-methylphenol	5.588	0.20	5	0	112	25 - 121	5.471	2.13	30	
4-Nitrophenol	4.79	1.0	5	0	95.8	30 - 130	4.842	1.07	20	
Acenaphthene	3.978	0.10	5	0	79.6	45 - 120	4.398	10	20	
Acenaphthylene	4.124	0.10	5	0	82.5	47 - 120	4.139	0.368	20	
Anthracene	4.564	0.10	5	0	91.3	45 - 120	4.609	0.985	20	
Benz(a)anthracene	4.402	0.10	5	0	88.0	40 - 120	4.284	2.73	20	
Benzo(a)pyrene	4.633	0.10	5	0	92.7	45 - 120	4.099	12.2	20	
Bis(2-chloroethoxy)methane	4.431	0.20	5	0	88.6	45 - 120	4.466	0.783	20	
Bis(2-ethylhexyl)phthalate	4.39	0.20	5	0	87.8	40 - 139	4.202	4.38	20	
Chrysene	4.79	0.10	5	0	95.8	43 - 120	4.745	0.959	20	
Dibenzofuran	4.261	0.10	5	0	85.2	50 - 120	4.287	0.602	20	
Di-n-butyl phthalate	4.792	0.20	5	0	95.8	45 - 123	4.89	2.02	20	
Fluoranthene	4.792	0.10	5	0	95.8	45 - 125	4.866	1.52	20	
Fluorene	4.279	0.10	5	0	85.6	49 - 120	4.384	2.42	20	
Naphthalene	4.282	0.10	5	0	85.6	45 - 120	4.293	0.27	20	
Nitrobenzene	4.617	0.20	5	0	92.3	44 - 120	4.568	1.08	20	
N-Nitrosodiphenylamine	4.664	0.20	5	0	93.3	40 - 125	4.806	2.99	20	
Pentachlorophenol	4.065	0.20	5	0	81.3	19 - 121	3.923	3.53	20	
Phenanthrene	4.457	0.10	5	0	89.1	45 - 121	4.546	1.98	20	
Phenol	4.533	0.20	5	0	90.7	20 - 124	4.568	0.761	20	
Pyrene	4.594	0.10	5	0	91.9	40 - 130	4.445	3.3	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.334</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.7</i>	<i>34 - 129</i>	<i>4.171</i>	<i>3.83</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.596</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.9</i>	<i>40 - 125</i>	<i>4.609</i>	<i>0.292</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.641</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.8</i>	<i>20 - 120</i>	<i>4.848</i>	<i>4.37</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.77</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.4</i>	<i>40 - 135</i>	<i>4.705</i>	<i>1.38</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>4.549</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.0</i>	<i>41 - 120</i>	<i>4.603</i>	<i>1.19</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.75</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.0</i>	<i>20 - 120</i>	<i>4.591</i>	<i>3.41</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010442-03    HS20010442-04    HS20010442-05    HS20010442-06  
 HS20010442-07    HS20010442-08

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

<b>Batch ID:</b> R354396 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200114</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 14:53</b>				
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435103</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.0</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200214</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 14:04</b>				
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435102</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18	1.0	20	0	90.0	70 - 124			
Benzene	17.81	1.0	20	0	89.1	74 - 120			
Chlorobenzene	18.74	1.0	20	0	93.7	76 - 113			
Ethylbenzene	18.45	1.0	20	0	92.3	77 - 117			
Methylene chloride	18.59	2.0	20	0	93.0	70 - 127			
Toluene	19.5	1.0	20	0	97.5	77 - 118			
Vinyl chloride	17.32	1.0	20	0	86.6	70 - 130			
Xylenes, Total	56.46	1.0	60	0	94.1	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.68</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.44</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>48.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.5</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

Batch ID: R354396 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>	Sample ID: <b>HS20010411-10MS</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 18:13</b>					
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435111</b>		PrepDate:		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.57	1.0	20	0	82.9	70 - 127				
Benzene	18.27	1.0	20	0.9903	86.4	70 - 127				
Chlorobenzene	17.64	1.0	20	0	88.2	70 - 114				
Ethylbenzene	22.73	1.0	20	5.13	88.0	70 - 124				
Methylene chloride	16.6	2.0	20	0	83.0	70 - 128				
Toluene	19.16	1.0	20	0	95.8	70 - 123				
Vinyl chloride	16.93	1.0	20	0	84.6	70 - 130				
Xylenes, Total	344.6	1.0	60	319.7	41.5	70 - 130				SO
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.2</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>52.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.53</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>47.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>82 - 127</i>				

<b>MSD</b>	Sample ID: <b>HS20010411-10MSD</b>	Units: <b>ug/L</b>			Analysis Date: <b>14-Jan-2020 18:37</b>					
Client ID:	Run ID: <b>VOA2_354396</b>	SeqNo: <b>5435112</b>		PrepDate:		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.84	1.0	20	0	84.2	70 - 127	16.57	1.58	20	
Benzene	17.96	1.0	20	0.9903	84.9	70 - 127	18.27	1.68	20	
Chlorobenzene	17.63	1.0	20	0	88.1	70 - 114	17.64	0.0493	20	
Ethylbenzene	22.1	1.0	20	5.13	84.8	70 - 124	22.73	2.82	20	
Methylene chloride	16.24	2.0	20	0	81.2	70 - 128	16.6	2.19	20	
Toluene	18.83	1.0	20	0	94.2	70 - 123	19.16	1.72	20	
Vinyl chloride	16.95	1.0	20	0	84.8	70 - 130	16.93	0.151	20	
Xylenes, Total	335.7	1.0	60	319.7	26.5	70 - 130	344.6	2.64	20	SO
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.9</i>	<i>70 - 126</i>	<i>48.09</i>	<i>1.33</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>52.3</i>	<i>1.73</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.0</i>	<i>77 - 123</i>	<i>48.53</i>	<i>1.09</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.6</i>	<i>82 - 127</i>	<i>47.83</i>	<i>0.978</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010442-02

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

<b>Batch ID:</b> R354484 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200115</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jan-2020 12:01</b>				
Client ID:	Run ID: <b>VOA4_354484</b>	SeqNo: <b>5436851</b>		PrepDate:			DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>59.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>119</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.5</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>54.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>108</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200115</b>	Units: <b>ug/L</b>			Analysis Date: <b>15-Jan-2020 11:12</b>				
Client ID:	Run ID: <b>VOA4_354484</b>	SeqNo: <b>5436850</b>		PrepDate:			DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	22.86	1.0	20	0	114	70 - 124			
Benzene	20.41	1.0	20	0	102	74 - 120			
Chlorobenzene	18.46	1.0	20	0	92.3	76 - 113			
Ethylbenzene	18.52	1.0	20	0	92.6	77 - 117			
Methylene chloride	22.58	2.0	20	0	113	70 - 127			
Toluene	20.42	1.0	20	0	102	77 - 118			
Vinyl chloride	22.6	1.0	20	0	113	70 - 130			
Xylenes, Total	58.07	1.0	60	0	96.8	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>58.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>118</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>54.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>110</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.15</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 120</i>			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QC BATCH REPORT**

**Batch ID:** R354484 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010445-02MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jan-2020 14:30</b>			
Client ID:		Run ID: <b>VOA4_354484</b>			SeqNo: <b>5436856</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	22.11	1.0	20	0	111	70 - 127				
Benzene	19.85	1.0	20	0	99.3	70 - 127				
Chlorobenzene	17.99	1.0	20	0	90.0	70 - 114				
Ethylbenzene	18.05	1.0	20	0	90.3	70 - 124				
Methylene chloride	19.71	2.0	20	0	98.5	70 - 128				
Toluene	20.34	1.0	20	0	102	70 - 123				
Vinyl chloride	24.78	1.0	20	0	124	70 - 130				
Xylenes, Total	59.57	1.0	60	0	99.3	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>61.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>124</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.65</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>57.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>115</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010445-02MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Jan-2020 14:55</b>			
Client ID:		Run ID: <b>VOA4_354484</b>			SeqNo: <b>5436857</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	22.53	1.0	20	0	113	70 - 127	22.11	1.85	20	
Benzene	19.22	1.0	20	0	96.1	70 - 127	19.85	3.24	20	
Chlorobenzene	17.29	1.0	20	0	86.5	70 - 114	17.99	3.97	20	
Ethylbenzene	18.2	1.0	20	0	91.0	70 - 124	18.05	0.8	20	
Methylene chloride	18.57	2.0	20	0	92.9	70 - 128	19.71	5.91	20	
Toluene	19.32	1.0	20	0	96.6	70 - 123	20.34	5.14	20	
Vinyl chloride	23.69	1.0	20	0	118	70 - 130	24.78	4.49	20	
Xylenes, Total	57.75	1.0	60	0	96.3	70 - 130	59.57	3.11	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>60.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>120</i>	<i>70 - 126</i>	<i>61.92</i>	<i>3.12</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.71</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>81 - 113</i>	<i>48.65</i>	<i>0.125</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>55.62</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>111</i>	<i>77 - 123</i>	<i>57.28</i>	<i>2.95</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>82 - 127</i>	<i>50.46</i>	<i>1.46</i>	<i>20</i>	

The following samples were analyzed in this batch: 

HS20010442-03	HS20010442-04	HS20010442-05	HS20010442-06
HS20010442-07	HS20010442-08		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010442

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010442

Date/Time Received: 10-Jan-2020 15:11
Received by: PMG

Checklist completed by: Asad Chaudhry
eSignature
Date: 11-Jan-2020

Reviewed by: Corey Grandits
eSignature
Date: 13-Jan-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [ ] No [checked]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:206384

Temperature(s)/Thermometer(s): 1.7c, 1.1c U/C IR 25
Cooler(s)/Kit(s): 24593, 25283
Date/Time sample(s) sent to storage: 01/10/2020 19:00
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by:

Login Notes: Sx WG-1620 FB 0420200110 time differs: COC= 10:30, Label= 10:00. Logged in per COC.

Client Contacted: Date Contacted: Person Contacted:
Contacted By: Regarding:
Comments:
Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Everett, WA  
+1 425 356 2600

Fort Collins, CO  
+1 970 490 1511

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page 1 of 1

COC ID: 206384

HS20010442

Golder Associates Inc.  
Houston TX-Wood Preserving Works

n, WV

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ALS Project Manager:



Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A
Work Order		Project Number	1620-07-Rev0 SR 92688	B
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	F
Phone	(512) 671-3434	Phone		G
Fax	(512) 671-3446	Fax		H
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		I
				J

8260\_LL\_W (5632528 Volatile Organics Site Specific)  
8260\_LL\_W (5632528 VOC Site Specific + V.C.)  
8270\_LOW\_W (5632532 SemiVolatiles Site specific)  
ICP\_TW (5636002 5652646 Metals - As)

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG 1620 TBO 20200110			Water	1	2		X									
2	WG 1620 TSCG 1719-74 20200110	1-10-2020	14:00	X	1	2		X									
3	WG 1620 TBC 121719-89 20200110	1-10-2020	14:00	X	1	2		X									
4	WG 1620 FDO 1 20200110	1-10-2020	-	X	1	6	X	X	X	X							
5	WG 1620 MW 3573 20200110	1-10-2020	12:45	X	1	6	X	X	X	X							
6	WG 1620 MW 3574 20200110	1-10-2020	11:20	X	1	6	X	X	X	X							
7	WG 1620 FB 04 20200110	1-10-2020	10:30	X	1	6	X	X	X	X							
8	WG 1620 MW 83B 20200110	1-10-2020	10:00	X	1	6	X	X	X	X							
9	WG 1620 MW 83C 20200110	1-10-2020	9:00	X	1	6	X	X	X	X							
10																	

Sampler(s) Please Print & Sign  
*Tim McSpadden T. McSpadden*  
 Requisitioned by: *T. McSpadden* Date: *1/10/20* Time: *15:11*  
 Relinquished by: *T. McSpadden* Date: *1/10/20* Time: *15:11*  
 Shipment Method: \_\_\_\_\_  
 Required Turnaround Time: (Check Box)  STD 10 Wk Days  5 Wk Days  2 Wk Days  24-hour  
 Results Due Date: \_\_\_\_\_  
 Notes: UPRR Houston MWPW  
 Cooler ID: *24593* Cooler Temp.: *1.70*  
*25283* *1.10*  
 QC Package: (Check One Box Below)  
 Level II Std OC  TRRP Check st  
 Level III Std OC/Raw Data  TRRP Level IV  
 Level IV SW614/CLP  
 Other

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

January 23, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010596**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 16 sample(s) on Jan 15, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "Dane J. Wacasey".

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey



Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: aaa			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010596			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149706,149707,149749,R354616,R354637,R354652,R354723,R354799			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			2
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group				LRC Date: aaa			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010596			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149706,149707,149749,R354616,R354637,R354652,R354723,R354799			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: aaa
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS20010596
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 149706,149707,149749,R354616,R354637,R354652,R354723,R354799
ER# <sup>5</sup>	Description	
1	Semivolatile Organics Method SW8270, samples WG-1620-MW40B-20200113, WG-1620-MW20A-20200114, the surrogate recoveries could not be determined due to dilution below the calibration range.  Semivolatile Organics Method SW8270, samples WG-1620-TW41B-20200113, WG-1620-MW12A-20200113 surrogates: 2,4,6-Tribromophenol, 4-Terphenyl-d14 due to a dilution required for sample analysis.	
2	Batch 149706, Semivolatile Organics Method SW8270: LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.  Batch 149707, Semivolatile Organics Method SW8270: LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.                      O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);                      NA = Not Applicable;                      NR = Not Reviewed;                      R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010596

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010596-01	WG-1620-MW40B-20200113	Water		13-Jan-2020 12:10	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-02	WG-1620-MW42B-20200113	Water		13-Jan-2020 13:10	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-03	WG-1620-TW41B-20200113	Water		13-Jan-2020 14:00	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-04	WG-1620-MW05-20200113	Water		13-Jan-2020 14:55	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-05	WG-1620-MW12C-20200113	Water		13-Jan-2020 15:55	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-06	WG-1620-MW12A-20200113	Water		13-Jan-2020 16:45	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-07	WG-1620-MW39B-20200113	Water		13-Jan-2020 17:35	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-08	WG-1620-FB05-20200113	Water		13-Jan-2020 17:50	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-09	WG-1620-MW13-20200114	Water		14-Jan-2020 08:10	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-10	WG-1620-MW09-20200114	Water		14-Jan-2020 09:00	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-11	WG-1620-MW21C-20200114	Water		14-Jan-2020 10:00	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-12	WG-1620-FD02-20200114	Water		14-Jan-2020 10:00	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-13	WG-1620-P11-20200114	Water		14-Jan-2020 10:55	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-14	WG-1620-MW88C-20200114	Water		14-Jan-2020 11:50	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-15	WG-1620-MW20A-20200114	Water		14-Jan-2020 12:55	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010596-16	WG-1620-TB03-20200114	Water		14-Jan-2020 00:00	15-Jan-2020 11:50	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW40B-20200113  
 Collection Date: 13-Jan-2020 12:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 06:51
<b>Benzene</b>	<b>0.011</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 06:51
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 06:51
<b>Ethylbenzene</b>	<b>0.089</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 06:51
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 06:51
<b>Toluene</b>	<b>0.022</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 06:51
<b>Xylenes, Total</b>	<b>0.16</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 06:51
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>120</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 06:51</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>112</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 06:51</i>
<i>Surr: Dibromofluoromethane</i>	<i>111</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 06:51</i>
<i>Surr: Toluene-d8</i>	<i>99.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 06:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW40B-20200113  
 Collection Date: 13-Jan-2020 12:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 19:44
<b>2,4-Dimethylphenol</b>	<b>0.0013</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 19:44
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 19:44
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 19:44
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 19:44
<b>2-Methylnaphthalene</b>	<b>0.34</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 19:44
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 19:44
<b>Acenaphthene</b>	<b>0.39</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
<b>Acenaphthylene</b>	<b>0.0021</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:44
<b>Anthracene</b>	<b>0.040</b>		<b>0.0014</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
<b>Benz(a)anthracene</b>	<b>0.000094</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:44
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 19:44
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 19:44
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 19:44
<b>Chrysene</b>	<b>0.000051</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:44
<b>Dibenzofuran</b>	<b>0.29</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 19:44
<b>Fluoranthene</b>	<b>0.017</b>		<b>0.0010</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
<b>Fluorene</b>	<b>0.30</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
<b>Naphthalene</b>	<b>7.9</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	21-Jan-2020 19:00
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 19:44
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 19:44
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 19:44
<b>Phenanthrene</b>	<b>0.22</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	21-Jan-2020 18:41
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 19:44
<b>Pyrene</b>	<b>0.0037</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 19:44
<i>Surr: 2,4,6-Tribromophenol</i>	<i>72.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:44</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>69.8</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>21-Jan-2020 18:41</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>1000</i>	<i>21-Jan-2020 19:00</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>68.1</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>21-Jan-2020 18:41</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>1000</i>	<i>21-Jan-2020 19:00</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>67.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>96.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>54.9</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>21-Jan-2020 18:41</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>1000</i>	<i>21-Jan-2020 19:00</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>115</i>	<i>J</i>		<i>40-135</i>	<i>%REC</i>	<i>100</i>	<i>21-Jan-2020 18:41</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>88.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 19:44</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>1000</i>	<i>21-Jan-2020 19:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW40B-20200113  
 Collection Date: 13-Jan-2020 12:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	114	J		41-120	%REC	100	21-Jan-2020 18:41
Surr: Nitrobenzene-d5	91.7			41-120	%REC	1	17-Jan-2020 19:44
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	21-Jan-2020 19:00
Surr: Phenol-d6	0	JS		20-120	%REC	1000	21-Jan-2020 19:00
Surr: Phenol-d6	72.7			20-120	%REC	1	17-Jan-2020 19:44
Surr: Phenol-d6	117	J		20-120	%REC	100	21-Jan-2020 18:41
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
Arsenic	0.0523		0.000400	0.00200	mg/L	1	20-Jan-2020 15:19

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW42B-20200113  
 Collection Date: 13-Jan-2020 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	20-Jan-2020 16:46
Benzene	U		0.00020	0.0010	mg/L	1	20-Jan-2020 16:46
Chlorobenzene	U		0.00030	0.0010	mg/L	1	20-Jan-2020 16:46
Ethylbenzene	U		0.00030	0.0010	mg/L	1	20-Jan-2020 16:46
Methylene chloride	U		0.0010	0.0020	mg/L	1	20-Jan-2020 16:46
Toluene	U		0.00020	0.0010	mg/L	1	20-Jan-2020 16:46
Xylenes, Total	U		0.00030	0.0010	mg/L	1	20-Jan-2020 16:46
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 16:46</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 16:46</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 16:46</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 16:46</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW42B-20200113  
 Collection Date: 13-Jan-2020 13:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 20:04
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 20:04
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 20:04
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 20:04
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 20:04
<b>2-Methylnaphthalene</b>	<b>0.000069</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 20:04
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 20:04
<b>Acenaphthene</b>	<b>0.00012</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 20:04
<b>Anthracene</b>	<b>0.000049</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 20:04
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 20:04
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 20:04
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00013</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 20:04
<b>Dibenzofuran</b>	<b>0.000065</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 20:04
<b>Fluoranthene</b>	<b>0.00018</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
<b>Fluorene</b>	<b>0.000067</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
<b>Naphthalene</b>	<b>0.00076</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 20:04
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 20:04
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 20:04
<b>Phenanthrene</b>	<b>0.00012</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 20:04
<b>Pyrene</b>	<b>0.00013</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:04
<i>Surr: 2,4,6-Tribromophenol</i>	67.8			34-129	%REC	1	17-Jan-2020 20:04
<i>Surr: 2-Fluorobiphenyl</i>	65.5			40-125	%REC	1	17-Jan-2020 20:04
<i>Surr: 2-Fluorophenol</i>	60.6			20-120	%REC	1	17-Jan-2020 20:04
<i>Surr: 4-Terphenyl-d14</i>	92.4			40-135	%REC	1	17-Jan-2020 20:04
<i>Surr: Nitrobenzene-d5</i>	63.9			41-120	%REC	1	17-Jan-2020 20:04
<i>Surr: Phenol-d6</i>	65.2			20-120	%REC	1	17-Jan-2020 20:04
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00133</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Jan-2020 16:18

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TW41B-20200113  
 Collection Date: 13-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	20-Jan-2020 14:15
<b>Benzene</b>	<b>0.00039</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	20-Jan-2020 14:15
Chlorobenzene	U		0.00030	0.0010	mg/L	1	20-Jan-2020 14:15
Ethylbenzene	U		0.00030	0.0010	mg/L	1	20-Jan-2020 14:15
Methylene chloride	U		0.0010	0.0020	mg/L	1	20-Jan-2020 14:15
Toluene	U		0.00020	0.0010	mg/L	1	20-Jan-2020 14:15
<b>Xylenes, Total</b>	<b>0.0021</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	20-Jan-2020 14:15
<i>Surr: 1,2-Dichloroethane-d4</i>	96.7			70-126	%REC	1	20-Jan-2020 14:15
<i>Surr: 4-Bromofluorobenzene</i>	97.2			81-113	%REC	1	20-Jan-2020 14:15
<i>Surr: Dibromofluoromethane</i>	99.8			77-123	%REC	1	20-Jan-2020 14:15
<i>Surr: Toluene-d8</i>	99.6			82-127	%REC	1	20-Jan-2020 14:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TW41B-20200113  
 Collection Date: 13-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 20:24
<b>2,4-Dimethylphenol</b>	<b>0.00017</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 20:24
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 20:24
<b>2-Chloronaphthalene</b>	<b>0.00012</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
<b>2-Methylnaphthalene</b>	<b>0.0071</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 20:24
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 20:24
<b>Acenaphthene</b>	<b>0.12</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	21-Jan-2020 19:19
<b>Acenaphthylene</b>	<b>0.0011</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
<b>Anthracene</b>	<b>0.0031</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 20:24
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 20:24
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 20:24
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 20:24
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 20:24
<b>Dibenzofuran</b>	<b>0.055</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	21-Jan-2020 17:42
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 20:24
<b>Fluoranthene</b>	<b>0.0021</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
<b>Fluorene</b>	<b>0.078</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	21-Jan-2020 17:42
<b>Naphthalene</b>	<b>0.058</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	21-Jan-2020 17:42
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 20:24
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 20:24
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 20:24
<b>Phenanthrene</b>	<b>0.0084</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 20:24
<b>Pyrene</b>	<b>0.00082</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:24
<i>Surr: 2,4,6-Tribromophenol</i>	<i>156</i>	<i>S</i>		<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>21-Jan-2020 17:42</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>60.3</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 19:19</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>83.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>68.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>118</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>21-Jan-2020 17:42</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>118</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 19:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>92.0</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>21-Jan-2020 17:42</i>
<i>Surr: 2-Fluorophenol</i>	<i>85.3</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 19:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>66.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:24</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:24</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>155</i>	<i>S</i>		<i>40-135</i>	<i>%REC</i>	<i>10</i>	<i>21-Jan-2020 17:42</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>101</i>			<i>40-135</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 19:19</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TW41B-20200113  
 Collection Date: 13-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	82.3			41-120	%REC	10	21-Jan-2020 17:42
Surr: Nitrobenzene-d5	83.5			41-120	%REC	20	21-Jan-2020 19:19
Surr: Nitrobenzene-d5	81.4			41-120	%REC	1	17-Jan-2020 20:24
Surr: Phenol-d6	71.7			20-120	%REC	1	17-Jan-2020 20:24
Surr: Phenol-d6	83.3			20-120	%REC	10	21-Jan-2020 17:42
Surr: Phenol-d6	91.0			20-120	%REC	20	21-Jan-2020 19:19
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
Arsenic	0.0557		0.000400	0.00200	mg/L	1	20-Jan-2020 16:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW05-20200113  
 Collection Date: 13-Jan-2020 14:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 07:40
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 07:40
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 07:40
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 07:40
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 07:40
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 07:40
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 07:40
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>116</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 07:40</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 07:40</i>
<i>Surr: Dibromofluoromethane</i>	<i>111</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 07:40</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 07:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW05-20200113  
 Collection Date: 13-Jan-2020 14:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 20:43
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 20:43
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 20:43
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 20:43
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 20:43
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 20:43
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 20:43
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 20:43
Acenaphthene	U		0.000027	0.00010	mg/L	1	17-Jan-2020 20:43
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 20:43
Anthracene	U		0.000014	0.00010	mg/L	1	17-Jan-2020 20:43
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 20:43
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 20:43
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 20:43
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 20:43
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 20:43
Dibenzofuran	U		0.000020	0.00010	mg/L	1	17-Jan-2020 20:43
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 20:43
Fluoranthene	U		0.000010	0.00010	mg/L	1	17-Jan-2020 20:43
Fluorene	U		0.000030	0.00010	mg/L	1	17-Jan-2020 20:43
<b>Naphthalene</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 20:43
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 20:43
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 20:43
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 20:43
Phenanthrene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 20:43
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 20:43
Pyrene	U		0.000019	0.00010	mg/L	1	17-Jan-2020 20:43
<i>Surr: 2,4,6-Tribromophenol</i>	<i>77.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:43</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:43</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:43</i>
<i>Surr: Nitrobenzene-d5</i>	<i>72.7</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:43</i>
<i>Surr: Phenol-d6</i>	<i>68.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 20:43</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00146</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jan-2020 16:46</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12C-20200113  
 Collection Date: 13-Jan-2020 15:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 08:04
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 08:04
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 08:04
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 08:04
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 08:04
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 08:04
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 08:04
<i>Surr: 1,2-Dichloroethane-d4</i>		126		70-126	%REC	1	17-Jan-2020 08:04
<i>Surr: 4-Bromofluorobenzene</i>		103		81-113	%REC	1	17-Jan-2020 08:04
<i>Surr: Dibromofluoromethane</i>		110		77-123	%REC	1	17-Jan-2020 08:04
<i>Surr: Toluene-d8</i>		100.0		82-127	%REC	1	17-Jan-2020 08:04

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12C-20200113  
 Collection Date: 13-Jan-2020 15:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 21:03
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 21:03
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 21:03
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 21:03
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 21:03
<b>2-Methylnaphthalene</b>	<b>0.00018</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 21:03
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 21:03
<b>Acenaphthene</b>	<b>0.00015</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 21:03
Anthracene	U		0.000014	0.00010	mg/L	1	17-Jan-2020 21:03
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 21:03
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 21:03
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 21:03
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 21:03
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 21:03
<b>Dibenzofuran</b>	<b>0.00012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 21:03
<b>Fluoranthene</b>	<b>0.000091</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
<b>Fluorene</b>	<b>0.00019</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
<b>Naphthalene</b>	<b>0.0017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 21:03
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 21:03
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 21:03
<b>Phenanthrene</b>	<b>0.00015</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
<b>Phenol</b>	<b>0.00012</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
<b>Pyrene</b>	<b>0.000068</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:03
<i>Surr: 2,4,6-Tribromophenol</i>	75.4			34-129	%REC	1	17-Jan-2020 21:03
<i>Surr: 2-Fluorobiphenyl</i>	62.8			40-125	%REC	1	17-Jan-2020 21:03
<i>Surr: 2-Fluorophenol</i>	59.3			20-120	%REC	1	17-Jan-2020 21:03
<i>Surr: 4-Terphenyl-d14</i>	85.7			40-135	%REC	1	17-Jan-2020 21:03
<i>Surr: Nitrobenzene-d5</i>	61.5			41-120	%REC	1	17-Jan-2020 21:03
<i>Surr: Phenol-d6</i>	61.6			20-120	%REC	1	17-Jan-2020 21:03
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00195</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Jan-2020 16:39

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12A-20200113  
 Collection Date: 13-Jan-2020 16:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 08:29
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 08:29
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 08:29
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 08:29
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 08:29
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 08:29
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 08:29
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>119</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 08:29</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 08:29</i>
<i>Surr: Dibromofluoromethane</i>	<i>113</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 08:29</i>
<i>Surr: Toluene-d8</i>	<i>99.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 08:29</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12A-20200113  
 Collection Date: 13-Jan-2020 16:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 21:23
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 21:23
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 21:23
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 21:23
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 21:23
<b>2-Methylnaphthalene</b>	<b>0.0023</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 21:23
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 21:23
<b>Acenaphthene</b>	<b>0.28</b>		<b>0.0011</b>	<b>0.0040</b>	<b>mg/L</b>	40	21-Jan-2020 19:39
<b>Acenaphthylene</b>	<b>0.0012</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
<b>Anthracene</b>	<b>0.0097</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
<b>Benz(a)anthracene</b>	<b>0.00015</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 21:23
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 21:23
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
<b>Chrysene</b>	<b>0.00015</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
<b>Dibenzofuran</b>	<b>0.23</b>		<b>0.00080</b>	<b>0.0040</b>	<b>mg/L</b>	40	21-Jan-2020 19:39
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 21:23
<b>Fluoranthene</b>	<b>0.0067</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
<b>Fluorene</b>	<b>0.28</b>		<b>0.0012</b>	<b>0.0040</b>	<b>mg/L</b>	40	21-Jan-2020 19:39
<b>Naphthalene</b>	<b>0.00051</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 21:23
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 21:23
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 21:23
<b>Phenanthrene</b>	<b>0.082</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	21-Jan-2020 18:02
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 21:23
<b>Pyrene</b>	<b>0.0031</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:23
<i>Surr: 2,4,6-Tribromophenol</i>	<i>74.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:23</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>135</i>	<i>S</i>		<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 18:02</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>93.5</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>40</i>	<i>21-Jan-2020 19:39</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>122</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 18:02</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>115</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>40</i>	<i>21-Jan-2020 19:39</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>62.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:23</i>
<i>Surr: 2-Fluorophenol</i>	<i>60.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:23</i>
<i>Surr: 2-Fluorophenol</i>	<i>88.7</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 18:02</i>
<i>Surr: 2-Fluorophenol</i>	<i>99.4</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>40</i>	<i>21-Jan-2020 19:39</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>165</i>	<i>S</i>		<i>40-135</i>	<i>%REC</i>	<i>20</i>	<i>21-Jan-2020 18:02</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>106</i>	<i>J</i>		<i>40-135</i>	<i>%REC</i>	<i>40</i>	<i>21-Jan-2020 19:39</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:23</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12A-20200113  
 Collection Date: 13-Jan-2020 16:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	72.0			41-120	%REC	1	17-Jan-2020 21:23
Surr: Nitrobenzene-d5	91.6			41-120	%REC	20	21-Jan-2020 18:02
Surr: Nitrobenzene-d5	89.2	J		41-120	%REC	40	21-Jan-2020 19:39
Surr: Phenol-d6	95.1			20-120	%REC	20	21-Jan-2020 18:02
Surr: Phenol-d6	79.5	J		20-120	%REC	40	21-Jan-2020 19:39
Surr: Phenol-d6	62.0			20-120	%REC	1	17-Jan-2020 21:23
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
Arsenic	0.00134	J	0.000400	0.00200	mg/L	1	20-Jan-2020 16:49

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW39B-20200113  
 Collection Date: 13-Jan-2020 17:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 13:19
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 13:19
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 13:19
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 13:19
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 13:19
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 13:19
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 13:19
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>92.0</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:19</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>95.0</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:19</i>
<i>Surr: Dibromofluoromethane</i>		<i>98.4</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:19</i>
<i>Surr: Toluene-d8</i>		<i>99.2</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:19</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW39B-20200113  
 Collection Date: 13-Jan-2020 17:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: GEY	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	17-Jan-2020 21:43
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	17-Jan-2020 21:43
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	17-Jan-2020 21:43
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	17-Jan-2020 21:43
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	17-Jan-2020 21:43
<b>2-Methylnaphthalene</b>	<b>0.00021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	17-Jan-2020 21:43
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	17-Jan-2020 21:43
<b>Acenaphthene</b>	<b>0.0012</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
Acenaphthylene	U		0.000015	0.00010	mg/L	1	17-Jan-2020 21:43
<b>Anthracene</b>	<b>0.00016</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	17-Jan-2020 21:43
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	17-Jan-2020 21:43
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	17-Jan-2020 21:43
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	17-Jan-2020 21:43
Chrysene	U		0.000021	0.00010	mg/L	1	17-Jan-2020 21:43
<b>Dibenzofuran</b>	<b>0.00019</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	17-Jan-2020 21:43
<b>Fluoranthene</b>	<b>0.00023</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
<b>Fluorene</b>	<b>0.00032</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
<b>Naphthalene</b>	<b>0.0023</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
Nitrobenzene	U		0.000024	0.00020	mg/L	1	17-Jan-2020 21:43
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	17-Jan-2020 21:43
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	17-Jan-2020 21:43
<b>Phenanthrene</b>	<b>0.00019</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
Phenol	U		0.000035	0.00020	mg/L	1	17-Jan-2020 21:43
<b>Pyrene</b>	<b>0.00022</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Jan-2020 21:43
<i>Surr: 2,4,6-Tribromophenol</i>	<i>81.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:43</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>81.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>76.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:43</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:43</i>
<i>Surr: Nitrobenzene-d5</i>	<i>81.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:43</i>
<i>Surr: Phenol-d6</i>	<i>75.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 21:43</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00248</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Jan-2020 16:51

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB05-20200113  
 Collection Date: 13-Jan-2020 17:50

**ANALYTICAL REPORT**

WorkOrder:HS20010596  
 Lab ID:HS20010596-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jan-2020 11:25
Benzene	U		0.00020	0.0010	mg/L	1	18-Jan-2020 11:25
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jan-2020 11:25
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jan-2020 11:25
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jan-2020 11:25
Toluene	U		0.00020	0.0010	mg/L	1	18-Jan-2020 11:25
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jan-2020 11:25
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 11:25</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 11:25</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 11:25</i>
<i>Surr: Toluene-d8</i>	<i>99.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 11:25</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB05-20200113  
 Collection Date: 13-Jan-2020 17:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 13:09
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 13:09
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 13:09
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 13:09
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 13:09
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 13:09
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 13:09
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 13:09
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 13:09
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 13:09
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 13:09
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 13:09
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 13:09
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 13:09
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 13:09
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 13:09
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 13:09
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 13:09
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 13:09
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 13:09
<b>Naphthalene</b>		<b>0.00033</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 13:09
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 13:09
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 13:09
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 13:09
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 13:09
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 13:09
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 13:09
<i>Surr: 2,4,6-Tribromophenol</i>		79.2		34-129	%REC	1	21-Jan-2020 13:09
<i>Surr: 2-Fluorobiphenyl</i>		79.4		40-125	%REC	1	21-Jan-2020 13:09
<i>Surr: 2-Fluorophenol</i>		67.5		20-120	%REC	1	21-Jan-2020 13:09
<i>Surr: 4-Terphenyl-d14</i>		87.4		40-135	%REC	1	21-Jan-2020 13:09
<i>Surr: Nitrobenzene-d5</i>		83.5		41-120	%REC	1	21-Jan-2020 13:09
<i>Surr: Phenol-d6</i>		66.4		20-120	%REC	1	21-Jan-2020 13:09
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	20-Jan-2020 16:53

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW13-20200114  
 Collection Date: 14-Jan-2020 08:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jan-2020 11:49
Benzene	U		0.00020	0.0010	mg/L	1	18-Jan-2020 11:49
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jan-2020 11:49
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jan-2020 11:49
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jan-2020 11:49
Toluene	U		0.00020	0.0010	mg/L	1	18-Jan-2020 11:49
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jan-2020 11:49
<i>Surr: 1,2-Dichloroethane-d4</i>		96.7		70-126	%REC	1	18-Jan-2020 11:49
<i>Surr: 4-Bromofluorobenzene</i>		95.4		81-113	%REC	1	18-Jan-2020 11:49
<i>Surr: Dibromofluoromethane</i>		99.8		77-123	%REC	1	18-Jan-2020 11:49
<i>Surr: Toluene-d8</i>		98.9		82-127	%REC	1	18-Jan-2020 11:49

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW13-20200114  
 Collection Date: 14-Jan-2020 08:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 13:29
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 13:29
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 13:29
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 13:29
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 13:29
<b>2-Methylnaphthalene</b>	<b>0.000054</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 13:29</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 13:29
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 13:29
<b>Acenaphthene</b>	<b>0.00022</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 13:29</b>
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 13:29
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 13:29
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 13:29
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 13:29
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 13:29
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 13:29
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 13:29
<b>Dibenzofuran</b>	<b>0.00016</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 13:29</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 13:29
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 13:29
<b>Fluorene</b>	<b>0.00016</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 13:29</b>
<b>Naphthalene</b>	<b>0.00026</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 13:29</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 13:29
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 13:29
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 13:29
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 13:29
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 13:29
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 13:29
<i>Surr: 2,4,6-Tribromophenol</i>	<i>99.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:29</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>77.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:29</i>
<i>Surr: 2-Fluorophenol</i>	<i>58.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:29</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:29</i>
<i>Surr: Nitrobenzene-d5</i>	<i>86.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:29</i>
<i>Surr: Phenol-d6</i>	<i>52.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:29</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.0642</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jan-2020 16:55</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW09-20200114  
 Collection Date: 14-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	18-Jan-2020 13:02
Benzene	U		0.00020	0.0010	mg/L	1	18-Jan-2020 13:02
Chlorobenzene	U		0.00030	0.0010	mg/L	1	18-Jan-2020 13:02
Ethylbenzene	U		0.00030	0.0010	mg/L	1	18-Jan-2020 13:02
Methylene chloride	U		0.0010	0.0020	mg/L	1	18-Jan-2020 13:02
Toluene	U		0.00020	0.0010	mg/L	1	18-Jan-2020 13:02
Xylenes, Total	U		0.00030	0.0010	mg/L	1	18-Jan-2020 13:02
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 13:02</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 13:02</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 13:02</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>18-Jan-2020 13:02</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW09-20200114  
 Collection Date: 14-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jan-2020 09:25
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	20-Jan-2020 09:25
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jan-2020 09:25
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jan-2020 09:25
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jan-2020 09:25
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	20-Jan-2020 09:25
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jan-2020 09:25
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jan-2020 09:25
Acenaphthene	U		0.000027	0.00010	mg/L	1	20-Jan-2020 09:25
Acenaphthylene	U		0.000015	0.00010	mg/L	1	20-Jan-2020 09:25
Anthracene	U		0.000014	0.00010	mg/L	1	20-Jan-2020 09:25
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jan-2020 09:25
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jan-2020 09:25
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jan-2020 09:25
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jan-2020 09:25
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jan-2020 09:25
Dibenzofuran	U		0.000020	0.00010	mg/L	1	20-Jan-2020 09:25
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jan-2020 09:25
Fluoranthene	U		0.000010	0.00010	mg/L	1	20-Jan-2020 09:25
Fluorene	U		0.000030	0.00010	mg/L	1	20-Jan-2020 09:25
<b>Naphthalene</b>	<b>0.00011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 09:25
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jan-2020 09:25
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jan-2020 09:25
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jan-2020 09:25
Phenanthrene	U		0.000021	0.00010	mg/L	1	20-Jan-2020 09:25
Phenol	U		0.000035	0.00020	mg/L	1	20-Jan-2020 09:25
Pyrene	U		0.000019	0.00010	mg/L	1	20-Jan-2020 09:25
<i>Surr: 2,4,6-Tribromophenol</i>	<i>82.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 09:25</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 09:25</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 09:25</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>98.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 09:25</i>
<i>Surr: Nitrobenzene-d5</i>	<i>76.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 09:25</i>
<i>Surr: Phenol-d6</i>	<i>71.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 09:25</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00430</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Jan-2020 16:58

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW21C-20200114  
 Collection Date: 14-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 02:20
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 02:20
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 02:20
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 02:20
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 02:20
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 02:20
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 02:20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:20</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:20</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:20</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW21C-20200114  
 Collection Date: 14-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 13:48
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 13:48
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 13:48
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 13:48
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 13:48
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 13:48
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 13:48
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 13:48
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 13:48
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 13:48
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 13:48
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 13:48
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 13:48
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 13:48
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00035</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 13:48
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 13:48
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 13:48
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 13:48
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 13:48
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 13:48
<b>Naphthalene</b>	<b>0.00011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 13:48
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 13:48
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 13:48
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 13:48
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 13:48
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 13:48
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 13:48
<i>Surr: 2,4,6-Tribromophenol</i>	<i>92.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:48</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>81.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:48</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:48</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>104</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:48</i>
<i>Surr: Nitrobenzene-d5</i>	<i>87.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:48</i>
<i>Surr: Phenol-d6</i>	<i>64.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 13:48</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00109</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jan-2020 17:00</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD02-20200114  
 Collection Date: 14-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 02:44
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 02:44
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 02:44
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 02:44
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 02:44
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 02:44
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 02:44
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>96.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:44</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:44</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:44</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 02:44</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD02-20200114  
 Collection Date: 14-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:08
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 14:08
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 14:08
<b>2,6-Dinitrotoluene</b>	<b>0.0030</b>		<b>0.000042</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:08
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:08
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:08
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 14:08
<b>Acenaphthene</b>	<b>0.00020</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 14:08
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 14:08
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 14:08
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:08
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 14:08
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000065</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 14:08
<b>Dibenzofuran</b>	<b>0.00017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:08
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 14:08
<b>Fluorene</b>	<b>0.00018</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
<b>Naphthalene</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 14:08
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 14:08
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 14:08
<b>Phenanthrene</b>	<b>0.000061</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 14:08
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 14:08
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:08
<i>Surr: 2,4,6-Tribromophenol</i>	<i>83.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:08</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:08</i>
<i>Surr: 2-Fluorophenol</i>	<i>65.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:08</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>107</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:08</i>
<i>Surr: Nitrobenzene-d5</i>	<i>76.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:08</i>
<i>Surr: Phenol-d6</i>	<i>66.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:08</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00109</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Jan-2020 17:02

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-P11-20200114  
 Collection Date: 14-Jan-2020 10:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:09
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:09
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:09
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:09
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 03:09
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:09
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:09
<i>Surr: 1,2-Dichloroethane-d4</i>		96.2		70-126	%REC	1	17-Jan-2020 03:09
<i>Surr: 4-Bromofluorobenzene</i>		94.7		81-113	%REC	1	17-Jan-2020 03:09
<i>Surr: Dibromofluoromethane</i>		99.1		77-123	%REC	1	17-Jan-2020 03:09
<i>Surr: Toluene-d8</i>		99.9		82-127	%REC	1	17-Jan-2020 03:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-P11-20200114  
 Collection Date: 14-Jan-2020 10:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:27
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 14:27
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 14:27
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 14:27
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:27
<b>2-Methylnaphthalene</b>	<b>0.000076</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:27
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 14:27
<b>Acenaphthene</b>	<b>0.0046</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 14:27
<b>Anthracene</b>	<b>0.000081</b>	<b>J</b>	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 14:27
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:27
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 14:27
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 14:27
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 14:27
<b>Dibenzofuran</b>	<b>0.00016</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:27
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 14:27
<b>Fluorene</b>	<b>0.0013</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
<b>Naphthalene</b>	<b>0.00048</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 14:27
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 14:27
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 14:27
<b>Phenanthrene</b>	<b>0.00035</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:27</b>
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 14:27
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:27
<i>Surr: 2,4,6-Tribromophenol</i>	<i>116</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:27</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:27</i>
<i>Surr: 2-Fluorophenol</i>	<i>58.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:27</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>120</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:27</i>
<i>Surr: Nitrobenzene-d5</i>	<i>82.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:27</i>
<i>Surr: Phenol-d6</i>	<i>58.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:27</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.0360</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>20-Jan-2020 17:19</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88C-20200114  
 Collection Date: 14-Jan-2020 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:33
Benzene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:33
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:33
Ethylbenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:33
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 03:33
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:33
Xylenes, Total	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:33
<i>Surr: 1,2-Dichloroethane-d4</i>	96.9			70-126	%REC	1	17-Jan-2020 03:33
<i>Surr: 4-Bromofluorobenzene</i>	93.0			81-113	%REC	1	17-Jan-2020 03:33
<i>Surr: Dibromofluoromethane</i>	100			77-123	%REC	1	17-Jan-2020 03:33
<i>Surr: Toluene-d8</i>	99.8			82-127	%REC	1	17-Jan-2020 03:33

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88C-20200114  
 Collection Date: 14-Jan-2020 11:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:46
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 14:46
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 14:46
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 14:46
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:46
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:46
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:46
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 14:46
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 14:46
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 14:46
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 14:46
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 14:46
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:46
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 14:46
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000064</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 14:46
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 14:46
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:46
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:46
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 14:46
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 14:46
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:46
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 14:46
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 14:46
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 14:46
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 14:46
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 14:46
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:46
<i>Surr: 2,4,6-Tribromophenol</i>	<i>110</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:46</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>87.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:46</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:46</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>114</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:46</i>
<i>Surr: Nitrobenzene-d5</i>	<i>81.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:46</i>
<i>Surr: Phenol-d6</i>	<i>73.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:46</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.000862</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Jan-2020 17:22

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW20A-20200114  
 Collection Date: 14-Jan-2020 12:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:58
<b>Benzene</b>	<b>0.020</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 03:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	17-Jan-2020 03:58
<b>Ethylbenzene</b>	<b>0.025</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 03:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	17-Jan-2020 03:58
Toluene	U		0.00020	0.0010	mg/L	1	17-Jan-2020 03:58
<b>Xylenes, Total</b>	<b>0.024</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	17-Jan-2020 03:58
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 03:58</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 03:58</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 03:58</i>
<i>Surr: Toluene-d8</i>	<i>98.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>17-Jan-2020 03:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW20A-20200114  
 Collection Date: 14-Jan-2020 12:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	20-Jan-2020 18:24
<b>2,4-Dimethylphenol</b>	<b>0.0018</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	20-Jan-2020 18:24
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	20-Jan-2020 18:24
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	20-Jan-2020 18:24
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	20-Jan-2020 18:24
<b>2-Methylnaphthalene</b>	<b>0.075</b>		<b>0.00048</b>	<b>0.0025</b>	<b>mg/L</b>	25	21-Jan-2020 18:21
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	20-Jan-2020 18:24
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	20-Jan-2020 18:24
<b>Acenaphthene</b>	<b>0.11</b>		<b>0.00068</b>	<b>0.0025</b>	<b>mg/L</b>	25	21-Jan-2020 18:21
<b>Acenaphthylene</b>	<b>0.00092</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:24
<b>Anthracene</b>	<b>0.0075</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:24
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	20-Jan-2020 18:24
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	20-Jan-2020 18:24
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	20-Jan-2020 18:24
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jan-2020 18:24
Chrysene	U		0.000021	0.00010	mg/L	1	20-Jan-2020 18:24
<b>Dibenzofuran</b>	<b>0.079</b>		<b>0.00050</b>	<b>0.0025</b>	<b>mg/L</b>	25	21-Jan-2020 18:21
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jan-2020 18:24
<b>Fluoranthene</b>	<b>0.00067</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:24
<b>Fluorene</b>	<b>0.080</b>		<b>0.00075</b>	<b>0.0025</b>	<b>mg/L</b>	25	21-Jan-2020 18:21
<b>Naphthalene</b>	<b>2.0</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	500	21-Jan-2020 19:58
Nitrobenzene	U		0.000024	0.00020	mg/L	1	20-Jan-2020 18:24
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	20-Jan-2020 18:24
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	20-Jan-2020 18:24
<b>Phenanthrene</b>	<b>0.028</b>		<b>0.00052</b>	<b>0.0025</b>	<b>mg/L</b>	25	21-Jan-2020 18:21
Phenol	U		0.000035	0.00020	mg/L	1	20-Jan-2020 18:24
<b>Pyrene</b>	<b>0.00030</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:24
Surr: 2,4,6-Tribromophenol	116			34-129	%REC	1	20-Jan-2020 18:24
Surr: 2,4,6-Tribromophenol	114			34-129	%REC	25	21-Jan-2020 18:21
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	500	21-Jan-2020 19:58
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	500	21-Jan-2020 19:58
Surr: 2-Fluorobiphenyl	116			40-125	%REC	25	21-Jan-2020 18:21
Surr: 2-Fluorobiphenyl	109			40-125	%REC	1	20-Jan-2020 18:24
Surr: 2-Fluorophenol	110			20-120	%REC	1	20-Jan-2020 18:24
Surr: 2-Fluorophenol	112			20-120	%REC	25	21-Jan-2020 18:21
Surr: 2-Fluorophenol	0	JS		20-120	%REC	500	21-Jan-2020 19:58
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	500	21-Jan-2020 19:58
Surr: 4-Terphenyl-d14	124			40-135	%REC	25	21-Jan-2020 18:21
Surr: 4-Terphenyl-d14	112			40-135	%REC	1	20-Jan-2020 18:24

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW20A-20200114  
 Collection Date: 14-Jan-2020 12:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	113			41-120	%REC	1	20-Jan-2020 18:24
Surr: Nitrobenzene-d5	107			41-120	%REC	25	21-Jan-2020 18:21
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	500	21-Jan-2020 19:58
Surr: Phenol-d6	99.1	J		20-120	%REC	25	21-Jan-2020 18:21
Surr: Phenol-d6	0	JS		20-120	%REC	500	21-Jan-2020 19:58
Surr: Phenol-d6	107			20-120	%REC	1	20-Jan-2020 18:24
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Jan-2020		Analyst: JHD	
Arsenic	0.00808		0.000400	0.00200	mg/L	1	20-Jan-2020 17:24

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-TB03-20200114  
 Collection Date: 14-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010596  
 Lab ID:HS20010596-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>					Analyst: AKP
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	16-Jan-2020 23:53
Benzene	U		0.00020	0.0010	mg/L	1	16-Jan-2020 23:53
Chlorobenzene	U		0.00030	0.0010	mg/L	1	16-Jan-2020 23:53
Ethylbenzene	U		0.00030	0.0010	mg/L	1	16-Jan-2020 23:53
Methylene chloride	U		0.0010	0.0020	mg/L	1	16-Jan-2020 23:53
Toluene	U		0.00020	0.0010	mg/L	1	16-Jan-2020 23:53
Xylenes, Total	U		0.00030	0.0010	mg/L	1	16-Jan-2020 23:53
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>16-Jan-2020 23:53</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>16-Jan-2020 23:53</i>
<i>Surr: Dibromofluoromethane</i>	<i>100.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>16-Jan-2020 23:53</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>16-Jan-2020 23:53</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**Batch ID:** 149706      **Start Date:** 16 Jan 2020 07:00      **End Date:** 16 Jan 2020 11:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010596-01	1	1000 (mL)	1 (mL)	0.001
HS20010596-02	1	1000 (mL)	1 (mL)	0.001
HS20010596-03	1	1000 (mL)	1 (mL)	0.001
HS20010596-04	1	1000 (mL)	1 (mL)	0.001
HS20010596-05	1	1000 (mL)	1 (mL)	0.001
HS20010596-06	1	1000 (mL)	1 (mL)	0.001
HS20010596-07	1	1000 (mL)	1 (mL)	0.001
HS20010596-08	1	1000 (mL)	1 (mL)	0.001
HS20010596-09	1	1000 (mL)	1 (mL)	0.001
HS20010596-10	1	1000 (mL)	1 (mL)	0.001
HS20010596-11	1	1000 (mL)	1 (mL)	0.001
HS20010596-12	1	1000 (mL)	1 (mL)	0.001
HS20010596-13	1	1000 (mL)	1 (mL)	0.001
HS20010596-14	1	1000 (mL)	1 (mL)	0.001

**Batch ID:** 149707      **Start Date:** 16 Jan 2020 11:00      **End Date:** 16 Jan 2020 13:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010596-15	1	1000 (mL)	1 (mL)	0.001

**Batch ID:** 149749      **Start Date:** 17 Jan 2020 08:00      **End Date:** 17 Jan 2020 12:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010596-01		10 (mL)	10 (mL)	1
HS20010596-02		10 (mL)	10 (mL)	1
HS20010596-03		10 (mL)	10 (mL)	1
HS20010596-04		10 (mL)	10 (mL)	1
HS20010596-05		10 (mL)	10 (mL)	1
HS20010596-06		10 (mL)	10 (mL)	1
HS20010596-07		10 (mL)	10 (mL)	1
HS20010596-08		10 (mL)	10 (mL)	1
HS20010596-09		10 (mL)	10 (mL)	1
HS20010596-10		10 (mL)	10 (mL)	1
HS20010596-11		10 (mL)	10 (mL)	1
HS20010596-12		10 (mL)	10 (mL)	1
HS20010596-13		10 (mL)	10 (mL)	1
HS20010596-14		10 (mL)	10 (mL)	1
HS20010596-15		10 (mL)	10 (mL)	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149706 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010596-01	WG-1620-MW40B-20200113	13 Jan 2020 12:10		16 Jan 2020 10:59	21 Jan 2020 19:00	1000
HS20010596-01	WG-1620-MW40B-20200113	13 Jan 2020 12:10		16 Jan 2020 10:59	21 Jan 2020 18:41	100
HS20010596-01	WG-1620-MW40B-20200113	13 Jan 2020 12:10		16 Jan 2020 10:59	17 Jan 2020 19:44	1
HS20010596-02	WG-1620-MW42B-20200113	13 Jan 2020 13:10		16 Jan 2020 10:59	17 Jan 2020 20:04	1
HS20010596-03	WG-1620-TW41B-20200113	13 Jan 2020 14:00		16 Jan 2020 10:59	21 Jan 2020 19:19	20
HS20010596-03	WG-1620-TW41B-20200113	13 Jan 2020 14:00		16 Jan 2020 10:59	21 Jan 2020 17:42	10
HS20010596-03	WG-1620-TW41B-20200113	13 Jan 2020 14:00		16 Jan 2020 10:59	17 Jan 2020 20:24	1
HS20010596-04	WG-1620-MW05-20200113	13 Jan 2020 14:55		16 Jan 2020 10:59	17 Jan 2020 20:43	1
HS20010596-05	WG-1620-MW12C-20200113	13 Jan 2020 15:55		16 Jan 2020 10:59	17 Jan 2020 21:03	1
HS20010596-06	WG-1620-MW12A-20200113	13 Jan 2020 16:45		16 Jan 2020 10:59	21 Jan 2020 19:39	40
HS20010596-06	WG-1620-MW12A-20200113	13 Jan 2020 16:45		16 Jan 2020 10:59	21 Jan 2020 18:02	20
HS20010596-06	WG-1620-MW12A-20200113	13 Jan 2020 16:45		16 Jan 2020 10:59	17 Jan 2020 21:23	1
HS20010596-07	WG-1620-MW39B-20200113	13 Jan 2020 17:35		16 Jan 2020 10:59	17 Jan 2020 21:43	1
HS20010596-08	WG-1620-FB05-20200113	13 Jan 2020 17:50		16 Jan 2020 10:59	21 Jan 2020 13:09	1
HS20010596-09	WG-1620-MW13-20200114	14 Jan 2020 08:10		16 Jan 2020 10:59	21 Jan 2020 13:29	1
HS20010596-10	WG-1620-MW09-20200114	14 Jan 2020 09:00		16 Jan 2020 10:59	20 Jan 2020 09:25	1
HS20010596-11	WG-1620-MW21C-20200114	14 Jan 2020 10:00		16 Jan 2020 10:59	21 Jan 2020 13:48	1
HS20010596-12	WG-1620-FD02-20200114	14 Jan 2020 10:00		16 Jan 2020 10:59	21 Jan 2020 14:08	1
HS20010596-13	WG-1620-P11-20200114	14 Jan 2020 10:55		16 Jan 2020 10:59	21 Jan 2020 14:27	1
HS20010596-14	WG-1620-MW88C-20200114	14 Jan 2020 11:50		16 Jan 2020 10:59	21 Jan 2020 14:46	1
<b>Batch ID: 149707 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010596-15	WG-1620-MW20A-20200114	14 Jan 2020 12:55		16 Jan 2020 11:00	21 Jan 2020 19:58	500
HS20010596-15	WG-1620-MW20A-20200114	14 Jan 2020 12:55		16 Jan 2020 11:00	21 Jan 2020 18:21	25
HS20010596-15	WG-1620-MW20A-20200114	14 Jan 2020 12:55		16 Jan 2020 11:00	20 Jan 2020 18:24	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID:</b> 149749 ( 0 )			<b>Test Name :</b> ICP-MS METALS BY SW6020A		<b>Matrix:</b> Water	
HS20010596-01	WG-1620-MW40B-20200113	13 Jan 2020 12:10		17 Jan 2020 12:00	20 Jan 2020 15:19	1
HS20010596-02	WG-1620-MW42B-20200113	13 Jan 2020 13:10		17 Jan 2020 12:00	20 Jan 2020 16:18	1
HS20010596-03	WG-1620-TW41B-20200113	13 Jan 2020 14:00		17 Jan 2020 12:00	20 Jan 2020 16:20	1
HS20010596-04	WG-1620-MW05-20200113	13 Jan 2020 14:55		17 Jan 2020 12:00	20 Jan 2020 16:46	1
HS20010596-05	WG-1620-MW12C-20200113	13 Jan 2020 15:55		17 Jan 2020 12:00	20 Jan 2020 16:39	1
HS20010596-06	WG-1620-MW12A-20200113	13 Jan 2020 16:45		17 Jan 2020 12:00	20 Jan 2020 16:49	1
HS20010596-07	WG-1620-MW39B-20200113	13 Jan 2020 17:35		17 Jan 2020 12:00	20 Jan 2020 16:51	1
HS20010596-08	WG-1620-FB05-20200113	13 Jan 2020 17:50		17 Jan 2020 12:00	20 Jan 2020 16:53	1
HS20010596-09	WG-1620-MW13-20200114	14 Jan 2020 08:10		17 Jan 2020 12:00	20 Jan 2020 16:55	1
HS20010596-10	WG-1620-MW09-20200114	14 Jan 2020 09:00		17 Jan 2020 12:00	20 Jan 2020 16:58	1
HS20010596-11	WG-1620-MW21C-20200114	14 Jan 2020 10:00		17 Jan 2020 12:00	20 Jan 2020 17:00	1
HS20010596-12	WG-1620-FD02-20200114	14 Jan 2020 10:00		17 Jan 2020 12:00	20 Jan 2020 17:02	1
HS20010596-13	WG-1620-P11-20200114	14 Jan 2020 10:55		17 Jan 2020 12:00	20 Jan 2020 17:19	1
HS20010596-14	WG-1620-MW88C-20200114	14 Jan 2020 11:50		17 Jan 2020 12:00	20 Jan 2020 17:22	1
HS20010596-15	WG-1620-MW20A-20200114	14 Jan 2020 12:55		17 Jan 2020 12:00	20 Jan 2020 17:24	1
<b>Batch ID:</b> R354616 ( 0 )			<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C		<b>Matrix:</b> Water	
HS20010596-11	WG-1620-MW21C-20200114	14 Jan 2020 10:00			17 Jan 2020 02:20	1
HS20010596-12	WG-1620-FD02-20200114	14 Jan 2020 10:00			17 Jan 2020 02:44	1
HS20010596-13	WG-1620-P11-20200114	14 Jan 2020 10:55			17 Jan 2020 03:09	1
HS20010596-14	WG-1620-MW88C-20200114	14 Jan 2020 11:50			17 Jan 2020 03:33	1
HS20010596-15	WG-1620-MW20A-20200114	14 Jan 2020 12:55			17 Jan 2020 03:58	1
HS20010596-16	WG-1620-TB03-20200114	14 Jan 2020 00:00			16 Jan 2020 23:53	1
<b>Batch ID:</b> R354637 ( 0 )			<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C		<b>Matrix:</b> Water	
HS20010596-01	WG-1620-MW40B-20200113	13 Jan 2020 12:10			17 Jan 2020 06:51	1
HS20010596-04	WG-1620-MW05-20200113	13 Jan 2020 14:55			17 Jan 2020 07:40	1
HS20010596-05	WG-1620-MW12C-20200113	13 Jan 2020 15:55			17 Jan 2020 08:04	1
HS20010596-06	WG-1620-MW12A-20200113	13 Jan 2020 16:45			17 Jan 2020 08:29	1
<b>Batch ID:</b> R354652 ( 0 )			<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C		<b>Matrix:</b> Water	
HS20010596-08	WG-1620-FB05-20200113	13 Jan 2020 17:50			18 Jan 2020 11:25	1
HS20010596-09	WG-1620-MW13-20200114	14 Jan 2020 08:10			18 Jan 2020 11:49	1
HS20010596-10	WG-1620-MW09-20200114	14 Jan 2020 09:00			18 Jan 2020 13:02	1
<b>Batch ID:</b> R354723 ( 0 )			<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C		<b>Matrix:</b> Water	
HS20010596-02	WG-1620-MW42B-20200113	13 Jan 2020 13:10			20 Jan 2020 16:46	1
HS20010596-03	WG-1620-TW41B-20200113	13 Jan 2020 14:00			20 Jan 2020 14:15	1
<b>Batch ID:</b> R354799 ( 0 )			<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C		<b>Matrix:</b> Water	
HS20010596-07	WG-1620-MW39B-20200113	13 Jan 2020 17:35			21 Jan 2020 13:19	1

WorkOrder: HS20010596  
InstrumentID: ICPMS05  
Test Code: ICP\_TW  
Test Number: SW6020  
Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20010596  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010596  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS20010596  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149749 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-149749</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Jan-2020 15:14</b>					
Client ID:		Run ID: <b>ICPMS05_354648</b>	SeqNo: <b>5441099</b>	PrepDate: <b>17-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-149749</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Jan-2020 15:16</b>					
Client ID:		Run ID: <b>ICPMS05_354648</b>	SeqNo: <b>5441100</b>	PrepDate: <b>17-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.04712	0.00200	0.05	0	94.2	80 - 120			
<b>MS</b>	Sample ID: <b>HS20010596-01MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Jan-2020 15:23</b>					
Client ID: <b>WG-1620-MW40B-20200113</b>		Run ID: <b>ICPMS05_354648</b>	SeqNo: <b>5441103</b>	PrepDate: <b>17-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.1024	0.00200	0.05	0.05228	100	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20010596-01MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Jan-2020 15:25</b>					
Client ID: <b>WG-1620-MW40B-20200113</b>		Run ID: <b>ICPMS05_354648</b>	SeqNo: <b>5441104</b>	PrepDate: <b>17-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.09984	0.00200	0.05	0.05228	95.1	80 - 120	0.1024	2.56	20
<b>PDS</b>	Sample ID: <b>HS20010596-01PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Jan-2020 15:27</b>					
Client ID: <b>WG-1620-MW40B-20200113</b>		Run ID: <b>ICPMS05_354648</b>	SeqNo: <b>5441105</b>	PrepDate: <b>17-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.1486	0.00200	0.1	0.05228	96.4	75 - 125			
<b>SD</b>	Sample ID: <b>HS20010596-01SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Jan-2020 15:21</b>					
Client ID: <b>WG-1620-MW40B-20200113</b>		Run ID: <b>ICPMS05_354648</b>	SeqNo: <b>5441102</b>	PrepDate: <b>17-Jan-2020</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	0.05319	0.0100					0.05228	1.75	10

The following samples were analyzed in this batch:

HS20010596-01	HS20010596-02	HS20010596-03	HS20010596-04
HS20010596-05	HS20010596-06	HS20010596-07	HS20010596-08
HS20010596-09	HS20010596-10	HS20010596-11	HS20010596-12
HS20010596-13	HS20010596-14	HS20010596-15	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149706 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149706	Units: ug/L			Analysis Date: 17-Jan-2020 11:10					
Client ID:	Run ID: SV-7_354580	SeqNo: 5438870	PrepDate: 16-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.801</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.834</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.7</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.028</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.043</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.556</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.1</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.987</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.7</i>	<i>20 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149706 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149706	Units: ug/L			Analysis Date: 17-Jan-2020 11:29					
Client ID:	Run ID: SV-7_354580	SeqNo: 5438871	PrepDate: 16-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.272	0.20	5	0	105	39 - 127				
2,4-Dimethylphenol	4.201	0.20	5	0	84.0	35 - 120				
2,4-Dinitrotoluene	4.666	0.20	5	0	93.3	50 - 122				
2,6-Dinitrotoluene	4.42	0.20	5	0	88.4	50 - 120				
2-Chloronaphthalene	4.636	0.20	5	0	92.7	50 - 120				
2-Methylnaphthalene	4.065	0.10	5	0	81.3	50 - 120				
4,6-Dinitro-2-methylphenol	5.471	0.20	5	0	109	25 - 121				
4-Nitrophenol	4.842	1.0	5	0	96.8	30 - 130				
Acenaphthene	4.398	0.10	5	0	88.0	45 - 120				
Acenaphthylene	4.139	0.10	5	0	82.8	47 - 120				
Anthracene	4.609	0.10	5	0	92.2	45 - 120				
Benz(a)anthracene	4.284	0.10	5	0	85.7	40 - 120				
Benzo(a)pyrene	4.099	0.10	5	0	82.0	45 - 120				
Bis(2-chloroethoxy)methane	4.466	0.20	5	0	89.3	45 - 120				
Bis(2-ethylhexyl)phthalate	4.202	0.20	5	0	84.0	40 - 139				
Chrysene	4.745	0.10	5	0	94.9	43 - 120				
Dibenzofuran	4.287	0.10	5	0	85.7	50 - 120				
Di-n-butyl phthalate	4.89	0.20	5	0	97.8	45 - 123				
Fluoranthene	4.866	0.10	5	0	97.3	45 - 125				
Fluorene	4.384	0.10	5	0	87.7	49 - 120				
Naphthalene	4.293	0.10	5	0	85.9	45 - 120				
Nitrobenzene	4.568	0.20	5	0	91.4	44 - 120				
N-Nitrosodiphenylamine	4.806	0.20	5	0	96.1	40 - 125				
Pentachlorophenol	3.923	0.20	5	0	78.5	19 - 121				
Phenanthrene	4.546	0.10	5	0	90.9	45 - 121				
Phenol	4.568	0.20	5	0	91.4	20 - 124				
Pyrene	4.445	0.10	5	0	88.9	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.171</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.4</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.609</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.848</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.705</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.1</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.603</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.1</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.591</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149706 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-149706		Units: ug/L		Analysis Date: 17-Jan-2020 11:49				
Client ID:		Run ID: SV-7_354580		SeqNo: 5438872		PrepDate: 16-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.285	0.20	5	0	106	39 - 127	5.272	0.234	20	
2,4-Dimethylphenol	4.2	0.20	5	0	84.0	35 - 120	4.201	0.0176	20	
2,4-Dinitrotoluene	4.635	0.20	5	0	92.7	50 - 122	4.666	0.65	20	
2,6-Dinitrotoluene	4.325	0.20	5	0	86.5	50 - 120	4.42	2.17	20	
2-Chloronaphthalene	4.49	0.20	5	0	89.8	50 - 120	4.636	3.19	20	
2-Methylnaphthalene	4.163	0.10	5	0	83.3	50 - 120	4.065	2.37	20	
4,6-Dinitro-2-methylphenol	5.588	0.20	5	0	112	25 - 121	5.471	2.13	30	
4-Nitrophenol	4.79	1.0	5	0	95.8	30 - 130	4.842	1.07	20	
Acenaphthene	3.978	0.10	5	0	79.6	45 - 120	4.398	10	20	
Acenaphthylene	4.124	0.10	5	0	82.5	47 - 120	4.139	0.368	20	
Anthracene	4.564	0.10	5	0	91.3	45 - 120	4.609	0.985	20	
Benz(a)anthracene	4.402	0.10	5	0	88.0	40 - 120	4.284	2.73	20	
Benzo(a)pyrene	4.633	0.10	5	0	92.7	45 - 120	4.099	12.2	20	
Bis(2-chloroethoxy)methane	4.431	0.20	5	0	88.6	45 - 120	4.466	0.783	20	
Bis(2-ethylhexyl)phthalate	4.39	0.20	5	0	87.8	40 - 139	4.202	4.38	20	
Chrysene	4.79	0.10	5	0	95.8	43 - 120	4.745	0.959	20	
Dibenzofuran	4.261	0.10	5	0	85.2	50 - 120	4.287	0.602	20	
Di-n-butyl phthalate	4.792	0.20	5	0	95.8	45 - 123	4.89	2.02	20	
Fluoranthene	4.792	0.10	5	0	95.8	45 - 125	4.866	1.52	20	
Fluorene	4.279	0.10	5	0	85.6	49 - 120	4.384	2.42	20	
Naphthalene	4.282	0.10	5	0	85.6	45 - 120	4.293	0.27	20	
Nitrobenzene	4.617	0.20	5	0	92.3	44 - 120	4.568	1.08	20	
N-Nitrosodiphenylamine	4.664	0.20	5	0	93.3	40 - 125	4.806	2.99	20	
Pentachlorophenol	4.065	0.20	5	0	81.3	19 - 121	3.923	3.53	20	
Phenanthrene	4.457	0.10	5	0	89.1	45 - 121	4.546	1.98	20	
Phenol	4.533	0.20	5	0	90.7	20 - 124	4.568	0.761	20	
Pyrene	4.594	0.10	5	0	91.9	40 - 130	4.445	3.3	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.334</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.7</i>	<i>34 - 129</i>	<i>4.171</i>	<i>3.83</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.596</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.9</i>	<i>40 - 125</i>	<i>4.609</i>	<i>0.292</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>4.641</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.8</i>	<i>20 - 120</i>	<i>4.848</i>	<i>4.37</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>4.77</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.4</i>	<i>40 - 135</i>	<i>4.705</i>	<i>1.38</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>4.549</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.0</i>	<i>41 - 120</i>	<i>4.603</i>	<i>1.19</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.75</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.0</i>	<i>20 - 120</i>	<i>4.591</i>	<i>3.41</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

<b>Batch ID:</b> 149706 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS20010596-01	HS20010596-02	HS20010596-03	HS20010596-04
HS20010596-05	HS20010596-06	HS20010596-07	HS20010596-08
HS20010596-09	HS20010596-10	HS20010596-11	HS20010596-12
HS20010596-13	HS20010596-14		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149707 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149707	Units: ug/L			Analysis Date: 20-Jan-2020 08:28					
Client ID:	Run ID: SV-7_354654	SeqNo: 5440530		PrepDate: 16-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.747</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.9</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.366</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>107</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.041</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.519</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>110</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.119</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>100</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149707 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149707	Units: ug/L			Analysis Date: 20-Jan-2020 08:47					
Client ID:	Run ID: SV-7_354654	SeqNo: 5440531		PrepDate: 16-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	5.534	0.20	5	0	111	39 - 127				
2,4-Dimethylphenol	4.512	0.20	5	0	90.2	35 - 120				
2,4-Dinitrotoluene	4.88	0.20	5	0	97.6	50 - 122				
2,6-Dinitrotoluene	4.738	0.20	5	0	94.8	50 - 120				
2-Chloronaphthalene	4.954	0.20	5	0	99.1	50 - 120				
2-Methylnaphthalene	4.718	0.10	5	0	94.4	50 - 120				
4,6-Dinitro-2-methylphenol	5.655	0.20	5	0	113	25 - 121				
4-Nitrophenol	5.049	1.0	5	0	101	30 - 130				
Acenaphthene	4.562	0.10	5	0	91.2	45 - 120				
Acenaphthylene	4.507	0.10	5	0	90.1	47 - 120				
Anthracene	4.822	0.10	5	0	96.4	45 - 120				
Benz(a)anthracene	4.637	0.10	5	0	92.7	40 - 120				
Benzo(a)pyrene	5.218	0.10	5	0	104	45 - 120				
Bis(2-chloroethoxy)methane	5.043	0.20	5	0	101	45 - 120				
Bis(2-ethylhexyl)phthalate	4.719	0.20	5	0	94.4	40 - 139				
Chrysene	5.07	0.10	5	0	101	43 - 120				
Dibenzofuran	4.702	0.10	5	0	94.0	50 - 120				
Di-n-butyl phthalate	5.079	0.20	5	0	102	45 - 123				
Fluoranthene	5.109	0.10	5	0	102	45 - 125				
Fluorene	4.754	0.10	5	0	95.1	49 - 120				
Naphthalene	4.92	0.10	5	0	98.4	45 - 120				
Nitrobenzene	5.121	0.20	5	0	102	44 - 120				
N-Nitrosodiphenylamine	4.699	0.20	5	0	94.0	40 - 125				
Pentachlorophenol	4.518	0.20	5	0	90.4	19 - 121				
Phenanthrene	4.712	0.10	5	0	94.2	45 - 121				
Phenol	4.916	0.20	5	0	98.3	20 - 124				
Pyrene	4.925	0.10	5	0	98.5	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.598</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.126</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>5.058</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.147</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>5.089</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>5.146</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

Batch ID: 149707 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-149707		Units: ug/L		Analysis Date: 20-Jan-2020 09:06				
Client ID:		Run ID: SV-7_354654		SeqNo: 5440532		PrepDate: 16-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	5.629	0.20	5	0	113	39 - 127	5.534	1.7	20	
2,4-Dimethylphenol	4.912	0.20	5	0	98.2	35 - 120	4.512	8.47	20	
2,4-Dinitrotoluene	5.416	0.20	5	0	108	50 - 122	4.88	10.4	20	
2,6-Dinitrotoluene	5.437	0.20	5	0	109	50 - 120	4.738	13.7	20	
2-Chloronaphthalene	5.688	0.20	5	0	114	50 - 120	4.954	13.8	20	
2-Methylnaphthalene	5.246	0.10	5	0	105	50 - 120	4.718	10.6	20	
4,6-Dinitro-2-methylphenol	5.603	0.20	5	0	112	25 - 121	5.655	0.93	30	
4-Nitrophenol	5.664	1.0	5	0	113	30 - 130	5.049	11.5	20	
Acenaphthene	4.935	0.10	5	0	98.7	45 - 120	4.562	7.85	20	
Acenaphthylene	5.162	0.10	5	0	103	47 - 120	4.507	13.6	20	
Anthracene	5.621	0.10	5	0	112	45 - 120	4.822	15.3	20	
Benz(a)anthracene	5.332	0.10	5	0	107	40 - 120	4.637	13.9	20	
Benzo(a)pyrene	5.606	0.10	5	0	112	45 - 120	5.218	7.16	20	
Bis(2-chloroethoxy)methane	5.487	0.20	5	0	110	45 - 120	5.043	8.43	20	
Bis(2-ethylhexyl)phthalate	5.4	0.20	5	0	108	40 - 139	4.719	13.5	20	
Chrysene	5.245	0.10	5	0	105	43 - 120	5.07	3.39	20	
Dibenzofuran	5.301	0.10	5	0	106	50 - 120	4.702	12	20	
Di-n-butyl phthalate	5.877	0.20	5	0	118	45 - 123	5.079	14.6	20	
Fluoranthene	5.965	0.10	5	0	119	45 - 125	5.109	15.4	20	
Fluorene	5.457	0.10	5	0	109	49 - 120	4.754	13.8	20	
Naphthalene	5.483	0.10	5	0	110	45 - 120	4.92	10.8	20	
Nitrobenzene	5.637	0.20	5	0	113	44 - 120	5.121	9.6	20	
N-Nitrosodiphenylamine	5.74	0.20	5	0	115	40 - 125	4.699	19.9	20	
Pentachlorophenol	4.914	0.20	5	0	98.3	19 - 121	4.518	8.4	20	
Phenanthrene	5.503	0.10	5	0	110	45 - 121	4.712	15.5	20	
Phenol	5.438	0.20	5	0	109	20 - 124	4.916	10.1	20	
Pyrene	5.793	0.10	5	0	116	40 - 130	4.925	16.2	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.247</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>34 - 129</i>	<i>4.598</i>	<i>13.2</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.627</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>113</i>	<i>40 - 125</i>	<i>5.126</i>	<i>9.31</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>5.548</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>20 - 120</i>	<i>5.058</i>	<i>9.25</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>5.667</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>113</i>	<i>40 - 135</i>	<i>5.147</i>	<i>9.62</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>5.474</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>109</i>	<i>41 - 120</i>	<i>5.089</i>	<i>7.3</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>5.535</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>20 - 120</i>	<i>5.146</i>	<i>7.28</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010596-15

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

<b>Batch ID:</b> R354616 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200116</b>	Units: <b>ug/L</b>			Analysis Date: <b>16-Jan-2020 23:29</b>				
Client ID:	Run ID: <b>VOA2_354616</b>	SeqNo: <b>5439503</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	48.3	1.0	50	0	96.6	70 - 123			
<i>Surr: 4-Bromofluorobenzene</i>	48.06	1.0	50	0	96.1	82 - 115			
<i>Surr: Dibromofluoromethane</i>	50.7	1.0	50	0	101	73 - 126			
<i>Surr: Toluene-d8</i>	50.27	1.0	50	0	101	81 - 120			

<b>LCS</b>	Sample ID: <b>VLCSW-200116</b>	Units: <b>ug/L</b>			Analysis Date: <b>16-Jan-2020 22:40</b>				
Client ID:	Run ID: <b>VOA2_354616</b>	SeqNo: <b>5439502</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	17.44	1.0	20	0	87.2	70 - 124			
Benzene	17.15	1.0	20	0	85.8	74 - 120			
Chlorobenzene	17.99	1.0	20	0	89.9	76 - 113			
Ethylbenzene	17.67	1.0	20	0	88.3	77 - 117			
Methylene chloride	16.86	2.0	20	0	84.3	70 - 127			
Toluene	18.73	1.0	20	0	93.6	77 - 118			
Xylenes, Total	53.91	1.0	60	0	89.8	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	47.73	1.0	50	0	95.5	70 - 130			
<i>Surr: 4-Bromofluorobenzene</i>	48.52	1.0	50	0	97.0	82 - 115			
<i>Surr: Dibromofluoromethane</i>	49.26	1.0	50	0	98.5	73 - 126			
<i>Surr: Toluene-d8</i>	48.43	1.0	50	0	96.9	81 - 120			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354616 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010635-14MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>17-Jan-2020 00:42</b>			
Client ID:		Run ID: <b>VOA2_354616</b>			SeqNo: <b>5439506</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.11	1.0	20	0	85.6	70 - 127				
Benzene	19.13	1.0	20	2.986	80.7	70 - 127				
Chlorobenzene	17.69	1.0	20	0	88.4	70 - 114				
Ethylbenzene	20.6	1.0	20	3.085	87.6	70 - 124				
Methylene chloride	16.17	2.0	20	0	80.8	70 - 128				
Toluene	18.76	1.0	20	0	93.8	70 - 123				
Xylenes, Total	53.49	1.0	60	0	89.2	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>53.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>107</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.18</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.4</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>47.62</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.2</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010635-14MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>17-Jan-2020 01:06</b>			
Client ID:		Run ID: <b>VOA2_354616</b>			SeqNo: <b>5439507</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.92	1.0	20	0	84.6	70 - 127	17.11	1.15	20	
Benzene	18.68	1.0	20	2.986	78.5	70 - 127	19.13	2.35	20	
Chlorobenzene	17.71	1.0	20	0	88.5	70 - 114	17.69	0.0902	20	
Ethylbenzene	20.29	1.0	20	3.085	86.0	70 - 124	20.6	1.51	20	
Methylene chloride	16.17	2.0	20	0	80.9	70 - 128	16.17	0.0223	20	
Toluene	18.59	1.0	20	0	92.9	70 - 123	18.76	0.918	20	
Xylenes, Total	52.95	1.0	60	0	88.2	70 - 130	53.49	1.02	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>70 - 126</i>	<i>48.95</i>	<i>2.31</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>53.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>107</i>	<i>81 - 113</i>	<i>53.28</i>	<i>0.0339</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.53</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.1</i>	<i>77 - 123</i>	<i>48.18</i>	<i>1.36</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>82 - 127</i>	<i>47.62</i>	<i>1.57</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010596-11    HS20010596-12    HS20010596-13    HS20010596-14  
 HS20010596-15    HS20010596-16



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

<b>Batch ID:</b> R354637 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200116</b>	Units: <b>ug/L</b>			Analysis Date: <b>16-Jan-2020 23:37</b>				
Client ID:	Run ID: <b>VOA4_354637</b>	SeqNo: <b>5440160</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	60.85	1.0	50	0	122	70 - 123			
<i>Surr: 4-Bromofluorobenzene</i>	49.27	1.0	50	0	98.5	82 - 115			
<i>Surr: Dibromofluoromethane</i>	55.4	1.0	50	0	111	73 - 126			
<i>Surr: Toluene-d8</i>	49.67	1.0	50	0	99.3	81 - 120			

<b>LCS</b>	Sample ID: <b>VLCSW-200116</b>	Units: <b>ug/L</b>			Analysis Date: <b>16-Jan-2020 22:48</b>				
Client ID:	Run ID: <b>VOA4_354637</b>	SeqNo: <b>5440159</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	23.82	1.0	20	0	119	70 - 124			
Benzene	19.98	1.0	20	0	99.9	74 - 120			
Chlorobenzene	17.86	1.0	20	0	89.3	76 - 113			
Ethylbenzene	18.9	1.0	20	0	94.5	77 - 117			
Methylene chloride	21.28	2.0	20	0	106	70 - 127			
Toluene	19.92	1.0	20	0	99.6	77 - 118			
Xylenes, Total	59.83	1.0	60	0	99.7	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	59.5	1.0	50	0	119	70 - 130			
<i>Surr: 4-Bromofluorobenzene</i>	52.1	1.0	50	0	104	82 - 115			
<i>Surr: Dibromofluoromethane</i>	58.76	1.0	50	0	118	73 - 126			
<i>Surr: Toluene-d8</i>	49.69	1.0	50	0	99.4	81 - 120			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354637 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010637-09MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>17-Jan-2020 01:15</b>			
Client ID:		Run ID: <b>VOA4_354637</b>			SeqNo: <b>5440164</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	20.59	1.0	20	0	103	70 - 127				
Benzene	18.49	1.0	20	0	92.4	70 - 127				
Chlorobenzene	16.6	1.0	20	0	83.0	70 - 114				
Ethylbenzene	17.67	1.0	20	0	88.3	70 - 124				
Methylene chloride	18.28	2.0	20	0	91.4	70 - 128				
Toluene	18.89	1.0	20	0	94.5	70 - 123				
Xylenes, Total	56.99	1.0	60	0	95.0	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>58.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>117</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>54.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>109</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>56.34</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>113</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>48.7</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010637-09MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>17-Jan-2020 01:40</b>			
Client ID:		Run ID: <b>VOA4_354637</b>			SeqNo: <b>5440165</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	20.81	1.0	20	0	104	70 - 127	20.59	1.1	20	
Benzene	17.85	1.0	20	0	89.2	70 - 127	18.49	3.52	20	
Chlorobenzene	16.04	1.0	20	0	80.2	70 - 114	16.6	3.42	20	
Ethylbenzene	16.79	1.0	20	0	83.9	70 - 124	17.67	5.11	20	
Methylene chloride	17.07	2.0	20	0	85.3	70 - 128	18.28	6.86	20	
Toluene	18.48	1.0	20	0	92.4	70 - 123	18.89	2.22	20	
Xylenes, Total	53.81	1.0	60	0	89.7	70 - 130	56.99	5.73	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>60.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>121</i>	<i>70 - 126</i>	<i>58.31</i>	<i>3.71</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 113</i>	<i>54.73</i>	<i>8.93</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>58.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>116</i>	<i>77 - 123</i>	<i>56.34</i>	<i>3.13</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.27</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>48.7</i>	<i>3.17</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010596-01    HS20010596-04    HS20010596-05    HS20010596-06

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354652 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200118</b>			Units: <b>ug/L</b>		Analysis Date: <b>18-Jan-2020 11:00</b>			
Client ID:		Run ID: <b>VOA2_354652</b>			SeqNo: <b>5440487</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.76</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.5</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.12</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200218</b>			Units: <b>ug/L</b>		Analysis Date: <b>18-Jan-2020 10:11</b>			
Client ID:		Run ID: <b>VOA2_354652</b>			SeqNo: <b>5440486</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.58	1.0	20	0	92.9	70 - 124				
Benzene	18.58	1.0	20	0	92.9	74 - 120				
Chlorobenzene	19.59	1.0	20	0	98.0	76 - 113				
Ethylbenzene	19.02	1.0	20	0	95.1	77 - 117				
Methylene chloride	18.17	2.0	20	0	90.9	70 - 127				
Toluene	20.48	1.0	20	0	102	77 - 118				
Xylenes, Total	58.37	1.0	60	0	97.3	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.85</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>48.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354652 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010596-09MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>18-Jan-2020 12:14</b>			
Client ID: <b>WG-1620-MW13-20200114</b>		Run ID: <b>VOA2_354652</b>			SeqNo: <b>5440490</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.12	1.0	20	0	85.6	70 - 127				
Benzene	17.46	1.0	20	0	87.3	70 - 127				
Chlorobenzene	18.21	1.0	20	0	91.0	70 - 114				
Ethylbenzene	18.38	1.0	20	0	91.9	70 - 124				
Methylene chloride	16.38	2.0	20	0	81.9	70 - 128				
Toluene	19.35	1.0	20	0	96.7	70 - 123				
Xylenes, Total	54.94	1.0	60	0	91.6	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.62</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010596-09MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>18-Jan-2020 12:38</b>			
Client ID: <b>WG-1620-MW13-20200114</b>		Run ID: <b>VOA2_354652</b>			SeqNo: <b>5440491</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.21	1.0	20	0	81.0	70 - 127	17.12	5.49	20	
Benzene	16.36	1.0	20	0	81.8	70 - 127	17.46	6.51	20	
Chlorobenzene	17.12	1.0	20	0	85.6	70 - 114	18.21	6.16	20	
Ethylbenzene	16.92	1.0	20	0	84.6	70 - 124	18.38	8.25	20	
Methylene chloride	15.64	2.0	20	0	78.2	70 - 128	16.38	4.6	20	
Toluene	17.97	1.0	20	0	89.8	70 - 123	19.35	7.39	20	
Xylenes, Total	51.5	1.0	60	0	85.8	70 - 130	54.94	6.46	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>70 - 126</i>	<i>49.23</i>	<i>1.71</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.19</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.4</i>	<i>81 - 113</i>	<i>48.62</i>	<i>0.878</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.8</i>	<i>77 - 123</i>	<i>49.52</i>	<i>1.23</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>82 - 127</i>	<i>49.04</i>	<i>0.665</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010596-08      HS20010596-09      HS20010596-10

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

<b>Batch ID:</b> R354723 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200120</b>	Units: <b>ug/L</b>			Analysis Date: <b>20-Jan-2020 11:48</b>				
Client ID:	Run ID: <b>VOA2_354723</b>	SeqNo: <b>5441718</b>		PrepDate:			DF: <b>1</b>		
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.8</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.6</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.2</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200220</b>	Units: <b>ug/L</b>			Analysis Date: <b>20-Jan-2020 10:59</b>				
Client ID:	Run ID: <b>VOA2_354723</b>	SeqNo: <b>5441717</b>		PrepDate:			DF: <b>1</b>		
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	18.96	1.0	20	0	94.8	70 - 124			
Benzene	18.33	1.0	20	0	91.6	74 - 120			
Chlorobenzene	19.14	1.0	20	0	95.7	76 - 113			
Ethylbenzene	19.07	1.0	20	0	95.3	77 - 117			
Methylene chloride	17.72	2.0	20	0	88.6	70 - 127			
Toluene	20.46	1.0	20	0	102	77 - 118			
Xylenes, Total	57.84	1.0	60	0	96.4	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.54</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.6</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354723 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010752-03MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>20-Jan-2020 15:04</b>			
Client ID:		Run ID: <b>VOA2_354723</b>			SeqNo: <b>5441726</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.27	1.0	20	0	81.4	70 - 127				
Benzene	16.74	1.0	20	0	83.7	70 - 127				
Chlorobenzene	17.35	1.0	20	0	86.8	70 - 114				
Ethylbenzene	17.58	1.0	20	0	87.9	70 - 124				
Methylene chloride	18.31	2.0	20	0	91.6	70 - 128				
Toluene	18.62	1.0	20	0	93.1	70 - 123				
Xylenes, Total	52.79	1.0	60	0	88.0	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.4</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.4</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.5</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>48.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.8</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010752-03MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>20-Jan-2020 15:28</b>			
Client ID:		Run ID: <b>VOA2_354723</b>			SeqNo: <b>5441727</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.28	1.0	20	0	81.4	70 - 127	16.27	0.0341	20	
Benzene	16.22	1.0	20	0	81.1	70 - 127	16.74	3.16	20	
Chlorobenzene	16.88	1.0	20	0	84.4	70 - 114	17.35	2.76	20	
Ethylbenzene	17	1.0	20	0	85.0	70 - 124	17.58	3.35	20	
Methylene chloride	17.83	2.0	20	0	89.1	70 - 128	18.31	2.68	20	
Toluene	17.86	1.0	20	0	89.3	70 - 123	18.62	4.19	20	
Xylenes, Total	51.27	1.0	60	0	85.4	70 - 130	52.79	2.92	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.12</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.2</i>	<i>70 - 126</i>	<i>47.72</i>	<i>1.28</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.62</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>81 - 113</i>	<i>48.21</i>	<i>0.842</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.87</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>77 - 123</i>	<i>48.74</i>	<i>0.256</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.42</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>82 - 127</i>	<i>48.92</i>	<i>1.01</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010596-02      HS20010596-03

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354799 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200121</b>			Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 12:55</b>			
Client ID:		Run ID: <b>VOA2_354799</b>			SeqNo: <b>5443190</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.22</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>90.4</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.91</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.8</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200221</b>			Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 12:06</b>			
Client ID:		Run ID: <b>VOA2_354799</b>			SeqNo: <b>5443189</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.66	1.0	20	0	88.3	70 - 124				
Benzene	17.2	1.0	20	0	86.0	74 - 120				
Chlorobenzene	18.44	1.0	20	0	92.2	76 - 113				
Ethylbenzene	17.94	1.0	20	0	89.7	77 - 117				
Methylene chloride	17.49	2.0	20	0	87.4	70 - 127				
Toluene	19.3	1.0	20	0	96.5	77 - 118				
Xylenes, Total	54.82	1.0	60	0	91.4	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.76</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QC BATCH REPORT**

**Batch ID:** R354799 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010767-05MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 17:48</b>			
Client ID:		Run ID: <b>VOA2_354799</b>			SeqNo: <b>5443202</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.36	1.0	20	0	81.8	70 - 127				
Benzene	24.4	1.0	20	6.383	90.1	70 - 127				
Chlorobenzene	17.96	1.0	20	0	89.8	70 - 114				
Ethylbenzene	28.72	1.0	20	9.361	96.8	70 - 124				
Methylene chloride	15.89	2.0	20	0	79.4	70 - 128				
Toluene	20.5	1.0	20	0	102	70 - 123				
Xylenes, Total	57.71	1.0	60	3.181	90.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.1</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.0</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.4</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010767-05MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 18:12</b>			
Client ID:		Run ID: <b>VOA2_354799</b>			SeqNo: <b>5443203</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.51	1.0	20	0	82.6	70 - 127	16.36	0.908	20	
Benzene	24.24	1.0	20	6.383	89.3	70 - 127	24.4	0.678	20	
Chlorobenzene	17.68	1.0	20	0	88.4	70 - 114	17.96	1.54	20	
Ethylbenzene	28.32	1.0	20	9.361	94.8	70 - 124	28.72	1.41	20	
Methylene chloride	15.24	2.0	20	0	76.2	70 - 128	15.89	4.15	20	
Toluene	20.18	1.0	20	0	101	70 - 123	20.5	1.56	20	
Xylenes, Total	56.81	1.0	60	3.181	89.4	70 - 130	57.71	1.57	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.11</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.2</i>	<i>70 - 126</i>	<i>46.04</i>	<i>0.147</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>81 - 113</i>	<i>48.8</i>	<i>0.867</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>77 - 123</i>	<i>47.49</i>	<i>0.191</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>82 - 127</i>	<i>49.2</i>	<i>0.339</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010596-07



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010596

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010596

Date/Time Received: 15-Jan-2020 11:50
Received by: AC

Checklist completed by: Paris Frazier
eSignature
Date: 15-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 21-Jan-2020

Matrices: WATER

Carrier name: Client

- Shipping container/cooler in good condition?
Custody seals intact on shipping container/cooler?
Custody seals intact on sample bottles?
VOA/TX1005/TX1006 Solids in hermetically sealed vials?
Chain of custody present?
Chain of custody signed when relinquished and received?
Samplers name present on COC?
Chain of custody agrees with sample labels?
Samples in proper container/bottle?
Sample containers intact?
Sufficient sample volume for indicated test?
All samples received within holding time?
Container/Temp Blank temperature in compliance?

- Yes/No checkboxes for each item in the list above.

Not Present checkboxes
2 Page(s)
COC IDs:206396/206383

Temperature(s)/Thermometer(s): 1.3C/1.3C ,3.1C/3.1C 1.8C/1.8C. 2.5C/2.5C UC/C IR25
Cooler(s)/Kit(s): 45442/45146/44319/45562
Date/Time sample(s) sent to storage: 01.15.2020 11:50

- Water - VOA vials have zero headspace?
Water - pH acceptable upon receipt?
pH adjusted?
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments: [Empty box]

Corrective Action: [Empty box]



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# Chain of Custody Form

Page 1 of 2

COC ID: 206396

## HS20010596

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TB0-202001</del>			Water	*	*		*									
2	WG-1620-MW40B-20200113	1-13-20	1210	W		6	X		X	X							
3	WG-1620-MW42B-20200113		1310	W		6	X		X	X							
4	WG-1620- <del>MW42B</del> -TW41B-20200113		1400	W		6	X		X	X							
5	WG-1620-MW05-20200113		1455	W		6	X		X	X							
6	WG-1620-MW12C-20200113		1555	W		6	X		X	X							
7	WG-1620-MW12A-20200113		1645	W		6	X		X	X							
8	WG-1620-MW39B-20200113		1735	W		6	X		X	X							
9	WG-1620-FB05-20200113		1750	W		6	X		X	X							
10	WG-1620-MW13-20200114	1-14-20	0810	W		6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>	Shipment Method <b>HAND DELIVERED</b>	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Deys <input type="checkbox"/> 5 Wk Deys <input type="checkbox"/> 2 Wk Deys <input type="checkbox"/> 24-hour	Other <input type="checkbox"/>	Results Due Date:
Relinquished by: <b>John Bz</b>	Date: <b>1-15-20</b>	Time: <b>11:50</b>	Received by: <b>AC</b>	Notes: <b>UPRR Houston MWPW</b>
Relinquished by: <b>John Bz</b>	Date: <b>1-15-2020</b>	Time: <b>11:50</b>	Received by (Laboratory): <b>AC</b>	QC Package: (Check One Box Below) <input type="checkbox"/> Level II Std QC <input type="checkbox"/> Level III Std QC/Raw Data <input type="checkbox"/> Level IV SV/843/CLP <input checked="" type="checkbox"/> TRRP Check st <input type="checkbox"/> TRRP Level IV
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	Other

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

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# Chain of Custody Form

Page 2 of 2

COC ID: 206383

## HS20010596

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:											
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W(5632528 Volatile Organics Site Specific)										
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W(5632528 VOC Site Specific + V.C.)										
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W(5632532 SemiVolatiles Site specific)										
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW(5636002 5652646 Metals - As)										
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E											
	Suite 4004		Stop 0750	F											
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G											
Phone	(512) 671-3434	Phone		H											
Fax	(512) 671-3446	Fax		I											
e-Mail Address	eric.matzner@pbwillc.com	e-Mail Address		J											

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TB01-20200114</del>			<del>Water</del>	<del>*</del>	<del>2</del>		<del>*</del>									
2	WG-1620-MW09-20200114	1-14-20	0900	W		6	X		X	X							
3	WG-1620-MW21C-20200114		1000	W		6	X		X	X							
4	WG-1620-FD02-20200114		1000	W		6	X		X	X							
5	WG-1620-P11-20200114		1055	W		6	X		X	X							
6	WG-1620-MWBBC-20200114		1150	W		6	X		X	X							
7	WG-1620-MW20A-20200114		1255	W		6	X		X	X							
8	WG-1620-TB03-20200114		-	W		2	X										
9																	
10																	

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 hour				Results Due Date:			
Relinquished by: <b>John Bray</b>		Date: 1-15-20	Time: 11:50	Received by:				Notes: UPRR Houston MWPW			
Relinquished by:		Date: 1-15-2020	Time: 11:50	Received by (Laboratory): <b>AC</b>				QC Package: (Check One Box Below)			
Logged by (Laboratory):		Date:	Time:	Checked by (Laboratory):				<input type="checkbox"/> Level II Std OC <input type="checkbox"/> Level III Std QC/Raw Data <input type="checkbox"/> Level IV SW614/CLP <input checked="" type="checkbox"/> TRRP Check st <input type="checkbox"/> TRRP Level IV			

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
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January 29, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010713**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 15 sample(s) on Jan 16, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: DAYNA.FISHER  
Dane J. Wacasey

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey



**Laboratory Review Checklist: Reportable Data**

Laboratory Name: ALS Laboratory Group			LRC Date: 01/29/2020				
Project Name: Houston TX-Wood Preserving Works			Laboratory Job Number: HS20010713				
Reviewer Name: Dane Wacasey			Prep Batch Number(s): 149790, 149807, 149808, R354726, R354799, R355099, R355193				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?		X			2
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

**Laboratory Review Checklist: Supporting Data**

Laboratory Name: ALS Laboratory Group		LRC Date: 01/29/2020					
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS20010713					
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 149790, 149807, 149808, R354726, R354799, R355099, R355193					
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section</b>					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods (NELAC Chap 5 or ISO/IEC 17025 Section 5)</b>					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 01/29/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20010713
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 149790, 149807, 149808, R354726, R354799, R355099, R355193

ER# <sup>5</sup>	Description
1	Semivolatiles by Method SW8270, Sample WG-1620-MW15A-20200114, Surrogate 4-Terphenyl-d14 recovered above upper control limits due to possible matrix effect.  Semivolatiles by Method SW8270, Samples WG-1620-MW46C-20200115, WG-1620-MW25C-20200115, surrogate recoveries could not be determined due to dilution below the calibration range.
2	Batch 149790, Semivolatiles by Method SW8270, Sample WG-1620-MW67B-20200115, MS/MSD RPD recovered above limits for 4,6-Dinitro-2-methylphenol due to possible matrix effect.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
 O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
 NA = Not Applicable;  
 NR = Not Reviewed;  
 R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010713

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010713-01	WG-1620-MW15A-20200114	Water		14-Jan-2020 13:45	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-02	WG-1620-MW15B-20200114	Water		14-Jan-2020 14:30	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-03	WG-1620-MW15C-20200114	Water		14-Jan-2020 15:20	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-04	WG-1620-MW14-20200114	Water		14-Jan-2020 16:15	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-05	WG-1620-MW17-20200114	Water		14-Jan-2020 17:10	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-06	WG-1620-FB06-20200114	Water		14-Jan-2020 17:25	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-07	WG-1620-MW17C-20200115	Water		15-Jan-2020 08:15	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-08	WG-1620-MW67B-20200115	Water		15-Jan-2020 09:30	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-09	WG-1620-MW46C-20200115	Water		15-Jan-2020 10:25	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-10	WG-1620-MW25A-20200115	Water		15-Jan-2020 11:20	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-11	WG-1620-MW44C-20200115	Water		15-Jan-2020 13:15	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-12	WG-1620-MW71B-20200115	Water		15-Jan-2020 14:00	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-13	WG-1620-MW33BR-20200115	Water		15-Jan-2020 15:00	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-14	WG-1620-MW25C-20200115	Water		15-Jan-2020 15:50	16-Jan-2020 12:40	<input type="checkbox"/>
HS20010713-15	WQ-1620-TB04-20200115	Water		15-Jan-2020 00:00	16-Jan-2020 12:40	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15A-20200114  
 Collection Date: 14-Jan-2020 13:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 15:46
<b>Benzene</b>	<b>0.00034</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 15:46
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 15:46
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 15:46
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 15:46
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 15:46
<b>Xylenes, Total</b>	<b>0.00093</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 15:46
<i>Surr: 1,2-Dichloroethane-d4</i>	93.1			70-126	%REC	1	21-Jan-2020 15:46
<i>Surr: 4-Bromofluorobenzene</i>	95.3			81-113	%REC	1	21-Jan-2020 15:46
<i>Surr: Dibromofluoromethane</i>	99.8			77-123	%REC	1	21-Jan-2020 15:46
<i>Surr: Toluene-d8</i>	99.6			82-127	%REC	1	21-Jan-2020 15:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15A-20200114  
 Collection Date: 14-Jan-2020 13:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
<b>1,2-Diphenylhydrazine</b>	<b>0.00036</b>		<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 20:32
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 20:32
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 20:32
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 20:32
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 20:32
<b>2-Methylnaphthalene</b>	<b>0.018</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:19
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 20:32
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 20:32
<b>Acenaphthene</b>	<b>0.14</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:19
<b>Acenaphthylene</b>	<b>0.00066</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:32
<b>Anthracene</b>	<b>0.0048</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:32
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 20:32
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 20:32
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 20:32
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 20:32
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 20:32
<b>Dibenzofuran</b>	<b>0.036</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:19
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 20:32
<b>Fluoranthene</b>	<b>0.0031</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:32
<b>Fluorene</b>	<b>0.065</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:19
<b>Naphthalene</b>	<b>0.00022</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:32
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 20:32
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 20:32
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 20:32
<b>Phenanthrene</b>	<b>0.025</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:19
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 20:32
<b>Pyrene</b>	<b>0.0013</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:32
Surr: 2,4,6-Tribromophenol	113			34-129	%REC	20	22-Jan-2020 13:19
Surr: 2,4,6-Tribromophenol	107			34-129	%REC	1	21-Jan-2020 20:32
Surr: 2-Fluorobiphenyl	72.9			40-125	%REC	1	21-Jan-2020 20:32
Surr: 2-Fluorobiphenyl	93.8			40-125	%REC	20	22-Jan-2020 13:19
Surr: 2-Fluorophenol	70.7	J		20-120	%REC	20	22-Jan-2020 13:19
Surr: 2-Fluorophenol	51.9			20-120	%REC	1	21-Jan-2020 20:32
Surr: 4-Terphenyl-d14	128			40-135	%REC	1	21-Jan-2020 20:32
Surr: 4-Terphenyl-d14	163	S		40-135	%REC	20	22-Jan-2020 13:19
Surr: Nitrobenzene-d5	58.0			41-120	%REC	1	21-Jan-2020 20:32
Surr: Nitrobenzene-d5	69.9	J		41-120	%REC	20	22-Jan-2020 13:19
Surr: Phenol-d6	71.8	J		20-120	%REC	20	22-Jan-2020 13:19
Surr: Phenol-d6	63.1			20-120	%REC	1	21-Jan-2020 20:32

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15A-20200114  
 Collection Date: 14-Jan-2020 13:45

**ANALYTICAL REPORT**

WorkOrder:HS20010713  
 Lab ID:HS20010713-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>	<b>Method:SW6020</b>			Prep:SW3010A / 20-Jan-2020		Analyst: JC	
Arsenic	0.0441		0.000400	0.00200	mg/L	1	22-Jan-2020 16:10

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15B-20200114  
 Collection Date: 14-Jan-2020 14:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:09
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:09
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:09
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:09
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 02:09
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:09
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:09
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:09</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:09</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:09</i>
<i>Surr: Toluene-d8</i>	<i>99.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:09</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15B-20200114  
 Collection Date: 14-Jan-2020 14:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 20:51
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 20:51
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 20:51
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 20:51
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 20:51
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 20:51
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 20:51
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 20:51
<b>Acenaphthene</b>	<b>0.000073</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 20:51
<b>Anthracene</b>	<b>0.00011</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 20:51
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 20:51
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 20:51
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000055</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 20:51
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 20:51
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 20:51
<b>Fluoranthene</b>	<b>0.00011</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
<b>Fluorene</b>	<b>0.000062</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 20:51
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 20:51
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 20:51
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 20:51
<b>Phenanthrene</b>	<b>0.000050</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 20:51
<b>Pyrene</b>	<b>0.000060</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 20:51
<i>Surr: 2,4,6-Tribromophenol</i>	<i>95.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:51</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>84.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:51</i>
<i>Surr: 2-Fluorophenol</i>	<i>52.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:51</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>110</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:51</i>
<i>Surr: Nitrobenzene-d5</i>	<i>58.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:51</i>
<i>Surr: Phenol-d6</i>	<i>66.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:51</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00343</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	22-Jan-2020 16:12

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15C-20200114  
 Collection Date: 14-Jan-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:33
<b>Benzene</b>	<b>0.00059</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 02:33
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:33
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:33
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 02:33
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:33
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:33
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:33</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:33</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:33</i>
<i>Surr: Toluene-d8</i>	<i>99.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 02:33</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15C-20200114  
 Collection Date: 14-Jan-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 21:10
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 21:10
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 21:10
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 21:10
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 21:10
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 21:10
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 21:10
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 21:10
<b>Acenaphthene</b>	<b>0.016</b>		<b>0.00011</b>	<b>0.00040</b>	<b>mg/L</b>	<b>4</b>	<b>22-Jan-2020 13:38</b>
<b>Acenaphthylene</b>	<b>0.00089</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
<b>Anthracene</b>	<b>0.00057</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 21:10
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:10
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 21:10
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 21:10
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 21:10
<b>Dibenzofuran</b>	<b>0.0045</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 21:10
<b>Fluoranthene</b>	<b>0.00066</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
<b>Fluorene</b>	<b>0.00051</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
<b>Naphthalene</b>	<b>0.00024</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 21:10
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 21:10
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 21:10
<b>Phenanthrene</b>	<b>0.000080</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 21:10
<b>Pyrene</b>	<b>0.00038</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 21:10</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>69.8</i>			<i>34-129</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 13:38</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>55.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:10</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>53.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:10</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>64.5</i>			<i>40-125</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 13:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>66.0</i>			<i>20-120</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 13:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>47.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:10</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>90.3</i>			<i>40-135</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 13:38</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>78.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:10</i>
<i>Surr: Nitrobenzene-d5</i>	<i>51.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:10</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.1</i>			<i>41-120</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 13:38</i>
<i>Surr: Phenol-d6</i>	<i>62.6</i>			<i>20-120</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 13:38</i>
<i>Surr: Phenol-d6</i>	<i>47.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:10</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW15C-20200114  
 Collection Date: 14-Jan-2020 15:20

**ANALYTICAL REPORT**

WorkOrder:HS20010713  
 Lab ID:HS20010713-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
Arsenic	0.000773	J	0.000400	0.00200	mg/L	1	21-Jan-2020 15:27

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW14-20200114  
 Collection Date: 14-Jan-2020 16:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:58
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:58
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 02:58
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 02:58
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 02:58
<i>Surr: 1,2-Dichloroethane-d4</i>	93.3			70-126	%REC	1	21-Jan-2020 02:58
<i>Surr: 4-Bromofluorobenzene</i>	94.3			81-113	%REC	1	21-Jan-2020 02:58
<i>Surr: Dibromofluoromethane</i>	98.7			77-123	%REC	1	21-Jan-2020 02:58
<i>Surr: Toluene-d8</i>	99.8			82-127	%REC	1	21-Jan-2020 02:58

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW14-20200114  
 Collection Date: 14-Jan-2020 16:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 21:29
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 21:29
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 21:29
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 21:29
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 21:29
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 21:29
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 21:29
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 21:29
<b>Acenaphthene</b>	<b>0.000091</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 21:29
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 21:29
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 21:29
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 21:29
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:29
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 21:29
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 21:29
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 21:29
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:29
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 21:29
<b>Fluoranthene</b>	<b>0.000064</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 21:29
<b>Fluorene</b>	<b>0.000057</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 21:29
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:29
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 21:29
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 21:29
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 21:29
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 21:29
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 21:29
<b>Pyrene</b>	<b>0.000060</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 21:29
<i>Surr: 2,4,6-Tribromophenol</i>	63.5			34-129	%REC	1	21-Jan-2020 21:29
<i>Surr: 2-Fluorobiphenyl</i>	86.0			40-125	%REC	1	21-Jan-2020 21:29
<i>Surr: 2-Fluorophenol</i>	71.4			20-120	%REC	1	21-Jan-2020 21:29
<i>Surr: 4-Terphenyl-d14</i>	88.9			40-135	%REC	1	21-Jan-2020 21:29
<i>Surr: Nitrobenzene-d5</i>	81.4			41-120	%REC	1	21-Jan-2020 21:29
<i>Surr: Phenol-d6</i>	74.5			20-120	%REC	1	21-Jan-2020 21:29
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00185</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	21-Jan-2020 15:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17-20200114  
 Collection Date: 14-Jan-2020 17:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 08:16
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 08:16
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 08:16
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 08:16
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 08:16
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 08:16
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 08:16
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 08:16</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 08:16</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 08:16</i>
<i>Surr: Toluene-d8</i>	<i>99.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 08:16</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17-20200114  
 Collection Date: 14-Jan-2020 17:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 21:48
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 21:48
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 21:48
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 21:48
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 21:48
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 21:48
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 21:48
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 21:48
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 21:48
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 21:48
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 21:48
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 21:48
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:48
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 21:48
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 21:48
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 21:48
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:48
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 21:48
<b>Fluoranthene</b>	<b>0.00011</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 21:48
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 21:48
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 21:48
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 21:48
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 21:48
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 21:48
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 21:48
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 21:48
<b>Pyrene</b>	<b>0.000066</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 21:48
<i>Surr: 2,4,6-Tribromophenol</i>	67.6			34-129	%REC	1	21-Jan-2020 21:48
<i>Surr: 2-Fluorobiphenyl</i>	96.0			40-125	%REC	1	21-Jan-2020 21:48
<i>Surr: 2-Fluorophenol</i>	79.7			20-120	%REC	1	21-Jan-2020 21:48
<i>Surr: 4-Terphenyl-d14</i>	92.6			40-135	%REC	1	21-Jan-2020 21:48
<i>Surr: Nitrobenzene-d5</i>	83.2			41-120	%REC	1	21-Jan-2020 21:48
<i>Surr: Phenol-d6</i>	79.0			20-120	%REC	1	21-Jan-2020 21:48
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00154</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	21-Jan-2020 15:39

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB06-20200114  
 Collection Date: 14-Jan-2020 17:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	27-Jan-2020 12:30
Benzene	U		0.00020	0.0010	mg/L	1	27-Jan-2020 12:30
Chlorobenzene	U		0.00030	0.0010	mg/L	1	27-Jan-2020 12:30
Ethylbenzene	U		0.00030	0.0010	mg/L	1	27-Jan-2020 12:30
Methylene chloride	U		0.0010	0.0020	mg/L	1	27-Jan-2020 12:30
Toluene	U		0.00020	0.0010	mg/L	1	27-Jan-2020 12:30
Vinyl chloride	U		0.00020	0.0010	mg/L	1	27-Jan-2020 12:30
Xylenes, Total	U		0.00030	0.0010	mg/L	1	27-Jan-2020 12:30
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>106</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 12:30</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>96.0</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 12:30</i>
<i>Surr: Dibromofluoromethane</i>		<i>104</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 12:30</i>
<i>Surr: Toluene-d8</i>		<i>101</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 12:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB06-20200114  
 Collection Date: 14-Jan-2020 17:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 22:07
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 22:07
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 22:07
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 22:07
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 22:07
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 22:07
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 22:07
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 22:07
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 22:07
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 22:07
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 22:07
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 22:07
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 22:07
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 22:07
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 22:07
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 22:07
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 22:07
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 22:07
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 22:07
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 22:07
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 22:07
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 22:07
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 22:07
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 22:07
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 22:07
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 22:07
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 22:07
<i>Surr: 2,4,6-Tribromophenol</i>		64.6		34-129	%REC	1	21-Jan-2020 22:07
<i>Surr: 2-Fluorobiphenyl</i>		97.3		40-125	%REC	1	21-Jan-2020 22:07
<i>Surr: 2-Fluorophenol</i>		74.6		20-120	%REC	1	21-Jan-2020 22:07
<i>Surr: 4-Terphenyl-d14</i>		95.7		40-135	%REC	1	21-Jan-2020 22:07
<i>Surr: Nitrobenzene-d5</i>		85.3		41-120	%REC	1	21-Jan-2020 22:07
<i>Surr: Phenol-d6</i>		77.5		20-120	%REC	1	21-Jan-2020 22:07
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
Arsenic	U		0.000400	0.00200	mg/L	1	21-Jan-2020 15:41

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17C-20200115  
 Collection Date: 15-Jan-2020 08:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 03:22
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 03:22
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 03:22
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 03:22
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 03:22
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 03:22
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 03:22
<i>Surr: 1,2-Dichloroethane-d4</i>		93.4		70-126	%REC	1	21-Jan-2020 03:22
<i>Surr: 4-Bromofluorobenzene</i>		95.0		81-113	%REC	1	21-Jan-2020 03:22
<i>Surr: Dibromofluoromethane</i>		97.3		77-123	%REC	1	21-Jan-2020 03:22
<i>Surr: Toluene-d8</i>		98.9		82-127	%REC	1	21-Jan-2020 03:22

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW17C-20200115  
 Collection Date: 15-Jan-2020 08:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 22:26
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 22:26
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 22:26
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 22:26
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 22:26
<b>2-Methylnaphthalene</b>	<b>0.000073</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 22:26
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 22:26
<b>Acenaphthene</b>	<b>0.00023</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 22:26
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 22:26
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 22:26
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 22:26
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 22:26
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 22:26
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 22:26
<b>Dibenzofuran</b>	<b>0.000065</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 22:26
<b>Fluoranthene</b>	<b>0.00034</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
<b>Fluorene</b>	<b>0.000061</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
<b>Naphthalene</b>	<b>0.00016</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 22:26
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 22:26
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 22:26
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 22:26
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 22:26
<b>Pyrene</b>	<b>0.00021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:26
<i>Surr: 2,4,6-Tribromophenol</i>	<i>68.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:26</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>101</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:26</i>
<i>Surr: 2-Fluorophenol</i>	<i>79.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:26</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>101</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:26</i>
<i>Surr: Nitrobenzene-d5</i>	<i>87.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:26</i>
<i>Surr: Phenol-d6</i>	<i>82.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:26</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00184</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	21-Jan-2020 15:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW67B-20200115  
 Collection Date: 15-Jan-2020 09:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 00:55
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 00:55
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 00:55
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 00:55
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 00:55
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 00:55
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 00:55
<i>Surr: 1,2-Dichloroethane-d4</i>		94.4		70-126	%REC	1	21-Jan-2020 00:55
<i>Surr: 4-Bromofluorobenzene</i>		96.4		81-113	%REC	1	21-Jan-2020 00:55
<i>Surr: Dibromofluoromethane</i>		99.5		77-123	%REC	1	21-Jan-2020 00:55
<i>Surr: Toluene-d8</i>		98.6		82-127	%REC	1	21-Jan-2020 00:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW67B-20200115  
 Collection Date: 15-Jan-2020 09:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:49
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 14:49
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 14:49
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 14:49
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 14:49
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:49
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:49
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 14:49
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 14:49
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 14:49
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 14:49
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	21-Jan-2020 14:49
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:49
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 14:49
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000090</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 14:49</b>
Chrysene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 14:49
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:49
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 14:49
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 14:49
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 14:49
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 14:49
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 14:49
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 14:49
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 14:49
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 14:49
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 14:49
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 14:49
<i>Surr: 2,4,6-Tribromophenol</i>	<i>66.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:49</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>78.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:49</i>
<i>Surr: 2-Fluorophenol</i>	<i>59.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:49</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>100</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:49</i>
<i>Surr: Nitrobenzene-d5</i>	<i>71.9</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:49</i>
<i>Surr: Phenol-d6</i>	<i>70.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 14:49</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.000467</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 15:16</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW46C-20200115  
 Collection Date: 15-Jan-2020 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 07:27
<b>Benzene</b>	<b>0.028</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 07:27
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 07:27
<b>Ethylbenzene</b>	<b>0.051</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 07:27
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 07:27
<b>Toluene</b>	<b>0.0047</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 07:27
<b>Xylenes, Total</b>	<b>0.11</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 07:27
<i>Surr: 1,2-Dichloroethane-d4</i>	92.2			70-126	%REC	1	21-Jan-2020 07:27
<i>Surr: 4-Bromofluorobenzene</i>	98.8			81-113	%REC	1	21-Jan-2020 07:27
<i>Surr: Dibromofluoromethane</i>	98.1			77-123	%REC	1	21-Jan-2020 07:27
<i>Surr: Toluene-d8</i>	98.3			82-127	%REC	1	21-Jan-2020 07:27

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW46C-20200115  
 Collection Date: 15-Jan-2020 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 22:45
<b>2,4-Dimethylphenol</b>	<b>0.00016</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 22:45
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 22:45
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 22:45
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 22:45
<b>2-Methylnaphthalene</b>	<b>0.20</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	22-Jan-2020 15:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 22:45
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 22:45
<b>Acenaphthene</b>	<b>0.16</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:57
<b>Acenaphthylene</b>	<b>0.0016</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:45
<b>Anthracene</b>	<b>0.073</b>		<b>0.00028</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:57
<b>Benz(a)anthracene</b>	<b>0.0071</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:45
<b>Benzo(a)pyrene</b>	<b>0.0021</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:45
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 22:45
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0011</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 22:45
<b>Chrysene</b>	<b>0.0068</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 22:45
<b>Dibenzofuran</b>	<b>0.16</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:57
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 22:45
<b>Fluoranthene</b>	<b>0.084</b>		<b>0.00020</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:57
<b>Fluorene</b>	<b>0.12</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:57
<b>Naphthalene</b>	<b>1.8</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	22-Jan-2020 15:37
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 22:45
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 22:45
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 22:45
<b>Phenanthrene</b>	<b>0.24</b>		<b>0.0042</b>	<b>0.020</b>	<b>mg/L</b>	200	22-Jan-2020 15:37
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 22:45
<b>Pyrene</b>	<b>0.053</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 13:57
<i>Surr: 2,4,6-Tribromophenol</i>	<i>104</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 13:57</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:37</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>68.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:45</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>53.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:45</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>95.2</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 13:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>98.2</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 13:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>58.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:45</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>94.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 22:45</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>122</i>			<i>40-135</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 13:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>0</i>	<i>JS</i>		<i>40-135</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:37</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW46C-20200115  
 Collection Date: 15-Jan-2020 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	87.1			41-120	%REC	20	22-Jan-2020 13:57
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	22-Jan-2020 15:37
Surr: Nitrobenzene-d5	46.7			41-120	%REC	1	21-Jan-2020 22:45
Surr: Phenol-d6	56.3			20-120	%REC	1	21-Jan-2020 22:45
Surr: Phenol-d6	83.5			20-120	%REC	20	22-Jan-2020 13:57
Surr: Phenol-d6	0	JS		20-120	%REC	200	22-Jan-2020 15:37
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
Arsenic	0.00272		0.000400	0.00200	mg/L	1	21-Jan-2020 15:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25A-20200115  
 Collection Date: 15-Jan-2020 11:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:10
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:10
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 16:10
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 16:10
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 16:10
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:10
Vinyl chloride	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:10
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 16:10
<i>Surr: 1,2-Dichloroethane-d4</i>		93.0		70-126	%REC	1	21-Jan-2020 16:10
<i>Surr: 4-Bromofluorobenzene</i>		93.2		81-113	%REC	1	21-Jan-2020 16:10
<i>Surr: Dibromofluoromethane</i>		98.8		77-123	%REC	1	21-Jan-2020 16:10
<i>Surr: Toluene-d8</i>		99.1		82-127	%REC	1	21-Jan-2020 16:10

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25A-20200115  
 Collection Date: 15-Jan-2020 11:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 23:04
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 23:04
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 23:04
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 23:04
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 23:04
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 23:04
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 23:04
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 23:04
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 23:04
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 23:04
<b>Anthracene</b>	<b>0.000089</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
<b>Benz(a)anthracene</b>	<b>0.00012</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
<b>Benzo(a)pyrene</b>	<b>0.000038</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 23:04
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 23:04
<b>Chrysene</b>	<b>0.000097</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 23:04
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 23:04
<b>Fluoranthene</b>	<b>0.00045</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 23:04
<b>Naphthalene</b>	<b>0.00010</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 23:04
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 23:04
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 23:04
<b>Phenanthrene</b>	<b>0.00017</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 23:04
<b>Pyrene</b>	<b>0.00035</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:04
<i>Surr: 2,4,6-Tribromophenol</i>	46.9			34-129	%REC	1	21-Jan-2020 23:04
<i>Surr: 2-Fluorobiphenyl</i>	68.3			40-125	%REC	1	21-Jan-2020 23:04
<i>Surr: 2-Fluorophenol</i>	48.4			20-120	%REC	1	21-Jan-2020 23:04
<i>Surr: 4-Terphenyl-d14</i>	97.4			40-135	%REC	1	21-Jan-2020 23:04
<i>Surr: Nitrobenzene-d5</i>	60.2			41-120	%REC	1	21-Jan-2020 23:04
<i>Surr: Phenol-d6</i>	61.0			20-120	%REC	1	21-Jan-2020 23:04
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00216</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	21-Jan-2020 15:48

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44C-20200115  
 Collection Date: 15-Jan-2020 13:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 07:51
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 07:51
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 07:51
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 07:51
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 07:51
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 07:51
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 07:51
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>89.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 07:51</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 07:51</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 07:51</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 07:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW44C-20200115  
 Collection Date: 15-Jan-2020 13:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	21-Jan-2020 23:23
<b>2,4-Dimethylphenol</b>	<b>0.00011</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	21-Jan-2020 23:23
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	21-Jan-2020 23:23
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	21-Jan-2020 23:23
<b>2-Methylnaphthalene</b>	<b>0.00012</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	21-Jan-2020 23:23
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	21-Jan-2020 23:23
<b>Acenaphthene</b>	<b>0.00045</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Acenaphthylene</b>	<b>0.00065</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Anthracene</b>	<b>0.0086</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Benz(a)anthracene</b>	<b>0.0018</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Benzo(a)pyrene</b>	<b>0.00065</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	21-Jan-2020 23:23
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0012</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Chrysene</b>	<b>0.0022</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Dibenzofuran</b>	<b>0.00034</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Di-n-butyl phthalate</b>	<b>0.0016</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Fluoranthene</b>	<b>0.0095</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Fluorene</b>	<b>0.00067</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<b>Naphthalene</b>	<b>0.000097</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
Nitrobenzene		U	0.000024	0.00020	mg/L	1	21-Jan-2020 23:23
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	21-Jan-2020 23:23
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	21-Jan-2020 23:23
<b>Phenanthrene</b>	<b>0.0077</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
Phenol		U	0.000035	0.00020	mg/L	1	21-Jan-2020 23:23
<b>Pyrene</b>	<b>0.0060</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:23
<i>Surr: 2,4,6-Tribromophenol</i>	<i>77.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:23</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>72.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:23</i>
<i>Surr: 2-Fluorophenol</i>	<i>59.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:23</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>90.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:23</i>
<i>Surr: Nitrobenzene-d5</i>	<i>61.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:23</i>
<i>Surr: Phenol-d6</i>	<i>61.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:23</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00314</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	21-Jan-2020 15:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW71B-20200115  
 Collection Date: 15-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 04:11
<b>Benzene</b>	<b>0.0021</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 04:11
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 04:11
<b>Ethylbenzene</b>	<b>0.0013</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 04:11
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 04:11
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 04:11
<b>Xylenes, Total</b>	<b>0.0030</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 04:11
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>94.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 04:11</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 04:11</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 04:11</i>
<i>Surr: Toluene-d8</i>	<i>98.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 04:11</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW71B-20200115  
 Collection Date: 15-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	21-Jan-2020 23:42
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	21-Jan-2020 23:42
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	21-Jan-2020 23:42
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	21-Jan-2020 23:42
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	21-Jan-2020 23:42
<b>2-Methylnaphthalene</b>	<b>0.0060</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	21-Jan-2020 23:42
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	21-Jan-2020 23:42
<b>Acenaphthene</b>	<b>0.0053</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Acenaphthylene</b>	<b>0.00015</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Anthracene</b>	<b>0.0039</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Benz(a)anthracene</b>	<b>0.0011</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Benzo(a)pyrene</b>	<b>0.00052</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	21-Jan-2020 23:42
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00063</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Chrysene</b>	<b>0.0011</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Dibenzofuran</b>	<b>0.0049</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Di-n-butyl phthalate</b>	<b>0.00098</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Fluoranthene</b>	<b>0.0057</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Fluorene</b>	<b>0.0038</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<b>Naphthalene</b>	<b>0.0064</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
Nitrobenzene	U		0.000024	0.00020	mg/L	1	21-Jan-2020 23:42
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	21-Jan-2020 23:42
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	21-Jan-2020 23:42
<b>Phenanthrene</b>	<b>0.017</b>		<b>0.000084</b>	<b>0.00040</b>	<b>mg/L</b>	4	22-Jan-2020 14:16
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 23:42
<b>Pyrene</b>	<b>0.0042</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 23:42
<i>Surr: 2,4,6-Tribromophenol</i>	<i>88.6</i>			<i>34-129</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 14:16</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>71.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:42</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>81.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:42</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>93.3</i>			<i>40-125</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 14:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.3</i>			<i>20-120</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 14:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>62.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:42</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>93.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:42</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>108</i>			<i>40-135</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 14:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>80.2</i>			<i>41-120</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 14:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>67.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:42</i>
<i>Surr: Phenol-d6</i>	<i>66.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:42</i>
<i>Surr: Phenol-d6</i>	<i>74.8</i>			<i>20-120</i>	<i>%REC</i>	<i>4</i>	<i>22-Jan-2020 14:16</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW71B-20200115  
 Collection Date: 15-Jan-2020 14:00

**ANALYTICAL REPORT**

WorkOrder:HS20010713  
 Lab ID:HS20010713-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
Arsenic	0.00279		0.000400	0.00200	mg/L	1	21-Jan-2020 15:53

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33BR-20200115  
 Collection Date: 15-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:34
<b>Benzene</b>	<b>0.12</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:34
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 16:34
<b>Ethylbenzene</b>	<b>0.043</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:34
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 16:34
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:34
Vinyl chloride	U		0.00020	0.0010	mg/L	1	21-Jan-2020 16:34
<b>Xylenes, Total</b>	<b>0.011</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:34
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>92.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:34</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:34</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:34</i>
<i>Surr: Toluene-d8</i>	<i>99.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:34</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33BR-20200115  
 Collection Date: 15-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	22-Jan-2020 00:01
<b>2,4-Dimethylphenol</b>	<b>0.00018</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	22-Jan-2020 00:01
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	22-Jan-2020 00:01
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	22-Jan-2020 00:01
<b>2-Methylnaphthalene</b>	<b>0.00025</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	22-Jan-2020 00:01
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	22-Jan-2020 00:01
<b>Acenaphthene</b>	<b>0.0051</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Acenaphthylene</b>	<b>0.00015</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Anthracene</b>	<b>0.00046</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Benz(a)anthracene</b>	<b>0.00010</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Benzo(a)pyrene</b>	<b>0.000039</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	22-Jan-2020 00:01
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00067</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Chrysene</b>	<b>0.00011</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Dibenzofuran</b>	<b>0.0039</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Di-n-butyl phthalate</b>	<b>0.00099</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Fluoranthene</b>	<b>0.00082</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Fluorene</b>	<b>0.0018</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Naphthalene</b>	<b>0.0051</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
Nitrobenzene		U	0.000024	0.00020	mg/L	1	22-Jan-2020 00:01
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	22-Jan-2020 00:01
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	22-Jan-2020 00:01
<b>Phenanthrene</b>	<b>0.0013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Phenol</b>	<b>0.00018</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<b>Pyrene</b>	<b>0.00057</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:01
<i>Surr: 2,4,6-Tribromophenol</i>	<i>74.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:01</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>78.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:01</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:01</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>93.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:01</i>
<i>Surr: Nitrobenzene-d5</i>	<i>75.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:01</i>
<i>Surr: Phenol-d6</i>	<i>68.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:01</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.000877</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	21-Jan-2020 15:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25C-20200115  
 Collection Date: 15-Jan-2020 15:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	21-Jan-2020 16:59
<b>Benzene</b>	<b>0.0017</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:59
Chlorobenzene		U	0.00030	0.0010	mg/L	1	21-Jan-2020 16:59
<b>Ethylbenzene</b>	<b>0.033</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:59
Methylene chloride		U	0.0010	0.0020	mg/L	1	21-Jan-2020 16:59
<b>Toluene</b>	<b>0.010</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:59
Vinyl chloride		U	0.00020	0.0010	mg/L	1	21-Jan-2020 16:59
<b>Xylenes, Total</b>	<b>0.22</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 16:59
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:59</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:59</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 16:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25C-20200115  
 Collection Date: 15-Jan-2020 15:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	22-Jan-2020 00:20
<b>2,4-Dimethylphenol</b>	<b>0.0062</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	22-Jan-2020 00:20
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	22-Jan-2020 00:20
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	22-Jan-2020 00:20
<b>2-Methylnaphthalene</b>	<b>0.96</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	22-Jan-2020 15:56
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	22-Jan-2020 00:20
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	22-Jan-2020 00:20
<b>Acenaphthene</b>	<b>0.36</b>		<b>0.0054</b>	<b>0.020</b>	<b>mg/L</b>	200	22-Jan-2020 15:56
<b>Acenaphthylene</b>	<b>0.0025</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<b>Anthracene</b>	<b>0.027</b>		<b>0.00028</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 14:35
<b>Benz(a)anthracene</b>	<b>0.00057</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<b>Benzo(a)pyrene</b>	<b>0.00019</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	22-Jan-2020 00:20
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00066</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<b>Chrysene</b>	<b>0.00056</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<b>Dibenzofuran</b>	<b>0.34</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	22-Jan-2020 15:56
<b>Di-n-butyl phthalate</b>	<b>0.00068</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<b>Fluoranthene</b>	<b>0.0058</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<b>Fluorene</b>	<b>0.15</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 14:35
<b>Naphthalene</b>	<b>7.5</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	22-Jan-2020 16:15
Nitrobenzene	U		0.000024	0.00020	mg/L	1	22-Jan-2020 00:20
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	22-Jan-2020 00:20
<b>Pentachlorophenol</b>	<b>0.17</b>		<b>0.0016</b>	<b>0.0040</b>	<b>mg/L</b>	20	22-Jan-2020 14:35
<b>Phenanthrene</b>	<b>0.18</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	22-Jan-2020 14:35
Phenol	U		0.000035	0.00020	mg/L	1	22-Jan-2020 00:20
<b>Pyrene</b>	<b>0.0049</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	22-Jan-2020 00:20
<i>Surr: 2,4,6-Tribromophenol</i>	<i>56.9</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:20</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>104</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 14:35</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:56</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>22-Jan-2020 16:15</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>22-Jan-2020 16:15</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>84.3</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 14:35</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:56</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>60.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>65.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>98.3</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>22-Jan-2020 14:35</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>22-Jan-2020 15:56</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>22-Jan-2020 16:15</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW25C-20200115  
 Collection Date: 15-Jan-2020 15:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	22-Jan-2020 16:15
Surr: 4-Terphenyl-d14	87.3			40-135	%REC	1	22-Jan-2020 00:20
Surr: 4-Terphenyl-d14	111			40-135	%REC	20	22-Jan-2020 14:35
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	22-Jan-2020 15:56
Surr: Nitrobenzene-d5	99.3			41-120	%REC	20	22-Jan-2020 14:35
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	22-Jan-2020 15:56
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	22-Jan-2020 16:15
Surr: Nitrobenzene-d5	49.3			41-120	%REC	1	22-Jan-2020 00:20
Surr: Phenol-d6	65.7			20-120	%REC	1	22-Jan-2020 00:20
Surr: Phenol-d6	0	JS		20-120	%REC	200	22-Jan-2020 15:56
Surr: Phenol-d6	0	JS		20-120	%REC	2000	22-Jan-2020 16:15
Surr: Phenol-d6	110			20-120	%REC	20	22-Jan-2020 14:35
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 20-Jan-2020		Analyst: JC	
Arsenic	0.00391		0.000400	0.00200	mg/L	1	21-Jan-2020 16:16

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB04-20200115  
 Collection Date: 15-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010713  
 Lab ID:HS20010713-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 00:06
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 00:06
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 00:06
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 00:06
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 00:06
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 00:06
Vinyl chloride	U		0.00020	0.0010	mg/L	1	28-Jan-2020 15:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 00:06
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>92.6</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 00:06</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>100</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:35</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 00:06</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:35</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 00:06</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:35</i>
<i>Surr: Toluene-d8</i>	<i>99.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 00:06</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:35</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

<b>Batch ID:</b> 149790	<b>Start Date:</b> 19 Jan 2020 09:00	<b>End Date:</b> 19 Jan 2020 16:00
<b>Method:</b> SV AQ SEP FUN EXTRACT-LOWLEV - 3510C		<b>Prep Code:</b> 3510_B_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010713-01	1	1000 (mL)	1 (mL)	0.001
HS20010713-02	1	1000 (mL)	1 (mL)	0.001
HS20010713-03	1	1000 (mL)	1 (mL)	0.001
HS20010713-04	1	1000 (mL)	1 (mL)	0.001
HS20010713-05	1	1000 (mL)	1 (mL)	0.001
HS20010713-06	1	1000 (mL)	1 (mL)	0.001
HS20010713-07	1	1000 (mL)	1 (mL)	0.001
HS20010713-08	1	1000 (mL)	1 (mL)	0.001
HS20010713-09	1	1000 (mL)	1 (mL)	0.001
HS20010713-10	1	1000 (mL)	1 (mL)	0.001
HS20010713-11	1	1000 (mL)	1 (mL)	0.001
HS20010713-12	1	1000 (mL)	1 (mL)	0.001
HS20010713-13	1	1000 (mL)	1 (mL)	0.001
HS20010713-14	1	1000 (mL)	1 (mL)	0.001

<b>Batch ID:</b> 149807	<b>Start Date:</b> 20 Jan 2020 10:00	<b>End Date:</b> 20 Jan 2020 14:00
<b>Method:</b> WATER - SW3010A		<b>Prep Code:</b> 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010713-01		10 (mL)	10 (mL)	1
HS20010713-02		10 (mL)	10 (mL)	1

<b>Batch ID:</b> 149808	<b>Start Date:</b> 20 Jan 2020 10:00	<b>End Date:</b> 20 Jan 2020 14:00
<b>Method:</b> WATER - SW3010A		<b>Prep Code:</b> 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010713-03		10 (mL)	10 (mL)	1
HS20010713-04		10 (mL)	10 (mL)	1
HS20010713-05		10 (mL)	10 (mL)	1
HS20010713-06		10 (mL)	10 (mL)	1
HS20010713-07		10 (mL)	10 (mL)	1
HS20010713-08		10 (mL)	10 (mL)	1
HS20010713-09		10 (mL)	10 (mL)	1
HS20010713-10		10 (mL)	10 (mL)	1
HS20010713-11		10 (mL)	10 (mL)	1
HS20010713-12		10 (mL)	10 (mL)	1
HS20010713-13		10 (mL)	10 (mL)	1
HS20010713-14		10 (mL)	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149790 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010713-01	WG-1620-MW15A-20200114	14 Jan 2020 13:45		19 Jan 2020 15:03	22 Jan 2020 13:19	20
HS20010713-01	WG-1620-MW15A-20200114	14 Jan 2020 13:45		19 Jan 2020 15:03	21 Jan 2020 20:32	1
HS20010713-02	WG-1620-MW15B-20200114	14 Jan 2020 14:30		19 Jan 2020 15:03	21 Jan 2020 20:51	1
HS20010713-03	WG-1620-MW15C-20200114	14 Jan 2020 15:20		19 Jan 2020 15:03	22 Jan 2020 13:38	4
HS20010713-03	WG-1620-MW15C-20200114	14 Jan 2020 15:20		19 Jan 2020 15:03	21 Jan 2020 21:10	1
HS20010713-04	WG-1620-MW14-20200114	14 Jan 2020 16:15		19 Jan 2020 15:03	21 Jan 2020 21:29	1
HS20010713-05	WG-1620-MW17-20200114	14 Jan 2020 17:10		19 Jan 2020 15:03	21 Jan 2020 21:48	1
HS20010713-06	WG-1620-FB06-20200114	14 Jan 2020 17:25		19 Jan 2020 15:03	21 Jan 2020 22:07	1
HS20010713-07	WG-1620-MW17C-20200115	15 Jan 2020 08:15		19 Jan 2020 15:03	21 Jan 2020 22:26	1
HS20010713-08	WG-1620-MW67B-20200115	15 Jan 2020 09:30		19 Jan 2020 15:03	21 Jan 2020 14:49	1
HS20010713-09	WG-1620-MW46C-20200115	15 Jan 2020 10:25		19 Jan 2020 15:03	22 Jan 2020 15:37	200
HS20010713-09	WG-1620-MW46C-20200115	15 Jan 2020 10:25		19 Jan 2020 15:03	22 Jan 2020 13:57	20
HS20010713-09	WG-1620-MW46C-20200115	15 Jan 2020 10:25		19 Jan 2020 15:03	21 Jan 2020 22:45	1
HS20010713-10	WG-1620-MW25A-20200115	15 Jan 2020 11:20		19 Jan 2020 15:03	21 Jan 2020 23:04	1
HS20010713-11	WG-1620-MW44C-20200115	15 Jan 2020 13:15		19 Jan 2020 15:03	21 Jan 2020 23:23	1
HS20010713-12	WG-1620-MW71B-20200115	15 Jan 2020 14:00		19 Jan 2020 15:03	22 Jan 2020 14:16	4
HS20010713-12	WG-1620-MW71B-20200115	15 Jan 2020 14:00		19 Jan 2020 15:03	21 Jan 2020 23:42	1
HS20010713-13	WG-1620-MW33BR-20200115	15 Jan 2020 15:00		19 Jan 2020 15:03	22 Jan 2020 00:01	1
HS20010713-14	WG-1620-MW25C-20200115	15 Jan 2020 15:50		19 Jan 2020 15:03	22 Jan 2020 16:15	2000
HS20010713-14	WG-1620-MW25C-20200115	15 Jan 2020 15:50		19 Jan 2020 15:03	22 Jan 2020 15:56	200
HS20010713-14	WG-1620-MW25C-20200115	15 Jan 2020 15:50		19 Jan 2020 15:03	22 Jan 2020 14:35	20
HS20010713-14	WG-1620-MW25C-20200115	15 Jan 2020 15:50		19 Jan 2020 15:03	22 Jan 2020 00:20	1
<b>Batch ID: 149807 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010713-01	WG-1620-MW15A-20200114	14 Jan 2020 13:45		20 Jan 2020 14:00	22 Jan 2020 16:10	1
HS20010713-02	WG-1620-MW15B-20200114	14 Jan 2020 14:30		20 Jan 2020 14:00	22 Jan 2020 16:12	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149808 ( 0 )</b>			<b>Test Name : ICP-MS METALS BY SW6020A</b>		<b>Matrix: Water</b>	
HS20010713-03	WG-1620-MW15C-20200114	14 Jan 2020 15:20		20 Jan 2020 14:00	21 Jan 2020 15:27	1
HS20010713-04	WG-1620-MW14-20200114	14 Jan 2020 16:15		20 Jan 2020 14:00	21 Jan 2020 15:37	1
HS20010713-05	WG-1620-MW17-20200114	14 Jan 2020 17:10		20 Jan 2020 14:00	21 Jan 2020 15:39	1
HS20010713-06	WG-1620-FB06-20200114	14 Jan 2020 17:25		20 Jan 2020 14:00	21 Jan 2020 15:41	1
HS20010713-07	WG-1620-MW17C-20200115	15 Jan 2020 08:15		20 Jan 2020 14:00	21 Jan 2020 15:44	1
HS20010713-08	WG-1620-MW67B-20200115	15 Jan 2020 09:30		20 Jan 2020 14:00	21 Jan 2020 15:16	1
HS20010713-09	WG-1620-MW46C-20200115	15 Jan 2020 10:25		20 Jan 2020 14:00	21 Jan 2020 15:46	1
HS20010713-10	WG-1620-MW25A-20200115	15 Jan 2020 11:20		20 Jan 2020 14:00	21 Jan 2020 15:48	1
HS20010713-11	WG-1620-MW44C-20200115	15 Jan 2020 13:15		20 Jan 2020 14:00	21 Jan 2020 15:50	1
HS20010713-12	WG-1620-MW71B-20200115	15 Jan 2020 14:00		20 Jan 2020 14:00	21 Jan 2020 15:53	1
HS20010713-13	WG-1620-MW33BR- 20200115	15 Jan 2020 15:00		20 Jan 2020 14:00	21 Jan 2020 15:55	1
HS20010713-14	WG-1620-MW25C-20200115	15 Jan 2020 15:50		20 Jan 2020 14:00	21 Jan 2020 16:16	1
<b>Batch ID: R354726 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS20010713-02	WG-1620-MW15B-20200114	14 Jan 2020 14:30			21 Jan 2020 02:09	1
HS20010713-03	WG-1620-MW15C-20200114	14 Jan 2020 15:20			21 Jan 2020 02:33	1
HS20010713-04	WG-1620-MW14-20200114	14 Jan 2020 16:15			21 Jan 2020 02:58	1
HS20010713-05	WG-1620-MW17-20200114	14 Jan 2020 17:10			21 Jan 2020 08:16	1
HS20010713-07	WG-1620-MW17C-20200115	15 Jan 2020 08:15			21 Jan 2020 03:22	1
HS20010713-08	WG-1620-MW67B-20200115	15 Jan 2020 09:30			21 Jan 2020 00:55	1
HS20010713-09	WG-1620-MW46C-20200115	15 Jan 2020 10:25			21 Jan 2020 07:27	1
HS20010713-11	WG-1620-MW44C-20200115	15 Jan 2020 13:15			21 Jan 2020 07:51	1
HS20010713-12	WG-1620-MW71B-20200115	15 Jan 2020 14:00			21 Jan 2020 04:11	1
HS20010713-15	WQ-1620-TB04-20200115	15 Jan 2020 00:00			21 Jan 2020 00:06	1
<b>Batch ID: R354799 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS20010713-01	WG-1620-MW15A-20200114	14 Jan 2020 13:45			21 Jan 2020 15:46	1
HS20010713-10	WG-1620-MW25A-20200115	15 Jan 2020 11:20			21 Jan 2020 16:10	1
HS20010713-13	WG-1620-MW33BR- 20200115	15 Jan 2020 15:00			21 Jan 2020 16:34	1
HS20010713-14	WG-1620-MW25C-20200115	15 Jan 2020 15:50			21 Jan 2020 16:59	1
<b>Batch ID: R355099 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS20010713-06	WG-1620-FB06-20200114	14 Jan 2020 17:25			27 Jan 2020 12:30	1
<b>Batch ID: R355193 ( 0 )</b>			<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>		<b>Matrix: Water</b>	
HS20010713-15	WQ-1620-TB04-20200115	15 Jan 2020 00:00			28 Jan 2020 15:35	1

WorkOrder: HS20010713  
 InstrumentID: ICPMS04  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000456	0.000400	0.00200

WorkOrder: HS20010713  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20010713  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010713  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS20010713  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00045	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

<b>Batch ID:</b> 149807 ( 0 )	<b>Instrument:</b> ICPMS04	<b>Method:</b> ICP-MS METALS BY SW6020A
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<b>MBLK</b>	Sample ID: <b>MBLK-149807</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 14:20</b>							
Client ID:	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5443749</b>	PrepDate: <b>20-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	RPD Qual

Arsenic U 0.00200

<b>LCS</b>	Sample ID: <b>LCS-149807</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 14:22</b>							
Client ID:	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5443750</b>	PrepDate: <b>20-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	RPD Qual

Arsenic 0.0482 0.00200 0.05 0 96.4 80 - 120

<b>MS</b>	Sample ID: <b>HS20010683-01MS</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 14:31</b>							
Client ID:	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5443754</b>	PrepDate: <b>20-Jan-2020</b> DF: <b>10</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	RPD Qual

Arsenic 0.05208 0.0200 0.05 0.00263 98.9 80 - 120

<b>MSD</b>	Sample ID: <b>HS20010683-01MSD</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 16:28</b>							
Client ID:	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444073</b>	PrepDate: <b>20-Jan-2020</b> DF: <b>10</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	RPD Qual

Arsenic 0.05405 0.0200 0.05 0 108 80 - 120 0.05208 3.72 20

<b>PDS</b>	Sample ID: <b>HS20010683-01PDS</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 14:36</b>							
Client ID:	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5443756</b>	PrepDate: <b>20-Jan-2020</b> DF: <b>10</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	RPD Qual

Arsenic 1.031 0.0200 1 0 103 75 - 125

<b>SD</b>	Sample ID: <b>HS20010683-01SD</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 14:29</b>							
Client ID:	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5443753</b>	PrepDate: <b>20-Jan-2020</b> DF: <b>50</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	RPD Qual

Arsenic U 0.100 0.00263 0 10

The following samples were analyzed in this batch: HS20010713-01 HS20010713-02

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

**Batch ID:** 149808 ( 0)      **Instrument:** ICPMS05      **Method:** ICP-MS METALS BY SW6020A

**MBLK**      Sample ID: **MBLK-149808**      Units: **mg/L**      Analysis Date: **21-Jan-2020 15:12**  
 Client ID:      Run ID: **ICPMS05\_354727**      SeqNo: **5442639**      PrepDate: **20-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      U      0.00200

**LCS**      Sample ID: **LCS-149808**      Units: **mg/L**      Analysis Date: **21-Jan-2020 15:14**  
 Client ID:      Run ID: **ICPMS05\_354727**      SeqNo: **5442640**      PrepDate: **20-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.04553      0.00200      0.05      0      91.1      80 - 120

**MS**      Sample ID: **HS20010713-08MS**      Units: **mg/L**      Analysis Date: **21-Jan-2020 15:20**  
 Client ID: **WG-1620-MW67B-20200115**      Run ID: **ICPMS05\_354727**      SeqNo: **5442643**      PrepDate: **20-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.04636      0.00200      0.05      0.000467      91.8      80 - 120

**MSD**      Sample ID: **HS20010713-08MSD**      Units: **mg/L**      Analysis Date: **21-Jan-2020 15:23**  
 Client ID: **WG-1620-MW67B-20200115**      Run ID: **ICPMS05\_354727**      SeqNo: **5442644**      PrepDate: **20-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.04707      0.00200      0.05      0.000467      93.2      80 - 120      0.04636      1.51      20

**PDS**      Sample ID: **HS20010713-08PDS**      Units: **mg/L**      Analysis Date: **21-Jan-2020 15:25**  
 Client ID: **WG-1620-MW67B-20200115**      Run ID: **ICPMS05\_354727**      SeqNo: **5442645**      PrepDate: **20-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.09734      0.00200      0.1      0.000467      96.9      75 - 125

**SD**      Sample ID: **HS20010713-08SD**      Units: **mg/L**      Analysis Date: **21-Jan-2020 15:18**  
 Client ID: **WG-1620-MW67B-20200115**      Run ID: **ICPMS05\_354727**      SeqNo: **5442642**      PrepDate: **20-Jan-2020**      DF: **5**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %D      %D Limit Qual

Arsenic      U      0.0100      0.000467      0      10

**The following samples were analyzed in this batch:**

HS20010713-03	HS20010713-04	HS20010713-05	HS20010713-06
HS20010713-07	HS20010713-08	HS20010713-09	HS20010713-10
HS20010713-11	HS20010713-12	HS20010713-13	HS20010713-14



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

Batch ID: 149790 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149790	Units: ug/L			Analysis Date: 21-Jan-2020 14:11					
Client ID:	Run ID: SV-6_354783	SeqNo: 5442894	PrepDate: 19-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	3.802	0.20	5	0	76.0	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.693	0.20	5	0	93.9	40 - 125				
<i>Surr: 2-Fluorophenol</i>	4.333	0.20	5	0	86.7	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.586	0.20	5	0	91.7	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	3.975	0.20	5	0	79.5	41 - 120				
<i>Surr: Phenol-d6</i>	4.137	0.20	5	0	82.7	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

Batch ID: 149790 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149790	Units: ug/L			Analysis Date: 21-Jan-2020 14:30					
Client ID:	Run ID: SV-6_354783	SeqNo: 5442895		PrepDate: 19-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.591	0.20	5	0	91.8	39 - 127				
2,4-Dimethylphenol	4.25	0.20	5	0	85.0	35 - 120				
2,4-Dinitrotoluene	5.089	0.20	5	0	102	50 - 122				
2,6-Dinitrotoluene	5.095	0.20	5	0	102	50 - 120				
2-Chloronaphthalene	4.451	0.20	5	0	89.0	50 - 120				
2-Methylnaphthalene	4.393	0.10	5	0	87.9	50 - 120				
4,6-Dinitro-2-methylphenol	5.82	0.20	5	0	116	25 - 121				
4-Nitrophenol	4.859	1.0	5	0	97.2	30 - 130				
Acenaphthene	4.701	0.10	5	0	94.0	45 - 120				
Acenaphthylene	4.377	0.10	5	0	87.5	47 - 120				
Anthracene	5.109	0.10	5	0	102	45 - 120				
Benz(a)anthracene	4.959	0.10	5	0	99.2	40 - 120				
Benzo(a)pyrene	4.774	0.10	5	0	95.5	45 - 120				
Bis(2-chloroethoxy)methane	4.305	0.20	5	0	86.1	45 - 120				
Bis(2-ethylhexyl)phthalate	6.051	0.20	5	0	121	40 - 139				
Chrysene	4.991	0.10	5	0	99.8	43 - 120				
Dibenzofuran	4.274	0.10	5	0	85.5	50 - 120				
Di-n-butyl phthalate	5.708	0.20	5	0	114	45 - 123				
Fluoranthene	4.84	0.10	5	0	96.8	45 - 125				
Fluorene	4.308	0.10	5	0	86.2	49 - 120				
Naphthalene	4.351	0.10	5	0	87.0	45 - 120				
Nitrobenzene	4.173	0.20	5	0	83.5	44 - 120				
N-Nitrosodiphenylamine	4.822	0.20	5	0	96.4	40 - 125				
Pentachlorophenol	4.568	0.20	5	0	91.4	19 - 121				
Phenanthrene	4.864	0.10	5	0	97.3	45 - 121				
Phenol	4.143	0.20	5	0	82.9	20 - 124				
Pyrene	5.11	0.10	5	0	102	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.644</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.9</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.364</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.709</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.2</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.728</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.07</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>81.4</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.948</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>79.0</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

Batch ID: 149790 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS20010713-08MS		Units: ug/L		Analysis Date: 21-Jan-2020 15:08				
Client ID: WG-1620-MW67B-20200115		Run ID: SV-6_354783		SeqNo: 5442897		PrepDate: 19-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.019	0.20	5	0	80.4	39 - 127				
2,4-Dimethylphenol	2.936	0.20	5	0	58.7	35 - 120				
2,4-Dinitrotoluene	4.832	0.20	5	0	96.6	50 - 122				
2,6-Dinitrotoluene	4.816	0.20	5	0	96.3	50 - 120				
2-Chloronaphthalene	3.838	0.20	5	0	76.8	50 - 120				
2-Methylnaphthalene	3.517	0.10	5	0	70.3	50 - 120				
4,6-Dinitro-2-methylphenol	5.71	0.20	5	0	114	25 - 121				
4-Nitrophenol	4.92	1.0	5	0	98.4	30 - 130				
Acenaphthene	4.264	0.10	5	0	85.3	45 - 120				
Acenaphthylene	3.808	0.10	5	0	76.2	47 - 120				
Anthracene	4.668	0.10	5	0	93.4	45 - 120				
Benz(a)anthracene	4.683	0.10	5	0	93.7	40 - 120				
Benzo(a)pyrene	4.504	0.10	5	0	90.1	45 - 120				
Bis(2-chloroethoxy)methane	3.413	0.20	5	0	68.3	45 - 120				
Bis(2-ethylhexyl)phthalate	5.894	0.20	5	0.08957	116	40 - 139				
Chrysene	4.496	0.10	5	0	89.9	43 - 120				
Dibenzofuran	3.835	0.10	5	0	76.7	50 - 120				
Di-n-butyl phthalate	5.43	0.20	5	0	109	45 - 123				
Fluoranthene	4.612	0.10	5	0	92.2	45 - 125				
Fluorene	4.02	0.10	5	0	80.4	49 - 120				
Naphthalene	3.538	0.10	5	0	70.8	45 - 120				
Nitrobenzene	3.265	0.20	5	0	65.3	44 - 120				
N-Nitrosodiphenylamine	4.782	0.20	5	0	95.6	40 - 125				
Pentachlorophenol	4.303	0.20	5	0	86.1	19 - 121				
Phenanthrene	4.398	0.10	5	0	88.0	45 - 121				
Phenol	3.368	0.20	5	0	67.4	20 - 124				
Pyrene	4.822	0.10	5	0	96.4	40 - 130				
Surr: 2,4,6-Tribromophenol	3.785	0.20	5	0	75.7	34 - 129				
Surr: 2-Fluorobiphenyl	3.928	0.20	5	0	78.6	40 - 125				
Surr: 2-Fluorophenol	3.33	0.20	5	0	66.6	20 - 120				
Surr: 4-Terphenyl-d14	4.641	0.20	5	0	92.8	40 - 135				
Surr: Nitrobenzene-d5	3.323	0.20	5	0	66.5	41 - 120				
Surr: Phenol-d6	3.289	0.20	5	0	65.8	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

Batch ID: 149790 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS20010713-08MSD			Units: ug/L		Analysis Date: 21-Jan-2020 15:27			
Client ID: WG-1620-MW67B-20200115		Run ID: SV-6_354783		SeqNo: 5442898		PrepDate: 19-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.253	0.20	5	0	85.1	39 - 127	4.019	5.65	20	
2,4-Dimethylphenol	2.99	0.20	5	0	59.8	35 - 120	2.936	1.82	20	
2,4-Dinitrotoluene	5.313	0.20	5	0	106	50 - 122	4.832	9.49	20	
2,6-Dinitrotoluene	4.866	0.20	5	0	97.3	50 - 120	4.816	1.02	20	
2-Chloronaphthalene	3.489	0.20	5	0	69.8	50 - 120	3.838	9.55	20	
2-Methylnaphthalene	3.303	0.10	5	0	66.1	50 - 120	3.517	6.29	20	
4,6-Dinitro-2-methylphenol	3.819	0.20	5	0	76.4	25 - 121	5.71	39.7	30	R
4-Nitrophenol	5.384	1.0	5	0	108	30 - 130	4.92	9.01	20	
Acenaphthene	3.975	0.10	5	0	79.5	45 - 120	4.264	7.01	20	
Acenaphthylene	3.616	0.10	5	0	72.3	47 - 120	3.808	5.17	20	
Anthracene	5.25	0.10	5	0	105	45 - 120	4.668	11.7	20	
Benz(a)anthracene	5.359	0.10	5	0	107	40 - 120	4.683	13.4	20	
Benzo(a)pyrene	5.276	0.10	5	0	106	45 - 120	4.504	15.8	20	
Bis(2-chloroethoxy)methane	3.093	0.20	5	0	61.9	45 - 120	3.413	9.83	20	
Bis(2-ethylhexyl)phthalate	6.076	0.20	5	0.08957	120	40 - 139	5.894	3.04	20	
Chrysene	5.125	0.10	5	0	102	43 - 120	4.496	13.1	20	
Dibenzofuran	3.691	0.10	5	0	73.8	50 - 120	3.835	3.84	20	
Di-n-butyl phthalate	6.04	0.20	5	0	121	45 - 123	5.43	10.6	20	
Fluoranthene	5.332	0.10	5	0	107	45 - 125	4.612	14.5	20	
Fluorene	4.082	0.10	5	0	81.6	49 - 120	4.02	1.53	20	
Naphthalene	3.323	0.10	5	0	66.5	45 - 120	3.538	6.29	20	
Nitrobenzene	3.067	0.20	5	0	61.3	44 - 120	3.265	6.27	20	
N-Nitrosodiphenylamine	5.371	0.20	5	0	107	40 - 125	4.782	11.6	20	
Pentachlorophenol	5.167	0.20	5	0	103	19 - 121	4.303	18.3	20	
Phenanthrene	4.947	0.10	5	0	98.9	45 - 121	4.398	11.7	20	
Phenol	3.164	0.20	5	0	63.3	20 - 124	3.368	6.23	20	
Pyrene	5.367	0.10	5	0	107	40 - 130	4.822	10.7	20	
Surr: 2,4,6-Tribromophenol	4.064	0.20	5	0	81.3	34 - 129	3.785	7.11	20	
Surr: 2-Fluorobiphenyl	3.411	0.20	5	0	68.2	40 - 125	3.928	14.1	20	
Surr: 2-Fluorophenol	2.871	0.20	5	0	57.4	20 - 120	3.33	14.8	20	
Surr: 4-Terphenyl-d14	5.181	0.20	5	0	104	40 - 135	4.641	11	20	
Surr: Nitrobenzene-d5	3.015	0.20	5	0	60.3	41 - 120	3.323	9.7	20	
Surr: Phenol-d6	3.041	0.20	5	0	60.8	20 - 120	3.289	7.86	20	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

**Batch ID:** 149790 ( 0 )      **Instrument:** SV-6      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

The following samples were analyzed in this batch:

HS20010713-01	HS20010713-02	HS20010713-03	HS20010713-04
HS20010713-05	HS20010713-06	HS20010713-07	HS20010713-08
HS20010713-09	HS20010713-10	HS20010713-11	HS20010713-12
HS20010713-13	HS20010713-14		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

<b>Batch ID:</b> R354726 ( 0 )	<b>Instrument:</b> VOA2	<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C
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<b>MBLK</b>		Sample ID: <b>VBLKW-200120</b>			Units: <b>ug/L</b>		Analysis Date: <b>20-Jan-2020 23:42</b>			
Client ID:		Run ID: <b>VOA2_354726</b>			SeqNo: <b>5441781</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.99</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.0</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.9</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200120</b>			Units: <b>ug/L</b>		Analysis Date: <b>20-Jan-2020 22:53</b>			
Client ID:		Run ID: <b>VOA2_354726</b>			SeqNo: <b>5441780</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.43	1.0	20	0	87.1	70 - 124				
Benzene	17.19	1.0	20	0	86.0	74 - 120				
Chlorobenzene	18.11	1.0	20	0	90.5	76 - 113				
Ethylbenzene	17.79	1.0	20	0	89.0	77 - 117				
Methylene chloride	17.05	2.0	20	0	85.3	70 - 127				
Toluene	18.88	1.0	20	0	94.4	77 - 118				
Xylenes, Total	54.29	1.0	60	0	90.5	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.0</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.41</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.15</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

**Batch ID:** R354726 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010713-08MS			Units: ug/L		Analysis Date: 21-Jan-2020 01:20			
Client ID: WG-1620-MW67B-20200115		Run ID: VOA2_354726			SeqNo: 5441843		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.86	1.0	20	0	84.3	70 - 127				
Benzene	17.23	1.0	20	0	86.2	70 - 127				
Chlorobenzene	17.78	1.0	20	0	88.9	70 - 114				
Ethylbenzene	17.83	1.0	20	0	89.1	70 - 124				
Methylene chloride	16.49	2.0	20	0	82.5	70 - 128				
Toluene	18.93	1.0	20	0	94.7	70 - 123				
Xylenes, Total	53.55	1.0	60	0	89.2	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.3</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.0</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.19</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.4</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>48.27</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.5</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010713-08MSD			Units: ug/L		Analysis Date: 21-Jan-2020 01:44			
Client ID: WG-1620-MW67B-20200115		Run ID: VOA2_354726			SeqNo: 5441844		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.39	1.0	20	0	82.0	70 - 127	16.86	2.79	20	
Benzene	16.52	1.0	20	0	82.6	70 - 127	17.23	4.21	20	
Chlorobenzene	17.4	1.0	20	0	87.0	70 - 114	17.78	2.16	20	
Ethylbenzene	17.35	1.0	20	0	86.8	70 - 124	17.83	2.72	20	
Methylene chloride	15.73	2.0	20	0	78.7	70 - 128	16.49	4.72	20	
Toluene	18.13	1.0	20	0	90.6	70 - 123	18.93	4.35	20	
Xylenes, Total	52.34	1.0	60	0	87.2	70 - 130	53.55	2.27	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>70 - 126</i>	<i>47.67</i>	<i>0.57</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>81 - 113</i>	<i>49.02</i>	<i>1.3</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.16</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.3</i>	<i>77 - 123</i>	<i>49.19</i>	<i>2.12</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.0</i>	<i>82 - 127</i>	<i>48.27</i>	<i>1.52</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS20010713-02	HS20010713-03	HS20010713-04	HS20010713-05
HS20010713-07	HS20010713-08	HS20010713-09	HS20010713-11
HS20010713-12	HS20010713-15		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

<b>Batch ID:</b> R354799 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200121</b>	Units: <b>ug/L</b>			Analysis Date: <b>21-Jan-2020 12:55</b>				
Client ID:	Run ID: <b>VOA2_354799</b>	SeqNo: <b>5443190</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.22</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>90.4</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.91</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200221</b>	Units: <b>ug/L</b>			Analysis Date: <b>21-Jan-2020 12:06</b>				
Client ID:	Run ID: <b>VOA2_354799</b>	SeqNo: <b>5443189</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	17.66	1.0	20	0	88.3	70 - 124			
Benzene	17.2	1.0	20	0	86.0	74 - 120			
Chlorobenzene	18.44	1.0	20	0	92.2	76 - 113			
Ethylbenzene	17.94	1.0	20	0	89.7	77 - 117			
Methylene chloride	17.49	2.0	20	0	87.4	70 - 127			
Toluene	19.3	1.0	20	0	96.5	77 - 118			
Vinyl chloride	15.17	1.0	20	0	75.8	70 - 130			
Xylenes, Total	54.82	1.0	60	0	91.4	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.76</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>81 - 120</i>			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

**Batch ID:** R354799 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010767-05MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 17:48</b>			
Client ID:		Run ID: <b>VOA2_354799</b>			SeqNo: <b>5443202</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.36	1.0	20	0	81.8	70 - 127				
Benzene	24.4	1.0	20	6.383	90.1	70 - 127				
Chlorobenzene	17.96	1.0	20	0	89.8	70 - 114				
Ethylbenzene	28.72	1.0	20	9.361	96.8	70 - 124				
Methylene chloride	15.89	2.0	20	0	79.4	70 - 128				
Toluene	20.5	1.0	20	0	102	70 - 123				
Vinyl chloride	17.82	1.0	20	0	89.1	70 - 130				
Xylenes, Total	57.71	1.0	60	3.181	90.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.1</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.0</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.4</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010767-05MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 18:12</b>			
Client ID:		Run ID: <b>VOA2_354799</b>			SeqNo: <b>5443203</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.51	1.0	20	0	82.6	70 - 127	16.36	0.908	20	
Benzene	24.24	1.0	20	6.383	89.3	70 - 127	24.4	0.678	20	
Chlorobenzene	17.68	1.0	20	0	88.4	70 - 114	17.96	1.54	20	
Ethylbenzene	28.32	1.0	20	9.361	94.8	70 - 124	28.72	1.41	20	
Methylene chloride	15.24	2.0	20	0	76.2	70 - 128	15.89	4.15	20	
Toluene	20.18	1.0	20	0	101	70 - 123	20.5	1.56	20	
Vinyl chloride	16.8	1.0	20	0	84.0	70 - 130	17.82	5.89	20	
Xylenes, Total	56.81	1.0	60	3.181	89.4	70 - 130	57.71	1.57	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.11</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.2</i>	<i>70 - 126</i>	<i>46.04</i>	<i>0.147</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>81 - 113</i>	<i>48.8</i>	<i>0.867</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>77 - 123</i>	<i>47.49</i>	<i>0.191</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>82 - 127</i>	<i>49.2</i>	<i>0.339</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010713-01      HS20010713-10      HS20010713-13      HS20010713-14

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

<b>Batch ID:</b> R355099 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200127</b>	Units: <b>ug/L</b>			Analysis Date: <b>27-Jan-2020 11:41</b>				
Client ID:	Run ID: <b>VOA4_355099</b>	SeqNo: <b>5448362</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>52.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.5</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>52.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>106</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200127</b>	Units: <b>ug/L</b>			Analysis Date: <b>27-Jan-2020 10:51</b>				
Client ID:	Run ID: <b>VOA4_355099</b>	SeqNo: <b>5448361</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18.52	1.0	20	0	92.6	70 - 124			
Benzene	19.58	1.0	20	0	97.9	74 - 120			
Chlorobenzene	19.39	1.0	20	0	97.0	76 - 113			
Ethylbenzene	19.72	1.0	20	0	98.6	77 - 117			
Methylene chloride	20.33	2.0	20	0	102	70 - 127			
Toluene	19.39	1.0	20	0	97.0	77 - 118			
Vinyl chloride	18.36	1.0	20	0	91.8	70 - 130			
Xylenes, Total	59.29	1.0	60	0	98.8	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.77</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.28</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.71</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.35</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.7</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

**Batch ID:** R355099 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010947-02MS			Units: ug/L		Analysis Date: 27-Jan-2020 14:11			
Client ID:		Run ID: VOA4_355099			SeqNo: 5448368		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.79	1.0	20	0	84.0	70 - 127				
Benzene	18.52	1.0	20	0	92.6	70 - 127				
Chlorobenzene	18.12	1.0	20	0	90.6	70 - 114				
Ethylbenzene	18.26	1.0	20	0	91.3	70 - 124				
Methylene chloride	19.39	2.0	20	0	97.0	70 - 128				
Toluene	19.1	1.0	20	0	95.5	70 - 123				
Vinyl chloride	18.48	1.0	20	0	92.4	70 - 130				
Xylenes, Total	55.63	1.0	60	0	92.7	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.93</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>52.13</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010947-02MSD			Units: ug/L		Analysis Date: 27-Jan-2020 14:36			
Client ID:		Run ID: VOA4_355099			SeqNo: 5448394		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.27	1.0	20	0	81.4	70 - 127	16.79	3.15	20	
Benzene	18.05	1.0	20	0	90.3	70 - 127	18.52	2.54	20	
Chlorobenzene	18	1.0	20	0	90.0	70 - 114	18.12	0.656	20	
Ethylbenzene	18.17	1.0	20	0	90.9	70 - 124	18.26	0.491	20	
Methylene chloride	18.89	2.0	20	0	94.4	70 - 128	19.39	2.63	20	
Toluene	18.46	1.0	20	0	92.3	70 - 123	19.1	3.39	20	
Vinyl chloride	17.75	1.0	20	0	88.7	70 - 130	18.48	4.06	20	
Xylenes, Total	54.43	1.0	60	0	90.7	70 - 130	55.63	2.19	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>	<i>50.93</i>	<i>0.53</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>	<i>50.89</i>	<i>0.119</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>51.77</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>77 - 123</i>	<i>52.13</i>	<i>0.708</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>82 - 127</i>	<i>50.6</i>	<i>1.23</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010713-06

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

Batch ID: R355193 ( 0 )		Instrument: VOA4		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MBLK</b>	Sample ID: <b>VBLKW-200128</b>	Units: <b>ug/L</b>			Analysis Date: <b>28-Jan-2020 13:04</b>					
Client ID:	Run ID: <b>VOA4_355193</b>	SeqNo: <b>5450213</b>		PrepDate:			DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
Vinyl chloride	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	50.07	1.0	50	0	100	70 - 123				
<i>Surr: 4-Bromofluorobenzene</i>	48.31	1.0	50	0	96.6	82 - 115				
<i>Surr: Dibromofluoromethane</i>	48.67	1.0	50	0	97.3	73 - 126				
<i>Surr: Toluene-d8</i>	50.19	1.0	50	0	100	81 - 120				
<b>LCS</b>	Sample ID: <b>VLCSW-200128</b>	Units: <b>ug/L</b>			Analysis Date: <b>28-Jan-2020 12:15</b>					
Client ID:	Run ID: <b>VOA4_355193</b>	SeqNo: <b>5450212</b>		PrepDate:			DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
Vinyl chloride	18.75	1.0	20	0	93.7	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	49.15	1.0	50	0	98.3	70 - 130				
<i>Surr: 4-Bromofluorobenzene</i>	50	1.0	50	0	100.0	82 - 115				
<i>Surr: Dibromofluoromethane</i>	49.33	1.0	50	0	98.7	73 - 126				
<i>Surr: Toluene-d8</i>	49.88	1.0	50	0	99.8	81 - 120				
<b>MS</b>	Sample ID: <b>HS20011178-01MS</b>	Units: <b>ug/L</b>			Analysis Date: <b>28-Jan-2020 14:18</b>					
Client ID:	Run ID: <b>VOA4_355193</b>	SeqNo: <b>5450501</b>		PrepDate:			DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
Vinyl chloride	17.86	1.0	20	0	89.3	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	49.02	1.0	50	0	98.0	70 - 126				
<i>Surr: 4-Bromofluorobenzene</i>	50.12	1.0	50	0	100	81 - 113				
<i>Surr: Dibromofluoromethane</i>	49.15	1.0	50	0	98.3	77 - 123				
<i>Surr: Toluene-d8</i>	48.65	1.0	50	0	97.3	82 - 127				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QC BATCH REPORT**

**Batch ID:** R355193 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MSD</b>		Sample ID: <b>HS20011178-01MSD</b>		Units: <b>ug/L</b>		Analysis Date: <b>28-Jan-2020 14:43</b>			
Client ID:		Run ID: <b>VOA4_355193</b>		SeqNo: <b>5450502</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
Vinyl chloride	16.89	1.0	20	0	84.4	70 - 130	17.86	5.6	20
<i>Surr: 1,2-Dichloroethane-d4</i>	49	1.0	50	0	98.0	70 - 126	49.02	0.0445	20
<i>Surr: 4-Bromofluorobenzene</i>	50.74	1.0	50	0	101	81 - 113	50.12	1.24	20
<i>Surr: Dibromofluoromethane</i>	49.79	1.0	50	0	99.6	77 - 123	49.15	1.29	20
<i>Surr: Toluene-d8</i>	49.76	1.0	50	0	99.5	82 - 127	48.65	2.25	20

The following samples were analyzed in this batch: HS20010713-15

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010713

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010713

Date/Time Received: 16-Jan-2020 12:40
Received by: NDD

Checklist completed by: Nelson D. Dusara
eSignature
Date: 17-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 21-Jan-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

Temperature(s)/Thermometer(s): 0.5,0.7,0.8,1.2 c UC/c IR 11

Cooler(s)/Kit(s): 45645,25638,45624,45665

Date/Time sample(s) sent to storage: Jan/17/2020 08:45

Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]

Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]

pH adjusted? Yes [ ] No [checked] N/A [ ]

pH adjusted by: [ ]

Login Notes: WG-1620-MW25C-20200115 time differs: COC =15:50 Label=12:15; logged per COC.

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments: [ ]

Corrective Action: [ ]





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# Chain of Custody Form

Page 1 of 2

COC ID: 206377

HS20010713

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As)
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E	
				F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TBO-202001</del>			<del>Water</del>	1	2	X	X	X								
2	WG-1620-MWISA-20200114	1-14-20	1345	W		6	X		X	X							
3	WG-1620-MWISB-20200114		1430	W		6	X		X	X							
4	WG-1620-MWISC-20200114		1520	W		6	X		X	X							
5	WG-1620-MWI4-20200114		1615	W		6	X		X	X							
6	WG-1620-MWI7-20200114		1710	W		6	X		X	X							
7	WG-1620-FB06-20200114		1725	W		6	X		X	X							
8	WG-1620-MWI7C-20200115	1-15-20	0815	W		6	X		X	X							
9	WG-1620-MWI67B-20200115		0930	W		6	X		X	X							
10	WG-1620-MWI67BMS-20200115		0930	W		6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour			Results Due Date:	
Relinquished by: <b>John Br</b>	Date: <b>1-16-20</b>	Time:	Received by:	Notes: <b>UPRR Houston MWPPW</b>		QC Package: (Check One Box Below)		
Relinquished by: <b>John Br</b>	Date: <b>Jan 16/20</b>	Time: <b>12:40</b>	Received by (Laboratory): <b>M. Nelson</b>	Cooler ID: <b>45645</b>	Cooler Temp. <b>4°C</b>	<input type="checkbox"/> Level II Std OC	<input checked="" type="checkbox"/> TRRP Check st	
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	<b>45638</b>	<b>0.5</b>	<input type="checkbox"/> Level III Std OC/Row Date	<input type="checkbox"/> TRRP Level IV	
				<b>45624</b>	<b>0.8</b>	<input type="checkbox"/> Level IV SV/843/CLP		
				<b>45665</b>	<b>1.2</b>	<input type="checkbox"/> Other		

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

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# Chain of Custody Form

Page 2 of 2

COC ID: 206378

HS20010713

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A 8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B 8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C 8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D ICP_TVW (5636002 5652646 Metals - As)
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	F
Phone	(512) 671-3434	Phone		G
Fax	(512) 671-3446	Fax		H
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		I
				J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TB0-202001			Water		2		*									
2	WG-1620-MW67BMSD-20200115	1-15-20	0930	W		6	X		X	X							
3	WG-1620-MW46C-20200115		1025	W		6	X		X	X							
4	WG-1620-MW25A-20200115		1120	W		6		X	X	X							
5	WG-1620-MW44C-20200115		1315	W		6	X		X	X							
6	WG-1620-MW71B-20200115		1400	W		6	X		X	X							
7	WG-1620-MW33BR-20200115		1500	W		6		X	X	X							
8	WG-1620-MW25C-20200115		1550	W		6		X	X	X							
9	WG-1620-TB04-20200115		-	W		2	X										
10																	

Sampler(s) Please Print & Sign <b>JOHN BEAUBON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Deys <input type="checkbox"/> 5 Wk Deys <input type="checkbox"/> 2 Wk Deys <input type="checkbox"/> 24 Hour				Results Due Date:			
Relinquished by: <b>John Beaubon</b>	Date: 1-16-20	Time:	Received by:	Notes: <b>UPRR Houston MWPW</b>							
Relinquished by: <b>John Beaubon</b>	Date: Jan 16/20	Time: 12:40	Received by (Laboratory): <b>Nelson</b>	Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)					
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	<input type="checkbox"/> Level II Std QC	<input checked="" type="checkbox"/> TRRP Check st						
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				<input type="checkbox"/> Level III Std QC/Raw Date	<input type="checkbox"/> TRRP Level IV						
				<input type="checkbox"/> Level IV SV483/CLP	<input type="checkbox"/> Other						

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January 28, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010767**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 29 sample(s) on Jan 17, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read 'Dane J. Wacasey'.

Generated By: DANE.WACASEY  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

---

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

**Laboratory Review Checklist: Reportable Data**

Laboratory Name: ALS Laboratory Group			LRC Date: 01/28/2020				
Project Name: Houston TX-Wood Preserving Works			Laboratory Job Number: HS20010767				
Reviewer Name: Dane Wacasey			Prep Batch Number(s): 149797,149831,149832,149881,149912,149913,149936,R354794,R354799,R354902,R354859,R355066				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSS included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group			LRC Date: 01/28/2020				
Project Name: Houston TX-Wood Preserving Works			Laboratory Job Number: HS20010767				
Reviewer Name: Dane Wacasey			Prep Batch Number(s): 149797,149831,149832,149881,149912,149913,149936,R354794,R354799,R354902,R354859,R355066				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 01/28/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20010767
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 149797,149831,149832,149881,149912,149913,149936,R354794,R354799,R354902,R354859,R355066

ER# <sup>5</sup>	Description
1	Semivolatile Organics Method SW8270, samples WG-1620-MW75B-20200117, WG-1620-MW79A-20200117, WG-1620-MW74B-20200117, WG-1620-MW78A-20200117, the surrogate recoveries could not be determined due to dilution below the calibration range.  Semivolatile Organics Method SW8270, sample WG-1620-MW75B-20200117, surrogate Phenol-d6 recovered above the upper control limit. This was due to a dilution required for sample analysis.
2	Batch 149881, Metals Method SW6020, sample HS20010752-08, MS was performed on unrelated sample.  Batch 149912, Semivolatile Organics Method SW8270, sample WG-1620-MW28C-20200116, MS recovered above the control limits for 4,6-Dinitro-2-methylphenol and Di-n-butyl phthalate due to suspect matrix effect.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
 O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
 NA = Not Applicable;  
 NR = Not Reviewed;  
 R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010767

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010767-01	WG-1620-MW28A-20200116	Water		16-Jan-2020 09:10	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-02	WG-1620-MW28C-20200116	Water		16-Jan-2020 10:15	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-03	WG-1620-FB08-20200116	Water		16-Jan-2020 12:00	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-04	WG-1620-MW89B-20200116	Water		16-Jan-2020 12:45	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-05	WG-1620-MW84B-20200116	Water		16-Jan-2020 13:45	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-06	WG-1620-MW63B-20200116	Water		16-Jan-2020 14:45	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-07	WG-1620-MW26A-20200116	Water		16-Jan-2020 16:00	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-08	WG-1620-MW68C-20200117	Water		17-Jan-2020 09:10	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-09	WG-1620-MW68A-20200117	Water		17-Jan-2020 10:15	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-10	WG-1620-FB09-20200117	Water		17-Jan-2020 10:00	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-11	WG-1620-MW27A-20200115	Water		15-Jan-2020 16:40	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-12	WG-1620-MW27C-20200115	Water		15-Jan-2020 17:20	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-13	WG-1620-MW59A-20200116	Water		16-Jan-2020 08:05	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-14	WG-1620-MW59B-20200116	Water		16-Jan-2020 09:00	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-15	WG-1620-MW59D-20200116	Water		16-Jan-2020 13:50	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-16	WG-1620-FD03-20200116	Water		16-Jan-2020 13:50	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-17	WG-1620-MW66D-20200116	Water		16-Jan-2020 15:00	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-18	WG-1620-MW65D-20200116	Water		16-Jan-2020 16:20	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-19	WG-1620-MW36D-20200116	Water		16-Jan-2020 17:25	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-20	WG-1620-MW86C-20200117	Water		17-Jan-2020 07:25	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-21	WG-1620-FD04-20200117	Water		17-Jan-2020 07:25	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-22	WG-1620-MW75B-20200117	Water		17-Jan-2020 08:15	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-23	WG-1620-MW79A-20200117	Water		17-Jan-2020 09:20	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-24	WG-1620-MW74B-20200117	Water		17-Jan-2020 10:20	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-25	WG-1620-MW78A-20200117	Water		17-Jan-2020 11:25	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-26	WQ-1620-TB05-2020117	Water		17-Jan-2020 00:00	17-Jan-2020 13:11	<input type="checkbox"/>

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010767

**SAMPLE SUMMARY**

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Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010767-27	WG-1620-MW48C-20200116	Water		16-Jan-2020 10:20	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-28	WG-1620-MW47C-20200116	Water		16-Jan-2020 11:05	17-Jan-2020 13:11	<input type="checkbox"/>
HS20010767-29	WG-1620-MW61A-20200116	Water		16-Jan-2020 12:00	17-Jan-2020 13:11	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28A-20200116  
 Collection Date: 16-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 00:55
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 00:55
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 00:55
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 00:55
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 00:55
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 00:55
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 00:55
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:55</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:55</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:55</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 00:55</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28A-20200116  
 Collection Date: 16-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
<b>1,2-Diphenylhydrazine</b>	<b>0.000034</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
2,4-Dimethylphenol		U	0.000040	0.00020	mg/L	1	25-Jan-2020 18:43
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	25-Jan-2020 18:43
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	25-Jan-2020 18:43
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	25-Jan-2020 18:43
2-Methylnaphthalene		U	0.000019	0.00010	mg/L	1	25-Jan-2020 18:43
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	25-Jan-2020 18:43
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	25-Jan-2020 18:43
<b>Acenaphthene</b>	<b>0.00014</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
Acenaphthylene		U	0.000015	0.00010	mg/L	1	25-Jan-2020 18:43
<b>Anthracene</b>	<b>0.00013</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Benz(a)anthracene</b>	<b>0.000073</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Benzo(a)pyrene</b>	<b>0.000057</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	25-Jan-2020 18:43
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000090</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Chrysene</b>	<b>0.000065</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Dibenzofuran</b>	<b>0.00010</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Di-n-butyl phthalate</b>	<b>0.000032</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Fluoranthene</b>	<b>0.00030</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Fluorene</b>	<b>0.00014</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<b>Naphthalene</b>	<b>0.00060</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
Nitrobenzene		U	0.000024	0.00020	mg/L	1	25-Jan-2020 18:43
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	25-Jan-2020 18:43
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	25-Jan-2020 18:43
<b>Phenanthrene</b>	<b>0.00048</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
Phenol		U	0.000035	0.00020	mg/L	1	25-Jan-2020 18:43
<b>Pyrene</b>	<b>0.00021</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 18:43
<i>Surr: 2,4,6-Tribromophenol</i>	69.8			34-129	%REC	1	25-Jan-2020 18:43
<i>Surr: 2-Fluorobiphenyl</i>	78.4			40-125	%REC	1	25-Jan-2020 18:43
<i>Surr: 2-Fluorophenol</i>	61.6			20-120	%REC	1	25-Jan-2020 18:43
<i>Surr: 4-Terphenyl-d14</i>	87.8			40-135	%REC	1	25-Jan-2020 18:43
<i>Surr: Nitrobenzene-d5</i>	72.2			41-120	%REC	1	25-Jan-2020 18:43
<i>Surr: Phenol-d6</i>	66.5			20-120	%REC	1	25-Jan-2020 18:43
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 21-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00664</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	22-Jan-2020 23:52

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28C-20200116  
 Collection Date: 16-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 23:41
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 23:41
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 23:41
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 23:41
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 23:41
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 23:41
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 23:41
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:41</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:41</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:41</i>
<i>Surr: Toluene-d8</i>	<i>97.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 23:41</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW28C-20200116  
 Collection Date: 16-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 12:03
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 12:03
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 12:03
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 12:03
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 12:03
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	25-Jan-2020 12:03
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 12:03
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 12:03
Acenaphthene	U		0.000027	0.00010	mg/L	1	25-Jan-2020 12:03
Acenaphthylene	U		0.000015	0.00010	mg/L	1	25-Jan-2020 12:03
Anthracene	U		0.000014	0.00010	mg/L	1	25-Jan-2020 12:03
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 12:03
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	25-Jan-2020 12:03
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 12:03
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000072</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 12:03
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 12:03
Dibenzofuran	U		0.000020	0.00010	mg/L	1	25-Jan-2020 12:03
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	25-Jan-2020 12:03
Fluoranthene	U		0.000010	0.00010	mg/L	1	25-Jan-2020 12:03
Fluorene	U		0.000030	0.00010	mg/L	1	25-Jan-2020 12:03
<b>Naphthalene</b>	<b>0.00016</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 12:03
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 12:03
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 12:03
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 12:03
Phenanthrene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 12:03
<b>Phenol</b>	<b>0.000086</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 12:03
Pyrene	U		0.000019	0.00010	mg/L	1	25-Jan-2020 12:03
<i>Surr: 2,4,6-Tribromophenol</i>	<i>74.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 12:03</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>71.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 12:03</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 12:03</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>96.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 12:03</i>
<i>Surr: Nitrobenzene-d5</i>	<i>63.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 12:03</i>
<i>Surr: Phenol-d6</i>	<i>71.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 12:03</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 21-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.000937</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	22-Jan-2020 23:32

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB08-20200116  
 Collection Date: 16-Jan-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 02:58
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 02:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 02:58
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 02:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 02:58
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 02:58
Vinyl chloride	U		0.00020	0.0010	mg/L	1	22-Jan-2020 02:58
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 02:58
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>101</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 02:58</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.5</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 02:58</i>
<i>Surr: Dibromofluoromethane</i>		<i>98.2</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 02:58</i>
<i>Surr: Toluene-d8</i>		<i>99.9</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 02:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB08-20200116  
 Collection Date: 16-Jan-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 19:02
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 19:02
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 19:02
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 19:02
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 19:02
<b>2-Methylnaphthalene</b>	<b>0.000052</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 19:02
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 19:02
<b>Acenaphthene</b>	<b>0.000077</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
Acenaphthylene	U		0.000015	0.00010	mg/L	1	25-Jan-2020 19:02
<b>Anthracene</b>	<b>0.000027</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 19:02
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	25-Jan-2020 19:02
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 19:02
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	25-Jan-2020 19:02
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 19:02
<b>Dibenzofuran</b>	<b>0.000066</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	25-Jan-2020 19:02
<b>Fluoranthene</b>	<b>0.000033</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
<b>Fluorene</b>	<b>0.000058</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
<b>Naphthalene</b>	<b>0.00054</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 19:02
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 19:02
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 19:02
<b>Phenanthrene</b>	<b>0.000067</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 19:02
<b>Pyrene</b>	<b>0.000022</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:02
Surr: 2,4,6-Tribromophenol	61.1			34-129	%REC	1	25-Jan-2020 19:02
Surr: 2-Fluorobiphenyl	92.5			40-125	%REC	1	25-Jan-2020 19:02
Surr: 2-Fluorophenol	71.4			20-120	%REC	1	25-Jan-2020 19:02
Surr: 4-Terphenyl-d14	93.8			40-135	%REC	1	25-Jan-2020 19:02
Surr: Nitrobenzene-d5	91.0			41-120	%REC	1	25-Jan-2020 19:02
Surr: Phenol-d6	79.1			20-120	%REC	1	25-Jan-2020 19:02
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jan-2020		Analyst: JC	
Arsenic	U		0.000400	0.00200	mg/L	1	24-Jan-2020 16:10

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW89B-20200116  
 Collection Date: 16-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 18:37
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 18:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 18:37
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 18:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 18:37
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 18:37
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 18:37
<i>Surr: 1,2-Dichloroethane-d4</i>		89.2		70-126	%REC	1	21-Jan-2020 18:37
<i>Surr: 4-Bromofluorobenzene</i>		94.6		81-113	%REC	1	21-Jan-2020 18:37
<i>Surr: Dibromofluoromethane</i>		96.3		77-123	%REC	1	21-Jan-2020 18:37
<i>Surr: Toluene-d8</i>		99.4		82-127	%REC	1	21-Jan-2020 18:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW89B-20200116  
 Collection Date: 16-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 19:22
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 19:22
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 19:22
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 19:22
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 19:22
<b>2-Methylnaphthalene</b>	<b>0.000032</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 19:22
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 19:22
<b>Acenaphthene</b>	<b>0.000083</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Acenaphthylene</b>	<b>0.000045</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Anthracene</b>	<b>0.00020</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Benz(a)anthracene</b>	<b>0.00017</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Benzo(a)pyrene</b>	<b>0.000079</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 19:22
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000099</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Chrysene</b>	<b>0.00024</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Dibenzofuran</b>	<b>0.00011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Di-n-butyl phthalate</b>	<b>0.000034</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Fluoranthene</b>	<b>0.0012</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Fluorene</b>	<b>0.00018</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<b>Naphthalene</b>	<b>0.00029</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 19:22
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 19:22
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 19:22
<b>Phenanthrene</b>	<b>0.00083</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 19:22
<b>Pyrene</b>	<b>0.00081</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:22
<i>Surr: 2,4,6-Tribromophenol</i>	73.7			34-129	%REC	1	25-Jan-2020 19:22
<i>Surr: 2-Fluorobiphenyl</i>	74.9			40-125	%REC	1	25-Jan-2020 19:22
<i>Surr: 2-Fluorophenol</i>	61.8			20-120	%REC	1	25-Jan-2020 19:22
<i>Surr: 4-Terphenyl-d14</i>	90.7			40-135	%REC	1	25-Jan-2020 19:22
<i>Surr: Nitrobenzene-d5</i>	75.3			41-120	%REC	1	25-Jan-2020 19:22
<i>Surr: Phenol-d6</i>	72.4			20-120	%REC	1	25-Jan-2020 19:22
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 21-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.000463</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	22-Jan-2020 23:54

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84B-20200116  
 Collection Date: 16-Jan-2020 13:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 17:23
<b>Benzene</b>	<b>0.0064</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 17:23
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 17:23
<b>Ethylbenzene</b>	<b>0.0094</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 17:23
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 17:23
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 17:23
<b>Xylenes, Total</b>	<b>0.0032</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 17:23
<i>Surr: 1,2-Dichloroethane-d4</i>	88.3			70-126	%REC	1	21-Jan-2020 17:23
<i>Surr: 4-Bromofluorobenzene</i>	95.9			81-113	%REC	1	21-Jan-2020 17:23
<i>Surr: Dibromofluoromethane</i>	95.4			77-123	%REC	1	21-Jan-2020 17:23
<i>Surr: Toluene-d8</i>	99.6			82-127	%REC	1	21-Jan-2020 17:23

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84B-20200116  
 Collection Date: 16-Jan-2020 13:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	25-Jan-2020 19:41
<b>2,4-Dimethylphenol</b>	<b>0.00048</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	25-Jan-2020 19:41
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	25-Jan-2020 19:41
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	25-Jan-2020 19:41
<b>2-Methylnaphthalene</b>	<b>0.0065</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
<b>4,6-Dinitro-2-methylphenol</b>	<b>0.00035</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	25-Jan-2020 19:41
<b>Acenaphthene</b>	<b>0.025</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 16:43
<b>Acenaphthylene</b>	<b>0.00031</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
<b>Anthracene</b>	<b>0.00095</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	25-Jan-2020 19:41
Benzo(a)pyrene		U	0.000020	0.00010	mg/L	1	25-Jan-2020 19:41
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	25-Jan-2020 19:41
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	25-Jan-2020 19:41
Chrysene		U	0.000021	0.00010	mg/L	1	25-Jan-2020 19:41
<b>Dibenzofuran</b>	<b>0.015</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 16:43
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	25-Jan-2020 19:41
<b>Fluoranthene</b>	<b>0.00037</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
<b>Fluorene</b>	<b>0.0063</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
<b>Naphthalene</b>	<b>0.13</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 16:43
Nitrobenzene		U	0.000024	0.00020	mg/L	1	25-Jan-2020 19:41
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	25-Jan-2020 19:41
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	25-Jan-2020 19:41
<b>Phenanthrene</b>	<b>0.0065</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
Phenol		U	0.000035	0.00020	mg/L	1	25-Jan-2020 19:41
<b>Pyrene</b>	<b>0.00025</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 19:41
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.0</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 16:43</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>73.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:41</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>62.7</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:41</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.4</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 16:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>85.3</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 16:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>55.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:41</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>98.7</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:41</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>104</i>			<i>40-135</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 16:43</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.9</i>	<i>J</i>		<i>41-120</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 16:43</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:41</i>
<i>Surr: Phenol-d6</i>	<i>62.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:41</i>
<i>Surr: Phenol-d6</i>	<i>40.8</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 16:43</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84B-20200116  
 Collection Date: 16-Jan-2020 13:45

**ANALYTICAL REPORT**

WorkOrder:HS20010767  
 Lab ID:HS20010767-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jan-2020		Analyst: JC	
Arsenic	0.00363		0.000400	0.00200	mg/L	1	24-Jan-2020 16:13

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW63B-20200116  
 Collection Date: 16-Jan-2020 14:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:01
<b>Benzene</b>	<b>0.018</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 19:01
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 19:01
<b>Ethylbenzene</b>	<b>0.051</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 19:01
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 19:01
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:01
<b>Xylenes, Total</b>	<b>0.013</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 19:01
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:01</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:01</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:01</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:01</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW63B-20200116  
 Collection Date: 16-Jan-2020 14:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:00
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 20:00
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 20:00
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 20:00
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:00
<b>2-Methylnaphthalene</b>	<b>0.000056</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 20:00
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 20:00
<b>Acenaphthene</b>	<b>0.000059</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
<b>Acenaphthylene</b>	<b>0.000032</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
<b>Anthracene</b>	<b>0.000081</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 20:00
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	25-Jan-2020 20:00
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 20:00
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000065</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 20:00
<b>Dibenzofuran</b>	<b>0.000063</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
<b>Di-n-butyl phthalate</b>	<b>0.000034</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
<b>Fluoranthene</b>	<b>0.000023</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
<b>Fluorene</b>	<b>0.00012</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
<b>Naphthalene</b>	<b>0.00035</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 20:00
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 20:00
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 20:00
<b>Phenanthrene</b>	<b>0.000072</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 20:00
<b>Pyrene</b>	<b>0.000021</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:00
Surr: 2,4,6-Tribromophenol	63.5			34-129	%REC	1	25-Jan-2020 20:00
Surr: 2-Fluorobiphenyl	57.6			40-125	%REC	1	25-Jan-2020 20:00
Surr: 2-Fluorophenol	51.8			20-120	%REC	1	25-Jan-2020 20:00
Surr: 4-Terphenyl-d14	87.3			40-135	%REC	1	25-Jan-2020 20:00
Surr: Nitrobenzene-d5	57.4			41-120	%REC	1	25-Jan-2020 20:00
Surr: Phenol-d6	58.3			20-120	%REC	1	25-Jan-2020 20:00
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00204</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 16:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW26A-20200116  
 Collection Date: 16-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 09:17
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 09:17
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 09:17
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 09:17
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 09:17
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 09:17
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 09:17
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 09:17</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 09:17</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 09:17</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 09:17</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW26A-20200116  
 Collection Date: 16-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:19
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 20:19
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 20:19
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 20:19
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:19
<b>2-Methylnaphthalene</b>	<b>0.000042</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 20:19
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 20:19
<b>Acenaphthene</b>	<b>0.00017</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
Acenaphthylene	U		0.000015	0.00010	mg/L	1	25-Jan-2020 20:19
<b>Anthracene</b>	<b>0.000087</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 20:19
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	25-Jan-2020 20:19
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 20:19
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	25-Jan-2020 20:19
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 20:19
<b>Dibenzofuran</b>	<b>0.000044</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	25-Jan-2020 20:19
<b>Fluoranthene</b>	<b>0.00073</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
<b>Fluorene</b>	<b>0.00029</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
<b>Naphthalene</b>	<b>0.00017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 20:19
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 20:19
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 20:19
<b>Phenanthrene</b>	<b>0.000030</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 20:19
<b>Pyrene</b>	<b>0.00035</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:19
Surr: 2,4,6-Tribromophenol	77.9			34-129	%REC	1	25-Jan-2020 20:19
Surr: 2-Fluorobiphenyl	85.4			40-125	%REC	1	25-Jan-2020 20:19
Surr: 2-Fluorophenol	72.2			20-120	%REC	1	25-Jan-2020 20:19
Surr: 4-Terphenyl-d14	88.9			40-135	%REC	1	25-Jan-2020 20:19
Surr: Nitrobenzene-d5	78.4			41-120	%REC	1	25-Jan-2020 20:19
Surr: Phenol-d6	72.2			20-120	%REC	1	25-Jan-2020 20:19
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.0217</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 16:17

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68C-20200117  
 Collection Date: 17-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:26
<b>Benzene</b>	<b>0.00056</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 19:26
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 19:26
<b>Ethylbenzene</b>	<b>0.00047</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 19:26
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 19:26
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:26
<b>Xylenes, Total</b>	<b>0.00062</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 19:26
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:26</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:26</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:26</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:26</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68C-20200117  
 Collection Date: 17-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:38
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 20:38
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 20:38
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 20:38
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:38
<b>2-Methylnaphthalene</b>	<b>0.00033</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 20:38
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 20:38
<b>Acenaphthene</b>	<b>0.000036</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Acenaphthylene</b>	<b>0.000023</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Anthracene</b>	<b>0.000036</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Benz(a)anthracene</b>	<b>0.000055</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Benzo(a)pyrene</b>	<b>0.000063</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 20:38
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00029</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Chrysene</b>	<b>0.000054</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Dibenzofuran</b>	<b>0.000030</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	25-Jan-2020 20:38
<b>Fluoranthene</b>	<b>0.000074</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Fluorene</b>	<b>0.000036</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<b>Naphthalene</b>	<b>0.00052</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 20:38
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 20:38
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 20:38
<b>Phenanthrene</b>	<b>0.000042</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 20:38
<b>Pyrene</b>	<b>0.000092</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 20:38
<i>Surr: 2,4,6-Tribromophenol</i>	69.6			34-129	%REC	1	25-Jan-2020 20:38
<i>Surr: 2-Fluorobiphenyl</i>	71.4			40-125	%REC	1	25-Jan-2020 20:38
<i>Surr: 2-Fluorophenol</i>	61.9			20-120	%REC	1	25-Jan-2020 20:38
<i>Surr: 4-Terphenyl-d14</i>	96.7			40-135	%REC	1	25-Jan-2020 20:38
<i>Surr: Nitrobenzene-d5</i>	73.4			41-120	%REC	1	25-Jan-2020 20:38
<i>Surr: Phenol-d6</i>	65.8			20-120	%REC	1	25-Jan-2020 20:38
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 22-Jan-2020		Analyst: JC	
Arsenic	U		0.000400	0.00200	mg/L	1	24-Jan-2020 16:21

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68A-20200117  
 Collection Date: 17-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:50
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:50
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 19:50
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 19:50
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 19:50
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 19:50
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 19:50
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:50</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:50</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:50</i>
<i>Surr: Toluene-d8</i>	<i>98.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 19:50</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68A-20200117  
 Collection Date: 17-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:57
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 20:57
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 20:57
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 20:57
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 20:57
<b>2-Methylnaphthalene</b>	<b>0.000048</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 20:57
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 20:57
<b>Acenaphthene</b>	<b>0.00030</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
<b>Acenaphthylene</b>	<b>0.000054</b>	<b>J</b>	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
<b>Anthracene</b>	<b>0.000059</b>	<b>J</b>	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 20:57
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	25-Jan-2020 20:57
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 20:57
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000061</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 20:57
Dibenzofuran	U		0.000020	0.00010	mg/L	1	25-Jan-2020 20:57
<b>Di-n-butyl phthalate</b>	<b>0.000034</b>	<b>J</b>	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
<b>Fluoranthene</b>	<b>0.00013</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
<b>Fluorene</b>	<b>0.000069</b>	<b>J</b>	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
<b>Naphthalene</b>	<b>0.00022</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 20:57
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 20:57
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 20:57
<b>Phenanthrene</b>	<b>0.00016</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 20:57
<b>Pyrene</b>	<b>0.000068</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>25-Jan-2020 20:57</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>80.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>97.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>100</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:57</i>
<i>Surr: Nitrobenzene-d5</i>	<i>95.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:57</i>
<i>Surr: Phenol-d6</i>	<i>86.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:57</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.0423</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jan-2020 03:07</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB09-20200117  
 Collection Date: 17-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:08
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:08
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:08
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:08
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 01:08
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:08
Vinyl chloride	U		0.00020	0.0010	mg/L	1	22-Jan-2020 13:50
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:08
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 01:08</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>103</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 13:50</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 01:08</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 13:50</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.2</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 01:08</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 13:50</i>
<i>Surr: Toluene-d8</i>	<i>99.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 01:08</i>
<i>Surr: Toluene-d8</i>	<i>97.1</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 13:50</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB09-20200117  
 Collection Date: 17-Jan-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 21:16
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 21:16
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 21:16
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 21:16
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 21:16
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	25-Jan-2020 21:16
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 21:16
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 21:16
Acenaphthene	U		0.000027	0.00010	mg/L	1	25-Jan-2020 21:16
Acenaphthylene	U		0.000015	0.00010	mg/L	1	25-Jan-2020 21:16
Anthracene	U		0.000014	0.00010	mg/L	1	25-Jan-2020 21:16
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 21:16
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	25-Jan-2020 21:16
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 21:16
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000050</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 21:16
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 21:16
Dibenzofuran	U		0.000020	0.00010	mg/L	1	25-Jan-2020 21:16
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	25-Jan-2020 21:16
Fluoranthene	U		0.000010	0.00010	mg/L	1	25-Jan-2020 21:16
Fluorene	U		0.000030	0.00010	mg/L	1	25-Jan-2020 21:16
<b>Naphthalene</b>	<b>0.00017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 21:16
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 21:16
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 21:16
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 21:16
Phenanthrene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 21:16
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 21:16
Pyrene	U		0.000019	0.00010	mg/L	1	25-Jan-2020 21:16
<i>Surr: 2,4,6-Tribromophenol</i>	<i>65.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>93.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>79.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:16</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>98.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>93.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:16</i>
<i>Surr: Phenol-d6</i>	<i>86.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:16</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	U		0.000400	0.00200	mg/L	1	24-Jan-2020 03:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27A-20200115  
 Collection Date: 15-Jan-2020 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 20:14
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 20:14
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 20:14
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 20:14
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 20:14
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 20:14
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 20:14
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.5</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:14</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:14</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:14</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:14</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27A-20200115  
 Collection Date: 15-Jan-2020 16:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	25-Jan-2020 21:35
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	25-Jan-2020 21:35
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	25-Jan-2020 21:35
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	25-Jan-2020 21:35
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	25-Jan-2020 21:35
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	25-Jan-2020 21:35
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	25-Jan-2020 21:35
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	25-Jan-2020 21:35
Acenaphthene	U		0.000027	0.00010	mg/L	1	25-Jan-2020 21:35
Acenaphthylene	U		0.000015	0.00010	mg/L	1	25-Jan-2020 21:35
Anthracene	U		0.000014	0.00010	mg/L	1	25-Jan-2020 21:35
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	25-Jan-2020 21:35
<b>Benzo(a)pyrene</b>	<b>0.000052</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 21:35
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	25-Jan-2020 21:35
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000062</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	25-Jan-2020 21:35
Chrysene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 21:35
Dibenzofuran	U		0.000020	0.00010	mg/L	1	25-Jan-2020 21:35
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	25-Jan-2020 21:35
Fluoranthene	U		0.000010	0.00010	mg/L	1	25-Jan-2020 21:35
Fluorene	U		0.000030	0.00010	mg/L	1	25-Jan-2020 21:35
<b>Naphthalene</b>	<b>0.000053</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	25-Jan-2020 21:35
Nitrobenzene	U		0.000024	0.00020	mg/L	1	25-Jan-2020 21:35
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	25-Jan-2020 21:35
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	25-Jan-2020 21:35
Phenanthrene	U		0.000021	0.00010	mg/L	1	25-Jan-2020 21:35
Phenol	U		0.000035	0.00020	mg/L	1	25-Jan-2020 21:35
Pyrene	U		0.000019	0.00010	mg/L	1	25-Jan-2020 21:35
<i>Surr: 2,4,6-Tribromophenol</i>	<i>53.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:35</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>68.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:35</i>
<i>Surr: 2-Fluorophenol</i>	<i>57.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:35</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>99.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:35</i>
<i>Surr: Nitrobenzene-d5</i>	<i>68.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:35</i>
<i>Surr: Phenol-d6</i>	<i>65.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:35</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000859</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 03:11

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27C-20200115  
 Collection Date: 15-Jan-2020 17:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 20:39
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 20:39
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 20:39
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 20:39
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 20:39
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 20:39
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 20:39
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:39</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:39</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:39</i>
<i>Surr: Toluene-d8</i>	<i>99.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 20:39</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW27C-20200115  
 Collection Date: 15-Jan-2020 17:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-12  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 13:17
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	26-Jan-2020 13:17
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 13:17
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 13:17
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 13:17
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 13:17
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 13:17
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 13:17
Acenaphthene	U		0.000027	0.00010	mg/L	1	26-Jan-2020 13:17
Acenaphthylene	U		0.000015	0.00010	mg/L	1	26-Jan-2020 13:17
Anthracene	U		0.000014	0.00010	mg/L	1	26-Jan-2020 13:17
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	26-Jan-2020 13:17
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	26-Jan-2020 13:17
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 13:17
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	26-Jan-2020 13:17
Chrysene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 13:17
Dibenzofuran	U		0.000020	0.00010	mg/L	1	26-Jan-2020 13:17
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	26-Jan-2020 13:17
Fluoranthene	U		0.000010	0.00010	mg/L	1	26-Jan-2020 13:17
Fluorene	U		0.000030	0.00010	mg/L	1	26-Jan-2020 13:17
<b>Naphthalene</b>	<b>0.000055</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 13:17</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 13:17
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 13:17
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 13:17
Phenanthrene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 13:17
Phenol	U		0.000035	0.00020	mg/L	1	26-Jan-2020 13:17
Pyrene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 13:17
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 13:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>85.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 13:17</i>
<i>Surr: 2-Fluorophenol</i>	<i>66.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 13:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>94.8</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 13:17</i>
<i>Surr: Nitrobenzene-d5</i>	<i>74.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 13:17</i>
<i>Surr: Phenol-d6</i>	<i>71.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 13:17</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000623</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jan-2020 03:13</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59A-20200116  
 Collection Date: 16-Jan-2020 08:05

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:03
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:03
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:03
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:03
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 21:03
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:03
Vinyl chloride	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:03
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:03
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>90.7</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:03</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.6</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:03</i>
<i>Surr: Dibromofluoromethane</i>		<i>98.4</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:03</i>
<i>Surr: Toluene-d8</i>		<i>99.3</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:03</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59A-20200116  
 Collection Date: 16-Jan-2020 08:05

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 13:36
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	26-Jan-2020 13:36
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 13:36
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 13:36
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 13:36
<b>2-Methylnaphthalene</b>	<b>0.00016</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 13:36
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 13:36
<b>Acenaphthene</b>	<b>0.00029</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
Acenaphthylene	U		0.000015	0.00010	mg/L	1	26-Jan-2020 13:36
<b>Anthracene</b>	<b>0.000027</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	26-Jan-2020 13:36
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	26-Jan-2020 13:36
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 13:36
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	26-Jan-2020 13:36
Chrysene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 13:36
<b>Dibenzofuran</b>	<b>0.00016</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	26-Jan-2020 13:36
<b>Fluoranthene</b>	<b>0.000032</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
<b>Fluorene</b>	<b>0.00011</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
<b>Naphthalene</b>	<b>0.0012</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 13:36
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 13:36
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 13:36
<b>Phenanthrene</b>	<b>0.00012</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
Phenol	U		0.000035	0.00020	mg/L	1	26-Jan-2020 13:36
<b>Pyrene</b>	<b>0.000023</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 13:36
<i>Surr: 2,4,6-Tribromophenol</i>	59.7			34-129	%REC	1	26-Jan-2020 13:36
<i>Surr: 2-Fluorobiphenyl</i>	63.2			40-125	%REC	1	26-Jan-2020 13:36
<i>Surr: 2-Fluorophenol</i>	55.1			20-120	%REC	1	26-Jan-2020 13:36
<i>Surr: 4-Terphenyl-d14</i>	85.8			40-135	%REC	1	26-Jan-2020 13:36
<i>Surr: Nitrobenzene-d5</i>	57.0			41-120	%REC	1	26-Jan-2020 13:36
<i>Surr: Phenol-d6</i>	56.8			20-120	%REC	1	26-Jan-2020 13:36
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00368</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 03:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59B-20200116  
 Collection Date: 16-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:28
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:28
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:28
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:28
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 21:28
<b>Toluene</b>	<b>0.18</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	21-Jan-2020 21:28
Vinyl chloride	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:28
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:28
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:28</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.7</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:28</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:28</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:28</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59B-20200116  
 Collection Date: 16-Jan-2020 09:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-14  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	27-Jan-2020 08:31
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	27-Jan-2020 08:31
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	27-Jan-2020 08:31
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	27-Jan-2020 08:31
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	27-Jan-2020 08:31
<b>2-Methylnaphthalene</b>	<b>0.000038</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	27-Jan-2020 08:31
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	27-Jan-2020 08:31
<b>Acenaphthene</b>	<b>0.000051</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
Acenaphthylene	U		0.000015	0.00010	mg/L	1	27-Jan-2020 08:31
<b>Anthracene</b>	<b>0.000031</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	27-Jan-2020 08:31
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	27-Jan-2020 08:31
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	27-Jan-2020 08:31
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	27-Jan-2020 08:31
Chrysene	U		0.000021	0.00010	mg/L	1	27-Jan-2020 08:31
<b>Dibenzofuran</b>	<b>0.000050</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	27-Jan-2020 08:31
<b>Fluoranthene</b>	<b>0.000025</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
<b>Fluorene</b>	<b>0.000042</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
<b>Naphthalene</b>	<b>0.00024</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
Nitrobenzene	U		0.000024	0.00020	mg/L	1	27-Jan-2020 08:31
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	27-Jan-2020 08:31
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	27-Jan-2020 08:31
<b>Phenanthrene</b>	<b>0.000099</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
Phenol	U		0.000035	0.00020	mg/L	1	27-Jan-2020 08:31
<b>Pyrene</b>	<b>0.000021</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 08:31
Surr: 2,4,6-Tribromophenol	71.4			34-129	%REC	1	27-Jan-2020 08:31
Surr: 2-Fluorobiphenyl	74.1			40-125	%REC	1	27-Jan-2020 08:31
Surr: 2-Fluorophenol	65.1			20-120	%REC	1	27-Jan-2020 08:31
Surr: 4-Terphenyl-d14	92.4			40-135	%REC	1	27-Jan-2020 08:31
Surr: Nitrobenzene-d5	68.0			41-120	%REC	1	27-Jan-2020 08:31
Surr: Phenol-d6	68.0			20-120	%REC	1	27-Jan-2020 08:31
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000486</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 03:18

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59D-20200116  
 Collection Date: 16-Jan-2020 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:52
Benzene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:52
Chlorobenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:52
Ethylbenzene	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:52
Methylene chloride	U		0.0010	0.0020	mg/L	1	21-Jan-2020 21:52
Toluene	U		0.00020	0.0010	mg/L	1	21-Jan-2020 21:52
Xylenes, Total	U		0.00030	0.0010	mg/L	1	21-Jan-2020 21:52
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:52</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:52</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:52</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 21:52</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW59D-20200116  
 Collection Date: 16-Jan-2020 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-15  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 14:14
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	26-Jan-2020 14:14
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 14:14
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 14:14
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 14:14
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 14:14
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 14:14
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 14:14
<b>Acenaphthene</b>	<b>0.000036</b>	<b>J</b>	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 14:14</b>
Acenaphthylene	U		0.000015	0.00010	mg/L	1	26-Jan-2020 14:14
Anthracene	U		0.000014	0.00010	mg/L	1	26-Jan-2020 14:14
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	26-Jan-2020 14:14
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	26-Jan-2020 14:14
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 14:14
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	26-Jan-2020 14:14
Chrysene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 14:14
<b>Dibenzofuran</b>	<b>0.000028</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 14:14</b>
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	26-Jan-2020 14:14
Fluoranthene	U		0.000010	0.00010	mg/L	1	26-Jan-2020 14:14
<b>Fluorene</b>	<b>0.000031</b>	<b>J</b>	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 14:14</b>
<b>Naphthalene</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 14:14</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 14:14
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 14:14
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 14:14
<b>Phenanthrene</b>	<b>0.000043</b>	<b>J</b>	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 14:14</b>
Phenol	U		0.000035	0.00020	mg/L	1	26-Jan-2020 14:14
Pyrene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 14:14
<i>Surr: 2,4,6-Tribromophenol</i>	<i>60.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:14</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>67.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:14</i>
<i>Surr: 2-Fluorophenol</i>	<i>49.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:14</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>82.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:14</i>
<i>Surr: Nitrobenzene-d5</i>	<i>59.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:14</i>
<i>Surr: Phenol-d6</i>	<i>52.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:14</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000508</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jan-2020 03:20</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD03-20200116  
 Collection Date: 16-Jan-2020 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:10
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:10
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:10
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:10
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 03:10
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:10
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:10
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:10</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:10</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:10</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:10</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD03-20200116  
 Collection Date: 16-Jan-2020 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-16  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 14:34
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	26-Jan-2020 14:34
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 14:34
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 14:34
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 14:34
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 14:34
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 14:34
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 14:34
Acenaphthene	U		0.000027	0.00010	mg/L	1	26-Jan-2020 14:34
Acenaphthylene	U		0.000015	0.00010	mg/L	1	26-Jan-2020 14:34
Anthracene	U		0.000014	0.00010	mg/L	1	26-Jan-2020 14:34
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	26-Jan-2020 14:34
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	26-Jan-2020 14:34
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 14:34
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	26-Jan-2020 14:34
Chrysene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 14:34
Dibenzofuran	U		0.000020	0.00010	mg/L	1	26-Jan-2020 14:34
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	26-Jan-2020 14:34
Fluoranthene	U		0.000010	0.00010	mg/L	1	26-Jan-2020 14:34
Fluorene	U		0.000030	0.00010	mg/L	1	26-Jan-2020 14:34
Naphthalene	U		0.000020	0.00010	mg/L	1	26-Jan-2020 14:34
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 14:34
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 14:34
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 14:34
Phenanthrene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 14:34
Phenol	U		0.000035	0.00020	mg/L	1	26-Jan-2020 14:34
Pyrene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 14:34
<i>Surr: 2,4,6-Tribromophenol</i>	<i>43.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:34</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:34</i>
<i>Surr: 2-Fluorophenol</i>	<i>46.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:34</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>93.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:34</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:34</i>
<i>Surr: Phenol-d6</i>	<i>54.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 14:34</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000478</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jan-2020 03:22</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW66D-20200116  
 Collection Date: 16-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:35
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:35
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:35
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:35
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 03:35
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:35
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:35
<i>Surr: 1,2-Dichloroethane-d4</i>		91.9		70-126	%REC	1	22-Jan-2020 03:35
<i>Surr: 4-Bromofluorobenzene</i>		94.7		81-113	%REC	1	22-Jan-2020 03:35
<i>Surr: Dibromofluoromethane</i>		97.7		77-123	%REC	1	22-Jan-2020 03:35
<i>Surr: Toluene-d8</i>		98.3		82-127	%REC	1	22-Jan-2020 03:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW66D-20200116  
 Collection Date: 16-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-17  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	27-Jan-2020 08:50
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	27-Jan-2020 08:50
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	27-Jan-2020 08:50
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	27-Jan-2020 08:50
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	27-Jan-2020 08:50
<b>2-Methylnaphthalene</b>	<b>0.000095</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>27-Jan-2020 08:50</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	27-Jan-2020 08:50
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	27-Jan-2020 08:50
Acenaphthene	U		0.000027	0.00010	mg/L	1	27-Jan-2020 08:50
Acenaphthylene	U		0.000015	0.00010	mg/L	1	27-Jan-2020 08:50
Anthracene	U		0.000014	0.00010	mg/L	1	27-Jan-2020 08:50
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	27-Jan-2020 08:50
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	27-Jan-2020 08:50
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	27-Jan-2020 08:50
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000046</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>27-Jan-2020 08:50</b>
Chrysene	U		0.000021	0.00010	mg/L	1	27-Jan-2020 08:50
Dibenzofuran	U		0.000020	0.00010	mg/L	1	27-Jan-2020 08:50
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	27-Jan-2020 08:50
Fluoranthene	U		0.000010	0.00010	mg/L	1	27-Jan-2020 08:50
Fluorene	U		0.000030	0.00010	mg/L	1	27-Jan-2020 08:50
<b>Naphthalene</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>27-Jan-2020 08:50</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	27-Jan-2020 08:50
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	27-Jan-2020 08:50
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	27-Jan-2020 08:50
Phenanthrene	U		0.000021	0.00010	mg/L	1	27-Jan-2020 08:50
Phenol	U		0.000035	0.00020	mg/L	1	27-Jan-2020 08:50
Pyrene	U		0.000019	0.00010	mg/L	1	27-Jan-2020 08:50
<i>Surr: 2,4,6-Tribromophenol</i>	<i>66.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 08:50</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>85.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 08:50</i>
<i>Surr: 2-Fluorophenol</i>	<i>73.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 08:50</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.5</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 08:50</i>
<i>Surr: Nitrobenzene-d5</i>	<i>79.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 08:50</i>
<i>Surr: Phenol-d6</i>	<i>74.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 08:50</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00138</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jan-2020 03:24</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW65D-20200116  
 Collection Date: 16-Jan-2020 16:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:57
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:57
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:57
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:57
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 01:57
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:57
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:57
<i>Surr: 1,2-Dichloroethane-d4</i>		90.4		70-126	%REC	1	22-Jan-2020 01:57
<i>Surr: 4-Bromofluorobenzene</i>		93.4		81-113	%REC	1	22-Jan-2020 01:57
<i>Surr: Dibromofluoromethane</i>		96.5		77-123	%REC	1	22-Jan-2020 01:57
<i>Surr: Toluene-d8</i>		99.4		82-127	%REC	1	22-Jan-2020 01:57

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW65D-20200116  
 Collection Date: 16-Jan-2020 16:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-18  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 09:40
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jan-2020 09:40
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 09:40
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 09:40
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 09:40
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jan-2020 09:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 09:40
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 09:40
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jan-2020 09:40
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jan-2020 09:40
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jan-2020 09:40
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jan-2020 09:40
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jan-2020 09:40
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 09:40
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00026</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 09:40
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jan-2020 09:40
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jan-2020 09:40
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jan-2020 09:40
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jan-2020 09:40
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jan-2020 09:40
Naphthalene	U		0.000020	0.00010	mg/L	1	23-Jan-2020 09:40
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 09:40
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 09:40
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 09:40
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jan-2020 09:40
Phenol	U		0.000035	0.00020	mg/L	1	23-Jan-2020 09:40
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jan-2020 09:40
<i>Surr: 2,4,6-Tribromophenol</i>	<i>111</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 09:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>110</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 09:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 09:40</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>123</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 09:40</i>
<i>Surr: Nitrobenzene-d5</i>	<i>89.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 09:40</i>
<i>Surr: Phenol-d6</i>	<i>87.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 09:40</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 21-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.000507</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jan-2020 00:56</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36D-20200116  
 Collection Date: 16-Jan-2020 17:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:59
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:59
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:59
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:59
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 03:59
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 03:59
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 03:59
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:59</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>93.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:59</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:59</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 03:59</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW36D-20200116  
 Collection Date: 16-Jan-2020 17:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-19  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 15:12
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	26-Jan-2020 15:12
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 15:12
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 15:12
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 15:12
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 15:12
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 15:12
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 15:12
Acenaphthene	U		0.000027	0.00010	mg/L	1	26-Jan-2020 15:12
Acenaphthylene	U		0.000015	0.00010	mg/L	1	26-Jan-2020 15:12
<b>Anthracene</b>	<b>0.000033</b>	<b>J</b>	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 15:12</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	26-Jan-2020 15:12
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	26-Jan-2020 15:12
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 15:12
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	26-Jan-2020 15:12
Chrysene	U		0.000021	0.00010	mg/L	1	26-Jan-2020 15:12
Dibenzofuran	U		0.000020	0.00010	mg/L	1	26-Jan-2020 15:12
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	26-Jan-2020 15:12
Fluoranthene	U		0.000010	0.00010	mg/L	1	26-Jan-2020 15:12
Fluorene	U		0.000030	0.00010	mg/L	1	26-Jan-2020 15:12
<b>Naphthalene</b>	<b>0.000033</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 15:12</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 15:12
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 15:12
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 15:12
<b>Phenanthrene</b>	<b>0.000047</b>	<b>J</b>	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>26-Jan-2020 15:12</b>
Phenol	U		0.000035	0.00020	mg/L	1	26-Jan-2020 15:12
Pyrene	U		0.000019	0.00010	mg/L	1	26-Jan-2020 15:12
<i>Surr: 2,4,6-Tribromophenol</i>	<i>60.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 15:12</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>80.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 15:12</i>
<i>Surr: 2-Fluorophenol</i>	<i>71.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 15:12</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>88.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 15:12</i>
<i>Surr: Nitrobenzene-d5</i>	<i>76.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 15:12</i>
<i>Surr: Phenol-d6</i>	<i>68.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>26-Jan-2020 15:12</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.000533</b>	<b>J</b>	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>24-Jan-2020 03:27</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW86C-20200117  
 Collection Date: 17-Jan-2020 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 04:24
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 04:24
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 04:24
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 04:24
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 04:24
<b>Toluene</b>	<b>0.0039</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 04:24
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 04:24
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>91.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:24</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:24</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:24</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:24</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW86C-20200117  
 Collection Date: 17-Jan-2020 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:00
<b>2,4-Dimethylphenol</b>	<b>0.00018</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 18:00
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 18:00
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 18:00
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:00
<b>2-Methylnaphthalene</b>	<b>0.000061</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:00
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 18:00
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 18:00
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jan-2020 18:00
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jan-2020 18:00
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jan-2020 18:00
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jan-2020 18:00
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jan-2020 18:00
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 18:00
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	23-Jan-2020 18:00
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jan-2020 18:00
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jan-2020 18:00
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jan-2020 18:00
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jan-2020 18:00
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jan-2020 18:00
<b>Naphthalene</b>	<b>0.00075</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:00
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 18:00
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 18:00
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 18:00
Phenanthrene	U		0.000021	0.00010	mg/L	1	23-Jan-2020 18:00
Phenol	U		0.000035	0.00020	mg/L	1	23-Jan-2020 18:00
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jan-2020 18:00
<i>Surr: 2,4,6-Tribromophenol</i>	<i>86.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:00</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:00</i>
<i>Surr: 2-Fluorophenol</i>	<i>53.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:00</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>107</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:00</i>
<i>Surr: Nitrobenzene-d5</i>	<i>64.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:00</i>
<i>Surr: Phenol-d6</i>	<i>59.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW86C-20200117  
 Collection Date: 17-Jan-2020 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-20  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 20-Jan-2020		Analyst: MBG	
nC6 to nC12	U		0.19	0.48	mg/L	1	21-Jan-2020 05:19
>nC12 to nC28	U		0.19	0.48	mg/L	1	21-Jan-2020 05:19
>nC28 to nC35	U		0.19	0.48	mg/L	1	21-Jan-2020 05:19
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	21-Jan-2020 05:19
Surr: 2-Fluorobiphenyl	109			70-130	%REC	1	21-Jan-2020 05:19
Surr: Trifluoromethyl benzene	114			70-130	%REC	1	21-Jan-2020 05:19
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	0.000645	J	0.000400	0.00200	mg/L	1	24-Jan-2020 03:33

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD04-20200117  
 Collection Date: 17-Jan-2020 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 04:48
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 04:48
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 04:48
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 04:48
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 04:48
<b>Toluene</b>	<b>0.0045</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 04:48
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 04:48
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:48</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>93.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:48</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:48</i>
<i>Surr: Toluene-d8</i>	<i>99.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 04:48</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD04-20200117  
 Collection Date: 17-Jan-2020 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:19
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jan-2020 18:19
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 18:19
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 18:19
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:19
<b>2-Methylnaphthalene</b>	<b>0.000066</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jan-2020 18:19</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 18:19
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 18:19
Acenaphthene	U		0.000027	0.00010	mg/L	1	23-Jan-2020 18:19
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jan-2020 18:19
<b>Anthracene</b>	<b>0.00012</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jan-2020 18:19</b>
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jan-2020 18:19
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jan-2020 18:19
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 18:19
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000053</b>	<b>J</b>	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jan-2020 18:19</b>
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jan-2020 18:19
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jan-2020 18:19
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jan-2020 18:19
Fluoranthene	U		0.000010	0.00010	mg/L	1	23-Jan-2020 18:19
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jan-2020 18:19
<b>Naphthalene</b>	<b>0.00022</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jan-2020 18:19</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 18:19
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 18:19
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 18:19
<b>Phenanthrene</b>	<b>0.00036</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>23-Jan-2020 18:19</b>
Phenol	U		0.000035	0.00020	mg/L	1	23-Jan-2020 18:19
Pyrene	U		0.000019	0.00010	mg/L	1	23-Jan-2020 18:19
<i>Surr: 2,4,6-Tribromophenol</i>	<i>78.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:19</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:19</i>
<i>Surr: 2-Fluorophenol</i>	<i>61.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:19</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>101</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:19</i>
<i>Surr: Nitrobenzene-d5</i>	<i>63.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:19</i>
<i>Surr: Phenol-d6</i>	<i>58.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:19</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD04-20200117  
 Collection Date: 17-Jan-2020 07:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-21  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 20-Jan-2020		Analyst: MBG	
nC6 to nC12	U		0.19	0.48	mg/L	1	21-Jan-2020 05:48
>nC12 to nC28	U		0.19	0.48	mg/L	1	21-Jan-2020 05:48
>nC28 to nC35	U		0.19	0.48	mg/L	1	21-Jan-2020 05:48
Total Petroleum Hydrocarbon	U		0.19	0.48	mg/L	1	21-Jan-2020 05:48
Surr: 2-Fluorobiphenyl	106			70-130	%REC	1	21-Jan-2020 05:48
Surr: Trifluoromethyl benzene	113			70-130	%REC	1	21-Jan-2020 05:48
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	0.000699	J	0.000400	0.00200	mg/L	1	24-Jan-2020 03:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW75B-20200117  
 Collection Date: 17-Jan-2020 08:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 05:13
<b>Benzene</b>	<b>0.12</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:13
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 05:13
<b>Ethylbenzene</b>	<b>0.037</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:13
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 05:13
<b>Toluene</b>	<b>0.13</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:13
<b>Xylenes, Total</b>	<b>0.11</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:13
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>90.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:13</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>100</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:13</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.4</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:13</i>
<i>Surr: Toluene-d8</i>	<i>97.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW75B-20200117  
 Collection Date: 17-Jan-2020 08:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:38
<b>2,4-Dimethylphenol</b>	<b>0.86</b>		<b>0.0080</b>	<b>0.040</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 18:38
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 18:38
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:38
<b>2-Methylnaphthalene</b>	<b>0.81</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 18:38
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 18:38
<b>Acenaphthene</b>	<b>0.67</b>		<b>0.0054</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<b>Acenaphthylene</b>	<b>0.0086</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:38
<b>Anthracene</b>	<b>0.39</b>		<b>0.0028</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<b>Benz(a)anthracene</b>	<b>0.12</b>		<b>0.0010</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 09:28
<b>Benzo(a)pyrene</b>	<b>0.035</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 09:28
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 18:38
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0021</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 18:38
<b>Chrysene</b>	<b>0.095</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 09:28
<b>Dibenzofuran</b>	<b>0.54</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<b>Di-n-butyl phthalate</b>	<b>0.0016</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 18:38
<b>Fluoranthene</b>	<b>0.75</b>		<b>0.0020</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<b>Fluorene</b>	<b>0.61</b>		<b>0.0060</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<b>Naphthalene</b>	<b>7.5</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	25-Jan-2020 21:55
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 18:38
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 18:38
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 18:38
<b>Phenanthrene</b>	<b>1.9</b>		<b>0.0042</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<b>Phenol</b>	<b>0.0029</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 18:38
<b>Pyrene</b>	<b>0.51</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 09:47
<i>Surr: 2,4,6-Tribromophenol</i>	<i>64.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:38</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>25-Jan-2020 21:55</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>109</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 09:28</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>27-Jan-2020 09:47</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>120</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 09:28</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>27-Jan-2020 09:47</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>25-Jan-2020 21:55</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>61.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>95.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:38</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>25-Jan-2020 21:55</i>
<i>Surr: 2-Fluorophenol</i>	<i>111</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 09:28</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>27-Jan-2020 09:47</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW75B-20200117  
 Collection Date: 17-Jan-2020 08:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-22  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	111			40-135	%REC	20	27-Jan-2020 09:28
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	27-Jan-2020 09:47
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	25-Jan-2020 21:55
Surr: 4-Terphenyl-d14	115			40-135	%REC	1	23-Jan-2020 18:38
Surr: Nitrobenzene-d5	72.8			41-120	%REC	1	23-Jan-2020 18:38
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	25-Jan-2020 21:55
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	27-Jan-2020 09:47
Surr: Nitrobenzene-d5	100			41-120	%REC	20	27-Jan-2020 09:28
Surr: Phenol-d6	121	S		20-120	%REC	20	27-Jan-2020 09:28
Surr: Phenol-d6	0	JS		20-120	%REC	200	27-Jan-2020 09:47
Surr: Phenol-d6	0	JS		20-120	%REC	2000	25-Jan-2020 21:55
Surr: Phenol-d6	91.1			20-120	%REC	1	23-Jan-2020 18:38
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 20-Jan-2020		Analyst: MBG	
nC6 to nC12	13		0.19	0.49	mg/L	1	21-Jan-2020 06:16
>nC12 to nC28	21		0.19	0.49	mg/L	1	21-Jan-2020 06:16
>nC28 to nC35	0.91		0.19	0.49	mg/L	1	21-Jan-2020 06:16
Total Petroleum Hydrocarbon	34.9		0.19	0.49	mg/L	1	21-Jan-2020 06:16
Surr: 2-Fluorobiphenyl	111			70-130	%REC	1	21-Jan-2020 06:16
Surr: Trifluoromethyl benzene	111			70-130	%REC	1	21-Jan-2020 06:16
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	0.00167	J	0.000400	0.00200	mg/L	1	24-Jan-2020 03:38

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW79A-20200117  
 Collection Date: 17-Jan-2020 09:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 05:37
<b>Benzene</b>	<b>0.12</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 05:37
<b>Ethylbenzene</b>	<b>0.075</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 05:37
<b>Toluene</b>	<b>0.26</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	5	22-Jan-2020 16:23
<b>Xylenes, Total</b>	<b>0.20</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 05:37
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>86.8</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:37</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>107</i>			<i>70-126</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:23</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:37</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:23</i>
<i>Surr: Dibromofluoromethane</i>	<i>95.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:37</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.8</i>			<i>77-123</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:23</i>
<i>Surr: Toluene-d8</i>	<i>98.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 05:37</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:23</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW79A-20200117  
 Collection Date: 17-Jan-2020 09:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:57
<b>2,4-Dimethylphenol</b>	<b>6.7</b>		<b>0.080</b>	<b>0.40</b>	<b>mg/L</b>	2000	26-Jan-2020 18:43
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 18:57
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 18:57
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 18:57
<b>2-Methylnaphthalene</b>	<b>0.43</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	27-Jan-2020 10:24
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 18:57
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 18:57
<b>Acenaphthene</b>	<b>0.19</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 10:06
<b>Acenaphthylene</b>	<b>0.0036</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<b>Anthracene</b>	<b>0.0094</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<b>Benz(a)anthracene</b>	<b>0.00081</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<b>Benzo(a)pyrene</b>	<b>0.00029</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 18:57
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	23-Jan-2020 18:57
<b>Chrysene</b>	<b>0.00072</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<b>Dibenzofuran</b>	<b>0.14</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 10:06
<b>Di-n-butyl phthalate</b>	<b>0.00098</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<b>Fluoranthene</b>	<b>0.0076</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<b>Fluorene</b>	<b>0.094</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 10:06
<b>Naphthalene</b>	<b>9.0</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	26-Jan-2020 18:43
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 18:57
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 18:57
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 18:57
<b>Phenanthrene</b>	<b>0.081</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	27-Jan-2020 10:06
<b>Phenol</b>	<b>0.35</b>		<b>0.0070</b>	<b>0.040</b>	<b>mg/L</b>	200	27-Jan-2020 10:24
<b>Pyrene</b>	<b>0.0035</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 18:57
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:57</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>27-Jan-2020 10:24</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>26-Jan-2020 18:43</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>110</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 10:06</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>114</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 10:06</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>26-Jan-2020 18:43</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>27-Jan-2020 10:24</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>103</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>105</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>23-Jan-2020 18:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>27-Jan-2020 10:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>26-Jan-2020 18:43</i>
<i>Surr: 2-Fluorophenol</i>	<i>114</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>27-Jan-2020 10:06</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW79A-20200117  
 Collection Date: 17-Jan-2020 09:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-23  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	101			40-135	%REC	20	27-Jan-2020 10:06
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	26-Jan-2020 18:43
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	27-Jan-2020 10:24
Surr: 4-Terphenyl-d14	105			40-135	%REC	1	23-Jan-2020 18:57
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	26-Jan-2020 18:43
Surr: Nitrobenzene-d5	91.5			41-120	%REC	1	23-Jan-2020 18:57
Surr: Nitrobenzene-d5	72.4	J		41-120	%REC	20	27-Jan-2020 10:06
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	27-Jan-2020 10:24
Surr: Phenol-d6	86.6			20-120	%REC	20	27-Jan-2020 10:06
Surr: Phenol-d6	0	JS		20-120	%REC	200	27-Jan-2020 10:24
Surr: Phenol-d6	84.0			20-120	%REC	1	23-Jan-2020 18:57
Surr: Phenol-d6	0	JS		20-120	%REC	2000	26-Jan-2020 18:43
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 20-Jan-2020		Analyst: MBG	
nC6 to nC12	21		0.19	0.47	mg/L	1	21-Jan-2020 06:45
>nC12 to nC28	16		0.19	0.47	mg/L	1	21-Jan-2020 06:45
>nC28 to nC35	U		0.19	0.47	mg/L	1	21-Jan-2020 06:45
<b>Total Petroleum Hydrocarbon</b>	<b>37.0</b>		<b>0.19</b>	<b>0.47</b>	<b>mg/L</b>	<b>1</b>	<b>21-Jan-2020 06:45</b>
Surr: 2-Fluorobiphenyl	117			70-130	%REC	1	21-Jan-2020 06:45
Surr: Trifluoromethyl benzene	113			70-130	%REC	1	21-Jan-2020 06:45
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	0.00893		0.000400	0.00200	mg/L	1	24-Jan-2020 03:40

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW74B-20200117  
 Collection Date: 17-Jan-2020 10:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 06:01
<b>Benzene</b>	<b>0.12</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:01
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 06:01
<b>Ethylbenzene</b>	<b>0.076</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:01
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 06:01
<b>Toluene</b>	<b>0.29</b>		<b>0.0010</b>	<b>0.0050</b>	<b>mg/L</b>	5	22-Jan-2020 16:50
<b>Xylenes, Total</b>	<b>0.20</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:01
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>88.4</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 06:01</i>
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>103</i>			<i>70-126</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:50</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 06:01</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.3</i>			<i>81-113</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:50</i>
<i>Surr: Dibromofluoromethane</i>	<i>94.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 06:01</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:50</i>
<i>Surr: Toluene-d8</i>	<i>97.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>22-Jan-2020 06:01</i>
<i>Surr: Toluene-d8</i>	<i>99.9</i>			<i>82-127</i>	<i>%REC</i>	<i>5</i>	<i>22-Jan-2020 16:50</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW74B-20200117  
 Collection Date: 17-Jan-2020 10:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 19:17
<b>2,4-Dimethylphenol</b>	<b>0.21</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 19:17
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 19:17
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 19:17
<b>2-Methylnaphthalene</b>	<b>1.3</b>		<b>0.019</b>	<b>0.10</b>	<b>mg/L</b>	1000	26-Jan-2020 19:40
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 19:17
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 19:17
<b>Acenaphthene</b>	<b>0.98</b>		<b>0.0027</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<b>Acenaphthylene</b>	<b>0.018</b>		<b>0.00015</b>	<b>0.0010</b>	<b>mg/L</b>	10	27-Jan-2020 10:43
<b>Anthracene</b>	<b>0.42</b>		<b>0.0014</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<b>Benz(a)anthracene</b>	<b>0.13</b>		<b>0.0050</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<b>Benzo(a)pyrene</b>	<b>0.034</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	27-Jan-2020 10:43
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 19:17
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0014</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 19:17
<b>Chrysene</b>	<b>0.082</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	27-Jan-2020 10:43
<b>Dibenzofuran</b>	<b>0.72</b>		<b>0.0020</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<b>Di-n-butyl phthalate</b>	<b>0.0022</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 19:17
<b>Fluoranthene</b>	<b>0.67</b>		<b>0.0010</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<b>Fluorene</b>	<b>0.88</b>		<b>0.0030</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<b>Naphthalene</b>	<b>8.5</b>		<b>0.020</b>	<b>0.10</b>	<b>mg/L</b>	1000	26-Jan-2020 19:40
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 19:17
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 19:17
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 19:17
<b>Phenanthrene</b>	<b>2.3</b>		<b>0.021</b>	<b>0.10</b>	<b>mg/L</b>	1000	26-Jan-2020 19:40
<b>Phenol</b>	<b>0.048</b>		<b>0.00035</b>	<b>0.0020</b>	<b>mg/L</b>	10	27-Jan-2020 10:43
<b>Pyrene</b>	<b>0.44</b>		<b>0.0019</b>	<b>0.010</b>	<b>mg/L</b>	100	27-Jan-2020 11:02
<i>Surr: 2,4,6-Tribromophenol</i>	78.6			34-129	%REC	1	23-Jan-2020 19:17
<i>Surr: 2,4,6-Tribromophenol</i>	86.1			34-129	%REC	10	27-Jan-2020 10:43
<i>Surr: 2,4,6-Tribromophenol</i>	115	J		34-129	%REC	100	27-Jan-2020 11:02
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	1000	26-Jan-2020 19:40
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	1000	26-Jan-2020 19:40
<i>Surr: 2-Fluorobiphenyl</i>	94.5			40-125	%REC	10	27-Jan-2020 10:43
<i>Surr: 2-Fluorobiphenyl</i>	120	J		40-125	%REC	100	27-Jan-2020 11:02
<i>Surr: 2-Fluorobiphenyl</i>	64.4			40-125	%REC	1	23-Jan-2020 19:17
<i>Surr: 2-Fluorophenol</i>	70.1			20-120	%REC	1	23-Jan-2020 19:17
<i>Surr: 2-Fluorophenol</i>	105			20-120	%REC	10	27-Jan-2020 10:43
<i>Surr: 2-Fluorophenol</i>	110	J		20-120	%REC	100	27-Jan-2020 11:02
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	1000	26-Jan-2020 19:40

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW74B-20200117  
 Collection Date: 17-Jan-2020 10:20

**ANALYTICAL REPORT**

WorkOrder:HS20010767  
 Lab ID:HS20010767-24  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	26-Jan-2020 19:40
Surr: 4-Terphenyl-d14	115			40-135	%REC	1	23-Jan-2020 19:17
Surr: 4-Terphenyl-d14	104			40-135	%REC	10	27-Jan-2020 10:43
Surr: 4-Terphenyl-d14	87.9	J		40-135	%REC	100	27-Jan-2020 11:02
Surr: Nitrobenzene-d5	95.5			41-120	%REC	10	27-Jan-2020 10:43
Surr: Nitrobenzene-d5	108	J		41-120	%REC	100	27-Jan-2020 11:02
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	26-Jan-2020 19:40
Surr: Nitrobenzene-d5	118			41-120	%REC	1	23-Jan-2020 19:17
Surr: Phenol-d6	62.2			20-120	%REC	1	23-Jan-2020 19:17
Surr: Phenol-d6	0	JS		20-120	%REC	1000	26-Jan-2020 19:40
Surr: Phenol-d6	101			20-120	%REC	10	27-Jan-2020 10:43
Surr: Phenol-d6	111	J		20-120	%REC	100	27-Jan-2020 11:02
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 20-Jan-2020		Analyst: MBG	
nC6 to nC12	12		0.19	0.48	mg/L	1	21-Jan-2020 07:14
>nC12 to nC28	13		0.19	0.48	mg/L	1	21-Jan-2020 07:14
>nC28 to nC35	0.47	J	0.19	0.48	mg/L	1	21-Jan-2020 07:14
Total Petroleum Hydrocarbon	25.5		0.19	0.48	mg/L	1	21-Jan-2020 07:14
Surr: 2-Fluorobiphenyl	113			70-130	%REC	1	21-Jan-2020 07:14
Surr: Trifluoromethyl benzene	112			70-130	%REC	1	21-Jan-2020 07:14
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	0.000443	J	0.000400	0.00200	mg/L	1	24-Jan-2020 03:42

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW78A-20200117  
 Collection Date: 17-Jan-2020 11:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 06:26
<b>Benzene</b>	<b>0.018</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:26
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 06:26
<b>Ethylbenzene</b>	<b>0.011</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:26
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 06:26
<b>Toluene</b>	<b>0.033</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:26
<b>Xylenes, Total</b>	<b>0.028</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	22-Jan-2020 06:26
Surr: 1,2-Dichloroethane-d4	91.1			70-126	%REC	1	22-Jan-2020 06:26
Surr: 4-Bromofluorobenzene	100.0			81-113	%REC	1	22-Jan-2020 06:26
Surr: Dibromofluoromethane	96.6			77-123	%REC	1	22-Jan-2020 06:26
Surr: Toluene-d8	99.8			82-127	%REC	1	22-Jan-2020 06:26

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW78A-20200117  
 Collection Date: 17-Jan-2020 11:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 15:31
<b>2,4-Dimethylphenol</b>	<b>2.5</b>		<b>0.016</b>	<b>0.080</b>	<b>mg/L</b>	400	28-Jan-2020 09:03
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 15:31
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 15:31
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 15:31
<b>2-Methylnaphthalene</b>	<b>0.13</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	28-Jan-2020 08:25
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 15:31
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 15:31
<b>Acenaphthene</b>	<b>0.071</b>		<b>0.00054</b>	<b>0.0020</b>	<b>mg/L</b>	20	28-Jan-2020 08:25
<b>Acenaphthylene</b>	<b>0.0015</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Anthracene</b>	<b>0.0040</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Benz(a)anthracene</b>	<b>0.00043</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Benzo(a)pyrene</b>	<b>0.00014</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 15:31
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00092</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Chrysene</b>	<b>0.00039</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Dibenzofuran</b>	<b>0.055</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	20	28-Jan-2020 08:25
<b>Di-n-butyl phthalate</b>	<b>0.0016</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Fluoranthene</b>	<b>0.0029</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
<b>Fluorene</b>	<b>0.040</b>		<b>0.00060</b>	<b>0.0020</b>	<b>mg/L</b>	20	28-Jan-2020 08:25
<b>Naphthalene</b>	<b>3.1</b>		<b>0.0080</b>	<b>0.040</b>	<b>mg/L</b>	400	28-Jan-2020 09:03
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 15:31
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 15:31
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 15:31
<b>Phenanthrene</b>	<b>0.031</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	28-Jan-2020 08:25
<b>Phenol</b>	<b>0.13</b>		<b>0.00070</b>	<b>0.0040</b>	<b>mg/L</b>	20	28-Jan-2020 08:25
<b>Pyrene</b>	<b>0.0018</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:31
Surr: 2,4,6-Tribromophenol	96.7			34-129	%REC	20	28-Jan-2020 08:25
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	400	28-Jan-2020 09:03
Surr: 2,4,6-Tribromophenol	68.2			34-129	%REC	1	26-Jan-2020 15:31
Surr: 2-Fluorobiphenyl	70.7			40-125	%REC	1	26-Jan-2020 15:31
Surr: 2-Fluorobiphenyl	115			40-125	%REC	20	28-Jan-2020 08:25
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	400	28-Jan-2020 09:03
Surr: 2-Fluorophenol	87.6			20-120	%REC	20	28-Jan-2020 08:25
Surr: 2-Fluorophenol	0	JS		20-120	%REC	400	28-Jan-2020 09:03
Surr: 2-Fluorophenol	59.4			20-120	%REC	1	26-Jan-2020 15:31
Surr: 4-Terphenyl-d14	86.7			40-135	%REC	1	26-Jan-2020 15:31
Surr: 4-Terphenyl-d14	110			40-135	%REC	20	28-Jan-2020 08:25
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	400	28-Jan-2020 09:03

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW78A-20200117  
 Collection Date: 17-Jan-2020 11:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-25  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
Surr: Nitrobenzene-d5	66.7			41-120	%REC	1	26-Jan-2020 15:31
Surr: Nitrobenzene-d5	93.6			41-120	%REC	20	28-Jan-2020 08:25
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	400	28-Jan-2020 09:03
Surr: Phenol-d6	102			20-120	%REC	20	28-Jan-2020 08:25
Surr: Phenol-d6	0	JS		20-120	%REC	400	28-Jan-2020 09:03
Surr: Phenol-d6	71.1			20-120	%REC	1	26-Jan-2020 15:31
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 20-Jan-2020		Analyst: MBG	
nC6 to nC12	2.7		0.19	0.48	mg/L	1	21-Jan-2020 07:42
>nC12 to nC28	1.6		0.19	0.48	mg/L	1	21-Jan-2020 07:42
>nC28 to nC35	U		0.19	0.48	mg/L	1	21-Jan-2020 07:42
<b>Total Petroleum Hydrocarbon</b>	<b>4.30</b>		<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	21-Jan-2020 07:42
Surr: 2-Fluorobiphenyl	108			70-130	%REC	1	21-Jan-2020 07:42
Surr: Trifluoromethyl benzene	107			70-130	%REC	1	21-Jan-2020 07:42
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
Arsenic	0.00991		0.000400	0.00200	mg/L	1	24-Jan-2020 03:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB05-2020117  
 Collection Date: 17-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-26  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:32
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:32
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:32
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:32
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 01:32
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 01:32
Vinyl chloride	U		0.00020	0.0010	mg/L	1	22-Jan-2020 14:15
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 01:32
<i>Surr: 1,2-Dichloroethane-d4</i>		90.8		70-126	%REC	1	22-Jan-2020 01:32
<i>Surr: 1,2-Dichloroethane-d4</i>		103		70-126	%REC	1	22-Jan-2020 14:15
<i>Surr: 4-Bromofluorobenzene</i>		93.3		81-113	%REC	1	22-Jan-2020 01:32
<i>Surr: 4-Bromofluorobenzene</i>		100		81-113	%REC	1	22-Jan-2020 14:15
<i>Surr: Dibromofluoromethane</i>		97.6		77-123	%REC	1	22-Jan-2020 01:32
<i>Surr: Dibromofluoromethane</i>		100		77-123	%REC	1	22-Jan-2020 14:15
<i>Surr: Toluene-d8</i>		99.2		82-127	%REC	1	22-Jan-2020 01:32
<i>Surr: Toluene-d8</i>		99.1		82-127	%REC	1	22-Jan-2020 14:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW48C-20200116  
 Collection Date: 16-Jan-2020 10:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-27  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 06:50
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 06:50
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 06:50
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 06:50
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 06:50
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 06:50
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 06:50
<i>Surr: 1,2-Dichloroethane-d4</i>		89.4		70-126	%REC	1	22-Jan-2020 06:50
<i>Surr: 4-Bromofluorobenzene</i>		96.9		81-113	%REC	1	22-Jan-2020 06:50
<i>Surr: Dibromofluoromethane</i>		95.4		77-123	%REC	1	22-Jan-2020 06:50
<i>Surr: Toluene-d8</i>		98.7		82-127	%REC	1	22-Jan-2020 06:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW48C-20200116  
 Collection Date: 16-Jan-2020 10:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-27  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	26-Jan-2020 15:50
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	26-Jan-2020 15:50
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	26-Jan-2020 15:50
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	26-Jan-2020 15:50
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	26-Jan-2020 15:50
<b>2-Methylnaphthalene</b>	<b>0.000076</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	26-Jan-2020 15:50
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	26-Jan-2020 15:50
<b>Acenaphthene</b>	<b>0.000073</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
Acenaphthylene	U		0.000015	0.00010	mg/L	1	26-Jan-2020 15:50
Anthracene	U		0.000014	0.00010	mg/L	1	26-Jan-2020 15:50
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	26-Jan-2020 15:50
<b>Benzo(a)pyrene</b>	<b>0.000021</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	26-Jan-2020 15:50
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000039</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
<b>Chrysene</b>	<b>0.000029</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
<b>Dibenzofuran</b>	<b>0.000054</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	26-Jan-2020 15:50
<b>Fluoranthene</b>	<b>0.000061</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
<b>Fluorene</b>	<b>0.000050</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
<b>Naphthalene</b>	<b>0.00052</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
Nitrobenzene	U		0.000024	0.00020	mg/L	1	26-Jan-2020 15:50
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	26-Jan-2020 15:50
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	26-Jan-2020 15:50
<b>Phenanthrene</b>	<b>0.000052</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
Phenol	U		0.000035	0.00020	mg/L	1	26-Jan-2020 15:50
<b>Pyrene</b>	<b>0.000065</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	26-Jan-2020 15:50
<i>Surr: 2,4,6-Tribromophenol</i>	54.9			34-129	%REC	1	26-Jan-2020 15:50
<i>Surr: 2-Fluorobiphenyl</i>	73.6			40-125	%REC	1	26-Jan-2020 15:50
<i>Surr: 2-Fluorophenol</i>	54.3			20-120	%REC	1	26-Jan-2020 15:50
<i>Surr: 4-Terphenyl-d14</i>	89.8			40-135	%REC	1	26-Jan-2020 15:50
<i>Surr: Nitrobenzene-d5</i>	66.1			41-120	%REC	1	26-Jan-2020 15:50
<i>Surr: Phenol-d6</i>	64.4			20-120	%REC	1	26-Jan-2020 15:50
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00126</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 03:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW47C-20200116  
 Collection Date: 16-Jan-2020 11:05

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-28  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	22-Jan-2020 07:15
Benzene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 07:15
Chlorobenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 07:15
Ethylbenzene	U		0.00030	0.0010	mg/L	1	22-Jan-2020 07:15
Methylene chloride	U		0.0010	0.0020	mg/L	1	22-Jan-2020 07:15
Toluene	U		0.00020	0.0010	mg/L	1	22-Jan-2020 07:15
Xylenes, Total	U		0.00030	0.0010	mg/L	1	22-Jan-2020 07:15
<i>Surr: 1,2-Dichloroethane-d4</i>		89.5		70-126	%REC	1	22-Jan-2020 07:15
<i>Surr: 4-Bromofluorobenzene</i>		95.1		81-113	%REC	1	22-Jan-2020 07:15
<i>Surr: Dibromofluoromethane</i>		95.7		77-123	%REC	1	22-Jan-2020 07:15
<i>Surr: Toluene-d8</i>		99.5		82-127	%REC	1	22-Jan-2020 07:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW47C-20200116  
 Collection Date: 16-Jan-2020 11:05

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-28  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	27-Jan-2020 09:09
<b>2,4-Dimethylphenol</b>	<b>0.00020</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	27-Jan-2020 09:09
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	27-Jan-2020 09:09
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	27-Jan-2020 09:09
<b>2-Methylnaphthalene</b>	<b>0.00010</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	27-Jan-2020 09:09
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	27-Jan-2020 09:09
<b>Acenaphthene</b>	<b>0.000057</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
Acenaphthylene		U	0.000015	0.00010	mg/L	1	27-Jan-2020 09:09
Anthracene		U	0.000014	0.00010	mg/L	1	27-Jan-2020 09:09
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	27-Jan-2020 09:09
<b>Benzo(a)pyrene</b>	<b>0.000040</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	27-Jan-2020 09:09
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00079</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<b>Chrysene</b>	<b>0.000060</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<b>Dibenzofuran</b>	<b>0.000049</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<b>Di-n-butyl phthalate</b>	<b>0.0014</b>		<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<b>Fluoranthene</b>	<b>0.00011</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<b>Fluorene</b>	<b>0.000042</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<b>Naphthalene</b>	<b>0.00056</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
Nitrobenzene		U	0.000024	0.00020	mg/L	1	27-Jan-2020 09:09
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	27-Jan-2020 09:09
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	27-Jan-2020 09:09
<b>Phenanthrene</b>	<b>0.000088</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
Phenol		U	0.000035	0.00020	mg/L	1	27-Jan-2020 09:09
<b>Pyrene</b>	<b>0.00010</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	27-Jan-2020 09:09
<i>Surr: 2,4,6-Tribromophenol</i>	<i>69.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 09:09</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>72.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 09:09</i>
<i>Surr: 2-Fluorophenol</i>	<i>62.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 09:09</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 09:09</i>
<i>Surr: Nitrobenzene-d5</i>	<i>67.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 09:09</i>
<i>Surr: Phenol-d6</i>	<i>67.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>27-Jan-2020 09:09</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00234</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 03:49

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW61A-20200116  
 Collection Date: 16-Jan-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-29  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:37
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 12:37
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 12:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 12:37
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:37
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:37
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 12:37
<i>Surr: 1,2-Dichloroethane-d4</i>		99.1		70-126	%REC	1	25-Jan-2020 12:37
<i>Surr: 4-Bromofluorobenzene</i>		95.5		81-113	%REC	1	25-Jan-2020 12:37
<i>Surr: Dibromofluoromethane</i>		102		77-123	%REC	1	25-Jan-2020 12:37
<i>Surr: Toluene-d8</i>		102		82-127	%REC	1	25-Jan-2020 12:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW61A-20200116  
 Collection Date: 16-Jan-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010767  
 Lab ID:HS20010767-29  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 22-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	23-Jan-2020 10:37
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	23-Jan-2020 10:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	23-Jan-2020 10:37
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	23-Jan-2020 10:37
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	23-Jan-2020 10:37
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	23-Jan-2020 10:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	23-Jan-2020 10:37
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	23-Jan-2020 10:37
<b>Acenaphthene</b>	<b>0.000052</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 10:37
Acenaphthylene	U		0.000015	0.00010	mg/L	1	23-Jan-2020 10:37
Anthracene	U		0.000014	0.00010	mg/L	1	23-Jan-2020 10:37
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	23-Jan-2020 10:37
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	23-Jan-2020 10:37
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	23-Jan-2020 10:37
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000081</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	23-Jan-2020 10:37
Chrysene	U		0.000021	0.00010	mg/L	1	23-Jan-2020 10:37
Dibenzofuran	U		0.000020	0.00010	mg/L	1	23-Jan-2020 10:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	23-Jan-2020 10:37
<b>Fluoranthene</b>	<b>0.000060</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 10:37
Fluorene	U		0.000030	0.00010	mg/L	1	23-Jan-2020 10:37
<b>Naphthalene</b>	<b>0.00021</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 10:37
Nitrobenzene	U		0.000024	0.00020	mg/L	1	23-Jan-2020 10:37
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	23-Jan-2020 10:37
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	23-Jan-2020 10:37
<b>Phenanthrene</b>	<b>0.00010</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 10:37
Phenol	U		0.000035	0.00020	mg/L	1	23-Jan-2020 10:37
<b>Pyrene</b>	<b>0.000059</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	23-Jan-2020 10:37
Surr: 2,4,6-Tribromophenol	112			34-129	%REC	1	23-Jan-2020 10:37
Surr: 2-Fluorobiphenyl	100			40-125	%REC	1	23-Jan-2020 10:37
Surr: 2-Fluorophenol	67.9			20-120	%REC	1	23-Jan-2020 10:37
Surr: 4-Terphenyl-d14	110			40-135	%REC	1	23-Jan-2020 10:37
Surr: Nitrobenzene-d5	77.2			41-120	%REC	1	23-Jan-2020 10:37
Surr: Phenol-d6	80.4			20-120	%REC	1	23-Jan-2020 10:37
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 23-Jan-2020		Analyst: ALR	
<b>Arsenic</b>	<b>0.00107</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 02:51

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**Batch ID:** 149797      **Start Date:** 20 Jan 2020 12:00      **End Date:** 20 Jan 2020 16:30  
**Method:** TX 1005 PREP      **Prep Code:** TX 1005\_W PR

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-20	1	30.95 (g)	3 (mL)	0.09693
HS20010767-21	1	31.02 (g)	3 (mL)	0.09671
HS20010767-22	1	30.9 (g)	3 (mL)	0.09709
HS20010767-23	1	31.7 (g)	3 (mL)	0.09464
HS20010767-24	1	30.99 (g)	3 (mL)	0.09681
HS20010767-25	1	31.01 (g)	3 (mL)	0.09674

**Batch ID:** 149831      **Start Date:** 21 Jan 2020 08:30      **End Date:** 21 Jan 2020 12:30  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-18		10 (mL)	10 (mL)	1

**Batch ID:** 149832      **Start Date:** 21 Jan 2020 13:00      **End Date:** 21 Jan 2020 17:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-01		10 (mL)	10 (mL)	1
HS20010767-02		10 (mL)	10 (mL)	1
HS20010767-04		10 (mL)	10 (mL)	1

**Batch ID:** 149881      **Start Date:** 22 Jan 2020 09:00      **End Date:** 22 Jan 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-03		10 (mL)	10 (mL)	1
HS20010767-05		10 (mL)	10 (mL)	1
HS20010767-06		10 (mL)	10 (mL)	1
HS20010767-07		10 (mL)	10 (mL)	1
HS20010767-08		10 (mL)	10 (mL)	1

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**Batch ID:** 149912      **Start Date:** 22 Jan 2020 10:00      **End Date:** 22 Jan 2020 17:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-01	1	1000 (mL)	1 (mL)	0.001
HS20010767-02	1	1000 (mL)	1 (mL)	0.001
HS20010767-03	1	1000 (mL)	1 (mL)	0.001
HS20010767-04	1	1000 (mL)	1 (mL)	0.001
HS20010767-05	1	1000 (mL)	1 (mL)	0.001
HS20010767-06	1	1000 (mL)	1 (mL)	0.001
HS20010767-07	1	1000 (mL)	1 (mL)	0.001
HS20010767-08	1	1000 (mL)	1 (mL)	0.001
HS20010767-09	1	1000 (mL)	1 (mL)	0.001
HS20010767-10	1	1000 (mL)	1 (mL)	0.001
HS20010767-11	1	1000 (mL)	1 (mL)	0.001
HS20010767-12	1	1000 (mL)	1 (mL)	0.001
HS20010767-13	1	1000 (mL)	1 (mL)	0.001
HS20010767-14	1	1000 (mL)	1 (mL)	0.001
HS20010767-15	1	1000 (mL)	1 (mL)	0.001
HS20010767-16	1	1000 (mL)	1 (mL)	0.001
HS20010767-17	1	1000 (mL)	1 (mL)	0.001
HS20010767-18	1	1000 (mL)	1 (mL)	0.001
HS20010767-19	1	1000 (mL)	1 (mL)	0.001

**Batch ID:** 149913      **Start Date:** 22 Jan 2020 11:30      **End Date:** 22 Jan 2020 17:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-20	1	1000 (mL)	1 (mL)	0.001
HS20010767-21	1	1000 (mL)	1 (mL)	0.001
HS20010767-22	1	1000 (mL)	1 (mL)	0.001
HS20010767-23	1	1000 (mL)	1 (mL)	0.001
HS20010767-24	1	1000 (mL)	1 (mL)	0.001
HS20010767-25	1	1000 (mL)	1 (mL)	0.001
HS20010767-27	1	1000 (mL)	1 (mL)	0.001
HS20010767-28	1	1000 (mL)	1 (mL)	0.001
HS20010767-29	1	1000 (mL)	1 (mL)	0.001

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**Batch ID:** 149936      **Start Date:** 23 Jan 2020 08:00      **End Date:** 23 Jan 2020 12:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010767-09		10 (mL)	10 (mL)	1
HS20010767-10		10 (mL)	10 (mL)	1
HS20010767-11		10 (mL)	10 (mL)	1
HS20010767-12		10 (mL)	10 (mL)	1
HS20010767-13		10 (mL)	10 (mL)	1
HS20010767-14		10 (mL)	10 (mL)	1
HS20010767-15		10 (mL)	10 (mL)	1
HS20010767-16		10 (mL)	10 (mL)	1
HS20010767-17		10 (mL)	10 (mL)	1
HS20010767-19		10 (mL)	10 (mL)	1
HS20010767-20		10 (mL)	10 (mL)	1
HS20010767-21		10 (mL)	10 (mL)	1
HS20010767-22		10 (mL)	10 (mL)	1
HS20010767-23		10 (mL)	10 (mL)	1
HS20010767-24		10 (mL)	10 (mL)	1
HS20010767-25		10 (mL)	10 (mL)	1
HS20010767-27		10 (mL)	10 (mL)	1
HS20010767-28		10 (mL)	10 (mL)	1
HS20010767-29		10 (mL)	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149797 ( 0 )</b>		<b>Test Name : LOW-LEVEL TEXAS TPH BY TX1005</b>			<b>Matrix: Water</b>	
HS20010767-20	WG-1620-MW86C-20200117	17 Jan 2020 07:25		20 Jan 2020 12:00	21 Jan 2020 05:19	1
HS20010767-21	WG-1620-FD04-20200117	17 Jan 2020 07:25		20 Jan 2020 12:00	21 Jan 2020 05:48	1
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15		20 Jan 2020 12:00	21 Jan 2020 06:16	1
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20		20 Jan 2020 12:00	21 Jan 2020 06:45	1
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20		20 Jan 2020 12:00	21 Jan 2020 07:14	1
HS20010767-25	WG-1620-MW78A-20200117	17 Jan 2020 11:25		20 Jan 2020 12:00	21 Jan 2020 07:42	1
<b>Batch ID: 149831 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010767-18	WG-1620-MW65D-20200116	16 Jan 2020 16:20		21 Jan 2020 12:30	23 Jan 2020 00:56	1
<b>Batch ID: 149832 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010767-01	WG-1620-MW28A-20200116	16 Jan 2020 09:10		21 Jan 2020 17:00	22 Jan 2020 23:52	1
HS20010767-02	WG-1620-MW28C-20200116	16 Jan 2020 10:15		21 Jan 2020 17:00	22 Jan 2020 23:32	1
HS20010767-04	WG-1620-MW89B-20200116	16 Jan 2020 12:45		21 Jan 2020 17:00	22 Jan 2020 23:54	1
<b>Batch ID: 149881 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010767-03	WG-1620-FB08-20200116	16 Jan 2020 12:00		22 Jan 2020 13:00	24 Jan 2020 16:10	1
HS20010767-05	WG-1620-MW84B-20200116	16 Jan 2020 13:45		22 Jan 2020 13:00	24 Jan 2020 16:13	1
HS20010767-06	WG-1620-MW63B-20200116	16 Jan 2020 14:45		22 Jan 2020 13:00	24 Jan 2020 16:15	1
HS20010767-07	WG-1620-MW26A-20200116	16 Jan 2020 16:00		22 Jan 2020 13:00	24 Jan 2020 16:17	1
HS20010767-08	WG-1620-MW68C-20200117	17 Jan 2020 09:10		22 Jan 2020 13:00	24 Jan 2020 16:21	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149912 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010767-01	WG-1620-MW28A-20200116	16 Jan 2020 09:10		22 Jan 2020 10:00	25 Jan 2020 18:43	1
HS20010767-02	WG-1620-MW28C-20200116	16 Jan 2020 10:15		22 Jan 2020 10:00	25 Jan 2020 12:03	1
HS20010767-03	WG-1620-FB08-20200116	16 Jan 2020 12:00		22 Jan 2020 10:00	25 Jan 2020 19:02	1
HS20010767-04	WG-1620-MW89B-20200116	16 Jan 2020 12:45		22 Jan 2020 10:00	25 Jan 2020 19:22	1
HS20010767-05	WG-1620-MW84B-20200116	16 Jan 2020 13:45		22 Jan 2020 10:00	27 Jan 2020 16:43	20
HS20010767-05	WG-1620-MW84B-20200116	16 Jan 2020 13:45		22 Jan 2020 10:00	25 Jan 2020 19:41	1
HS20010767-06	WG-1620-MW63B-20200116	16 Jan 2020 14:45		22 Jan 2020 10:00	25 Jan 2020 20:00	1
HS20010767-07	WG-1620-MW26A-20200116	16 Jan 2020 16:00		22 Jan 2020 10:00	25 Jan 2020 20:19	1
HS20010767-08	WG-1620-MW68C-20200117	17 Jan 2020 09:10		22 Jan 2020 10:00	25 Jan 2020 20:38	1
HS20010767-09	WG-1620-MW68A-20200117	17 Jan 2020 10:15		22 Jan 2020 10:00	25 Jan 2020 20:57	1
HS20010767-10	WG-1620-FB09-20200117	17 Jan 2020 10:00		22 Jan 2020 10:00	25 Jan 2020 21:16	1
HS20010767-11	WG-1620-MW27A-20200115	15 Jan 2020 16:40		22 Jan 2020 10:00	25 Jan 2020 21:35	1
HS20010767-12	WG-1620-MW27C-20200115	15 Jan 2020 17:20		22 Jan 2020 10:00	26 Jan 2020 13:17	1
HS20010767-13	WG-1620-MW59A-20200116	16 Jan 2020 08:05		22 Jan 2020 10:00	26 Jan 2020 13:36	1
HS20010767-14	WG-1620-MW59B-20200116	16 Jan 2020 09:00		22 Jan 2020 10:00	27 Jan 2020 08:31	1
HS20010767-15	WG-1620-MW59D-20200116	16 Jan 2020 13:50		22 Jan 2020 10:00	26 Jan 2020 14:14	1
HS20010767-16	WG-1620-FD03-20200116	16 Jan 2020 13:50		22 Jan 2020 10:00	26 Jan 2020 14:34	1
HS20010767-17	WG-1620-MW66D-20200116	16 Jan 2020 15:00		22 Jan 2020 10:00	27 Jan 2020 08:50	1
HS20010767-18	WG-1620-MW65D-20200116	16 Jan 2020 16:20		22 Jan 2020 10:00	23 Jan 2020 09:40	1
HS20010767-19	WG-1620-MW36D-20200116	16 Jan 2020 17:25		22 Jan 2020 10:00	26 Jan 2020 15:12	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149913 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010767-20	WG-1620-MW86C-20200117	17 Jan 2020 07:25		22 Jan 2020 11:30	23 Jan 2020 18:00	1
HS20010767-21	WG-1620-FD04-20200117	17 Jan 2020 07:25		22 Jan 2020 11:30	23 Jan 2020 18:19	1
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15		22 Jan 2020 11:30	27 Jan 2020 09:47	200
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15		22 Jan 2020 11:30	27 Jan 2020 09:28	20
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15		22 Jan 2020 11:30	25 Jan 2020 21:55	2000
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15		22 Jan 2020 11:30	23 Jan 2020 18:38	1
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20		22 Jan 2020 11:30	27 Jan 2020 10:24	200
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20		22 Jan 2020 11:30	27 Jan 2020 10:06	20
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20		22 Jan 2020 11:30	26 Jan 2020 18:43	2000
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20		22 Jan 2020 11:30	23 Jan 2020 18:57	1
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20		22 Jan 2020 11:30	27 Jan 2020 10:43	10
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20		22 Jan 2020 11:30	27 Jan 2020 11:02	100
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20		22 Jan 2020 11:30	26 Jan 2020 19:40	1000
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20		22 Jan 2020 11:30	23 Jan 2020 19:17	1
HS20010767-25	WG-1620-MW78A-20200117	17 Jan 2020 11:25		22 Jan 2020 11:30	28 Jan 2020 09:03	400
HS20010767-25	WG-1620-MW78A-20200117	17 Jan 2020 11:25		22 Jan 2020 11:30	28 Jan 2020 08:25	20
HS20010767-25	WG-1620-MW78A-20200117	17 Jan 2020 11:25		22 Jan 2020 11:30	26 Jan 2020 15:31	1
HS20010767-27	WG-1620-MW48C-20200116	16 Jan 2020 10:20		22 Jan 2020 11:30	26 Jan 2020 15:50	1
HS20010767-28	WG-1620-MW47C-20200116	16 Jan 2020 11:05		22 Jan 2020 11:30	27 Jan 2020 09:09	1
HS20010767-29	WG-1620-MW61A-20200116	16 Jan 2020 12:00		22 Jan 2020 11:30	23 Jan 2020 10:37	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149936 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010767-09	WG-1620-MW68A-20200117	17 Jan 2020 10:15		23 Jan 2020 12:00	24 Jan 2020 03:07	1
HS20010767-10	WG-1620-FB09-20200117	17 Jan 2020 10:00		23 Jan 2020 12:00	24 Jan 2020 03:09	1
HS20010767-11	WG-1620-MW27A-20200115	15 Jan 2020 16:40		23 Jan 2020 12:00	24 Jan 2020 03:11	1
HS20010767-12	WG-1620-MW27C-20200115	15 Jan 2020 17:20		23 Jan 2020 12:00	24 Jan 2020 03:13	1
HS20010767-13	WG-1620-MW59A-20200116	16 Jan 2020 08:05		23 Jan 2020 12:00	24 Jan 2020 03:15	1
HS20010767-14	WG-1620-MW59B-20200116	16 Jan 2020 09:00		23 Jan 2020 12:00	24 Jan 2020 03:18	1
HS20010767-15	WG-1620-MW59D-20200116	16 Jan 2020 13:50		23 Jan 2020 12:00	24 Jan 2020 03:20	1
HS20010767-16	WG-1620-FD03-20200116	16 Jan 2020 13:50		23 Jan 2020 12:00	24 Jan 2020 03:22	1
HS20010767-17	WG-1620-MW66D-20200116	16 Jan 2020 15:00		23 Jan 2020 12:00	24 Jan 2020 03:24	1
HS20010767-19	WG-1620-MW36D-20200116	16 Jan 2020 17:25		23 Jan 2020 12:00	24 Jan 2020 03:27	1
HS20010767-20	WG-1620-MW86C-20200117	17 Jan 2020 07:25		23 Jan 2020 12:00	24 Jan 2020 03:33	1
HS20010767-21	WG-1620-FD04-20200117	17 Jan 2020 07:25		23 Jan 2020 12:00	24 Jan 2020 03:35	1
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15		23 Jan 2020 12:00	24 Jan 2020 03:38	1
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20		23 Jan 2020 12:00	24 Jan 2020 03:40	1
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20		23 Jan 2020 12:00	24 Jan 2020 03:42	1
HS20010767-25	WG-1620-MW78A-20200117	17 Jan 2020 11:25		23 Jan 2020 12:00	24 Jan 2020 03:44	1
HS20010767-27	WG-1620-MW48C-20200116	16 Jan 2020 10:20		23 Jan 2020 12:00	24 Jan 2020 03:46	1
HS20010767-28	WG-1620-MW47C-20200116	16 Jan 2020 11:05		23 Jan 2020 12:00	24 Jan 2020 03:49	1
HS20010767-29	WG-1620-MW61A-20200116	16 Jan 2020 12:00		23 Jan 2020 12:00	24 Jan 2020 02:51	1
<b>Batch ID: R354794 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010767-01	WG-1620-MW28A-20200116	16 Jan 2020 09:10			22 Jan 2020 00:55	1
HS20010767-02	WG-1620-MW28C-20200116	16 Jan 2020 10:15			21 Jan 2020 23:41	1
HS20010767-03	WG-1620-FB08-20200116	16 Jan 2020 12:00			22 Jan 2020 02:58	1
<b>Batch ID: R354799 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010767-04	WG-1620-MW89B-20200116	16 Jan 2020 12:45			21 Jan 2020 18:37	1
HS20010767-05	WG-1620-MW84B-20200116	16 Jan 2020 13:45			21 Jan 2020 17:23	1
HS20010767-06	WG-1620-MW63B-20200116	16 Jan 2020 14:45			21 Jan 2020 19:01	1
HS20010767-08	WG-1620-MW68C-20200117	17 Jan 2020 09:10			21 Jan 2020 19:26	1
HS20010767-09	WG-1620-MW68A-20200117	17 Jan 2020 10:15			21 Jan 2020 19:50	1
HS20010767-11	WG-1620-MW27A-20200115	15 Jan 2020 16:40			21 Jan 2020 20:14	1
HS20010767-12	WG-1620-MW27C-20200115	15 Jan 2020 17:20			21 Jan 2020 20:39	1
HS20010767-13	WG-1620-MW59A-20200116	16 Jan 2020 08:05			21 Jan 2020 21:03	1
HS20010767-14	WG-1620-MW59B-20200116	16 Jan 2020 09:00			21 Jan 2020 21:28	1
HS20010767-15	WG-1620-MW59D-20200116	16 Jan 2020 13:50			21 Jan 2020 21:52	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: R354802 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010767-07	WG-1620-MW26A-20200116	16 Jan 2020 16:00			22 Jan 2020 09:17	1
HS20010767-10	WG-1620-FB09-20200117	17 Jan 2020 10:00			22 Jan 2020 01:08	1
HS20010767-16	WG-1620-FD03-20200116	16 Jan 2020 13:50			22 Jan 2020 03:10	1
HS20010767-17	WG-1620-MW66D-20200116	16 Jan 2020 15:00			22 Jan 2020 03:35	1
HS20010767-18	WG-1620-MW65D-20200116	16 Jan 2020 16:20			22 Jan 2020 01:57	1
HS20010767-19	WG-1620-MW36D-20200116	16 Jan 2020 17:25			22 Jan 2020 03:59	1
HS20010767-20	WG-1620-MW86C-20200117	17 Jan 2020 07:25			22 Jan 2020 04:24	1
HS20010767-21	WG-1620-FD04-20200117	17 Jan 2020 07:25			22 Jan 2020 04:48	1
HS20010767-22	WG-1620-MW75B-20200117	17 Jan 2020 08:15			22 Jan 2020 05:13	1
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20			22 Jan 2020 05:37	1
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20			22 Jan 2020 06:01	1
HS20010767-25	WG-1620-MW78A-20200117	17 Jan 2020 11:25			22 Jan 2020 06:26	1
HS20010767-26	WQ-1620-TB05-2020117	17 Jan 2020 00:00			22 Jan 2020 01:32	1
HS20010767-27	WG-1620-MW48C-20200116	16 Jan 2020 10:20			22 Jan 2020 06:50	1
HS20010767-28	WG-1620-MW47C-20200116	16 Jan 2020 11:05			22 Jan 2020 07:15	1
<b>Batch ID: R354859 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010767-10	WG-1620-FB09-20200117	17 Jan 2020 10:00			22 Jan 2020 13:50	1
HS20010767-23	WG-1620-MW79A-20200117	17 Jan 2020 09:20			22 Jan 2020 16:23	5
HS20010767-24	WG-1620-MW74B-20200117	17 Jan 2020 10:20			22 Jan 2020 16:50	5
HS20010767-26	WQ-1620-TB05-2020117	17 Jan 2020 00:00			22 Jan 2020 14:15	1
<b>Batch ID: R355066 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010767-29	WG-1620-MW61A-20200116	16 Jan 2020 12:00			25 Jan 2020 12:37	1

WorkOrder: HS20010767  
 InstrumentID: FID-10  
 Test Code: TX1005\_W\_Low  
 Test Number: TX1005  
 Test Name: Low-level Texas TPH by TX1005

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	nC6 to nC12	TPH-1005-1	0.25	0.26	0.20	0.50
A	>nC12 to nC28	TPH-1005-2	0.25	0.31	0.20	0.50
A	>nC28 to nC35	TPH-1005-4	0.25	0.26	0.20	0.50
A	Total Petroleum Hydrocarbon	TPH	0.25	0.26	0.20	0.50
S	2-Fluorobiphenyl	321-60-8	0	0	0	0
S	Trifluoromethyl benzene	98-08-8	0	0	0	0

WorkOrder: HS20010767  
 InstrumentID: ICPMS04  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000456	0.000400	0.00200

WorkOrder: HS20010767  
InstrumentID: ICPMS05  
Test Code: ICP\_TW  
Test Number: SW6020  
Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20010767  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010767  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010767  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00045	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010



WorkOrder: HS20010767  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** 149797 ( 0 )      **Instrument:** FID-10      **Method:** LOW-LEVEL TEXAS TPH BY TX1005

<b>MBLK</b>		Sample ID: <b>MBLK-149797</b>		Units: <b>mg/L</b>		Analysis Date: <b>21-Jan-2020 04:50</b>			
Client ID:		Run ID: <b>FID-10_354806</b>		SeqNo: <b>5443281</b>		PrepDate: <b>20-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
nC6 to nC12	U	0.50							
>nC12 to nC28	U	0.50							
>nC28 to nC35	U	0.50							
Total Petroleum Hydrocarbon	U	0.50							
Surr: 2-Fluorobiphenyl	2.764	0	2.5	0	111	70 - 130			
Surr: Trifluoromethyl benzene	2.717	0	2.5	0	109	70 - 130			

<b>LCS</b>		Sample ID: <b>LCS-149797</b>		Units: <b>mg/L</b>		Analysis Date: <b>21-Jan-2020 05:19</b>			
Client ID:		Run ID: <b>FID-10_354806</b>		SeqNo: <b>5443282</b>		PrepDate: <b>20-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
nC6 to nC12	28.98	0.50	25	0	116	75 - 125			
>nC12 to nC28	25.19	0.50	25	0	101	75 - 125			
Surr: 2-Fluorobiphenyl	3.022	0	2.5	0	121	70 - 130			
Surr: Trifluoromethyl benzene	2.779	0	2.5	0	111	70 - 130			

<b>LCSD</b>		Sample ID: <b>LCSD-149797</b>		Units: <b>mg/L</b>		Analysis Date: <b>21-Jan-2020 05:48</b>			
Client ID:		Run ID: <b>FID-10_354806</b>		SeqNo: <b>5443283</b>		PrepDate: <b>20-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
nC6 to nC12	28.84	0.50	25	0	115	75 - 125	28.98	0.499	20
>nC12 to nC28	21.62	0.50	25	0	86.5	75 - 125	25.19	15.3	20
Surr: 2-Fluorobiphenyl	2.954	0	2.5	0	118	70 - 130	3.022	2.25	20
Surr: Trifluoromethyl benzene	2.626	0	2.5	0	105	70 - 130	2.779	5.68	20

<b>MS</b>		Sample ID: <b>HS20010754-08MS</b>		Units: <b>mg/L</b>		Analysis Date: <b>21-Jan-2020 06:45</b>			
Client ID:		Run ID: <b>FID-10_354806</b>		SeqNo: <b>5443285</b>		PrepDate: <b>20-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
nC6 to nC12	28.22	0.48	24.23	0	116	75 - 125			
>nC12 to nC28	24.36	0.48	24.23	0	101	75 - 125			
Surr: 2-Fluorobiphenyl	3.105	0	2.423	0	128	70 - 130			
Surr: Trifluoromethyl benzene	2.633	0	2.423	0	109	70 - 130			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** 149797 ( 0 )      **Instrument:** FID-10      **Method:** LOW-LEVEL TEXAS TPH BY TX1005

**MSD**      Sample ID: **HS20010754-08MSD**      Units: **mg/L**      Analysis Date: **21-Jan-2020 07:14**  
 Client ID:      Run ID: **FID-10\_354806**      SeqNo: **5443286**      PrepDate: **20-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

nC6 to nC12	28.04	0.48	23.96	0	117	75 - 125	28.22	0.644	20
>nC12 to nC28	22.82	0.48	23.96	0	95.2	75 - 125	24.36	6.52	20
Surr: 2-Fluorobiphenyl	2.928	0	2.396	0	122	70 - 130	3.105	5.88	20
Surr: Trifluoromethyl benzene	2.517	0	2.396	0	105	70 - 130	2.633	4.52	20

The following samples were analyzed in this batch: HS20010767-20    HS20010767-21    HS20010767-22    HS20010767-23  
 HS20010767-24    HS20010767-25

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149831 ( 0 )		Instrument: ICPMS06		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-149831</b>	Units: <b>mg/L</b>		Analysis Date: <b>23-Jan-2020 13:06</b>					
Client ID:		Run ID: <b>ICPMS06_354915</b>	SeqNo: <b>5445305</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-149831</b>	Units: <b>mg/L</b>		Analysis Date: <b>23-Jan-2020 00:54</b>					
Client ID:		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444295</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.04846	0.00200	0.05	0	96.9	80 - 120			
<b>MS</b>	Sample ID: <b>HS20010767-18MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>23-Jan-2020 01:01</b>					
Client ID: <b>WG-1620-MW65D-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444244</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.04904	0.00200	0.05	0.000507	97.1	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20010767-18MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>23-Jan-2020 01:03</b>					
Client ID: <b>WG-1620-MW65D-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444245</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.04827	0.00200	0.05	0.000507	95.5	80 - 120	0.04904	1.58	20
<b>PDS</b>	Sample ID: <b>HS20010767-18PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>23-Jan-2020 01:05</b>					
Client ID: <b>WG-1620-MW65D-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444246</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.09677	0.00200	0.1	0.000507	96.3	75 - 125			
<b>SD</b>	Sample ID: <b>HS20010767-18SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>23-Jan-2020 00:59</b>					
Client ID: <b>WG-1620-MW65D-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444297</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	U	0.0100					0.000507	0	10

The following samples were analyzed in this batch: HS20010767-18

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149832 ( 0 )		Instrument: ICPMS04		Method: ICP-MS METALS BY SW6020A						
<b>MBLK</b>	Sample ID: <b>MBLKF1-149832</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jan-2020 23:27</b>						
Client ID:		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444345</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	U	0.00200								
<b>MBLK</b>	Sample ID: <b>MBLK-149832</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jan-2020 23:25</b>						
Client ID:		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444228</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-149832</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jan-2020 23:30</b>						
Client ID:		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444230</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.0493	0.00200	0.05	0	98.6	80 - 120				
<b>MS</b>	Sample ID: <b>HS20010767-02MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jan-2020 23:36</b>						
Client ID: <b>WG-1620-MW28C-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444233</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.04794	0.00200	0.05	0.000937	94.0	80 - 120				
<b>MSD</b>	Sample ID: <b>HS20010767-02MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jan-2020 23:39</b>						
Client ID: <b>WG-1620-MW28C-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444234</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.05018	0.00200	0.05	0.000937	98.5	80 - 120	0.04794	4.57	20	
<b>PDS</b>	Sample ID: <b>HS20010767-02PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>22-Jan-2020 23:41</b>						
Client ID: <b>WG-1620-MW28C-20200116</b>		Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444235</b>	PrepDate: <b>21-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.09536	0.00200	0.1	0.000937	94.4	75 - 125				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** 149832 ( 0 )      **Instrument:** ICPMS04      **Method:** ICP-MS METALS BY SW6020A

<b>SD</b>	Sample ID: <b>HS20010767-02SD</b>	Units: <b>mg/L</b>	Analysis Date: <b>22-Jan-2020 23:34</b>						
Client ID: <b>WG-1620-MW28C-20200116</b>	Run ID: <b>ICPMS04_354801</b>	SeqNo: <b>5444232</b>	PrepDate: <b>21-Jan-2020</b> DF: <b>5</b>						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	Limit Qual

Arsenic	U	0.0100					0.000937	0	10
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The following samples were analyzed in this batch: 

HS20010767-01	HS20010767-02	HS20010767-04
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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149881 ( 0 )		Instrument: ICPMS05			Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLKF1-149881</b>	Units: <b>mg/L</b>			Analysis Date: <b>24-Jan-2020 14:21</b>					
Client ID:		Run ID: <b>ICPMS05_354996</b>			SeqNo: <b>5447636</b>		PrepDate: <b>22-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>MBLK</b>	Sample ID: <b>MBLK-149881</b>	Units: <b>mg/L</b>			Analysis Date: <b>24-Jan-2020 14:18</b>					
Client ID:		Run ID: <b>ICPMS05_354996</b>			SeqNo: <b>5447635</b>		PrepDate: <b>22-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-149881</b>	Units: <b>mg/L</b>			Analysis Date: <b>24-Jan-2020 14:23</b>					
Client ID:		Run ID: <b>ICPMS05_354996</b>			SeqNo: <b>5447637</b>		PrepDate: <b>22-Jan-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.04927	0.00200	0.05	0	98.6	80 - 120				
<b>MS</b>	Sample ID: <b>HS20010752-08MS</b>	Units: <b>mg/L</b>			Analysis Date: <b>24-Jan-2020 15:26</b>					
Client ID:		Run ID: <b>ICPMS05_354996</b>			SeqNo: <b>5447650</b>		PrepDate: <b>22-Jan-2020</b>		DF: <b>10</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.1068	0.0200	0.05	0.06877	76.0	80 - 120				S
<b>MSD</b>	Sample ID: <b>HS20010752-08MSD</b>	Units: <b>mg/L</b>			Analysis Date: <b>27-Jan-2020 14:34</b>					
Client ID:		Run ID: <b>ICPMS05_355087</b>			SeqNo: <b>5448409</b>		PrepDate: <b>22-Jan-2020</b>		DF: <b>10</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.1141	0.0200	0.05	0.06877	90.6	80 - 120	0.1068	6.58	20	
<b>PDS</b>	Sample ID: <b>HS20010752-08PDS</b>	Units: <b>mg/L</b>			Analysis Date: <b>27-Jan-2020 14:39</b>					
Client ID:		Run ID: <b>ICPMS05_355087</b>			SeqNo: <b>5448411</b>		PrepDate: <b>22-Jan-2020</b>		DF: <b>10</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	1.132	0.0200	1	0.06877	106	75 - 125				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> 149881 ( 0 )		<b>Instrument:</b> ICPMS05		<b>Method:</b> ICP-MS METALS BY SW6020A					
<b>SD</b>	Sample ID: <b>HS20010752-08SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 15:24</b>					
Client ID:	Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447649</b>	PrepDate: <b>22-Jan-2020</b>	DF: <b>50</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	Limit Qual

Arsenic	0.07175	0.100					0.06877	0 10	J
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The following samples were analyzed in this batch:

HS20010767-03	HS20010767-05	HS20010767-06	HS20010767-07
HS20010767-08			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> 149936 ( 0 )	<b>Instrument:</b> ICPMS05	<b>Method:</b> ICP-MS METALS BY SW6020A
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<b>MBLK</b>	Sample ID: <b>MBLK-149936</b>	Units: <b>mg/L</b>	Analysis Date: <b>24-Jan-2020 02:47</b>							
Client ID:	Run ID: <b>ICPMS05_354962</b>	SeqNo: <b>5446103</b>	PrepDate: <b>23-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic U 0.00200

<b>LCS</b>	Sample ID: <b>LCS-149936</b>	Units: <b>mg/L</b>	Analysis Date: <b>24-Jan-2020 02:49</b>							
Client ID:	Run ID: <b>ICPMS05_354962</b>	SeqNo: <b>5446104</b>	PrepDate: <b>23-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.04886 0.00200 0.05 0 97.7 80 - 120

<b>MS</b>	Sample ID: <b>HS20010767-29MS</b>	Units: <b>mg/L</b>	Analysis Date: <b>24-Jan-2020 02:56</b>							
Client ID: <b>WG-1620-MW61A-20200116</b>	Run ID: <b>ICPMS05_354962</b>	SeqNo: <b>5446107</b>	PrepDate: <b>23-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.05254 0.00200 0.05 0.00107 103 80 - 120

<b>MSD</b>	Sample ID: <b>HS20010767-29MSD</b>	Units: <b>mg/L</b>	Analysis Date: <b>24-Jan-2020 02:58</b>							
Client ID: <b>WG-1620-MW61A-20200116</b>	Run ID: <b>ICPMS05_354962</b>	SeqNo: <b>5446108</b>	PrepDate: <b>23-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.05469 0.00200 0.05 0.00107 107 80 - 120 0.05254 4 20

<b>PDS</b>	Sample ID: <b>HS20010767-29PDS</b>	Units: <b>mg/L</b>	Analysis Date: <b>24-Jan-2020 03:00</b>							
Client ID: <b>WG-1620-MW61A-20200116</b>	Run ID: <b>ICPMS05_354962</b>	SeqNo: <b>5446109</b>	PrepDate: <b>23-Jan-2020</b> DF: <b>1</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual

Arsenic 0.1055 0.00200 0.1 0.00107 104 75 - 125

<b>SD</b>	Sample ID: <b>HS20010767-29SD</b>	Units: <b>mg/L</b>	Analysis Date: <b>24-Jan-2020 02:53</b>							
Client ID: <b>WG-1620-MW61A-20200116</b>	Run ID: <b>ICPMS05_354962</b>	SeqNo: <b>5446106</b>	PrepDate: <b>23-Jan-2020</b> DF: <b>5</b>							
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual

Arsenic U 0.0100 0.00107 0 10

<b>The following samples were analyzed in this batch:</b>	HS20010767-09	HS20010767-10	HS20010767-11	HS20010767-12
	HS20010767-13	HS20010767-14	HS20010767-15	HS20010767-16
	HS20010767-17	HS20010767-19	HS20010767-20	HS20010767-21
	HS20010767-22	HS20010767-23	HS20010767-24	HS20010767-25
	HS20010767-27	HS20010767-28	HS20010767-29	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149912 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149912	Units: ug/L			Analysis Date: 23-Jan-2020 08:56					
Client ID:	Run ID: SV-7_354913	SeqNo: 5445204		PrepDate: 22-Jan-2020		DF: 1				
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.765</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>115</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.735</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>115</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.308</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.2</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>6.463</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>129</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.615</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.471</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.4</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149912 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149912	Units: ug/L			Analysis Date: 23-Jan-2020 09:15					
Client ID:	Run ID: SV-7_354913	SeqNo: 5445205		PrepDate: 22-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.941	0.20	5	0	98.8	39 - 127				
2,4-Dimethylphenol	4.975	0.20	5	0	99.5	35 - 120				
2,4-Dinitrotoluene	5.465	0.20	5	0	109	50 - 122				
2,6-Dinitrotoluene	5.315	0.20	5	0	106	50 - 120				
2-Chloronaphthalene	5.375	0.20	5	0	108	50 - 120				
2-Methylnaphthalene	5.452	0.10	5	0	109	50 - 120				
4,6-Dinitro-2-methylphenol	4.139	0.20	5	0	82.8	25 - 121				
4-Nitrophenol	4.795	1.0	5	0	95.9	30 - 130				
Acenaphthene	5.261	0.10	5	0	105	45 - 120				
Acenaphthylene	5.595	0.10	5	0	112	47 - 120				
Anthracene	5.443	0.10	5	0	109	45 - 120				
Benz(a)anthracene	5.163	0.10	5	0	103	40 - 120				
Benzo(a)pyrene	5.388	0.10	5	0	108	45 - 120				
Bis(2-chloroethoxy)methane	4.569	0.20	5	0	91.4	45 - 120				
Bis(2-ethylhexyl)phthalate	6.14	0.20	5	0	123	40 - 139				
Chrysene	5.073	0.10	5	0	101	43 - 120				
Dibenzofuran	5.456	0.10	5	0	109	50 - 120				
Di-n-butyl phthalate	6.076	0.20	5	0	122	45 - 123				
Fluoranthene	5.46	0.10	5	0	109	45 - 125				
Fluorene	5.591	0.10	5	0	112	49 - 120				
Naphthalene	5.915	0.10	5	0	118	45 - 120				
Nitrobenzene	4.369	0.20	5	0	87.4	44 - 120				
N-Nitrosodiphenylamine	5.411	0.20	5	0	108	40 - 125				
Pentachlorophenol	2.724	0.20	5	0	54.5	19 - 121				
Phenanthrene	5.334	0.10	5	0	107	45 - 121				
Phenol	4.829	0.20	5	0	96.6	20 - 124				
Pyrene	5.438	0.10	5	0	109	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>6.301</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>126</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>6.146</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>123</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.409</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>88.2</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>6.275</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>126</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.93</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.815</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.3</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149912 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS20010767-18MS		Units: ug/L		Analysis Date: 23-Jan-2020 09:59				
Client ID: WG-1620-MW65D-20200116		Run ID: SV-7_354913		SeqNo: 5445207		PrepDate: 22-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.323	0.20	5	0	86.5	39 - 127				
2,4-Dimethylphenol	4.078	0.20	5	0	81.6	35 - 120				
2,4-Dinitrotoluene	4.77	0.20	5	0	95.4	50 - 122				
2,6-Dinitrotoluene	4.895	0.20	5	0	97.9	50 - 120				
2-Chloronaphthalene	4.764	0.20	5	0	95.3	50 - 120				
2-Methylnaphthalene	4.519	0.10	5	0	90.4	50 - 120				
4,6-Dinitro-2-methylphenol	4.287	0.20	5	0	85.7	25 - 121				
4-Nitrophenol	4.37	1.0	5	0	87.4	30 - 130				
Acenaphthene	4.646	0.10	5	0	92.9	45 - 120				
Acenaphthylene	4.886	0.10	5	0	97.7	47 - 120				
Anthracene	4.867	0.10	5	0	97.3	45 - 120				
Benz(a)anthracene	4.722	0.10	5	0	94.4	40 - 120				
Benzo(a)pyrene	4.84	0.10	5	0	96.8	45 - 120				
Bis(2-chloroethoxy)methane	3.874	0.20	5	0	77.5	45 - 120				
Bis(2-ethylhexyl)phthalate	5.758	0.20	5	0.2648	110	40 - 139				
Chrysene	4.566	0.10	5	0	91.3	43 - 120				
Dibenzofuran	4.741	0.10	5	0	94.8	50 - 120				
Di-n-butyl phthalate	5.562	0.20	5	0	111	45 - 123				
Fluoranthene	4.981	0.10	5	0	99.6	45 - 125				
Fluorene	4.89	0.10	5	0	97.8	49 - 120				
Naphthalene	4.673	0.10	5	0	93.5	45 - 120				
Nitrobenzene	3.616	0.20	5	0	72.3	44 - 120				
N-Nitrosodiphenylamine	4.609	0.20	5	0	92.2	40 - 125				
Pentachlorophenol	3.229	0.20	5	0	64.6	19 - 121				
Phenanthrene	4.672	0.10	5	0	93.4	45 - 121				
Phenol	4.224	0.20	5	0	84.5	20 - 124				
Pyrene	4.975	0.10	5	0	99.5	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.909</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>118</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.365</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>107</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.067</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>81.3</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.888</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>118</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.346</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.9</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.567</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.3</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149912 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS20010767-02MS		Units: ug/L		Analysis Date: 25-Jan-2020 12:22				
Client ID: WG-1620-MW28C-20200116		Run ID: SV-6_355049		SeqNo: 5448627		PrepDate: 22-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.094	0.20	5	0	102	39 - 127				
2,4-Dimethylphenol	3.359	0.20	5	0	67.2	35 - 120				
2,4-Dinitrotoluene	5.668	0.20	5	0	113	50 - 122				
2,6-Dinitrotoluene	4.959	0.20	5	0	99.2	50 - 120				
2-Chloronaphthalene	4.264	0.20	5	0	85.3	50 - 120				
2-Methylnaphthalene	4.045	0.10	5	0	80.9	50 - 120				
4,6-Dinitro-2-methylphenol	6.467	0.20	5	0	129	25 - 121				S
4-Nitrophenol	5.942	1.0	5	0	119	30 - 130				
Acenaphthene	4.553	0.10	5	0	91.1	45 - 120				
Acenaphthylene	4.175	0.10	5	0	83.5	47 - 120				
Anthracene	5.523	0.10	5	0	110	45 - 120				
Benz(a)anthracene	5.628	0.10	5	0	113	40 - 120				
Benzo(a)pyrene	5.453	0.10	5	0	109	45 - 120				
Bis(2-chloroethoxy)methane	4.015	0.20	5	0	80.3	45 - 120				
Bis(2-ethylhexyl)phthalate	5.277	0.20	5	0.07179	104	40 - 139				
Chrysene	5.372	0.10	5	0	107	43 - 120				
Dibenzofuran	4.21	0.10	5	0	84.2	50 - 120				
Di-n-butyl phthalate	6.154	0.20	5	0	123	45 - 123				S
Fluoranthene	5.615	0.10	5	0	112	45 - 125				
Fluorene	4.421	0.10	5	0	88.4	49 - 120				
Naphthalene	3.715	0.10	5	0.1556	71.2	45 - 120				
Nitrobenzene	3.908	0.20	5	0	78.2	44 - 120				
N-Nitrosodiphenylamine	5.898	0.20	5	0	118	40 - 125				
Pentachlorophenol	5.126	0.20	5	0	103	19 - 121				
Phenanthrene	5.305	0.10	5	0	106	45 - 121				
Phenol	3.48	0.20	5	0.08592	67.9	20 - 124				
Pyrene	5.61	0.10	5	0	112	40 - 130				
Surr: 2,4,6-Tribromophenol	4.646	0.20	5	0	92.9	34 - 129				
Surr: 2-Fluorobiphenyl	3.968	0.20	5	0	79.4	40 - 125				
Surr: 2-Fluorophenol	2.894	0.20	5	0	57.9	20 - 120				
Surr: 4-Terphenyl-d14	5.42	0.20	5	0	108	40 - 135				
Surr: Nitrobenzene-d5	3.91	0.20	5	0	78.2	41 - 120				
Surr: Phenol-d6	3.403	0.20	5	0	68.1	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149912 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS20010767-18MSD			Units: ug/L		Analysis Date: 23-Jan-2020 10:18			
Client ID: WG-1620-MW65D-20200116		Run ID: SV-7_354913			SeqNo: 5445208		PrepDate: 22-Jan-2020		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.677	0.20	5	0	93.5	39 - 127	4.323	7.87	20	
2,4-Dimethylphenol	4.291	0.20	5	0	85.8	35 - 120	4.078	5.09	20	
2,4-Dinitrotoluene	4.91	0.20	5	0	98.2	50 - 122	4.77	2.9	20	
2,6-Dinitrotoluene	4.907	0.20	5	0	98.1	50 - 120	4.895	0.258	20	
2-Chloronaphthalene	4.818	0.20	5	0	96.4	50 - 120	4.764	1.14	20	
2-Methylnaphthalene	4.716	0.10	5	0	94.3	50 - 120	4.519	4.27	20	
4,6-Dinitro-2-methylphenol	4.275	0.20	5	0	85.5	25 - 121	4.287	0.282	30	
4-Nitrophenol	4.178	1.0	5	0	83.6	30 - 130	4.37	4.5	20	
Acenaphthene	4.732	0.10	5	0	94.6	45 - 120	4.646	1.83	20	
Acenaphthylene	5.037	0.10	5	0	101	47 - 120	4.886	3.04	20	
Anthracene	4.947	0.10	5	0	98.9	45 - 120	4.867	1.63	20	
Benz(a)anthracene	4.832	0.10	5	0	96.6	40 - 120	4.722	2.32	20	
Benzo(a)pyrene	4.934	0.10	5	0	98.7	45 - 120	4.84	1.92	20	
Bis(2-chloroethoxy)methane	4.109	0.20	5	0	82.2	45 - 120	3.874	5.88	20	
Bis(2-ethylhexyl)phthalate	5.819	0.20	5	0.2648	111	40 - 139	5.758	1.06	20	
Chrysene	4.668	0.10	5	0	93.4	43 - 120	4.566	2.2	20	
Dibenzofuran	4.871	0.10	5	0	97.4	50 - 120	4.741	2.7	20	
Di-n-butyl phthalate	5.586	0.20	5	0	112	45 - 123	5.562	0.444	20	
Fluoranthene	4.883	0.10	5	0	97.7	45 - 125	4.981	1.97	20	
Fluorene	4.983	0.10	5	0	99.7	49 - 120	4.89	1.88	20	
Naphthalene	4.894	0.10	5	0	97.9	45 - 120	4.673	4.61	20	
Nitrobenzene	3.889	0.20	5	0	77.8	44 - 120	3.616	7.28	20	
N-Nitrosodiphenylamine	4.885	0.20	5	0	97.7	40 - 125	4.609	5.82	20	
Pentachlorophenol	3.134	0.20	5	0	62.7	19 - 121	3.229	3	20	
Phenanthrene	4.996	0.10	5	0	99.9	45 - 121	4.672	6.69	20	
Phenol	4.351	0.20	5	0	87.0	20 - 124	4.224	2.96	20	
Pyrene	5.082	0.10	5	0	102	40 - 130	4.975	2.14	20	
Surr: 2,4,6-Tribromophenol	5.731	0.20	5	0	115	34 - 129	5.909	3.05	20	
Surr: 2-Fluorobiphenyl	5.483	0.20	5	0	110	40 - 125	5.365	2.17	20	
Surr: 2-Fluorophenol	4.007	0.20	5	0	80.1	20 - 120	4.067	1.48	20	
Surr: 4-Terphenyl-d14	5.938	0.20	5	0	119	40 - 135	5.888	0.844	20	
Surr: Nitrobenzene-d5	4.206	0.20	5	0	84.1	41 - 120	4.346	3.27	20	
Surr: Phenol-d6	4.422	0.20	5	0	88.4	20 - 120	4.567	3.23	20	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149912 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS20010767-02MSD			Units: ug/L		Analysis Date: 25-Jan-2020 12:41			
Client ID: WG-1620-MW28C-20200116		Run ID: SV-6_355049		SeqNo: 5448628		PrepDate: 22-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.722	0.20	5	0	94.4	39 - 127	5.094	7.57	20	
2,4-Dimethylphenol	3.021	0.20	5	0	60.4	35 - 120	3.359	10.6	20	
2,4-Dinitrotoluene	5.415	0.20	5	0	108	50 - 122	5.668	4.56	20	
2,6-Dinitrotoluene	5.001	0.20	5	0	100	50 - 120	4.959	0.848	20	
2-Chloronaphthalene	4.118	0.20	5	0	82.4	50 - 120	4.264	3.49	20	
2-Methylnaphthalene	3.849	0.10	5	0	77.0	50 - 120	4.045	4.97	20	
4,6-Dinitro-2-methylphenol	5.971	0.20	5	0	119	25 - 121	6.467	7.97	30	
4-Nitrophenol	5.634	1.0	5	0	113	30 - 130	5.942	5.33	20	
Acenaphthene	4.507	0.10	5	0	90.1	45 - 120	4.553	1.02	20	
Acenaphthylene	4.003	0.10	5	0	80.1	47 - 120	4.175	4.2	20	
Anthracene	5.093	0.10	5	0	102	45 - 120	5.523	8.1	20	
Benz(a)anthracene	5.227	0.10	5	0	105	40 - 120	5.628	7.39	20	
Benzo(a)pyrene	5.12	0.10	5	0	102	45 - 120	5.453	6.31	20	
Bis(2-chloroethoxy)methane	3.329	0.20	5	0	66.6	45 - 120	4.015	18.7	20	
Bis(2-ethylhexyl)phthalate	6.092	0.20	5	0.07179	120	40 - 139	5.277	14.3	20	
Chrysene	5.384	0.10	5	0	108	43 - 120	5.372	0.21	20	
Dibenzofuran	4.23	0.10	5	0	84.6	50 - 120	4.21	0.471	20	
Di-n-butyl phthalate	6.028	0.20	5	0	121	45 - 123	6.154	2.06	20	
Fluoranthene	5.168	0.10	5	0	103	45 - 125	5.615	8.29	20	
Fluorene	4.566	0.10	5	0	91.3	49 - 120	4.421	3.21	20	
Naphthalene	3.661	0.10	5	0.1556	70.1	45 - 120	3.715	1.46	20	
Nitrobenzene	3.397	0.20	5	0	67.9	44 - 120	3.908	14	20	
N-Nitrosodiphenylamine	5.397	0.20	5	0	108	40 - 125	5.898	8.86	20	
Pentachlorophenol	4.902	0.20	5	0	98.0	19 - 121	5.126	4.48	20	
Phenanthrene	4.949	0.10	5	0	99.0	45 - 121	5.305	6.95	20	
Phenol	3.956	0.20	5	0.08592	77.4	20 - 124	3.48	12.8	20	
Pyrene	5.27	0.10	5	0	105	40 - 130	5.61	6.25	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.371</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.4</i>	<i>34 - 129</i>	<i>4.646</i>	<i>6.08</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.867</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.3</i>	<i>40 - 125</i>	<i>3.968</i>	<i>2.56</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.454</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.1</i>	<i>20 - 120</i>	<i>2.894</i>	<i>17.6</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>5.15</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>40 - 135</i>	<i>5.42</i>	<i>5.1</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.342</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>66.8</i>	<i>41 - 120</i>	<i>3.91</i>	<i>15.7</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.686</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.7</i>	<i>20 - 120</i>	<i>3.403</i>	<i>8</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> 149912 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D
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The following samples were analyzed in this batch:

HS20010767-01	HS20010767-02	HS20010767-03	HS20010767-04
HS20010767-05	HS20010767-06	HS20010767-07	HS20010767-08
HS20010767-09	HS20010767-10	HS20010767-11	HS20010767-12
HS20010767-13	HS20010767-14	HS20010767-15	HS20010767-16
HS20010767-17	HS20010767-18	HS20010767-19	



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149913 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149913	Units: ug/L			Analysis Date: 23-Jan-2020 08:18					
Client ID:	Run ID: SV-7_354913	SeqNo: 5447234		PrepDate: 22-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	5.682	0.20	5	0	114	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	5.86	0.20	5	0	117	40 - 125				
<i>Surr: 2-Fluorophenol</i>	4.972	0.20	5	0	99.4	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	5.966	0.20	5	0	119	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	4.68	0.20	5	0	93.6	41 - 120				
<i>Surr: Phenol-d6</i>	4.665	0.20	5	0	93.3	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149913 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149913	Units: ug/L			Analysis Date: 23-Jan-2020 08:37					
Client ID:	Run ID: SV-7_354913	SeqNo: 5447235		PrepDate: 22-Jan-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.666	0.20	5	0	93.3	39 - 127				
2,4-Dimethylphenol	4.637	0.20	5	0	92.7	35 - 120				
2,4-Dinitrotoluene	4.853	0.20	5	0	97.1	50 - 122				
2,6-Dinitrotoluene	4.895	0.20	5	0	97.9	50 - 120				
2-Chloronaphthalene	5.237	0.20	5	0	105	50 - 120				
2-Methylnaphthalene	4.811	0.10	5	0	96.2	50 - 120				
4,6-Dinitro-2-methylphenol	3.833	0.20	5	0	76.7	25 - 121				
4-Nitrophenol	4.307	1.0	5	0	86.1	30 - 130				
Acenaphthene	4.824	0.10	5	0	96.5	45 - 120				
Acenaphthylene	4.998	0.10	5	0	100.0	47 - 120				
Anthracene	5.009	0.10	5	0	100	45 - 120				
Benz(a)anthracene	4.636	0.10	5	0	92.7	40 - 120				
Benzo(a)pyrene	4.899	0.10	5	0	98.0	45 - 120				
Bis(2-chloroethoxy)methane	4.26	0.20	5	0	85.2	45 - 120				
Bis(2-ethylhexyl)phthalate	5.545	0.20	5	0	111	40 - 139				
Chrysene	4.739	0.10	5	0	94.8	43 - 120				
Dibenzofuran	4.863	0.10	5	0	97.3	50 - 120				
Di-n-butyl phthalate	5.557	0.20	5	0	111	45 - 123				
Fluoranthene	4.941	0.10	5	0	98.8	45 - 125				
Fluorene	4.979	0.10	5	0	99.6	49 - 120				
Naphthalene	5.181	0.10	5	0	104	45 - 120				
Nitrobenzene	3.945	0.20	5	0	78.9	44 - 120				
N-Nitrosodiphenylamine	4.857	0.20	5	0	97.1	40 - 125				
Pentachlorophenol	2.362	0.20	5	0	47.2	19 - 121				
Phenanthrene	4.929	0.10	5	0	98.6	45 - 121				
Phenol	4.554	0.20	5	0	91.1	20 - 124				
Pyrene	4.944	0.10	5	0	98.9	40 - 130				
Surr: 2,4,6-Tribromophenol	5.524	0.20	5	0	110	34 - 129				
Surr: 2-Fluorobiphenyl	5.551	0.20	5	0	111	40 - 125				
Surr: 2-Fluorophenol	4.237	0.20	5	0	84.7	20 - 120				
Surr: 4-Terphenyl-d14	5.865	0.20	5	0	117	40 - 135				
Surr: Nitrobenzene-d5	4.474	0.20	5	0	89.5	41 - 120				
Surr: Phenol-d6	4.723	0.20	5	0	94.5	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149913 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS20010767-29MS		Units: ug/L		Analysis Date: 23-Jan-2020 10:57				
Client ID: WG-1620-MW61A-20200116		Run ID: SV-7_354913		SeqNo: 5447237		PrepDate: 22-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.127	0.20	5	0	82.5	39 - 127				
2,4-Dimethylphenol	3.82	0.20	5	0	76.4	35 - 120				
2,4-Dinitrotoluene	4.735	0.20	5	0	94.7	50 - 122				
2,6-Dinitrotoluene	4.441	0.20	5	0	88.8	50 - 120				
2-Chloronaphthalene	4.418	0.20	5	0	88.4	50 - 120				
2-Methylnaphthalene	4.114	0.10	5	0	82.3	50 - 120				
4,6-Dinitro-2-methylphenol	4.515	0.20	5	0	90.3	25 - 121				
4-Nitrophenol	4.885	1.0	5	0	97.7	30 - 130				
Acenaphthene	4.136	0.10	5	0.05203	81.7	45 - 120				
Acenaphthylene	4.391	0.10	5	0	87.8	47 - 120				
Anthracene	4.875	0.10	5	0	97.5	45 - 120				
Benz(a)anthracene	4.904	0.10	5	0	98.1	40 - 120				
Benzo(a)pyrene	4.989	0.10	5	0	99.8	45 - 120				
Bis(2-chloroethoxy)methane	3.628	0.20	5	0	72.6	45 - 120				
Bis(2-ethylhexyl)phthalate	5.934	0.20	5	0.08059	117	40 - 139				
Chrysene	4.795	0.10	5	0	95.9	43 - 120				
Dibenzofuran	4.243	0.10	5	0	84.9	50 - 120				
Di-n-butyl phthalate	5.583	0.20	5	0	112	45 - 123				
Fluoranthene	5.114	0.10	5	0.06004	101	45 - 125				
Fluorene	4.513	0.10	5	0	90.3	49 - 120				
Naphthalene	4.262	0.10	5	0.2136	81.0	45 - 120				
Nitrobenzene	3.349	0.20	5	0	67.0	44 - 120				
N-Nitrosodiphenylamine	4.286	0.20	5	0	85.7	40 - 125				
Pentachlorophenol	3.54	0.20	5	0	70.8	19 - 121				
Phenanthrene	4.755	0.10	5	0.1016	93.1	45 - 121				
Phenol	3.665	0.20	5	0	73.3	20 - 124				
Pyrene	5.074	0.10	5	0.0591	100	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.896</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>118</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.696</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.9</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.424</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>68.5</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>6.154</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>123</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.847</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.9</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.851</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.0</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

Batch ID: 149913 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS20010767-29MSD	Units: ug/L			Analysis Date: 23-Jan-2020 11:16					
Client ID: WG-1620-MW61A-20200116	Run ID: SV-7_354913	SeqNo: 5447238	PrepDate: 22-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.446	0.20	5	0	88.9	39 - 127	4.127	7.43	20	
2,4-Dimethylphenol	4.161	0.20	5	0	83.2	35 - 120	3.82	8.54	20	
2,4-Dinitrotoluene	4.907	0.20	5	0	98.1	50 - 122	4.735	3.57	20	
2,6-Dinitrotoluene	4.699	0.20	5	0	94.0	50 - 120	4.441	5.64	20	
2-Chloronaphthalene	4.799	0.20	5	0	96.0	50 - 120	4.418	8.26	20	
2-Methylnaphthalene	4.319	0.10	5	0	86.4	50 - 120	4.114	4.88	20	
4,6-Dinitro-2-methylphenol	4.43	0.20	5	0	88.6	25 - 121	4.515	1.91	30	
4-Nitrophenol	4.913	1.0	5	0	98.3	30 - 130	4.885	0.58	20	
Acenaphthene	4.492	0.10	5	0.05203	88.8	45 - 120	4.136	8.24	20	
Acenaphthylene	4.652	0.10	5	0	93.0	47 - 120	4.391	5.76	20	
Anthracene	4.969	0.10	5	0	99.4	45 - 120	4.875	1.9	20	
Benz(a)anthracene	5.055	0.10	5	0	101	40 - 120	4.904	3.05	20	
Benzo(a)pyrene	4.885	0.10	5	0	97.7	45 - 120	4.989	2.12	20	
Bis(2-chloroethoxy)methane	3.667	0.20	5	0	73.3	45 - 120	3.628	1.06	20	
Bis(2-ethylhexyl)phthalate	6.102	0.20	5	0.08059	120	40 - 139	5.934	2.79	20	
Chrysene	4.891	0.10	5	0	97.8	43 - 120	4.795	2	20	
Dibenzofuran	4.52	0.10	5	0	90.4	50 - 120	4.243	6.32	20	
Di-n-butyl phthalate	5.617	0.20	5	0	112	45 - 123	5.583	0.62	20	
Fluoranthene	5.088	0.10	5	0.06004	101	45 - 125	5.114	0.51	20	
Fluorene	4.732	0.10	5	0	94.6	49 - 120	4.513	4.73	20	
Naphthalene	4.399	0.10	5	0.2136	83.7	45 - 120	4.262	3.17	20	
Nitrobenzene	3.289	0.20	5	0	65.8	44 - 120	3.349	1.8	20	
N-Nitrosodiphenylamine	4.566	0.20	5	0	91.3	40 - 125	4.286	6.33	20	
Pentachlorophenol	3.743	0.20	5	0	74.9	19 - 121	3.54	5.57	20	
Phenanthrene	4.944	0.10	5	0.1016	96.8	45 - 121	4.755	3.89	20	
Phenol	3.887	0.20	5	0	77.7	20 - 124	3.665	5.87	20	
Pyrene	5.317	0.10	5	0.0591	105	40 - 130	5.074	4.68	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.221</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>34 - 129</i>	<i>5.896</i>	<i>12.2</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.822</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.4</i>	<i>40 - 125</i>	<i>4.696</i>	<i>2.65</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.562</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>71.2</i>	<i>20 - 120</i>	<i>3.424</i>	<i>3.95</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>6.108</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>122</i>	<i>40 - 135</i>	<i>6.154</i>	<i>0.75</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.608</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.2</i>	<i>41 - 120</i>	<i>3.847</i>	<i>6.42</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.046</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>80.9</i>	<i>20 - 120</i>	<i>3.851</i>	<i>4.95</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> 149913 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS20010767-20	HS20010767-21	HS20010767-22	HS20010767-23
HS20010767-24	HS20010767-25	HS20010767-27	HS20010767-28
HS20010767-29			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> R354794 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200121</b>	Units: <b>ug/L</b>			Analysis Date: <b>21-Jan-2020 21:37</b>				
Client ID:	Run ID: <b>VOA4_354794</b>	SeqNo: <b>5443118</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.87</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.61</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200121</b>	Units: <b>ug/L</b>			Analysis Date: <b>21-Jan-2020 20:48</b>				
Client ID:	Run ID: <b>VOA4_354794</b>	SeqNo: <b>5443117</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	19.81	1.0	20	0	99.0	70 - 124			
Benzene	20.07	1.0	20	0	100	74 - 120			
Chlorobenzene	19.38	1.0	20	0	96.9	76 - 113			
Ethylbenzene	19.76	1.0	20	0	98.8	77 - 117			
Methylene chloride	21.54	2.0	20	0	108	70 - 127			
Toluene	19.62	1.0	20	0	98.1	77 - 118			
Vinyl chloride	19.87	1.0	20	0	99.3	70 - 130			
Xylenes, Total	62.57	1.0	60	0	104	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.85</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.58</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> R354794 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MS</b>		Sample ID: <b>HS20010767-02MS</b>		Units: <b>ug/L</b>		Analysis Date: <b>22-Jan-2020 00:05</b>			
Client ID: <b>WG-1620-MW28C-20200116</b>		Run ID: <b>VOA4_354794</b>		SeqNo: <b>5443123</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	17.06	1.0	20	0	85.3	70 - 127			
Benzene	17.8	1.0	20	0	89.0	70 - 127			
Chlorobenzene	17.78	1.0	20	0	88.9	70 - 114			
Ethylbenzene	18.09	1.0	20	0	90.5	70 - 124			
Methylene chloride	16.05	2.0	20	0	80.2	70 - 128			
Toluene	17.93	1.0	20	0	89.6	70 - 123			
Vinyl chloride	22.17	1.0	20	0.477	108	70 - 130			
Xylenes, Total	56.44	1.0	60	0	94.1	70 - 130			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.2</i>	<i>70 - 126</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.59</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>81 - 113</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.87</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>77 - 123</i>			
<i>Surr: Toluene-d8</i>	<i>50.43</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>			

<b>MSD</b>		Sample ID: <b>HS20010767-02MSD</b>		Units: <b>ug/L</b>		Analysis Date: <b>22-Jan-2020 00:30</b>			
Client ID: <b>WG-1620-MW28C-20200116</b>		Run ID: <b>VOA4_354794</b>		SeqNo: <b>5443124</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	17.07	1.0	20	0	85.3	70 - 127	17.06	0.0484	20
Benzene	17.72	1.0	20	0	88.6	70 - 127	17.8	0.415	20
Chlorobenzene	17.36	1.0	20	0	86.8	70 - 114	17.78	2.36	20
Ethylbenzene	17.78	1.0	20	0	88.9	70 - 124	18.09	1.74	20
Methylene chloride	15.15	2.0	20	0	75.8	70 - 128	16.05	5.74	20
Toluene	17.64	1.0	20	0	88.2	70 - 123	17.93	1.64	20
Vinyl chloride	21.02	1.0	20	0.477	103	70 - 130	22.17	5.33	20
Xylenes, Total	55.22	1.0	60	0	92.0	70 - 130	56.44	2.18	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>70 - 126</i>	<i>49.09</i>	<i>1.85</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.69</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>81 - 113</i>	<i>49.59</i>	<i>0.216</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>49.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>77 - 123</i>	<i>49.87</i>	<i>0.535</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>50.93</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 127</i>	<i>50.43</i>	<i>0.976</i>	<i>20</i>

The following samples were analyzed in this batch: HS20010767-01      HS20010767-02      HS20010767-03

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> R354799 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200121</b>	Units: <b>ug/L</b>			Analysis Date: <b>21-Jan-2020 12:55</b>				
Client ID:	Run ID: <b>VOA2_354799</b>	SeqNo: <b>5443190</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.22</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>90.4</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.91</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200221</b>	Units: <b>ug/L</b>			Analysis Date: <b>21-Jan-2020 12:06</b>				
Client ID:	Run ID: <b>VOA2_354799</b>	SeqNo: <b>5443189</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	17.66	1.0	20	0	88.3	70 - 124			
Benzene	17.2	1.0	20	0	86.0	74 - 120			
Chlorobenzene	18.44	1.0	20	0	92.2	76 - 113			
Ethylbenzene	17.94	1.0	20	0	89.7	77 - 117			
Methylene chloride	17.49	2.0	20	0	87.4	70 - 127			
Toluene	19.3	1.0	20	0	96.5	77 - 118			
Vinyl chloride	15.17	1.0	20	0	75.8	70 - 130			
Xylenes, Total	54.82	1.0	60	0	91.4	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.76</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>81 - 120</i>			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** R354799 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010767-05MS			Units: ug/L		Analysis Date: 21-Jan-2020 17:48			
Client ID: WG-1620-MW84B-20200116		Run ID: VOA2_354799			SeqNo: 5443202		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.36	1.0	20	0	81.8	70 - 127				
Benzene	24.4	1.0	20	6.383	90.1	70 - 127				
Chlorobenzene	17.96	1.0	20	0	89.8	70 - 114				
Ethylbenzene	28.72	1.0	20	9.361	96.8	70 - 124				
Methylene chloride	15.89	2.0	20	0	79.4	70 - 128				
Toluene	20.5	1.0	20	0	102	70 - 123				
Vinyl chloride	17.82	1.0	20	0	89.1	70 - 130				
Xylenes, Total	57.71	1.0	60	3.181	90.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.1</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.0</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.2</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.4</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010767-05MSD			Units: ug/L		Analysis Date: 21-Jan-2020 18:12			
Client ID: WG-1620-MW84B-20200116		Run ID: VOA2_354799			SeqNo: 5443203		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.51	1.0	20	0	82.6	70 - 127	16.36	0.908	20	
Benzene	24.24	1.0	20	6.383	89.3	70 - 127	24.4	0.678	20	
Chlorobenzene	17.68	1.0	20	0	88.4	70 - 114	17.96	1.54	20	
Ethylbenzene	28.32	1.0	20	9.361	94.8	70 - 124	28.72	1.41	20	
Methylene chloride	15.24	2.0	20	0	76.2	70 - 128	15.89	4.15	20	
Toluene	20.18	1.0	20	0	101	70 - 123	20.5	1.56	20	
Vinyl chloride	16.8	1.0	20	0	84.0	70 - 130	17.82	5.89	20	
Xylenes, Total	56.81	1.0	60	3.181	89.4	70 - 130	57.71	1.57	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.11</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.2</i>	<i>70 - 126</i>	<i>46.04</i>	<i>0.147</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>81 - 113</i>	<i>48.8</i>	<i>0.867</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.8</i>	<i>77 - 123</i>	<i>47.49</i>	<i>0.191</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.1</i>	<i>82 - 127</i>	<i>49.2</i>	<i>0.339</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS20010767-04	HS20010767-05	HS20010767-06	HS20010767-08
HS20010767-09	HS20010767-11	HS20010767-12	HS20010767-13
HS20010767-14	HS20010767-15		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** R354802 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200121</b>		Units: <b>ug/L</b>		Analysis Date: <b>22-Jan-2020 00:44</b>			
Client ID:		Run ID: <b>VOA2_354802</b>		SeqNo: <b>5443239</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>45.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>90.0</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.5</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.3</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.91</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>81 - 120</i>			

<b>LCS</b>		Sample ID: <b>VLCSW-200121</b>		Units: <b>ug/L</b>		Analysis Date: <b>21-Jan-2020 23:55</b>			
Client ID:		Run ID: <b>VOA2_354802</b>		SeqNo: <b>5443238</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	17.23	1.0	20	0	86.2	70 - 124			
Benzene	17.15	1.0	20	0	85.8	74 - 120			
Chlorobenzene	18.1	1.0	20	0	90.5	76 - 113			
Ethylbenzene	17.55	1.0	20	0	87.7	77 - 117			
Methylene chloride	18.45	2.0	20	0	92.3	70 - 127			
Toluene	19.21	1.0	20	0	96.0	77 - 118			
Xylenes, Total	53.45	1.0	60	0	89.1	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.2</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>47.27</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.5</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.15</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** R354802 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010767-18MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>22-Jan-2020 02:21</b>			
Client ID: <b>WG-1620-MW65D-20200116</b>		Run ID: <b>VOA2_354802</b>			SeqNo: <b>5443243</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.51	1.0	20	0	82.6	70 - 127				
Benzene	16.52	1.0	20	0	82.6	70 - 127				
Chlorobenzene	17.7	1.0	20	0	88.5	70 - 114				
Ethylbenzene	17.86	1.0	20	0	89.3	70 - 124				
Methylene chloride	15.38	2.0	20	0	76.9	70 - 128				
Toluene	18.94	1.0	20	0	94.7	70 - 123				
Xylenes, Total	53.57	1.0	60	0	89.3	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.68</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.4</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010767-18MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>22-Jan-2020 02:46</b>			
Client ID: <b>WG-1620-MW65D-20200116</b>		Run ID: <b>VOA2_354802</b>			SeqNo: <b>5443244</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.43	1.0	20	0	82.2	70 - 127	16.51	0.508	20	
Benzene	16.5	1.0	20	0	82.5	70 - 127	16.52	0.127	20	
Chlorobenzene	17.09	1.0	20	0	85.5	70 - 114	17.7	3.52	20	
Ethylbenzene	17.06	1.0	20	0	85.3	70 - 124	17.86	4.58	20	
Methylene chloride	15.4	2.0	20	0	77.0	70 - 128	15.38	0.0821	20	
Toluene	18.34	1.0	20	0	91.7	70 - 123	18.94	3.23	20	
Xylenes, Total	51.58	1.0	60	0	86.0	70 - 130	53.57	3.78	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>46.38</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>92.8</i>	<i>70 - 126</i>	<i>46.68</i>	<i>0.626</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>81 - 113</i>	<i>49.26</i>	<i>0.814</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>47.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>77 - 123</i>	<i>48.04</i>	<i>0.511</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>48.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>82 - 127</i>	<i>49.45</i>	<i>1.73</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS20010767-07	HS20010767-10	HS20010767-16	HS20010767-17
HS20010767-18	HS20010767-19	HS20010767-20	HS20010767-21
HS20010767-22	HS20010767-23	HS20010767-24	HS20010767-25
HS20010767-26	HS20010767-27	HS20010767-28	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> R354859 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200122</b>	Units: <b>ug/L</b>			Analysis Date: <b>22-Jan-2020 13:01</b>				
Client ID:	Run ID: <b>VOA4_354859</b>	SeqNo: <b>5444319</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

Toluene	U	1.0							
Vinyl chloride	U	1.0							
Surr: 1,2-Dichloroethane-d4	49.62	1.0	50	0	99.2	70 - 123			
Surr: 4-Bromofluorobenzene	48.7	1.0	50	0	97.4	82 - 115			
Surr: Dibromofluoromethane	48.78	1.0	50	0	97.6	73 - 126			
Surr: Toluene-d8	51.64	1.0	50	0	103	81 - 120			

<b>LCS</b>	Sample ID: <b>VLCSW-200122</b>	Units: <b>ug/L</b>			Analysis Date: <b>22-Jan-2020 12:12</b>				
Client ID:	Run ID: <b>VOA4_354859</b>	SeqNo: <b>5444318</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
Toluene	20.43	1.0	20	0	102	77 - 118			
Vinyl chloride	20.42	1.0	20	0	102	70 - 130			
Surr: 1,2-Dichloroethane-d4	49.37	1.0	50	0	98.7	70 - 130			
Surr: 4-Bromofluorobenzene	50.11	1.0	50	0	100	82 - 115			
Surr: Dibromofluoromethane	49.33	1.0	50	0	98.7	73 - 126			
Surr: Toluene-d8	50.79	1.0	50	0	102	81 - 120			

<b>MS</b>	Sample ID: <b>HS20010767-29MS</b>	Units: <b>ug/L</b>			Analysis Date: <b>22-Jan-2020 15:31</b>				
Client ID: <b>WG-1620-MW61A-20200116</b>	Run ID: <b>VOA4_354859</b>	SeqNo: <b>5444325</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
Toluene	18.26	1.0	20	0	91.3	70 - 123			
Vinyl chloride	23.34	1.0	20	0	117	70 - 130			
Surr: 1,2-Dichloroethane-d4	48.8	1.0	50	0	97.6	70 - 126			
Surr: 4-Bromofluorobenzene	49.91	1.0	50	0	99.8	81 - 113			
Surr: Dibromofluoromethane	49.62	1.0	50	0	99.2	77 - 123			
Surr: Toluene-d8	51.16	1.0	50	0	102	82 - 127			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** R354859 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MSD</b>		Sample ID: <b>HS20010767-29MSD</b>		Units: <b>ug/L</b>		Analysis Date: <b>22-Jan-2020 15:56</b>			
Client ID: <b>WG-1620-MW61A-20200116</b>		Run ID: <b>VOA4_354859</b>		SeqNo: <b>5444326</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
Toluene	17.93	1.0	20	0	89.7	70 - 123	18.26	1.81	20
Vinyl chloride	22.63	1.0	20	0	113	70 - 130	23.34	3.09	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.43</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>70 - 126</i>	<i>48.8</i>	<i>1.28</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 113</i>	<i>49.91</i>	<i>0.267</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>49.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>77 - 123</i>	<i>49.62</i>	<i>0.325</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>51.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 127</i>	<i>51.16</i>	<i>0.159</i>	<i>20</i>

The following samples were analyzed in this batch: HS20010767-10    HS20010767-23    HS20010767-24    HS20010767-26

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

<b>Batch ID:</b> R355066 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200125</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 11:48</b>				
Client ID:	Run ID: <b>VOA2_355066</b>	SeqNo: <b>5447954</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>51.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200225</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 10:59</b>				
Client ID:	Run ID: <b>VOA2_355066</b>	SeqNo: <b>5447953</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18.57	1.0	20	0	92.9	70 - 124			
Benzene	18.24	1.0	20	0	91.2	74 - 120			
Chlorobenzene	18.75	1.0	20	0	93.8	76 - 113			
Ethylbenzene	18.83	1.0	20	0	94.1	77 - 117			
Methylene chloride	18.77	2.0	20	0	93.8	70 - 127			
Toluene	20.88	1.0	20	0	104	77 - 118			
Vinyl chloride	19.59	1.0	20	0	98.0	70 - 130			
Xylenes, Total	58.21	1.0	60	0	97.0	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.59</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QC BATCH REPORT**

**Batch ID:** R355066 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010767-29MS			Units: ug/L		Analysis Date: 25-Jan-2020 13:28			
Client ID: WG-1620-MW61A-20200116		Run ID: VOA2_355066			SeqNo: 5447958		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.72	1.0	20	0	88.6	70 - 127				
Benzene	17.87	1.0	20	0	89.3	70 - 127				
Chlorobenzene	18.14	1.0	20	0	90.7	70 - 114				
Ethylbenzene	18.64	1.0	20	0	93.2	70 - 124				
Methylene chloride	18.41	2.0	20	0	92.0	70 - 128				
Toluene	20.39	1.0	20	0	102	70 - 123				
Vinyl chloride	18.29	1.0	20	0	91.4	70 - 130				
Xylenes, Total	57.1	1.0	60	0	95.2	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.56</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.1</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010767-29MSD			Units: ug/L		Analysis Date: 25-Jan-2020 13:52			
Client ID: WG-1620-MW61A-20200116		Run ID: VOA2_355066			SeqNo: 5447959		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.67	1.0	20	0	88.4	70 - 127	17.72	0.267	20	
Benzene	17.9	1.0	20	0	89.5	70 - 127	17.87	0.197	20	
Chlorobenzene	18.23	1.0	20	0	91.2	70 - 114	18.14	0.504	20	
Ethylbenzene	18.75	1.0	20	0	93.8	70 - 124	18.64	0.579	20	
Methylene chloride	17.79	2.0	20	0	88.9	70 - 128	18.41	3.41	20	
Toluene	20.63	1.0	20	0	103	70 - 123	20.39	1.13	20	
Vinyl chloride	18.66	1.0	20	0	93.3	70 - 130	18.29	2	20	
Xylenes, Total	57.29	1.0	60	0	95.5	70 - 130	57.1	0.342	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>	<i>51.56</i>	<i>1.47</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>81 - 113</i>	<i>50.55</i>	<i>1.63</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>77 - 123</i>	<i>50.1</i>	<i>0.63</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>49.72</i>	<i>1.16</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010767-29

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010767

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter



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**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

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<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010767

**SAMPLE TRACKING**

---

Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS20010767-01	WG-1620-MW28A-20200116	Login	1/17/2020 2:56:40 PM	JRM	MET002
HS20010767-01	WG-1620-MW28A-20200116	Login	1/17/2020 2:56:40 PM	JRM	EXT090
HS20010767-01	WG-1620-MW28A-20200116	Login	1/17/2020 2:56:40 PM	JRM	VOA239

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010767

Date/Time Received: 17-Jan-2020 13:11
Received by: PMG

Checklist completed by: Jared R. Makan
eSignature
Date: 17-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 21-Jan-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition?
Custody seals intact on shipping container/cooler?
Custody seals intact on sample bottles?
VOA/TX1005/TX1006 Solids in hermetically sealed vials?
Chain of custody present?
Chain of custody signed when relinquished and received?
Samplers name present on COC?
Chain of custody agrees with sample labels?
Samples in proper container/bottle?
Sample containers intact?
Sufficient sample volume for indicated test?
All samples received within holding time?
Container/Temp Blank temperature in compliance?
Temperature(s)/Thermometer(s):

- Yes/No/Not Present checkboxes for each item in the list above.

4 Page(s)
COC IDs:206390, 215871, 215867, 215868

Table with 2 columns: Temperature(s)/Thermometer(s) and Corrected Temp. Values include 1.7°C, 1.8°C, 1.3°C, 2.0°C, 1.3°C, 1.4°C, 1.7°C and IR25.

Cooler(s)/Kit(s):

45153, 43016, 5236, 42700, 45137, 45594, 45067

Date/Time sample(s) sent to storage:

01/07/2020 19:30

Water - VOA vials have zero headspace?

- Yes/No checkboxes and No VOA vials submitted checkbox.

Water - pH acceptable upon receipt?

- Yes/No checkboxes and N/A checkbox.

pH adjusted?

- Yes/No checkboxes and N/A checkbox.

pH adjusted by:

[Empty text box]

Login Notes:

All VOCs received in cooler 45153 w/Trip Blank.
Samples received, not listed on COC:
MW48C - 01/16/20 10:20
MW47C - 01/16/20 11:05
MW61A (MS & MSD) - 01/16/20 12:00
Samples logged for analysis. IDs added to COC by ALS

Client Contacted:

Date Contacted:

Person Contacted:

Contacted By:

Regarding:

Comments:

[Empty text box for comments]

Corrective Action:

[Empty text box for corrective action]



Cincinnati, OH  
+1 513 733 5336

Everett, WA  
+1 425 356 2600

Fort Collins, CO  
+1 970 490 1511

Holland, MI  
+1 616 399 6070

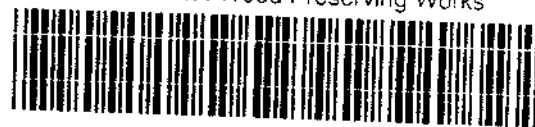
# Chain of Custody Form

Page 1 of 4

COC ID: 206390

HS20010767

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WAQ-1620-TBO-202001			Water	1	2		X									
2	WG-1620-MW28A-20200116	1-16-20	0910	W		6	X		X	X							
3	WG-1620-MW28C-20200116		1015	W		6	X		X	X							
4	WG-1620-MW28CMS-20200116		1015	W		6	X		X	X							
5	WG-1620-MW28MSD-20200116		1015	W		6	X		X	X							
6	WG-1620-FB08-20200116		1200	W		6	X		X	X							
7	WG-1620-MW89B-20200116		1245	W		6	X		X	X							
8	WG-1620-84B-20200116		1345	W		6	X		X	X							
9	WG-1620-63B-20200116		1445	W		6	X		X	X							
10	WG-1620-MW26A-20200116		1600	W		6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BEAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24-hour		Results Due Date:	
Relinquished by: <b>John Bry</b>	Date: <b>1-17-20</b>	Time: <b>1311</b>	Received by: <b>[Signature]</b>	Notes: <b>UPRR Houston MWPW</b>			
Relinquished by:	Date:	Time:	Received by (Laboratory): <b>[Signature]</b>	Cooler ID: <b>45236</b>	Cooler Temp: <b>1.3°</b>	QC Package: (Check One Box Below)	
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory): <b>[Signature]</b>	<b>43016</b>	<b>1.8°</b>	<input type="checkbox"/> Level II Std QC	<input checked="" type="checkbox"/> TRRP Check st
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				<b>45153</b>	<b>1.7°</b>	<input type="checkbox"/> Level III Str. QC/Raw Data	<input type="checkbox"/> TRRP Level IV
						<input type="checkbox"/> Level IV SW643/CLP	<input type="checkbox"/> Other

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
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# Chain of Custody Form

Page 2 of 4

COC ID: 215871

HS20010767

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:											
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A											
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 Volatile Organics Site Specific)										
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8260_LL_W (5632528 VOC Site Specific + V.C.)										
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	8270_LOW W (5632532 SemiVolatiles Site specific)										
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	ICP_TW (5636002 5652646 Metals - As)										
	Suite 4004		Stop 0750	F											
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G											
Phone	(512) 671-3434	Phone		H											
Fax	(512) 671-3446	Fax		I											
e-Mail Address	eric.matzner@pbwilc.com	e-Mail Address		J											

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TBO-202001			Water	1	2		X									
2	WG-1620-MW68C-20200117	1-17-20	0910	W		6	X		X	X							
3	WG-1620-MW68A-20200117	↓	1015	W		6	X		X	X							
4	WG-1620-PB09-20200117	↓	1000	W		6	X		X	X							
5	WG-1620-MW27A-20200115	1-15-20	1640	W		6	X		X	X							
6	WG-1620-MW27C-20200115	↓	1720	W		6	X		X	X							
7	WG-1620-MW59A-20200116	1-16-20	0805	W		6		X	X	X							
8	WG-1620-MW59B-20200116	↓	0900	W		6		X	X	X							
9	WG-1620-MW59D-20200116	↓	1350	W		6	X		X	X							
10	WG-1620-FD03-20200116	↓	1350	W		6	X		X	X							

Sampler(s) Please Print & Sign <b>JOHN BRAYTON John Br</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box)				Results Due Date:		
Relinquished by: <i>[Signature]</i>		Date: 1-17-20	Time: 1311	<input checked="" type="checkbox"/> STD 10 Wk Days		<input type="checkbox"/> 5 Wk Days		<input type="checkbox"/> 24 Hour		
Relinquished by:		Date:	Time:	Received by (Laboratory): <i>[Signature]</i>		Notes: UPRR Houston MWPW				
Logged by (Laboratory):		Date:	Time:	Checked by (Laboratory):		Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)		
						42700	2.00	<input type="checkbox"/> Level II Std CC	<input checked="" type="checkbox"/> TRRP Checklist	
						45137	1.30	<input type="checkbox"/> Level III Std QC/Raw Data	<input type="checkbox"/> TRRP Level IV	
						45594	1.40	<input type="checkbox"/> Level IV SIA618/CLP	<input type="checkbox"/> Other	

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

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# Chain of Custody Form

Page 3 of 4

COC ID: 215867

## HS20010767

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As)
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E	TPH TX 1005
				F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TBO-202001</del>			Water	1	2		X									
2	WG-1620-MW66D-20200116	1-16-20	1500	W		6	X		X	X							
3	WG-1620-MW65D-2020116		1620	W		6	X		X	X							
4	WG-1620-MW65DMS-2020116		1620	W		6	X		X	X							
5	WG-1620-MW65DMSD-20200116		1620	W		6	X		X	X							
6	WG-1620-MW36D-20200116		1725	W		6	X		X	X							
7	WG-1620-MW86C-20200117	1-17-20	0725	W		9	X		X	X	X						
8	WG-1620- <del>MW</del> FDO4-20200117		0725	W		9	X		X	X	X						
9	WG-1620-MW75B-20200117		0815	W		9	X		X	X	X						
10	WG-1620-MW79A-20200117		0920	W		9	X		X	X	X						

Sampler(s) Please Print & Sign <i>[Signature]</i>		Shipment Method HAND DELIVERED		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour			Results Due Date:	
Relinquished by: <i>[Signature]</i>		Date: 1-17-20	Time: 13:11	Received by: <i>[Signature]</i>		Notes: UPRR Houston MWPPW		
Relinquished by: <i>[Signature]</i>		Date:	Time:	Received by (Laboratory): <i>[Signature]</i>		Cooler ID:	Cooler Temp.:	QC Package: (Check One Box Below)
Logged by (Laboratory):		Date:	Time:	Checked by (Laboratory):		<input type="checkbox"/> Level II Std QC	<input checked="" type="checkbox"/> TRRP Checklist	
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035						<input type="checkbox"/> Level III Std QC/Raw Data	<input type="checkbox"/> TRRP Level IV	
						<input type="checkbox"/> Level IV SW846/CLP		
						<input type="checkbox"/> Other		

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# Chain of Custody Form

Page 4 of 4

COC ID: 215868

HS20010767

Golder Associates Inc.  
Houston TX-Wood Preserving Works



Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address	

A	8260_LL_W(5632528 Volatile Organics Site Specific)
B	8260_LL_W(5632528 VOC Site Specific + V.C.)
C	8270_LOW_W(5632532 SemiVolatiles Site specific)
D	ICP_TW(5636002 5652646 Metals - As)
E	TPH TX1005
F	MS   MSD
G	
H	
I	
J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	<del>WG-1620-TBO-202001</del>			Water	1	2		X									
2	WG-1620-MW74B-20200117	1-17-20	1020			9	X		X	X	X						
3	WG-1620-MW78A-20200117		1125			9	X		X	X	X						
4	WG-1620-TBOS-2020117					2	X										
5	WG-1620-MW61A-20200116	1-16-20	1200	Ground Water	1,2,8	18		X	X	X		X					
6	WG-1620-MW47C-20200116	1-16-20	1105	Ground Water	1,2,8	6	X		X	X							
7	WG-1620-MW48C-20200116	1-16-20	1020	Ground Water	1,2,8	6	X		X	X							
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>John Brauch</i>		Shipment Method HAND DELIVERED		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour				Results Due Date:			
Relinquished by: <i>John Brauch</i>	Date: 1-17-20	Time: 13:11	Received by:	Notes: UPRR Houston MWPW						QC Package: (Check One Box Below)	
Relinquished by:	Date:	Time:	Received by (Laboratory):	Cooler ID	Cooler Temp.	<input type="checkbox"/> Level II Std QC		<input checked="" type="checkbox"/> TRRF Checklist			
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):			<input type="checkbox"/> Level III Std QC/Raw Date		<input type="checkbox"/> TRRF Level IV			
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035						<input type="checkbox"/> Level IV SW846/CLP					

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January 29, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010930**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 10 sample(s) on Jan 21, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

---

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/29/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010930			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149963,149998,R355061,R355066,R355150			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?		X			3
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				4
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/29/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010930			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149963,149998,R355061,R355066,R355150			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section)					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 01/29/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20010930
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 149963,149998,R355061,R355066,R355150

ER# <sup>5</sup>	Description
1	Semivolatile Organics Method SW8270, sample WG-1620-MW32B-20200120, WG-1620-MW70B-20200120; the surrogate recoveries could not be determined due to dilution below the calibration range in the higher dilutions.
2	Batch 149998, Semivolatile Organics Method SW8260, LCS recovery was above the control limits for 2,4-Dinitrotoluene and 2,6-Dinitrotoluene. The analytes were not detected in the associated samples.
3	Batch 149998, Semivolatile Organics Method SW8260, sample HS20010910-03, MS/MSD RPD is for an unrelated sample.
4	Batch R355150, Volatile Organics Method SW8260, samples WG-1620-MW32B-20200120 and WG-1620-MW70B-20200120: Lowest practical dilution of 25x/50x performed due to high concentration of non-target analyte(s).  Batch 149998, Semivolatile Organics Method SW8270, samples WG-1620-MW32B-20200120, WG-1620-MW70B-20200120; The GCMS semi-volatile extract of the samples were run at a dilution due to a high level of matrix interference.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010930

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010930-01	WQ-1620-TB06-20200120	Water	CG-121719 -108	20-Jan-2020 00:00	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-02	WG-1620-MW32AR-20200120	Water		20-Jan-2020 09:10	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-03	WG-1620-MW32B-20200120	Water		20-Jan-2020 10:15	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-04	WG-1620-MW70B-20200120	Water		20-Jan-2020 11:30	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-05	WG-1620-MW33A-20200120	Water		20-Jan-2020 12:45	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-06	WG-1620-MW87C-20200120	Water		20-Jan-2020 14:00	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-07	WG-1620-MW90B-20200120	Water		20-Jan-2020 15:15	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-08	WG-1620-MW45C-20200120	Water		20-Jan-2020 16:30	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-09	WG-1620-FB10-20200120	Water		20-Jan-2020 09:30	21-Jan-2020 17:10	<input type="checkbox"/>
HS20010930-10	WG-1620-FD05-20200120	Water		20-Jan-2020 00:00	21-Jan-2020 17:10	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB06-20200120  
 Collection Date: 20-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:41
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:41
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 12:41
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 12:41
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 12:41
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:41
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 12:41
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 12:41
<i>Surr: 1,2-Dichloroethane-d4</i>		103		70-126	%REC	1	25-Jan-2020 12:41
<i>Surr: 4-Bromofluorobenzene</i>		97.3		81-113	%REC	1	25-Jan-2020 12:41
<i>Surr: Dibromofluoromethane</i>		102		77-123	%REC	1	25-Jan-2020 12:41
<i>Surr: Toluene-d8</i>		102		82-127	%REC	1	25-Jan-2020 12:41

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32AR-20200120  
 Collection Date: 20-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:22
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:22
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:22
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:22
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 19:22
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:22
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:22
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:22
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>97.3</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:22</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>96.7</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:22</i>
<i>Surr: Dibromofluoromethane</i>		<i>101</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:22</i>
<i>Surr: Toluene-d8</i>		<i>101</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 19:22</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32AR-20200120  
 Collection Date: 20-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	29-Jan-2020 00:59
2,4-Dimethylphenol		U	0.000040	0.00020	mg/L	1	29-Jan-2020 00:59
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	29-Jan-2020 00:59
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	29-Jan-2020 00:59
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	29-Jan-2020 00:59
<b>2-Methylnaphthalene</b>	<b>0.000028</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	29-Jan-2020 00:59
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	29-Jan-2020 00:59
<b>Acenaphthene</b>	<b>0.0014</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
Acenaphthylene		U	0.000015	0.00010	mg/L	1	29-Jan-2020 00:59
Anthracene		U	0.000014	0.00010	mg/L	1	29-Jan-2020 00:59
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	29-Jan-2020 00:59
<b>Benzo(a)pyrene</b>	<b>0.000081</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	29-Jan-2020 00:59
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000043</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
<b>Chrysene</b>	<b>0.000042</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
Dibenzofuran		U	0.000020	0.00010	mg/L	1	29-Jan-2020 00:59
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	29-Jan-2020 00:59
<b>Fluoranthene</b>	<b>0.00029</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
<b>Fluorene</b>	<b>0.000038</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
<b>Naphthalene</b>	<b>0.00014</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
Nitrobenzene		U	0.000024	0.00020	mg/L	1	29-Jan-2020 00:59
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	29-Jan-2020 00:59
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	29-Jan-2020 00:59
<b>Phenanthrene</b>	<b>0.000032</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
Phenol		U	0.000035	0.00020	mg/L	1	29-Jan-2020 00:59
<b>Pyrene</b>	<b>0.00042</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 00:59
<i>Surr: 2,4,6-Tribromophenol</i>	75.9			34-129	%REC	1	29-Jan-2020 00:59
<i>Surr: 2-Fluorobiphenyl</i>	86.4			40-125	%REC	1	29-Jan-2020 00:59
<i>Surr: 2-Fluorophenol</i>	72.7			20-120	%REC	1	29-Jan-2020 00:59
<i>Surr: 4-Terphenyl-d14</i>	99.3			40-135	%REC	1	29-Jan-2020 00:59
<i>Surr: Nitrobenzene-d5</i>	75.2			41-120	%REC	1	29-Jan-2020 00:59
<i>Surr: Phenol-d6</i>	70.0			20-120	%REC	1	29-Jan-2020 00:59
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00770</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 23:38

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32B-20200120  
 Collection Date: 20-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0050	0.025	mg/L	25	28-Jan-2020 08:52
<b>Benzene</b>	<b>1.8</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	28-Jan-2020 08:52
Chlorobenzene	U		0.0075	0.025	mg/L	25	28-Jan-2020 08:52
<b>Ethylbenzene</b>	<b>0.73</b>		<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	28-Jan-2020 08:52
Methylene chloride	U		0.025	0.050	mg/L	25	28-Jan-2020 08:52
<b>Toluene</b>	<b>2.3</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	28-Jan-2020 08:52
Vinyl chloride	U		0.0050	0.025	mg/L	25	28-Jan-2020 08:52
<b>Xylenes, Total</b>	<b>2.1</b>		<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	28-Jan-2020 08:52
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	25	<i>28-Jan-2020 08:52</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>104</i>			<i>81-113</i>	<i>%REC</i>	25	<i>28-Jan-2020 08:52</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.9</i>			<i>77-123</i>	<i>%REC</i>	25	<i>28-Jan-2020 08:52</i>
<i>Surr: Toluene-d8</i>	<i>95.4</i>			<i>82-127</i>	<i>%REC</i>	25	<i>28-Jan-2020 08:52</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32B-20200120  
 Collection Date: 20-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.00042	0.0040	mg/L	10	29-Jan-2020 13:29
<b>2,4-Dimethylphenol</b>	<b>26</b>		<b>0.16</b>	<b>0.80</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
2,4-Dinitrotoluene	U		0.0012	0.0040	mg/L	10	29-Jan-2020 13:29
2,6-Dinitrotoluene	U		0.00084	0.0040	mg/L	10	29-Jan-2020 13:29
2-Chloronaphthalene	U		0.00042	0.0040	mg/L	10	29-Jan-2020 13:29
<b>2-Methylnaphthalene</b>	<b>50</b>		<b>0.76</b>	<b>4.0</b>	<b>mg/L</b>	20000	29-Jan-2020 15:40
4,6-Dinitro-2-methylphenol	U		0.00040	0.0040	mg/L	10	29-Jan-2020 13:29
4-Nitrophenol	U		0.00094	0.020	mg/L	10	29-Jan-2020 13:29
<b>Acenaphthene</b>	<b>26</b>		<b>0.11</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
Acenaphthylene	U		0.00030	0.0020	mg/L	10	29-Jan-2020 13:29
<b>Anthracene</b>	<b>23</b>		<b>0.056</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
<b>Benz(a)anthracene</b>	<b>2.2</b>		<b>0.020</b>	<b>0.040</b>	<b>mg/L</b>	200	29-Jan-2020 14:44
<b>Benzo(a)pyrene</b>	<b>0.74</b>		<b>0.0080</b>	<b>0.040</b>	<b>mg/L</b>	200	29-Jan-2020 14:44
Bis(2-chloroethoxy)methane	U		0.00060	0.0040	mg/L	10	29-Jan-2020 13:29
Bis(2-ethylhexyl)phthalate	U		0.00074	0.0040	mg/L	10	29-Jan-2020 13:29
<b>Chrysene</b>	<b>2.4</b>		<b>0.0084</b>	<b>0.040</b>	<b>mg/L</b>	200	29-Jan-2020 14:44
<b>Dibenzofuran</b>	<b>28</b>		<b>0.080</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
Di-n-butyl phthalate	U		0.00040	0.0040	mg/L	10	29-Jan-2020 13:29
<b>Fluoranthene</b>	<b>21</b>		<b>0.040</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
<b>Fluorene</b>	<b>23</b>		<b>0.12</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
<b>Naphthalene</b>	<b>300</b>		<b>0.80</b>	<b>4.0</b>	<b>mg/L</b>	20000	29-Jan-2020 15:40
Nitrobenzene	U		0.00048	0.0040	mg/L	10	29-Jan-2020 13:29
N-Nitrosodiphenylamine	U		0.00050	0.0040	mg/L	10	29-Jan-2020 13:29
Pentachlorophenol	U		0.0016	0.0040	mg/L	10	29-Jan-2020 13:29
<b>Phenanthrene</b>	<b>69</b>		<b>0.84</b>	<b>4.0</b>	<b>mg/L</b>	20000	29-Jan-2020 15:40
<b>Phenol</b>	<b>17</b>		<b>0.14</b>	<b>0.80</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
<b>Pyrene</b>	<b>14</b>		<b>0.076</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:03
<i>Surr: 2,4,6-Tribromophenol</i>	<i>83.0</i>			<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>29-Jan-2020 13:29</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>29-Jan-2020 14:44</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>29-Jan-2020 15:03</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>20000</i>	<i>29-Jan-2020 15:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>112</i>			<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>29-Jan-2020 13:29</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>29-Jan-2020 14:44</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>29-Jan-2020 15:03</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>20000</i>	<i>29-Jan-2020 15:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>20000</i>	<i>29-Jan-2020 15:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>98.6</i>			<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>29-Jan-2020 13:29</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>29-Jan-2020 14:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>29-Jan-2020 15:03</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW32B-20200120  
 Collection Date: 20-Jan-2020 10:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	79.5	J		40-135	%REC	10	29-Jan-2020 13:29
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	29-Jan-2020 14:44
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	29-Jan-2020 15:03
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	20000	29-Jan-2020 15:40
Surr: Nitrobenzene-d5	84.4			41-120	%REC	10	29-Jan-2020 13:29
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	29-Jan-2020 14:44
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	29-Jan-2020 15:03
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	20000	29-Jan-2020 15:40
Surr: Phenol-d6	0	JS		20-120	%REC	2000	29-Jan-2020 15:03
Surr: Phenol-d6	0	JS		20-120	%REC	20000	29-Jan-2020 15:40
Surr: Phenol-d6	105			20-120	%REC	10	29-Jan-2020 13:29
Surr: Phenol-d6	0	JS		20-120	%REC	200	29-Jan-2020 14:44
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
Arsenic	0.00193	J	0.000400	0.00200	mg/L	1	27-Jan-2020 14:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW70B-20200120  
 Collection Date: 20-Jan-2020 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.010	0.050	mg/L	50	28-Jan-2020 09:19
<b>Benzene</b>	<b>1.9</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	50	28-Jan-2020 09:19
Chlorobenzene	U		0.015	0.050	mg/L	50	28-Jan-2020 09:19
<b>Ethylbenzene</b>	<b>0.57</b>		<b>0.015</b>	<b>0.050</b>	<b>mg/L</b>	50	28-Jan-2020 09:19
Methylene chloride	U		0.050	0.10	mg/L	50	28-Jan-2020 09:19
<b>Toluene</b>	<b>2.1</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	50	28-Jan-2020 09:19
Vinyl chloride	U		0.010	0.050	mg/L	50	28-Jan-2020 09:19
<b>Xylenes, Total</b>	<b>1.5</b>		<b>0.015</b>	<b>0.050</b>	<b>mg/L</b>	50	28-Jan-2020 09:19
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>103</i>			<i>70-126</i>	<i>%REC</i>	<i>50</i>	<i>28-Jan-2020 09:19</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>50</i>	<i>28-Jan-2020 09:19</i>
<i>Surr: Dibromofluoromethane</i>	<i>95.7</i>			<i>77-123</i>	<i>%REC</i>	<i>50</i>	<i>28-Jan-2020 09:19</i>
<i>Surr: Toluene-d8</i>	<i>97.6</i>			<i>82-127</i>	<i>%REC</i>	<i>50</i>	<i>28-Jan-2020 09:19</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW70B-20200120  
 Collection Date: 20-Jan-2020 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.00042	0.0040	mg/L	10	29-Jan-2020 14:06
<b>2,4-Dimethylphenol</b>	<b>49</b>		<b>0.64</b>	<b>3.2</b>	<b>mg/L</b>	8000	29-Jan-2020 15:59
2,4-Dinitrotoluene	U		0.0012	0.0040	mg/L	10	29-Jan-2020 14:06
2,6-Dinitrotoluene	U		0.00084	0.0040	mg/L	10	29-Jan-2020 14:06
2-Chloronaphthalene	U		0.00042	0.0040	mg/L	10	29-Jan-2020 14:06
<b>2-Methylnaphthalene</b>	<b>10</b>		<b>0.076</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
4,6-Dinitro-2-methylphenol	U		0.00040	0.0040	mg/L	10	29-Jan-2020 14:06
4-Nitrophenol	U		0.00094	0.020	mg/L	10	29-Jan-2020 14:06
<b>Acenaphthene</b>	<b>7.1</b>		<b>0.11</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
<b>Acenaphthylene</b>	<b>0.079</b>		<b>0.00030</b>	<b>0.0020</b>	<b>mg/L</b>	10	29-Jan-2020 14:06
<b>Anthracene</b>	<b>5.8</b>		<b>0.056</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
<b>Benz(a)anthracene</b>	<b>0.52</b>		<b>0.020</b>	<b>0.040</b>	<b>mg/L</b>	200	29-Jan-2020 14:25
<b>Benzo(a)pyrene</b>	<b>0.13</b>		<b>0.00040</b>	<b>0.0020</b>	<b>mg/L</b>	10	29-Jan-2020 14:06
Bis(2-chloroethoxy)methane	U		0.00060	0.0040	mg/L	10	29-Jan-2020 14:06
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0039</b>	J	<b>0.00074</b>	<b>0.0040</b>	<b>mg/L</b>	10	29-Jan-2020 14:06
<b>Chrysene</b>	<b>0.63</b>		<b>0.0084</b>	<b>0.040</b>	<b>mg/L</b>	200	29-Jan-2020 14:25
<b>Dibenzofuran</b>	<b>6.2</b>		<b>0.080</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
Di-n-butyl phthalate	U		0.00040	0.0040	mg/L	10	29-Jan-2020 14:06
<b>Fluoranthene</b>	<b>4.7</b>		<b>0.040</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
<b>Fluorene</b>	<b>6.7</b>		<b>0.12</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
<b>Naphthalene</b>	<b>71</b>		<b>0.32</b>	<b>1.6</b>	<b>mg/L</b>	8000	29-Jan-2020 15:59
Nitrobenzene	U		0.00048	0.0040	mg/L	10	29-Jan-2020 14:06
N-Nitrosodiphenylamine	U		0.00050	0.0040	mg/L	10	29-Jan-2020 14:06
Pentachlorophenol	U		0.0016	0.0040	mg/L	10	29-Jan-2020 14:06
<b>Phenanthrene</b>	<b>14</b>		<b>0.084</b>	<b>0.40</b>	<b>mg/L</b>	2000	29-Jan-2020 15:21
<b>Phenol</b>	<b>3.4</b>		<b>0.014</b>	<b>0.080</b>	<b>mg/L</b>	200	29-Jan-2020 14:25
<b>Pyrene</b>	<b>3.2</b>		<b>0.0076</b>	<b>0.040</b>	<b>mg/L</b>	200	29-Jan-2020 14:25
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	2000	29-Jan-2020 15:21
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	8000	29-Jan-2020 15:59
<i>Surr: 2,4,6-Tribromophenol</i>	82.9			34-129	%REC	10	29-Jan-2020 14:06
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	200	29-Jan-2020 14:25
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	200	29-Jan-2020 14:25
<i>Surr: 2-Fluorobiphenyl</i>	98.3			40-125	%REC	10	29-Jan-2020 14:06
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	2000	29-Jan-2020 15:21
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	8000	29-Jan-2020 15:59
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	2000	29-Jan-2020 15:21
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	8000	29-Jan-2020 15:59
<i>Surr: 2-Fluorophenol</i>	94.2			20-120	%REC	10	29-Jan-2020 14:06
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	200	29-Jan-2020 14:25

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW70B-20200120  
 Collection Date: 20-Jan-2020 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	29-Jan-2020 14:25
Surr: 4-Terphenyl-d14	77.9	J		40-135	%REC	10	29-Jan-2020 14:06
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	29-Jan-2020 15:21
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	8000	29-Jan-2020 15:59
Surr: Nitrobenzene-d5	82.7			41-120	%REC	10	29-Jan-2020 14:06
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	29-Jan-2020 14:25
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	29-Jan-2020 15:21
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	8000	29-Jan-2020 15:59
Surr: Phenol-d6	0	JS		20-120	%REC	200	29-Jan-2020 14:25
Surr: Phenol-d6	0	JS		20-120	%REC	2000	29-Jan-2020 15:21
Surr: Phenol-d6	0	JS		20-120	%REC	8000	29-Jan-2020 15:59
Surr: Phenol-d6	98.9			20-120	%REC	10	29-Jan-2020 14:06
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
Arsenic	0.00120	J	0.000400	0.00200	mg/L	1	27-Jan-2020 14:43

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33A-20200120  
 Collection Date: 20-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:47
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:47
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:47
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:47
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 19:47
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:47
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:47
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:47
<i>Surr: 1,2-Dichloroethane-d4</i>		96.7		70-126	%REC	1	25-Jan-2020 19:47
<i>Surr: 4-Bromofluorobenzene</i>		95.9		81-113	%REC	1	25-Jan-2020 19:47
<i>Surr: Dibromofluoromethane</i>		100		77-123	%REC	1	25-Jan-2020 19:47
<i>Surr: Toluene-d8</i>		101		82-127	%REC	1	25-Jan-2020 19:47

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW33A-20200120  
 Collection Date: 20-Jan-2020 12:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	29-Jan-2020 01:18
2,4-Dimethylphenol		U	0.000040	0.00020	mg/L	1	29-Jan-2020 01:18
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	29-Jan-2020 01:18
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	29-Jan-2020 01:18
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	29-Jan-2020 01:18
<b>2-Methylnaphthalene</b>	<b>0.000063</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	29-Jan-2020 01:18
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	29-Jan-2020 01:18
<b>Acenaphthene</b>	<b>0.00073</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Acenaphthylene</b>	<b>0.000049</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Anthracene</b>	<b>0.000057</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Benz(a)anthracene</b>	<b>0.00012</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Benzo(a)pyrene</b>	<b>0.000066</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	29-Jan-2020 01:18
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	29-Jan-2020 01:18
<b>Chrysene</b>	<b>0.000078</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Dibenzofuran</b>	<b>0.000096</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	29-Jan-2020 01:18
<b>Fluoranthene</b>	<b>0.00035</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Fluorene</b>	<b>0.000059</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<b>Naphthalene</b>	<b>0.00050</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
Nitrobenzene		U	0.000024	0.00020	mg/L	1	29-Jan-2020 01:18
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	29-Jan-2020 01:18
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	29-Jan-2020 01:18
<b>Phenanthrene</b>	<b>0.000056</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
Phenol		U	0.000035	0.00020	mg/L	1	29-Jan-2020 01:18
<b>Pyrene</b>	<b>0.00019</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:18
<i>Surr: 2,4,6-Tribromophenol</i>	<i>81.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 01:18</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>101</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 01:18</i>
<i>Surr: 2-Fluorophenol</i>	<i>75.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 01:18</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>100</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 01:18</i>
<i>Surr: Nitrobenzene-d5</i>	<i>86.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 01:18</i>
<i>Surr: Phenol-d6</i>	<i>73.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 01:18</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00810</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	24-Jan-2020 23:58

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW87C-20200120  
 Collection Date: 20-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:11
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:11
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:11
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:11
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 20:11
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:11
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:11
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:11
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:11</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:11</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:11</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:11</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW87C-20200120  
 Collection Date: 20-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	29-Jan-2020 01:37
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	29-Jan-2020 01:37
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	29-Jan-2020 01:37
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	29-Jan-2020 01:37
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	29-Jan-2020 01:37
<b>2-Methylnaphthalene</b>	<b>0.00023</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	29-Jan-2020 01:37
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	29-Jan-2020 01:37
<b>Acenaphthene</b>	<b>0.00011</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
Acenaphthylene	U		0.000015	0.00010	mg/L	1	29-Jan-2020 01:37
<b>Anthracene</b>	<b>0.000088</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	29-Jan-2020 01:37
<b>Benzo(a)pyrene</b>	<b>0.000066</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	29-Jan-2020 01:37
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	29-Jan-2020 01:37
Chrysene	U		0.000021	0.00010	mg/L	1	29-Jan-2020 01:37
<b>Dibenzofuran</b>	<b>0.00011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	29-Jan-2020 01:37
<b>Fluoranthene</b>	<b>0.000086</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
<b>Fluorene</b>	<b>0.000093</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
<b>Naphthalene</b>	<b>0.0011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
Nitrobenzene	U		0.000024	0.00020	mg/L	1	29-Jan-2020 01:37
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	29-Jan-2020 01:37
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	29-Jan-2020 01:37
<b>Phenanthrene</b>	<b>0.00028</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
<b>Phenol</b>	<b>0.00048</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
<b>Pyrene</b>	<b>0.000057</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:37
<i>Surr: 2,4,6-Tribromophenol</i>	69.4			34-129	%REC	1	29-Jan-2020 01:37
<i>Surr: 2-Fluorobiphenyl</i>	94.4			40-125	%REC	1	29-Jan-2020 01:37
<i>Surr: 2-Fluorophenol</i>	79.6			20-120	%REC	1	29-Jan-2020 01:37
<i>Surr: 4-Terphenyl-d14</i>	99.6			40-135	%REC	1	29-Jan-2020 01:37
<i>Surr: Nitrobenzene-d5</i>	85.2			41-120	%REC	1	29-Jan-2020 01:37
<i>Surr: Phenol-d6</i>	78.9			20-120	%REC	1	29-Jan-2020 01:37
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00135</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	25-Jan-2020 00:00

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW90B-20200120  
 Collection Date: 20-Jan-2020 15:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:36
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:36
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:36
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:36
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 20:36
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:36
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:36
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:36
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>96.4</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:36</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.7</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:36</i>
<i>Surr: Dibromofluoromethane</i>		<i>101</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:36</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:36</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW90B-20200120  
 Collection Date: 20-Jan-2020 15:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	29-Jan-2020 01:55
<b>2,4-Dimethylphenol</b>	<b>0.000075</b>	J	<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	29-Jan-2020 01:55
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	29-Jan-2020 01:55
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	29-Jan-2020 01:55
<b>2-Methylnaphthalene</b>	<b>0.00034</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	29-Jan-2020 01:55
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	29-Jan-2020 01:55
<b>Acenaphthene</b>	<b>0.00015</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
Acenaphthylene		U	0.000015	0.00010	mg/L	1	29-Jan-2020 01:55
<b>Anthracene</b>	<b>0.00014</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
Benz(a)anthracene		U	0.000050	0.00010	mg/L	1	29-Jan-2020 01:55
<b>Benzo(a)pyrene</b>	<b>0.000052</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	29-Jan-2020 01:55
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	29-Jan-2020 01:55
Chrysene		U	0.000021	0.00010	mg/L	1	29-Jan-2020 01:55
<b>Dibenzofuran</b>	<b>0.00011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	29-Jan-2020 01:55
<b>Fluoranthene</b>	<b>0.00012</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
<b>Fluorene</b>	<b>0.00013</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
<b>Naphthalene</b>	<b>0.0030</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
Nitrobenzene		U	0.000024	0.00020	mg/L	1	29-Jan-2020 01:55
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	29-Jan-2020 01:55
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	29-Jan-2020 01:55
<b>Phenanthrene</b>	<b>0.00040</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
<b>Phenol</b>	<b>0.00019</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
<b>Pyrene</b>	<b>0.000077</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 01:55
Surr: 2,4,6-Tribromophenol	67.0			34-129	%REC	1	29-Jan-2020 01:55
Surr: 2-Fluorobiphenyl	90.7			40-125	%REC	1	29-Jan-2020 01:55
Surr: 2-Fluorophenol	68.7			20-120	%REC	1	29-Jan-2020 01:55
Surr: 4-Terphenyl-d14	108			40-135	%REC	1	29-Jan-2020 01:55
Surr: Nitrobenzene-d5	79.6			41-120	%REC	1	29-Jan-2020 01:55
Surr: Phenol-d6	73.0			20-120	%REC	1	29-Jan-2020 01:55
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00290</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	25-Jan-2020 00:02

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW45C-20200120  
 Collection Date: 20-Jan-2020 16:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 21:00
<b>Benzene</b>	<b>0.00050</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	25-Jan-2020 21:00
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 21:00
<b>Ethylbenzene</b>	<b>0.00044</b>	J	<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	25-Jan-2020 21:00
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 21:00
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 21:00
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 21:00
<b>Xylenes, Total</b>	<b>0.0051</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	25-Jan-2020 21:00
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:00</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:00</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:00</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 21:00</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW45C-20200120  
 Collection Date: 20-Jan-2020 16:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-08  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	0.000085	J	0.000021	0.00020	mg/L	1	29-Jan-2020 13:10
2,4-Dimethylphenol		U	0.000040	0.00020	mg/L	1	29-Jan-2020 13:10
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	29-Jan-2020 13:10
2,6-Dinitrotoluene	0.0022		0.000042	0.00020	mg/L	1	29-Jan-2020 13:10
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	29-Jan-2020 13:10
2-Methylnaphthalene	0.00013		0.000019	0.00010	mg/L	1	29-Jan-2020 13:10
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	29-Jan-2020 13:10
4-Nitrophenol	0.000067	J	0.000047	0.0010	mg/L	1	29-Jan-2020 13:10
Acenaphthene	0.000084	J	0.000027	0.00010	mg/L	1	29-Jan-2020 13:10
Acenaphthylene	0.000037	J	0.000015	0.00010	mg/L	1	29-Jan-2020 13:10
Anthracene	0.00090		0.000014	0.00010	mg/L	1	29-Jan-2020 13:10
Benz(a)anthracene	0.00039		0.000050	0.00010	mg/L	1	29-Jan-2020 13:10
Benzo(a)pyrene	0.00042		0.000020	0.00010	mg/L	1	29-Jan-2020 13:10
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	29-Jan-2020 13:10
Bis(2-ethylhexyl)phthalate	0.000067	J	0.000037	0.00020	mg/L	1	29-Jan-2020 13:10
Chrysene	0.00062		0.000021	0.00010	mg/L	1	29-Jan-2020 13:10
Dibenzofuran	0.000075	J	0.000020	0.00010	mg/L	1	29-Jan-2020 13:10
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	29-Jan-2020 13:10
Fluoranthene	0.00062		0.000010	0.00010	mg/L	1	29-Jan-2020 13:10
Fluorene	0.00023		0.000030	0.00010	mg/L	1	29-Jan-2020 13:10
Naphthalene	0.00061		0.000020	0.00010	mg/L	1	29-Jan-2020 13:10
Nitrobenzene		U	0.000024	0.00020	mg/L	1	29-Jan-2020 13:10
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	29-Jan-2020 13:10
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	29-Jan-2020 13:10
Phenanthrene	0.00027		0.000021	0.00010	mg/L	1	29-Jan-2020 13:10
Phenol		U	0.000035	0.00020	mg/L	1	29-Jan-2020 13:10
Pyrene	0.00068		0.000019	0.00010	mg/L	1	29-Jan-2020 13:10
Surr: 2,4,6-Tribromophenol	85.3			34-129	%REC	1	29-Jan-2020 13:10
Surr: 2-Fluorobiphenyl	95.0			40-125	%REC	1	29-Jan-2020 13:10
Surr: 2-Fluorophenol	75.8			20-120	%REC	1	29-Jan-2020 13:10
Surr: 4-Terphenyl-d14	112			40-135	%REC	1	29-Jan-2020 13:10
Surr: Nitrobenzene-d5	91.7			41-120	%REC	1	29-Jan-2020 13:10
Surr: Phenol-d6	81.6			20-120	%REC	1	29-Jan-2020 13:10
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
Arsenic	0.000730	J	0.000400	0.00200	mg/L	1	25-Jan-2020 00:05

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB10-20200120  
 Collection Date: 20-Jan-2020 09:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	SQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 13:06
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 13:06
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 13:06
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 13:06
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 13:06
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 13:06
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 13:06
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 13:06
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>104</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 13:06</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>97.7</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 13:06</i>
<i>Surr: Dibromofluoromethane</i>		<i>99.2</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 13:06</i>
<i>Surr: Toluene-d8</i>		<i>101</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 13:06</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB10-20200120  
 Collection Date: 20-Jan-2020 09:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-09  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
<b>1,2-Diphenylhydrazine</b>	<b>0.000047</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 02:33
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	29-Jan-2020 02:33
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	29-Jan-2020 02:33
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	29-Jan-2020 02:33
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	29-Jan-2020 02:33
<b>2-Methylnaphthalene</b>	<b>0.000035</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:33
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	29-Jan-2020 02:33
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	29-Jan-2020 02:33
Acenaphthene	U		0.000027	0.00010	mg/L	1	29-Jan-2020 02:33
Acenaphthylene	U		0.000015	0.00010	mg/L	1	29-Jan-2020 02:33
Anthracene	U		0.000014	0.00010	mg/L	1	29-Jan-2020 02:33
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	29-Jan-2020 02:33
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	29-Jan-2020 02:33
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	29-Jan-2020 02:33
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	29-Jan-2020 02:33
Chrysene	U		0.000021	0.00010	mg/L	1	29-Jan-2020 02:33
Dibenzofuran	U		0.000020	0.00010	mg/L	1	29-Jan-2020 02:33
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	29-Jan-2020 02:33
<b>Fluoranthene</b>	<b>0.000060</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:33
Fluorene	U		0.000030	0.00010	mg/L	1	29-Jan-2020 02:33
<b>Naphthalene</b>	<b>0.00019</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:33
Nitrobenzene	U		0.000024	0.00020	mg/L	1	29-Jan-2020 02:33
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	29-Jan-2020 02:33
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	29-Jan-2020 02:33
<b>Phenanthrene</b>	<b>0.000066</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:33
Phenol	U		0.000035	0.00020	mg/L	1	29-Jan-2020 02:33
<b>Pyrene</b>	<b>0.000040</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:33
Surr: 2,4,6-Tribromophenol	68.3			34-129	%REC	1	29-Jan-2020 02:33
Surr: 2-Fluorobiphenyl	95.0			40-125	%REC	1	29-Jan-2020 02:33
Surr: 2-Fluorophenol	88.9			20-120	%REC	1	29-Jan-2020 02:33
Surr: 4-Terphenyl-d14	106			40-135	%REC	1	29-Jan-2020 02:33
Surr: Nitrobenzene-d5	92.6			41-120	%REC	1	29-Jan-2020 02:33
Surr: Phenol-d6	85.1			20-120	%REC	1	29-Jan-2020 02:33
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
Arsenic	U		0.000400	0.00200	mg/L	1	25-Jan-2020 00:07

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD05-20200120  
 Collection Date: 20-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:44
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:44
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:44
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:44
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 19:44
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:44
Vinyl chloride	U		0.00020	0.0010	mg/L	1	25-Jan-2020 19:44
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 19:44
<i>Surr: 1,2-Dichloroethane-d4</i>		106		70-126	%REC	1	25-Jan-2020 19:44
<i>Surr: 4-Bromofluorobenzene</i>		96.6		81-113	%REC	1	25-Jan-2020 19:44
<i>Surr: Dibromofluoromethane</i>		103		77-123	%REC	1	25-Jan-2020 19:44
<i>Surr: Toluene-d8</i>		101		82-127	%REC	1	25-Jan-2020 19:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD05-20200120  
 Collection Date: 20-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010930  
 Lab ID:HS20010930-10  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine		U	0.000021	0.00020	mg/L	1	29-Jan-2020 02:52
2,4-Dimethylphenol		U	0.000040	0.00020	mg/L	1	29-Jan-2020 02:52
2,4-Dinitrotoluene		U	0.000058	0.00020	mg/L	1	29-Jan-2020 02:52
2,6-Dinitrotoluene		U	0.000042	0.00020	mg/L	1	29-Jan-2020 02:52
2-Chloronaphthalene		U	0.000021	0.00020	mg/L	1	29-Jan-2020 02:52
<b>2-Methylnaphthalene</b>	<b>0.000095</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
4,6-Dinitro-2-methylphenol		U	0.000020	0.00020	mg/L	1	29-Jan-2020 02:52
4-Nitrophenol		U	0.000047	0.0010	mg/L	1	29-Jan-2020 02:52
<b>Acenaphthene</b>	<b>0.0011</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Acenaphthylene</b>	<b>0.000038</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Anthracene</b>	<b>0.000044</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Benz(a)anthracene</b>	<b>0.00012</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Benzo(a)pyrene</b>	<b>0.000072</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
Bis(2-chloroethoxy)methane		U	0.000030	0.00020	mg/L	1	29-Jan-2020 02:52
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	29-Jan-2020 02:52
<b>Chrysene</b>	<b>0.000093</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Dibenzofuran</b>	<b>0.000067</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
Di-n-butyl phthalate		U	0.000020	0.00020	mg/L	1	29-Jan-2020 02:52
<b>Fluoranthene</b>	<b>0.00057</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Fluorene</b>	<b>0.000072</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<b>Naphthalene</b>	<b>0.00066</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
Nitrobenzene		U	0.000024	0.00020	mg/L	1	29-Jan-2020 02:52
N-Nitrosodiphenylamine		U	0.000025	0.00020	mg/L	1	29-Jan-2020 02:52
Pentachlorophenol		U	0.000079	0.00020	mg/L	1	29-Jan-2020 02:52
<b>Phenanthrene</b>	<b>0.000046</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
Phenol		U	0.000035	0.00020	mg/L	1	29-Jan-2020 02:52
<b>Pyrene</b>	<b>0.00052</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 02:52
<i>Surr: 2,4,6-Tribromophenol</i>	63.6			34-129	%REC	1	29-Jan-2020 02:52
<i>Surr: 2-Fluorobiphenyl</i>	80.8			40-125	%REC	1	29-Jan-2020 02:52
<i>Surr: 2-Fluorophenol</i>	72.7			20-120	%REC	1	29-Jan-2020 02:52
<i>Surr: 4-Terphenyl-d14</i>	98.1			40-135	%REC	1	29-Jan-2020 02:52
<i>Surr: Nitrobenzene-d5</i>	86.5			41-120	%REC	1	29-Jan-2020 02:52
<i>Surr: Phenol-d6</i>	77.2			20-120	%REC	1	29-Jan-2020 02:52
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00755</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	25-Jan-2020 00:09

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**Batch ID:** 149963      **Start Date:** 24 Jan 2020 09:00      **End Date:** 24 Jan 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010930-02		10 (mL)	10 (mL)	1
HS20010930-03		10 (mL)	10 (mL)	1
HS20010930-04		10 (mL)	10 (mL)	1
HS20010930-05		10 (mL)	10 (mL)	1
HS20010930-06		10 (mL)	10 (mL)	1
HS20010930-07		10 (mL)	10 (mL)	1
HS20010930-08		10 (mL)	10 (mL)	1
HS20010930-09		10 (mL)	10 (mL)	1
HS20010930-10		10 (mL)	10 (mL)	1

**Batch ID:** 149998      **Start Date:** 27 Jan 2020 07:00      **End Date:** 27 Jan 2020 13:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010930-02	1	1000 (mL)	1 (mL)	0.001
HS20010930-03	1	1000 (mL)	2 (mL)	0.002
HS20010930-04	1	1000 (mL)	2 (mL)	0.002
HS20010930-05	1	1000 (mL)	1 (mL)	0.001
HS20010930-06	1	1000 (mL)	1 (mL)	0.001
HS20010930-07	1	1000 (mL)	1 (mL)	0.001
HS20010930-08	1	1000 (mL)	1 (mL)	0.001
HS20010930-09	1	1000 (mL)	1 (mL)	0.001
HS20010930-10	1	1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149963 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010930-02	WG-1620-MW32AR-20200120	20 Jan 2020 09:10		24 Jan 2020 13:00	24 Jan 2020 23:38	1
HS20010930-03	WG-1620-MW32B-20200120	20 Jan 2020 10:15		24 Jan 2020 13:00	27 Jan 2020 14:37	1
HS20010930-04	WG-1620-MW70B-20200120	20 Jan 2020 11:30		24 Jan 2020 13:00	27 Jan 2020 14:43	1
HS20010930-05	WG-1620-MW33A-20200120	20 Jan 2020 12:45		24 Jan 2020 13:00	24 Jan 2020 23:58	1
HS20010930-06	WG-1620-MW87C-20200120	20 Jan 2020 14:00		24 Jan 2020 13:00	25 Jan 2020 00:00	1
HS20010930-07	WG-1620-MW90B-20200120	20 Jan 2020 15:15		24 Jan 2020 13:00	25 Jan 2020 00:02	1
HS20010930-08	WG-1620-MW45C-20200120	20 Jan 2020 16:30		24 Jan 2020 13:00	25 Jan 2020 00:05	1
HS20010930-09	WG-1620-FB10-20200120	20 Jan 2020 09:30		24 Jan 2020 13:00	25 Jan 2020 00:07	1
HS20010930-10	WG-1620-FD05-20200120	20 Jan 2020 00:00		24 Jan 2020 13:00	25 Jan 2020 00:09	1
<b>Batch ID: 149998 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010930-02	WG-1620-MW32AR-20200120	20 Jan 2020 09:10		27 Jan 2020 10:49	29 Jan 2020 00:59	1
HS20010930-03	WG-1620-MW32B-20200120	20 Jan 2020 10:15		27 Jan 2020 10:49	29 Jan 2020 15:03	2000
HS20010930-03	WG-1620-MW32B-20200120	20 Jan 2020 10:15		27 Jan 2020 10:49	29 Jan 2020 15:40	2000 0
HS20010930-03	WG-1620-MW32B-20200120	20 Jan 2020 10:15		27 Jan 2020 10:49	29 Jan 2020 14:44	200
HS20010930-03	WG-1620-MW32B-20200120	20 Jan 2020 10:15		27 Jan 2020 10:49	29 Jan 2020 13:29	10
HS20010930-04	WG-1620-MW70B-20200120	20 Jan 2020 11:30		27 Jan 2020 10:49	29 Jan 2020 15:59	8000
HS20010930-04	WG-1620-MW70B-20200120	20 Jan 2020 11:30		27 Jan 2020 10:49	29 Jan 2020 15:21	2000
HS20010930-04	WG-1620-MW70B-20200120	20 Jan 2020 11:30		27 Jan 2020 10:49	29 Jan 2020 14:25	200
HS20010930-04	WG-1620-MW70B-20200120	20 Jan 2020 11:30		27 Jan 2020 10:49	29 Jan 2020 14:06	10
HS20010930-05	WG-1620-MW33A-20200120	20 Jan 2020 12:45		27 Jan 2020 10:49	29 Jan 2020 01:18	1
HS20010930-06	WG-1620-MW87C-20200120	20 Jan 2020 14:00		27 Jan 2020 10:49	29 Jan 2020 01:37	1
HS20010930-07	WG-1620-MW90B-20200120	20 Jan 2020 15:15		27 Jan 2020 10:49	29 Jan 2020 01:55	1
HS20010930-08	WG-1620-MW45C-20200120	20 Jan 2020 16:30		27 Jan 2020 10:49	29 Jan 2020 13:10	1
HS20010930-09	WG-1620-FB10-20200120	20 Jan 2020 09:30		27 Jan 2020 10:49	29 Jan 2020 02:33	1
HS20010930-10	WG-1620-FD05-20200120	20 Jan 2020 00:00		27 Jan 2020 10:49	29 Jan 2020 02:52	1
<b>Batch ID: R355061 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010930-01	WQ-1620-TB06-20200120	20 Jan 2020 00:00			25 Jan 2020 12:41	1
HS20010930-09	WG-1620-FB10-20200120	20 Jan 2020 09:30			25 Jan 2020 13:06	1
HS20010930-10	WG-1620-FD05-20200120	20 Jan 2020 00:00			25 Jan 2020 19:44	1
<b>Batch ID: R355066 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010930-02	WG-1620-MW32AR-20200120	20 Jan 2020 09:10			25 Jan 2020 19:22	1
HS20010930-05	WG-1620-MW33A-20200120	20 Jan 2020 12:45			25 Jan 2020 19:47	1
HS20010930-06	WG-1620-MW87C-20200120	20 Jan 2020 14:00			25 Jan 2020 20:11	1
HS20010930-07	WG-1620-MW90B-20200120	20 Jan 2020 15:15			25 Jan 2020 20:36	1
HS20010930-08	WG-1620-MW45C-20200120	20 Jan 2020 16:30			25 Jan 2020 21:00	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID:</b> R355150 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Water	
HS20010930-03	WG-1620-MW32B-20200120	20 Jan 2020 10:15			28 Jan 2020 08:52	25
HS20010930-04	WG-1620-MW70B-20200120	20 Jan 2020 11:30			28 Jan 2020 09:19	50

WorkOrder: HS20010930  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20010930  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020



WorkOrder: HS20010930  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00045	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

WorkOrder: HS20010930  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

Batch ID: 149963 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-149963</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 23:34</b>					
Client ID:		Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447747</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	U	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-149963</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 23:36</b>					
Client ID:		Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447748</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.043	0.00200	0.05	0	86.0	80 - 120			
<b>MS</b>	Sample ID: <b>HS20010930-02MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 23:42</b>					
Client ID: <b>WG-1620-MW32AR-20200120</b>		Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447751</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05468	0.00200	0.05	0.007701	94.0	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20010930-02MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 23:45</b>					
Client ID: <b>WG-1620-MW32AR-20200120</b>		Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447752</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.05348	0.00200	0.05	0.007701	91.6	80 - 120	0.05468	2.22	20
<b>PDS</b>	Sample ID: <b>HS20010930-02PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 23:47</b>					
Client ID: <b>WG-1620-MW32AR-20200120</b>		Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447753</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual
Arsenic	0.09766	0.00200	0.1	0.007701	90.0	75 - 125			
<b>SD</b>	Sample ID: <b>HS20010930-02SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>24-Jan-2020 23:40</b>					
Client ID: <b>WG-1620-MW32AR-20200120</b>		Run ID: <b>ICPMS05_354996</b>	SeqNo: <b>5447750</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual
Arsenic	0.007533	0.0100					0.007701	0 10	J

The following samples were analyzed in this batch:

HS20010930-02	HS20010930-03	HS20010930-04	HS20010930-05
HS20010930-06	HS20010930-07	HS20010930-08	HS20010930-09
HS20010930-10			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149998	Units: ug/L			Analysis Date: 28-Jan-2020 19:21					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452454	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.165</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.567</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.869</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.797</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>116</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.964</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.941</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149998	Units: ug/L			Analysis Date: 28-Jan-2020 19:40					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452455	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.783	0.20	5	0	116	39 - 127				
2,4-Dimethylphenol	5.012	0.20	5	0	100	35 - 120				
2,4-Dinitrotoluene	6.415	0.20	5	0	128	50 - 122				S
2,6-Dinitrotoluene	6.03	0.20	5	0	121	50 - 120				S
2-Chloronaphthalene	5.424	0.20	5	0	108	50 - 120				
2-Methylnaphthalene	5.383	0.10	5	0	108	50 - 120				
4,6-Dinitro-2-methylphenol	5.692	0.20	5	0	114	25 - 121				
4-Nitrophenol	5.835	1.0	5	0	117	30 - 130				
Acenaphthene	5.432	0.10	5	0	109	45 - 120				
Acenaphthylene	5.667	0.10	5	0	113	47 - 120				
Anthracene	5.847	0.10	5	0	117	45 - 120				
Benz(a)anthracene	5.919	0.10	5	0	118	40 - 120				
Benzo(a)pyrene	5.832	0.10	5	0	117	45 - 120				
Bis(2-chloroethoxy)methane	5.315	0.20	5	0	106	45 - 120				
Bis(2-ethylhexyl)phthalate	5.645	0.20	5	0	113	40 - 139				
Chrysene	5.825	0.10	5	0	117	43 - 120				
Dibenzofuran	5.475	0.10	5	0	109	50 - 120				
Di-n-butyl phthalate	5.996	0.20	5	0	120	45 - 123				
Fluoranthene	5.728	0.10	5	0	115	45 - 125				
Fluorene	5.748	0.10	5	0	115	49 - 120				
Naphthalene	5.43	0.10	5	0	109	45 - 120				
Nitrobenzene	5.209	0.20	5	0	104	44 - 120				
N-Nitrosodiphenylamine	5.868	0.20	5	0	117	40 - 125				
Pentachlorophenol	2.037	0.20	5	0	40.7	19 - 121				
Phenanthrene	5.866	0.10	5	0	117	45 - 121				
Phenol	5.563	0.20	5	0	111	20 - 124				
Pyrene	5.924	0.10	5	0	118	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.466</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.582</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>112</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.846</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.9</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.609</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>112</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.978</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.891</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS	Sample ID: HS20010910-03MS	Units: ug/L			Analysis Date: 28-Jan-2020 23:44					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452170	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.258	0.20	5	0	85.2	39 - 127				
2,4-Dimethylphenol	2.953	0.20	5	0.09533	57.1	35 - 120				
2,4-Dinitrotoluene	5.24	0.20	5	0	105	50 - 122				
2,6-Dinitrotoluene	4.544	0.20	5	0	90.9	50 - 120				
2-Chloronaphthalene	3.621	0.20	5	0	72.4	50 - 120				
2-Methylnaphthalene	3.197	0.10	5	0	63.9	50 - 120				
4,6-Dinitro-2-methylphenol	5.327	0.20	5	0	107	25 - 121				
4-Nitrophenol	5.515	1.0	5	0	110	30 - 130				
Acenaphthene	3.484	0.10	5	0	69.7	45 - 120				
Acenaphthylene	3.505	0.10	5	0	70.1	47 - 120				
Anthracene	5.162	0.10	5	0	103	45 - 120				
Benz(a)anthracene	5.438	0.10	5	0	109	40 - 120				
Benzo(a)pyrene	5.398	0.10	5	0	108	45 - 120				
Bis(2-chloroethoxy)methane	2.957	0.20	5	0	59.1	45 - 120				
Bis(2-ethylhexyl)phthalate	5.76	0.20	5	0	115	40 - 139				
Chrysene	5.195	0.10	5	0	104	43 - 120				
Dibenzofuran	3.564	0.10	5	0	71.3	50 - 120				
Di-n-butyl phthalate	6.041	0.20	5	0	121	45 - 123				
Fluoranthene	5.294	0.10	5	0.02986	105	45 - 125				
Fluorene	3.974	0.10	5	0	79.5	49 - 120				
Naphthalene	3.352	0.10	5	0.22	62.6	45 - 120				
Nitrobenzene	3.206	0.20	5	0	64.1	44 - 120				
N-Nitrosodiphenylamine	5.464	0.20	5	0	109	40 - 125				
Pentachlorophenol	3.381	0.20	5	0	67.6	19 - 121				
Phenanthrene	4.995	0.10	5	0	99.9	45 - 121				
Phenol	3.121	0.20	5	0	62.4	20 - 124				
Pyrene	5.526	0.10	5	0.02733	110	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.181</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.6</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.463</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.1</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>62.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.541</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.186</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.7</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.21</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>64.2</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS20010910-03MSD	Units: ug/L			Analysis Date: 29-Jan-2020 00:03					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452171	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.047	0.20	5	0	101	39 - 127	4.258	17	20	
2,4-Dimethylphenol	3.862	0.20	5	0.09533	75.3	35 - 120	2.953	26.7	20	R
2,4-Dinitrotoluene	5.46	0.20	5	0	109	50 - 122	5.24	4.12	20	
2,6-Dinitrotoluene	5.056	0.20	5	0	101	50 - 120	4.544	10.7	20	
2-Chloronaphthalene	4.215	0.20	5	0	84.3	50 - 120	3.621	15.2	20	
2-Methylnaphthalene	3.899	0.10	5	0	78.0	50 - 120	3.197	19.8	20	
4,6-Dinitro-2-methylphenol	5.595	0.20	5	0	112	25 - 121	5.327	4.9	30	
4-Nitrophenol	5.641	1.0	5	0	113	30 - 130	5.515	2.26	20	
Acenaphthene	4.245	0.10	5	0	84.9	45 - 120	3.484	19.7	20	
Acenaphthylene	4.266	0.10	5	0	85.3	47 - 120	3.505	19.6	20	
Anthracene	5.342	0.10	5	0	107	45 - 120	5.162	3.43	20	
Benz(a)anthracene	5.45	0.10	5	0	109	40 - 120	5.438	0.221	20	
Benzo(a)pyrene	5.228	0.10	5	0	105	45 - 120	5.398	3.2	20	
Bis(2-chloroethoxy)methane	3.686	0.20	5	0	73.7	45 - 120	2.957	21.9	20	R
Bis(2-ethylhexyl)phthalate	5.666	0.20	5	0	113	40 - 139	5.76	1.64	20	
Chrysene	5.136	0.10	5	0	103	43 - 120	5.195	1.14	20	
Dibenzofuran	4.314	0.10	5	0	86.3	50 - 120	3.564	19	20	
Di-n-butyl phthalate	6.065	0.20	5	0	121	45 - 123	6.041	0.393	20	
Fluoranthene	5.068	0.10	5	0.02986	101	45 - 125	5.294	4.37	20	
Fluorene	4.577	0.10	5	0	91.5	49 - 120	3.974	14.1	20	
Naphthalene	4.118	0.10	5	0.22	78.0	45 - 120	3.352	20.5	20	R
Nitrobenzene	3.792	0.20	5	0	75.8	44 - 120	3.206	16.7	20	
N-Nitrosodiphenylamine	5.62	0.20	5	0	112	40 - 125	5.464	2.8	20	
Pentachlorophenol	3.323	0.20	5	0	66.5	19 - 121	3.381	1.74	20	
Phenanthrene	5.142	0.10	5	0	103	45 - 121	4.995	2.9	20	
Phenol	3.883	0.20	5	0	77.7	20 - 124	3.121	21.7	20	R
Pyrene	5.443	0.10	5	0.02733	108	40 - 130	5.526	1.51	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.11</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>82.2</i>	<i>34 - 129</i>	<i>4.181</i>	<i>1.73</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.108</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>82.2</i>	<i>40 - 125</i>	<i>3.463</i>	<i>17</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.476</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.5</i>	<i>20 - 120</i>	<i>3.1</i>	<i>11.4</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>5.339</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>107</i>	<i>40 - 135</i>	<i>5.541</i>	<i>3.71</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.75</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>75.0</i>	<i>41 - 120</i>	<i>3.186</i>	<i>16.3</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.663</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.3</i>	<i>20 - 120</i>	<i>3.21</i>	<i>13.2</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

<b>Batch ID:</b> 149998 ( 0 )	<b>Instrument:</b> SV-6	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS20010930-02	HS20010930-03	HS20010930-04	HS20010930-05
HS20010930-06	HS20010930-07	HS20010930-08	HS20010930-09
HS20010930-10			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

<b>Batch ID:</b> R355061 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200125</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 11:52</b>				
Client ID:	Run ID: <b>VOA4_355061</b>	SeqNo: <b>5447903</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.99</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>48.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200125</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 11:02</b>				
Client ID:	Run ID: <b>VOA4_355061</b>	SeqNo: <b>5447902</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	18.78	1.0	20	0	93.9	70 - 124			
Benzene	19.4	1.0	20	0	97.0	74 - 120			
Chlorobenzene	19.61	1.0	20	0	98.0	76 - 113			
Ethylbenzene	19.57	1.0	20	0	97.8	77 - 117			
Methylene chloride	20.36	2.0	20	0	102	70 - 127			
Toluene	19.83	1.0	20	0	99.1	77 - 118			
Vinyl chloride	19.46	1.0	20	0	97.3	70 - 130			
Xylenes, Total	58.33	1.0	60	0	97.2	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

**Batch ID:** R355061 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010937-01MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>25-Jan-2020 15:09</b>			
Client ID:		Run ID: <b>VOA4_355061</b>			SeqNo: <b>5447911</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.24	1.0	20	0	81.2	70 - 127				
Benzene	17.71	1.0	20	0	88.5	70 - 127				
Chlorobenzene	17.17	1.0	20	0	85.8	70 - 114				
Ethylbenzene	17.53	1.0	20	0	87.6	70 - 124				
Methylene chloride	18.41	2.0	20	0	92.1	70 - 128				
Toluene	18.12	1.0	20	0	90.6	70 - 123				
Vinyl chloride	17.88	1.0	20	0	89.4	70 - 130				
Xylenes, Total	51.84	1.0	60	0	86.4	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>51.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010937-01MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>25-Jan-2020 15:34</b>			
Client ID:		Run ID: <b>VOA4_355061</b>			SeqNo: <b>5447912</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.9	1.0	20	0	79.5	70 - 127	16.24	2.11	20	
Benzene	17.13	1.0	20	0	85.6	70 - 127	17.71	3.33	20	
Chlorobenzene	17.02	1.0	20	0	85.1	70 - 114	17.17	0.844	20	
Ethylbenzene	16.92	1.0	20	0	84.6	70 - 124	17.53	3.5	20	
Methylene chloride	17.25	2.0	20	0	86.3	70 - 128	18.41	6.49	20	
Toluene	16.89	1.0	20	0	84.4	70 - 123	18.12	7.08	20	
Vinyl chloride	16.93	1.0	20	0	84.6	70 - 130	17.88	5.47	20	
Xylenes, Total	50.49	1.0	60	0	84.1	70 - 130	51.84	2.64	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>	<i>51.17</i>	<i>0.402</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.63</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>49.86</i>	<i>3.48</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>51.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>77 - 123</i>	<i>50.64</i>	<i>2.29</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.8</i>	<i>82 - 127</i>	<i>51.21</i>	<i>3.59</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010930-01      HS20010930-09      HS20010930-10

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

<b>Batch ID:</b> R355066 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200125</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 11:48</b>				
Client ID:	Run ID: <b>VOA2_355066</b>	SeqNo: <b>5447954</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.24</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>51.29</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200225</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 10:59</b>				
Client ID:	Run ID: <b>VOA2_355066</b>	SeqNo: <b>5447953</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18.57	1.0	20	0	92.9	70 - 124			
Benzene	18.24	1.0	20	0	91.2	74 - 120			
Chlorobenzene	18.75	1.0	20	0	93.8	76 - 113			
Ethylbenzene	18.83	1.0	20	0	94.1	77 - 117			
Methylene chloride	18.77	2.0	20	0	93.8	70 - 127			
Toluene	20.88	1.0	20	0	104	77 - 118			
Vinyl chloride	19.59	1.0	20	0	98.0	70 - 130			
Xylenes, Total	58.21	1.0	60	0	97.0	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.59</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

**Batch ID:** R355066 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010767-29MS			Units: ug/L		Analysis Date: 25-Jan-2020 13:28			
Client ID:		Run ID: VOA2_355066			SeqNo: 5447958		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.72	1.0	20	0	88.6	70 - 127				
Benzene	17.87	1.0	20	0	89.3	70 - 127				
Chlorobenzene	18.14	1.0	20	0	90.7	70 - 114				
Ethylbenzene	18.64	1.0	20	0	93.2	70 - 124				
Methylene chloride	18.41	2.0	20	0	92.0	70 - 128				
Toluene	20.39	1.0	20	0	102	70 - 123				
Vinyl chloride	18.29	1.0	20	0	91.4	70 - 130				
Xylenes, Total	57.1	1.0	60	0	95.2	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.56</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.1</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010767-29MSD			Units: ug/L		Analysis Date: 25-Jan-2020 13:52			
Client ID:		Run ID: VOA2_355066			SeqNo: 5447959		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.67	1.0	20	0	88.4	70 - 127	17.72	0.267	20	
Benzene	17.9	1.0	20	0	89.5	70 - 127	17.87	0.197	20	
Chlorobenzene	18.23	1.0	20	0	91.2	70 - 114	18.14	0.504	20	
Ethylbenzene	18.75	1.0	20	0	93.8	70 - 124	18.64	0.579	20	
Methylene chloride	17.79	2.0	20	0	88.9	70 - 128	18.41	3.41	20	
Toluene	20.63	1.0	20	0	103	70 - 123	20.39	1.13	20	
Vinyl chloride	18.66	1.0	20	0	93.3	70 - 130	18.29	2	20	
Xylenes, Total	57.29	1.0	60	0	95.5	70 - 130	57.1	0.342	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>	<i>51.56</i>	<i>1.47</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>81 - 113</i>	<i>50.55</i>	<i>1.63</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>77 - 123</i>	<i>50.1</i>	<i>0.63</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>49.72</i>	<i>1.16</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010930-02 HS20010930-05 HS20010930-06 HS20010930-07  
 HS20010930-08

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

<b>Batch ID:</b> R355150 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200127</b>	Units: <b>ug/L</b>			Analysis Date: <b>28-Jan-2020 00:04</b>				
Client ID:	Run ID: <b>VOA4_355150</b>	SeqNo: <b>5449458</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>53.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>107</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>51.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200127</b>	Units: <b>ug/L</b>			Analysis Date: <b>27-Jan-2020 23:14</b>				
Client ID:	Run ID: <b>VOA4_355150</b>	SeqNo: <b>5449457</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	17.77	1.0	20	0	88.9	70 - 124			
Benzene	18.82	1.0	20	0	94.1	74 - 120			
Chlorobenzene	18.67	1.0	20	0	93.4	76 - 113			
Ethylbenzene	18.47	1.0	20	0	92.3	77 - 117			
Methylene chloride	20	2.0	20	0	100.0	70 - 127			
Toluene	18.6	1.0	20	0	93.0	77 - 118			
Vinyl chloride	17.6	1.0	20	0	88.0	70 - 130			
Xylenes, Total	57.11	1.0	60	0	95.2	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>51.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.7</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QC BATCH REPORT**

**Batch ID:** R355150 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010958-02MS			Units: ug/L		Analysis Date: 28-Jan-2020 02:07			
Client ID:		Run ID: VOA4_355150			SeqNo: 5449463		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.91	1.0	20	0	84.5	70 - 127				
Benzene	18.82	1.0	20	0	94.1	70 - 127				
Chlorobenzene	17.88	1.0	20	0	89.4	70 - 114				
Ethylbenzene	17.99	1.0	20	0	89.9	70 - 124				
Methylene chloride	18.93	2.0	20	0	94.6	70 - 128				
Toluene	18.76	1.0	20	0	93.8	70 - 123				
Vinyl chloride	17.58	1.0	20	0	87.9	70 - 130				
Xylenes, Total	54.55	1.0	60	0	90.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>51.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010958-02MSD			Units: ug/L		Analysis Date: 28-Jan-2020 02:32			
Client ID:		Run ID: VOA4_355150			SeqNo: 5449464		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.29	1.0	20	0	81.4	70 - 127	16.91	3.76	20	
Benzene	17.94	1.0	20	0	89.7	70 - 127	18.82	4.77	20	
Chlorobenzene	17.53	1.0	20	0	87.7	70 - 114	17.88	1.94	20	
Ethylbenzene	17.71	1.0	20	0	88.5	70 - 124	17.99	1.57	20	
Methylene chloride	18.74	2.0	20	0	93.7	70 - 128	18.93	0.98	20	
Toluene	17.72	1.0	20	0	88.6	70 - 123	18.76	5.69	20	
Vinyl chloride	16.97	1.0	20	0	84.8	70 - 130	17.58	3.57	20	
Xylenes, Total	53.22	1.0	60	0	88.7	70 - 130	54.55	2.47	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>	<i>51.55</i>	<i>0.229</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>	<i>50.92</i>	<i>0.238</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>51.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>77 - 123</i>	<i>51.17</i>	<i>0.604</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>	<i>50.57</i>	<i>0.637</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010930-03      HS20010930-04

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010930

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010930

**SAMPLE TRACKING**

Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS20010930-01	WQ-1620-TB06-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA232
HS20010930-02	WG-1620-MW32AR-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-02	WG-1620-MW32AR-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT106
HS20010930-02	WG-1620-MW32AR-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-03	WG-1620-MW32B-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-03	WG-1620-MW32B-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT106
HS20010930-03	WG-1620-MW32B-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-04	WG-1620-MW70B-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-04	WG-1620-MW70B-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT106
HS20010930-04	WG-1620-MW70B-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-05	WG-1620-MW33A-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-05	WG-1620-MW33A-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT106
HS20010930-05	WG-1620-MW33A-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-06	WG-1620-MW87C-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-06	WG-1620-MW87C-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT106
HS20010930-06	WG-1620-MW87C-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-07	WG-1620-MW90B-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-07	WG-1620-MW90B-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT106
HS20010930-07	WG-1620-MW90B-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-08	WG-1620-MW45C-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-08	WG-1620-MW45C-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT111
HS20010930-08	WG-1620-MW45C-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-09	WG-1620-FB10-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-09	WG-1620-FB10-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT111
HS20010930-09	WG-1620-FB10-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252
HS20010930-10	WG-1620-FD05-20200120	Login	1/21/2020 6:41:02 PM	JRM	MET074
HS20010930-10	WG-1620-FD05-20200120	Login	1/21/2020 6:41:02 PM	JRM	EXT111
HS20010930-10	WG-1620-FD05-20200120	Login	1/21/2020 6:41:02 PM	JRM	VOA252

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010930

Date/Time Received: 21-Jan-2020 17:10
Received by: DDG

Checklist completed by: Jared R. Makan
eSignature
Date: 21-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 27-Jan-2020

Matrices: Water

Carrier name: ALS Courier

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:206397

Temperature(s)/Thermometer(s): 1.3°C/1.3°C, 2.0°C/2.0°C, 1.6°C/1.6°C UC/C IR25
Cooler(s)/Kit(s): 45675, 45543, 44845
Date/Time sample(s) sent to storage: 01/21/2020 18:52

- Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]

pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

## HS20010930

Golder Associates Inc.  
Houston TX-Wood Preserving Works

Page 1 of 1

COC ID: 206397



ALS Project Manager:

Customer Information		Project Information		ALS Project Manager:											
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)										
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)										
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)										
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As)										
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E											
	Suite 4004		Stop 0750	F											
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G											
Phone	(512) 671-3434	Phone		H											
Fax	(512) 671-3446	Fax		I											
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		J											

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TB005 20200121	1-20-20	9:10	Water	1	2		X									
2	WG 1620 MW 32AR 20200120	1-20-20	9:10	Water	1	6	X	X	X	X							
3	WG 1620 MW 32B 20200120	1-20-20	10:15	H <sub>2</sub> O	1	6	X	X	X	X							
4	WG 1620 MW 70B 20200120	1-20-20	11:30	H <sub>2</sub> O	1	6	X	X	X	X							
5	WG 1620 MW 33A 20200120	1-20-20	12:48	H <sub>2</sub> O	1	6	X	X	X	X							
6	WG 1620 MW 87C 20200120	1-20-20	14:00	H <sub>2</sub> O	1	6	X	X	X	X							
7	WG 1620 MW 90B 20200120	1-20-20	15:15	H <sub>2</sub> O	1	6	X	X	X	X							
8	WG 1620 MW 45C 20200120	1-20-20	16:30	H <sub>2</sub> O	1	6	X	X	X	X							
9	WG 1620 PB 10 20200120	1-20-20	9:30	H <sub>2</sub> O	1	6	X	X	X	X							
0	WG 1620 FD 05 20200120	1-20-20	-	H <sub>2</sub> O	1	6	X	X	X	X							

Sampler(s) Please Print & Sign: Tim McSpauld T. McSpauld

Shipment Method: \_\_\_\_\_ Required Turnaround Time: (Check Box)  STD 10 Wk Days  5 Wk Days  2 Wk Days  24-hour

Results Due Date: \_\_\_\_\_

Relinquished by: T. McSpauld Date: 1-20-20 Time: 10:20

Received by: D.S. Date: 1-21-20 Time: 17:10

Relinquished by: D.S. Date: 1-21-20 Time: 17:10

Received by (Laboratory): D.S.

Checked by (Laboratory): \_\_\_\_\_

Logged by (Laboratory): \_\_\_\_\_

Notes: UPRR Houston MWPW

Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)
45643	2.0	<input type="checkbox"/> Level II Std QC <input checked="" type="checkbox"/> TRRP Check 1st
44815	1.6	<input type="checkbox"/> Level III Std QC/Raw Date <input type="checkbox"/> TRRP Level IV
45675	1.3	<input type="checkbox"/> Level IV SW843/CLP <input type="checkbox"/> Other

Reservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Notes: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

Copyright 2011 by ALS Environmental.



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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

January 30, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010958**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 7 sample(s) on Jan 22, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read 'Dane J. Wacasey'.

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

---

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/30/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010958			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149970,149998,R355061,R355150			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?	X				
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?		X			1
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Were MS/MSD RPDs within laboratory QC limits?		X			3
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/30/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010958			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149970,149998,R355061,R355150			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?		X			4
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).



**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 01/30/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20010958
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 149970,149998,R355061,R355150

ER# <sup>5</sup>	Description
1	Batch 149998, Semivolatile Organics Method SW8270, LCS recovery was above the control limits for 2,4-Dinitrotoluene and 2,6-Dinitrotoluene. The analytes were not detected in the associated samples.
2	Batch 149998, Semivolatile Organics Method SW8270, sample WG-1620-MW82B-20200121, MS recovered outside the control limit for 4-Nitrophenol due to possible matrix effect.
3	Batch 149998, Semivolatile Organics Method SW8270, sample WG-1620-MW82B-20200121, MS/MSD RPD recovered above the RPD limits for select compounds due to possible matrix effect.
4	Batch 149970, Metals Method SW6020, sample WG-1620-MW82B-20200121, PDS recovered above the control limit for Arsenic.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010958

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010958-01	WQ-1620-TB07-20200121	Water	CG-121719 -114	21-Jan-2020 00:00	22-Jan-2020 11:30	<input type="checkbox"/>
HS20010958-02	WG-1620-MW82B-20200121	Water		21-Jan-2020 09:10	22-Jan-2020 11:30	<input type="checkbox"/>
HS20010958-03	WG-1620-MW38A-20200121	Water		21-Jan-2020 11:10	22-Jan-2020 11:30	<input type="checkbox"/>
HS20010958-04	WG-1620-MW38B-20200121	Water		21-Jan-2020 13:45	22-Jan-2020 11:30	<input type="checkbox"/>
HS20010958-05	WG-1620-MW03-20200121	Water		21-Jan-2020 15:00	22-Jan-2020 11:30	<input type="checkbox"/>
HS20010958-06	WG-1620-MW04-20200121	Water		21-Jan-2020 16:00	22-Jan-2020 11:30	<input type="checkbox"/>
HS20010958-07	WG-1620-FB11-20200121	Water		21-Jan-2020 15:30	22-Jan-2020 11:30	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB07-20200121  
 Collection Date: 21-Jan-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	28-Jan-2020 00:28
Benzene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 00:28
Chlorobenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 00:28
Ethylbenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 00:28
Methylene chloride	U		0.0010	0.0020	mg/L	1	28-Jan-2020 00:28
Toluene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 00:28
Xylenes, Total	U		0.00030	0.0010	mg/L	1	28-Jan-2020 00:28
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>108</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:28</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:28</i>
<i>Surr: Dibromofluoromethane</i>	<i>102</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:28</i>
<i>Surr: Toluene-d8</i>	<i>99.8</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:28</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW82B-20200121  
 Collection Date: 21-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	28-Jan-2020 01:42
Benzene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 01:42
Chlorobenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 01:42
Ethylbenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 01:42
Methylene chloride	U		0.0010	0.0020	mg/L	1	28-Jan-2020 01:42
Toluene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 01:42
Xylenes, Total	U		0.00030	0.0010	mg/L	1	28-Jan-2020 01:42
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>108</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 01:42</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.0</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 01:42</i>
<i>Surr: Dibromofluoromethane</i>		<i>105</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 01:42</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 01:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW82B-20200121  
 Collection Date: 21-Jan-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	29-Jan-2020 22:17
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	29-Jan-2020 22:17
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	29-Jan-2020 22:17
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	29-Jan-2020 22:17
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	29-Jan-2020 22:17
<b>2-Methylnaphthalene</b>	<b>0.00010</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	29-Jan-2020 22:17
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	29-Jan-2020 22:17
<b>Acenaphthene</b>	<b>0.000043</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
Acenaphthylene	U		0.000015	0.00010	mg/L	1	29-Jan-2020 22:17
<b>Anthracene</b>	<b>0.000031</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	29-Jan-2020 22:17
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	29-Jan-2020 22:17
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	29-Jan-2020 22:17
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	29-Jan-2020 22:17
Chrysene	U		0.000021	0.00010	mg/L	1	29-Jan-2020 22:17
<b>Dibenzofuran</b>	<b>0.000040</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	29-Jan-2020 22:17
Fluoranthene	U		0.000010	0.00010	mg/L	1	29-Jan-2020 22:17
<b>Fluorene</b>	<b>0.000037</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
<b>Naphthalene</b>	<b>0.00041</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
Nitrobenzene	U		0.000024	0.00020	mg/L	1	29-Jan-2020 22:17
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	29-Jan-2020 22:17
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	29-Jan-2020 22:17
<b>Phenanthrene</b>	<b>0.00012</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 22:17
Phenol	U		0.000035	0.00020	mg/L	1	29-Jan-2020 22:17
Pyrene	U		0.000019	0.00010	mg/L	1	29-Jan-2020 22:17
<i>Surr: 2,4,6-Tribromophenol</i>	<i>93.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 22:17</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>93.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 22:17</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 22:17</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>115</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 22:17</i>
<i>Surr: Nitrobenzene-d5</i>	<i>79.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 22:17</i>
<i>Surr: Phenol-d6</i>	<i>74.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 22:17</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00484</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Jan-2020 21:28

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38A-20200121  
 Collection Date: 21-Jan-2020 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:08
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:08
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:08
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:08
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 20:08
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:08
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:08
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>105</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:08</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.5</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:08</i>
<i>Surr: Dibromofluoromethane</i>		<i>101</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:08</i>
<i>Surr: Toluene-d8</i>		<i>100</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:08</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38A-20200121  
 Collection Date: 21-Jan-2020 11:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	29-Jan-2020 23:14
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	29-Jan-2020 23:14
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	29-Jan-2020 23:14
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	29-Jan-2020 23:14
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	29-Jan-2020 23:14
<b>2-Methylnaphthalene</b>	<b>0.00017</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	29-Jan-2020 23:14
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	29-Jan-2020 23:14
<b>Acenaphthene</b>	<b>0.00068</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Acenaphthylene	U		0.000015	0.00010	mg/L	1	29-Jan-2020 23:14
<b>Anthracene</b>	<b>0.00015</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	29-Jan-2020 23:14
<b>Benzo(a)pyrene</b>	<b>0.000032</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	29-Jan-2020 23:14
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000063</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
<b>Chrysene</b>	<b>0.000060</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
<b>Dibenzofuran</b>	<b>0.00036</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	29-Jan-2020 23:14
<b>Fluoranthene</b>	<b>0.00018</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
<b>Fluorene</b>	<b>0.00029</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
<b>Naphthalene</b>	<b>0.0011</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Nitrobenzene	U		0.000024	0.00020	mg/L	1	29-Jan-2020 23:14
<b>N-Nitrosodiphenylamine</b>	<b>0.00011</b>	J	<b>0.000025</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	29-Jan-2020 23:14
<b>Phenanthrene</b>	<b>0.00039</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Phenol	U		0.000035	0.00020	mg/L	1	29-Jan-2020 23:14
<b>Pyrene</b>	<b>0.00014</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:14
Surr: 2,4,6-Tribromophenol	68.4			34-129	%REC	1	29-Jan-2020 23:14
Surr: 2-Fluorobiphenyl	72.4			40-125	%REC	1	29-Jan-2020 23:14
Surr: 2-Fluorophenol	63.9			20-120	%REC	1	29-Jan-2020 23:14
Surr: 4-Terphenyl-d14	92.4			40-135	%REC	1	29-Jan-2020 23:14
Surr: Nitrobenzene-d5	67.8			41-120	%REC	1	29-Jan-2020 23:14
Surr: Phenol-d6	66.4			20-120	%REC	1	29-Jan-2020 23:14
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.0177</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Jan-2020 21:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38B-20200121  
 Collection Date: 21-Jan-2020 13:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:33
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:33
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:33
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:33
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 20:33
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:33
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:33
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>104</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:33</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>99.4</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:33</i>
<i>Surr: Dibromofluoromethane</i>		<i>98.8</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:33</i>
<i>Surr: Toluene-d8</i>		<i>96.9</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:33</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38B-20200121  
 Collection Date: 21-Jan-2020 13:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
<b>1,2-Diphenylhydrazine</b>	<b>0.00025</b>		<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	29-Jan-2020 23:33
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	29-Jan-2020 23:33
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	29-Jan-2020 23:33
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	29-Jan-2020 23:33
<b>2-Methylnaphthalene</b>	<b>0.00017</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	29-Jan-2020 23:33
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	29-Jan-2020 23:33
<b>Acenaphthene</b>	<b>0.054</b>		<b>0.00027</b>	<b>0.0010</b>	<b>mg/L</b>	10	30-Jan-2020 14:04
<b>Acenaphthylene</b>	<b>0.00045</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
<b>Anthracene</b>	<b>0.0012</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
<b>Benz(a)anthracene</b>	<b>0.000055</b>	J	<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	29-Jan-2020 23:33
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	29-Jan-2020 23:33
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	29-Jan-2020 23:33
<b>Chrysene</b>	<b>0.000057</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
<b>Dibenzofuran</b>	<b>0.0054</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
<b>Di-n-butyl phthalate</b>	<b>0.000036</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
<b>Fluoranthene</b>	<b>0.0025</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
<b>Fluorene</b>	<b>0.015</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	10	30-Jan-2020 14:04
<b>Naphthalene</b>	<b>0.00098</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
Nitrobenzene	U		0.000024	0.00020	mg/L	1	29-Jan-2020 23:33
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	29-Jan-2020 23:33
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	29-Jan-2020 23:33
<b>Phenanthrene</b>	<b>0.00095</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
Phenol	U		0.000035	0.00020	mg/L	1	29-Jan-2020 23:33
<b>Pyrene</b>	<b>0.0015</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:33
Surr: 2,4,6-Tribromophenol	80.6			34-129	%REC	1	29-Jan-2020 23:33
Surr: 2,4,6-Tribromophenol	93.5			34-129	%REC	10	30-Jan-2020 14:04
Surr: 2-Fluorobiphenyl	106			40-125	%REC	10	30-Jan-2020 14:04
Surr: 2-Fluorobiphenyl	87.7			40-125	%REC	1	29-Jan-2020 23:33
Surr: 2-Fluorophenol	71.8			20-120	%REC	1	29-Jan-2020 23:33
Surr: 2-Fluorophenol	74.8			20-120	%REC	10	30-Jan-2020 14:04
Surr: 4-Terphenyl-d14	117			40-135	%REC	10	30-Jan-2020 14:04
Surr: 4-Terphenyl-d14	95.1			40-135	%REC	1	29-Jan-2020 23:33
Surr: Nitrobenzene-d5	75.9			41-120	%REC	1	29-Jan-2020 23:33
Surr: Nitrobenzene-d5	94.3			41-120	%REC	10	30-Jan-2020 14:04
Surr: Phenol-d6	93.3			20-120	%REC	10	30-Jan-2020 14:04
Surr: Phenol-d6	79.3			20-120	%REC	1	29-Jan-2020 23:33

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW38B-20200121  
 Collection Date: 21-Jan-2020 13:45

**ANALYTICAL REPORT**

WorkOrder:HS20010958  
 Lab ID:HS20010958-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 24-Jan-2020		Analyst: JHD
Arsenic	0.0173		0.000400	0.00200	mg/L	1	27-Jan-2020 21:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW03-20200121  
 Collection Date: 21-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:58
Benzene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:58
Chlorobenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:58
Ethylbenzene	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:58
Methylene chloride	U		0.0010	0.0020	mg/L	1	25-Jan-2020 20:58
Toluene	U		0.00020	0.0010	mg/L	1	25-Jan-2020 20:58
Xylenes, Total	U		0.00030	0.0010	mg/L	1	25-Jan-2020 20:58
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>104</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:58</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>97.5</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:58</i>
<i>Surr: Dibromofluoromethane</i>		<i>100</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:58</i>
<i>Surr: Toluene-d8</i>		<i>99.2</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Jan-2020 20:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW03-20200121  
 Collection Date: 21-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	29-Jan-2020 23:52
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	29-Jan-2020 23:52
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	29-Jan-2020 23:52
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	29-Jan-2020 23:52
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	29-Jan-2020 23:52
<b>2-Methylnaphthalene</b>	<b>0.000045</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	29-Jan-2020 23:52
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	29-Jan-2020 23:52
<b>Acenaphthene</b>	<b>0.000061</b>	J	<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
Acenaphthylene	U		0.000015	0.00010	mg/L	1	29-Jan-2020 23:52
<b>Anthracene</b>	<b>0.000069</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	29-Jan-2020 23:52
<b>Benzo(a)pyrene</b>	<b>0.000024</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	29-Jan-2020 23:52
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
Chrysene	U		0.000021	0.00010	mg/L	1	29-Jan-2020 23:52
Dibenzofuran	U		0.000020	0.00010	mg/L	1	29-Jan-2020 23:52
<b>Di-n-butyl phthalate</b>	<b>0.000029</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
<b>Fluoranthene</b>	<b>0.000055</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
Fluorene	U		0.000030	0.00010	mg/L	1	29-Jan-2020 23:52
<b>Naphthalene</b>	<b>0.00024</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
Nitrobenzene	U		0.000024	0.00020	mg/L	1	29-Jan-2020 23:52
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	29-Jan-2020 23:52
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	29-Jan-2020 23:52
Phenanthrene	U		0.000021	0.00010	mg/L	1	29-Jan-2020 23:52
Phenol	U		0.000035	0.00020	mg/L	1	29-Jan-2020 23:52
<b>Pyrene</b>	<b>0.000036</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	29-Jan-2020 23:52
<i>Surr: 2,4,6-Tribromophenol</i>	68.7			34-129	%REC	1	29-Jan-2020 23:52
<i>Surr: 2-Fluorobiphenyl</i>	91.9			40-125	%REC	1	29-Jan-2020 23:52
<i>Surr: 2-Fluorophenol</i>	75.5			20-120	%REC	1	29-Jan-2020 23:52
<i>Surr: 4-Terphenyl-d14</i>	94.9			40-135	%REC	1	29-Jan-2020 23:52
<i>Surr: Nitrobenzene-d5</i>	80.4			41-120	%REC	1	29-Jan-2020 23:52
<i>Surr: Phenol-d6</i>	77.0			20-120	%REC	1	29-Jan-2020 23:52
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00207</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Jan-2020 21:48

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW04-20200121  
 Collection Date: 21-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	28-Jan-2020 02:56
Benzene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 02:56
Chlorobenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 02:56
Ethylbenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 02:56
Methylene chloride	U		0.0010	0.0020	mg/L	1	28-Jan-2020 02:56
Toluene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 02:56
Xylenes, Total	U		0.00030	0.0010	mg/L	1	28-Jan-2020 02:56
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>105</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 02:56</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>98.5</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 02:56</i>
<i>Surr: Dibromofluoromethane</i>		<i>103</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 02:56</i>
<i>Surr: Toluene-d8</i>		<i>99.1</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 02:56</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW04-20200121  
 Collection Date: 21-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-06  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	30-Jan-2020 00:11
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	30-Jan-2020 00:11
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	30-Jan-2020 00:11
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	30-Jan-2020 00:11
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	30-Jan-2020 00:11
<b>2-Methylnaphthalene</b>	<b>0.00026</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	30-Jan-2020 00:11
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	30-Jan-2020 00:11
<b>Acenaphthene</b>	<b>0.00015</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
Acenaphthylene	U		0.000015	0.00010	mg/L	1	30-Jan-2020 00:11
<b>Anthracene</b>	<b>0.00010</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	30-Jan-2020 00:11
<b>Benzo(a)pyrene</b>	<b>0.000033</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	30-Jan-2020 00:11
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0018</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
<b>Chrysene</b>	<b>0.000034</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
<b>Dibenzofuran</b>	<b>0.000097</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	30-Jan-2020 00:11
<b>Fluoranthene</b>	<b>0.00029</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
<b>Fluorene</b>	<b>0.000067</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
<b>Naphthalene</b>	<b>0.0013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
Nitrobenzene	U		0.000024	0.00020	mg/L	1	30-Jan-2020 00:11
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	30-Jan-2020 00:11
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	30-Jan-2020 00:11
<b>Phenanthrene</b>	<b>0.00017</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
Phenol	U		0.000035	0.00020	mg/L	1	30-Jan-2020 00:11
<b>Pyrene</b>	<b>0.00020</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:11
<i>Surr: 2,4,6-Tribromophenol</i>	70.9			34-129	%REC	1	30-Jan-2020 00:11
<i>Surr: 2-Fluorobiphenyl</i>	85.8			40-125	%REC	1	30-Jan-2020 00:11
<i>Surr: 2-Fluorophenol</i>	62.6			20-120	%REC	1	30-Jan-2020 00:11
<i>Surr: 4-Terphenyl-d14</i>	95.4			40-135	%REC	1	30-Jan-2020 00:11
<i>Surr: Nitrobenzene-d5</i>	71.1			41-120	%REC	1	30-Jan-2020 00:11
<i>Surr: Phenol-d6</i>	71.4			20-120	%REC	1	30-Jan-2020 00:11
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00223</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Jan-2020 21:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB11-20200121  
 Collection Date: 21-Jan-2020 15:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	28-Jan-2020 00:53
Benzene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 00:53
Chlorobenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 00:53
Ethylbenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 00:53
Methylene chloride	U		0.0010	0.0020	mg/L	1	28-Jan-2020 00:53
Toluene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 00:53
Xylenes, Total	U		0.00030	0.0010	mg/L	1	28-Jan-2020 00:53
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>108</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:53</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:53</i>
<i>Surr: Dibromofluoromethane</i>	<i>104</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:53</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 00:53</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB11-20200121  
 Collection Date: 21-Jan-2020 15:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010958  
 Lab ID:HS20010958-07  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 27-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	30-Jan-2020 00:30
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	30-Jan-2020 00:30
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	30-Jan-2020 00:30
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	30-Jan-2020 00:30
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	30-Jan-2020 00:30
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	30-Jan-2020 00:30
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	30-Jan-2020 00:30
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	30-Jan-2020 00:30
Acenaphthene	U		0.000027	0.00010	mg/L	1	30-Jan-2020 00:30
Acenaphthylene	U		0.000015	0.00010	mg/L	1	30-Jan-2020 00:30
Anthracene	U		0.000014	0.00010	mg/L	1	30-Jan-2020 00:30
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	30-Jan-2020 00:30
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	30-Jan-2020 00:30
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	30-Jan-2020 00:30
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	30-Jan-2020 00:30
Chrysene	U		0.000021	0.00010	mg/L	1	30-Jan-2020 00:30
Dibenzofuran	U		0.000020	0.00010	mg/L	1	30-Jan-2020 00:30
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	30-Jan-2020 00:30
Fluoranthene	U		0.000010	0.00010	mg/L	1	30-Jan-2020 00:30
Fluorene	U		0.000030	0.00010	mg/L	1	30-Jan-2020 00:30
<b>Naphthalene</b>	<b>0.00018</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 00:30
Nitrobenzene	U		0.000024	0.00020	mg/L	1	30-Jan-2020 00:30
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	30-Jan-2020 00:30
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	30-Jan-2020 00:30
Phenanthrene	U		0.000021	0.00010	mg/L	1	30-Jan-2020 00:30
Phenol	U		0.000035	0.00020	mg/L	1	30-Jan-2020 00:30
Pyrene	U		0.000019	0.00010	mg/L	1	30-Jan-2020 00:30
<i>Surr: 2,4,6-Tribromophenol</i>	<i>69.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 00:30</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>94.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 00:30</i>
<i>Surr: 2-Fluorophenol</i>	<i>80.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 00:30</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 00:30</i>
<i>Surr: Nitrobenzene-d5</i>	<i>86.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 00:30</i>
<i>Surr: Phenol-d6</i>	<i>81.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 00:30</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 24-Jan-2020		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	27-Jan-2020 21:52

Note: See Qualifiers Page for a list of qualifiers and their explanation.



## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**Batch ID:** 149970      **Start Date:** 24 Jan 2020 09:00      **End Date:** 24 Jan 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010958-02		10 (mL)	10 (mL)	1
HS20010958-03		10 (mL)	10 (mL)	1
HS20010958-04		10 (mL)	10 (mL)	1
HS20010958-05		10 (mL)	10 (mL)	1
HS20010958-06		10 (mL)	10 (mL)	1
HS20010958-07		10 (mL)	10 (mL)	1

**Batch ID:** 149998      **Start Date:** 27 Jan 2020 07:00      **End Date:** 27 Jan 2020 13:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010958-02	1	1000 (mL)	1 (mL)	0.001
HS20010958-03	1	1000 (mL)	1 (mL)	0.001
HS20010958-04	1	1000 (mL)	1 (mL)	0.001
HS20010958-05	1	1000 (mL)	1 (mL)	0.001
HS20010958-06	1	1000 (mL)	1 (mL)	0.001
HS20010958-07	1	1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149970 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20010958-02	WG-1620-MW82B-20200121	21 Jan 2020 09:10		24 Jan 2020 13:00	27 Jan 2020 21:28	1
HS20010958-03	WG-1620-MW38A-20200121	21 Jan 2020 11:10		24 Jan 2020 13:00	27 Jan 2020 21:44	1
HS20010958-04	WG-1620-MW38B-20200121	21 Jan 2020 13:45		24 Jan 2020 13:00	27 Jan 2020 21:46	1
HS20010958-05	WG-1620-MW03-20200121	21 Jan 2020 15:00		24 Jan 2020 13:00	27 Jan 2020 21:48	1
HS20010958-06	WG-1620-MW04-20200121	21 Jan 2020 16:00		24 Jan 2020 13:00	27 Jan 2020 21:50	1
HS20010958-07	WG-1620-FB11-20200121	21 Jan 2020 15:30		24 Jan 2020 13:00	27 Jan 2020 21:52	1
<b>Batch ID: 149998 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010958-02	WG-1620-MW82B-20200121	21 Jan 2020 09:10		27 Jan 2020 10:49	29 Jan 2020 22:17	1
HS20010958-03	WG-1620-MW38A-20200121	21 Jan 2020 11:10		27 Jan 2020 10:49	29 Jan 2020 23:14	1
HS20010958-04	WG-1620-MW38B-20200121	21 Jan 2020 13:45		27 Jan 2020 10:49	30 Jan 2020 14:04	10
HS20010958-04	WG-1620-MW38B-20200121	21 Jan 2020 13:45		27 Jan 2020 10:49	29 Jan 2020 23:33	1
HS20010958-05	WG-1620-MW03-20200121	21 Jan 2020 15:00		27 Jan 2020 10:49	29 Jan 2020 23:52	1
HS20010958-06	WG-1620-MW04-20200121	21 Jan 2020 16:00		27 Jan 2020 10:49	30 Jan 2020 00:11	1
HS20010958-07	WG-1620-FB11-20200121	21 Jan 2020 15:30		27 Jan 2020 10:49	30 Jan 2020 00:30	1
<b>Batch ID: R355061 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010958-03	WG-1620-MW38A-20200121	21 Jan 2020 11:10			25 Jan 2020 20:08	1
HS20010958-04	WG-1620-MW38B-20200121	21 Jan 2020 13:45			25 Jan 2020 20:33	1
HS20010958-05	WG-1620-MW03-20200121	21 Jan 2020 15:00			25 Jan 2020 20:58	1
<b>Batch ID: R355150 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20010958-01	WQ-1620-TB07-20200121	21 Jan 2020 00:00			28 Jan 2020 00:28	1
HS20010958-02	WG-1620-MW82B-20200121	21 Jan 2020 09:10			28 Jan 2020 01:42	1
HS20010958-06	WG-1620-MW04-20200121	21 Jan 2020 16:00			28 Jan 2020 02:56	1
HS20010958-07	WG-1620-FB11-20200121	21 Jan 2020 15:30			28 Jan 2020 00:53	1

WorkOrder: HS20010958  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20010958  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20010958  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149970 ( 0 )		Instrument: ICPMS05			Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-149970</b>	Units: <b>mg/L</b>			Analysis Date: <b>27-Jan-2020 21:24</b>					
Client ID:		Run ID: <b>ICPMS05_355087</b>	SeqNo: <b>5449127</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>MBLK</b>	Sample ID: <b>MBLK-149970</b>	Units: <b>mg/L</b>			Analysis Date: <b>28-Jan-2020 13:11</b>					
Client ID:		Run ID: <b>ICPMS05_355152</b>	SeqNo: <b>5450164</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-149970</b>	Units: <b>mg/L</b>			Analysis Date: <b>27-Jan-2020 21:26</b>					
Client ID:		Run ID: <b>ICPMS05_355087</b>	SeqNo: <b>5449128</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.04847	0.00200	0.05	0	96.9	80 - 120				
<b>MS</b>	Sample ID: <b>HS20010958-02MS</b>	Units: <b>mg/L</b>			Analysis Date: <b>27-Jan-2020 21:33</b>					
Client ID: <b>WG-1620-MW82B-20200121</b>		Run ID: <b>ICPMS05_355087</b>	SeqNo: <b>5449131</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.05447	0.00200	0.05	0.004841	99.3	80 - 120				
<b>MSD</b>	Sample ID: <b>HS20010958-02MSD</b>	Units: <b>mg/L</b>			Analysis Date: <b>27-Jan-2020 21:35</b>					
Client ID: <b>WG-1620-MW82B-20200121</b>		Run ID: <b>ICPMS05_355087</b>	SeqNo: <b>5449132</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.05461	0.00200	0.05	0.004841	99.5	80 - 120	0.05447	0.248	20	
<b>PDS</b>	Sample ID: <b>HS20010958-02PDS</b>	Units: <b>mg/L</b>			Analysis Date: <b>28-Jan-2020 13:18</b>					
Client ID: <b>WG-1620-MW82B-20200121</b>		Run ID: <b>ICPMS05_355152</b>	SeqNo: <b>5450167</b>	PrepDate: <b>24-Jan-2020</b>	DF: <b>20</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	1.996	0.0400	0.1	0.00472	1990	75 - 125				S

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

<b>Batch ID:</b> 149970 ( 0 )		<b>Instrument:</b> ICPMS05		<b>Method:</b> ICP-MS METALS BY SW6020A					
<b>PDS</b>	Sample ID: <b>HS20010958-02PDS</b>		Units: <b>mg/L</b>		Analysis Date: <b>27-Jan-2020 21:37</b>				
Client ID:	<b>WG-1620-MW82B-20200121</b>	Run ID:	<b>ICPMS05_355087</b>	SeqNo:	<b>5449133</b>	PrepDate:	<b>24-Jan-2020</b>	DF:	<b>1</b>
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
Arsenic	0.1045	0.00200	0.1	0.004841	99.6	75 - 125			

<b>SD</b>	Sample ID: <b>HS20010958-02SD</b>		Units: <b>mg/L</b>		Analysis Date: <b>27-Jan-2020 21:30</b>				
Client ID:	<b>WG-1620-MW82B-20200121</b>	Run ID:	<b>ICPMS05_355087</b>	SeqNo:	<b>5449130</b>	PrepDate:	<b>24-Jan-2020</b>	DF:	<b>5</b>
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	%D Limit Qual
Arsenic	0.005043	0.0100					0.004841	0	10 J

<b>SD</b>	Sample ID: <b>HS20010958-02SD</b>		Units: <b>mg/L</b>		Analysis Date: <b>28-Jan-2020 13:15</b>				
Client ID:	<b>WG-1620-MW82B-20200121</b>	Run ID:	<b>ICPMS05_355152</b>	SeqNo:	<b>5450166</b>	PrepDate:	<b>24-Jan-2020</b>	DF:	<b>100</b>
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	%D Limit Qual
Arsenic	U	0.200					0.00472	0	10

The following samples were analyzed in this batch:

HS20010958-02	HS20010958-03	HS20010958-04	HS20010958-05
HS20010958-06	HS20010958-07		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-149998	Units: ug/L			Analysis Date: 28-Jan-2020 19:21					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452454	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.165</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.567</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.869</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.797</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>116</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.964</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.941</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>98.8</i>	<i>20 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-149998	Units: ug/L			Analysis Date: 28-Jan-2020 19:40					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452455	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.783	0.20	5	0	116	39 - 127				
2,4-Dimethylphenol	5.012	0.20	5	0	100	35 - 120				
2,4-Dinitrotoluene	6.415	0.20	5	0	128	50 - 122				S
2,6-Dinitrotoluene	6.03	0.20	5	0	121	50 - 120				S
2-Chloronaphthalene	5.424	0.20	5	0	108	50 - 120				
2-Methylnaphthalene	5.383	0.10	5	0	108	50 - 120				
4,6-Dinitro-2-methylphenol	5.692	0.20	5	0	114	25 - 121				
4-Nitrophenol	5.835	1.0	5	0	117	30 - 130				
Acenaphthene	5.432	0.10	5	0	109	45 - 120				
Acenaphthylene	5.667	0.10	5	0	113	47 - 120				
Anthracene	5.847	0.10	5	0	117	45 - 120				
Benz(a)anthracene	5.919	0.10	5	0	118	40 - 120				
Benzo(a)pyrene	5.832	0.10	5	0	117	45 - 120				
Bis(2-chloroethoxy)methane	5.315	0.20	5	0	106	45 - 120				
Bis(2-ethylhexyl)phthalate	5.645	0.20	5	0	113	40 - 139				
Chrysene	5.825	0.10	5	0	117	43 - 120				
Dibenzofuran	5.475	0.10	5	0	109	50 - 120				
Di-n-butyl phthalate	5.996	0.20	5	0	120	45 - 123				
Fluoranthene	5.728	0.10	5	0	115	45 - 125				
Fluorene	5.748	0.10	5	0	115	49 - 120				
Naphthalene	5.43	0.10	5	0	109	45 - 120				
Nitrobenzene	5.209	0.20	5	0	104	44 - 120				
N-Nitrosodiphenylamine	5.868	0.20	5	0	117	40 - 125				
Pentachlorophenol	2.037	0.20	5	0	40.7	19 - 121				
Phenanthrene	5.866	0.10	5	0	117	45 - 121				
Phenol	5.563	0.20	5	0	111	20 - 124				
Pyrene	5.924	0.10	5	0	118	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.466</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.582</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>112</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.846</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.9</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.609</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>112</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.978</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.891</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>97.8</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS20010958-02MS		Units: ug/L		Analysis Date: 29-Jan-2020 22:36				
Client ID: WG-1620-MW82B-20200121		Run ID: SV-6_355351		SeqNo: 5453263		PrepDate: 27-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	5.036	0.20	5	0	101	39 - 127				
2,4-Dimethylphenol	3.403	0.20	5	0	68.1	35 - 120				
2,4-Dinitrotoluene	6.071	0.20	5	0	121	50 - 122				
2,6-Dinitrotoluene	5.246	0.20	5	0	105	50 - 120				
2-Chloronaphthalene	3.81	0.20	5	0	76.2	50 - 120				
2-Methylnaphthalene	3.528	0.10	5	0.1009	68.5	50 - 120				
4,6-Dinitro-2-methylphenol	4.25	0.20	5	0	85.0	25 - 121				
4-Nitrophenol	6.787	1.0	5	0	136	30 - 130			S	
Acenaphthene	4.374	0.10	5	0.04257	86.6	45 - 120				
Acenaphthylene	4.018	0.10	5	0	80.4	47 - 120				
Anthracene	5.906	0.10	5	0.03111	117	45 - 120				
Benz(a)anthracene	5.831	0.10	5	0	117	40 - 120				
Benzo(a)pyrene	5.639	0.10	5	0	113	45 - 120				
Bis(2-chloroethoxy)methane	3.489	0.20	5	0	69.8	45 - 120				
Bis(2-ethylhexyl)phthalate	3.59	0.20	5	0	71.8	40 - 139				
Chrysene	5.562	0.10	5	0	111	43 - 120				
Dibenzofuran	4.131	0.10	5	0.04005	81.8	50 - 120				
Di-n-butyl phthalate	4.49	0.20	5	0	89.8	45 - 123				
Fluoranthene	5.834	0.10	5	0	117	45 - 125				
Fluorene	4.534	0.10	5	0.03736	89.9	49 - 120				
Naphthalene	3.669	0.10	5	0.4092	65.2	45 - 120				
Nitrobenzene	3.406	0.20	5	0	68.1	44 - 120				
N-Nitrosodiphenylamine	5.951	0.20	5	0	119	40 - 125				
Pentachlorophenol	5.92	0.20	5	0	118	19 - 121				
Phenanthrene	5.624	0.10	5	0.1179	110	45 - 121				
Phenol	3.512	0.20	5	0	70.2	20 - 124				
Pyrene	5.934	0.10	5	0	119	40 - 130				
Surr: 2,4,6-Tribromophenol	4.8	0.20	5	0	96.0	34 - 129				
Surr: 2-Fluorobiphenyl	3.948	0.20	5	0	79.0	40 - 125				
Surr: 2-Fluorophenol	3.307	0.20	5	0	66.1	20 - 120				
Surr: 4-Terphenyl-d14	5.739	0.20	5	0	115	40 - 135				
Surr: Nitrobenzene-d5	3.377	0.20	5	0	67.5	41 - 120				
Surr: Phenol-d6	3.333	0.20	5	0	66.7	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS	Sample ID: HS20010910-03MS	Units: ug/L			Analysis Date: 28-Jan-2020 23:44					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452170	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.258	0.20	5	0	85.2	39 - 127				
2,4-Dimethylphenol	2.953	0.20	5	0.09533	57.1	35 - 120				
2,4-Dinitrotoluene	5.24	0.20	5	0	105	50 - 122				
2,6-Dinitrotoluene	4.544	0.20	5	0	90.9	50 - 120				
2-Chloronaphthalene	3.621	0.20	5	0	72.4	50 - 120				
2-Methylnaphthalene	3.197	0.10	5	0	63.9	50 - 120				
4,6-Dinitro-2-methylphenol	5.327	0.20	5	0	107	25 - 121				
4-Nitrophenol	5.515	1.0	5	0	110	30 - 130				
Acenaphthene	3.484	0.10	5	0	69.7	45 - 120				
Acenaphthylene	3.505	0.10	5	0	70.1	47 - 120				
Anthracene	5.162	0.10	5	0	103	45 - 120				
Benz(a)anthracene	5.438	0.10	5	0	109	40 - 120				
Benzo(a)pyrene	5.398	0.10	5	0	108	45 - 120				
Bis(2-chloroethoxy)methane	2.957	0.20	5	0	59.1	45 - 120				
Bis(2-ethylhexyl)phthalate	5.76	0.20	5	0	115	40 - 139				
Chrysene	5.195	0.10	5	0	104	43 - 120				
Dibenzofuran	3.564	0.10	5	0	71.3	50 - 120				
Di-n-butyl phthalate	6.041	0.20	5	0	121	45 - 123				
Fluoranthene	5.294	0.10	5	0.02986	105	45 - 125				
Fluorene	3.974	0.10	5	0	79.5	49 - 120				
Naphthalene	3.352	0.10	5	0.22	62.6	45 - 120				
Nitrobenzene	3.206	0.20	5	0	64.1	44 - 120				
N-Nitrosodiphenylamine	5.464	0.20	5	0	109	40 - 125				
Pentachlorophenol	3.381	0.20	5	0	67.6	19 - 121				
Phenanthrene	4.995	0.10	5	0	99.9	45 - 121				
Phenol	3.121	0.20	5	0	62.4	20 - 124				
Pyrene	5.526	0.10	5	0.02733	110	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.181</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.6</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.463</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.3</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.1</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>62.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.541</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>111</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.186</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.7</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.21</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>64.2</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS20010958-02MSD		Units: ug/L		Analysis Date: 29-Jan-2020 22:55				
Client ID: WG-1620-MW82B-20200121		Run ID: SV-6_355351		SeqNo: 5453264		PrepDate: 27-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.703	0.20	5	0	94.1	39 - 127	5.036	6.83	20	
2,4-Dimethylphenol	3.362	0.20	5	0	67.2	35 - 120	3.403	1.21	20	
2,4-Dinitrotoluene	5.237	0.20	5	0	105	50 - 122	6.071	14.8	20	
2,6-Dinitrotoluene	4.788	0.20	5	0	95.8	50 - 120	5.246	9.14	20	
2-Chloronaphthalene	3.704	0.20	5	0	74.1	50 - 120	3.81	2.82	20	
2-Methylnaphthalene	3.424	0.10	5	0.1009	66.5	50 - 120	3.528	2.98	20	
4,6-Dinitro-2-methylphenol	4.301	0.20	5	0	86.0	25 - 121	4.25	1.18	30	
4-Nitrophenol	4.754	1.0	5	0	95.1	30 - 130	6.787	35.2	20	R
Acenaphthene	4.092	0.10	5	0.04257	81.0	45 - 120	4.374	6.66	20	
Acenaphthylene	3.822	0.10	5	0	76.4	47 - 120	4.018	4.98	20	
Anthracene	5.18	0.10	5	0.03111	103	45 - 120	5.906	13.1	20	
Benz(a)anthracene	5.424	0.10	5	0	108	40 - 120	5.831	7.23	20	
Benzo(a)pyrene	5.25	0.10	5	0	105	45 - 120	5.639	7.13	20	
Bis(2-chloroethoxy)methane	3.359	0.20	5	0	67.2	45 - 120	3.489	3.81	20	
Bis(2-ethylhexyl)phthalate	6.024	0.20	5	0	120	40 - 139	3.59	50.6	20	R
Chrysene	4.985	0.10	5	0	99.7	43 - 120	5.562	10.9	20	
Dibenzofuran	3.858	0.10	5	0.04005	76.4	50 - 120	4.131	6.83	20	
Di-n-butyl phthalate	6.119	0.20	5	0	122	45 - 123	4.49	30.7	20	R
Fluoranthene	5.225	0.10	5	0	105	45 - 125	5.834	11	20	
Fluorene	4.231	0.10	5	0.03736	83.9	49 - 120	4.534	6.92	20	
Naphthalene	3.7	0.10	5	0.4092	65.8	45 - 120	3.669	0.855	20	
Nitrobenzene	3.342	0.20	5	0	66.8	44 - 120	3.406	1.89	20	
N-Nitrosodiphenylamine	5.416	0.20	5	0	108	40 - 125	5.951	9.43	20	
Pentachlorophenol	5.036	0.20	5	0	101	19 - 121	5.92	16.1	20	
Phenanthrene	5.126	0.10	5	0.1179	100	45 - 121	5.624	9.27	20	
Phenol	3.5	0.20	5	0	70.0	20 - 124	3.512	0.332	20	
Pyrene	5.407	0.10	5	0	108	40 - 130	5.934	9.3	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.064</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>81.3</i>	<i>34 - 129</i>	<i>4.8</i>	<i>16.6</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.604</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.1</i>	<i>40 - 125</i>	<i>3.948</i>	<i>9.1</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.185</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.7</i>	<i>20 - 120</i>	<i>3.307</i>	<i>3.77</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>5.088</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>102</i>	<i>40 - 135</i>	<i>5.739</i>	<i>12</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.252</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>65.0</i>	<i>41 - 120</i>	<i>3.377</i>	<i>3.77</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.158</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.2</i>	<i>20 - 120</i>	<i>3.333</i>	<i>5.39</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

Batch ID: 149998 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS20010910-03MSD	Units: ug/L			Analysis Date: 29-Jan-2020 00:03					
Client ID:	Run ID: SV-6_355238	SeqNo: 5452171	PrepDate: 27-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.047	0.20	5	0	101	39 - 127	4.258	17	20	
2,4-Dimethylphenol	3.862	0.20	5	0.09533	75.3	35 - 120	2.953	26.7	20	R
2,4-Dinitrotoluene	5.46	0.20	5	0	109	50 - 122	5.24	4.12	20	
2,6-Dinitrotoluene	5.056	0.20	5	0	101	50 - 120	4.544	10.7	20	
2-Chloronaphthalene	4.215	0.20	5	0	84.3	50 - 120	3.621	15.2	20	
2-Methylnaphthalene	3.899	0.10	5	0	78.0	50 - 120	3.197	19.8	20	
4,6-Dinitro-2-methylphenol	5.595	0.20	5	0	112	25 - 121	5.327	4.9	30	
4-Nitrophenol	5.641	1.0	5	0	113	30 - 130	5.515	2.26	20	
Acenaphthene	4.245	0.10	5	0	84.9	45 - 120	3.484	19.7	20	
Acenaphthylene	4.266	0.10	5	0	85.3	47 - 120	3.505	19.6	20	
Anthracene	5.342	0.10	5	0	107	45 - 120	5.162	3.43	20	
Benz(a)anthracene	5.45	0.10	5	0	109	40 - 120	5.438	0.221	20	
Benzo(a)pyrene	5.228	0.10	5	0	105	45 - 120	5.398	3.2	20	
Bis(2-chloroethoxy)methane	3.686	0.20	5	0	73.7	45 - 120	2.957	21.9	20	R
Bis(2-ethylhexyl)phthalate	5.666	0.20	5	0	113	40 - 139	5.76	1.64	20	
Chrysene	5.136	0.10	5	0	103	43 - 120	5.195	1.14	20	
Dibenzofuran	4.314	0.10	5	0	86.3	50 - 120	3.564	19	20	
Di-n-butyl phthalate	6.065	0.20	5	0	121	45 - 123	6.041	0.393	20	
Fluoranthene	5.068	0.10	5	0.02986	101	45 - 125	5.294	4.37	20	
Fluorene	4.577	0.10	5	0	91.5	49 - 120	3.974	14.1	20	
Naphthalene	4.118	0.10	5	0.22	78.0	45 - 120	3.352	20.5	20	R
Nitrobenzene	3.792	0.20	5	0	75.8	44 - 120	3.206	16.7	20	
N-Nitrosodiphenylamine	5.62	0.20	5	0	112	40 - 125	5.464	2.8	20	
Pentachlorophenol	3.323	0.20	5	0	66.5	19 - 121	3.381	1.74	20	
Phenanthrene	5.142	0.10	5	0	103	45 - 121	4.995	2.9	20	
Phenol	3.883	0.20	5	0	77.7	20 - 124	3.121	21.7	20	R
Pyrene	5.443	0.10	5	0.02733	108	40 - 130	5.526	1.51	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.11</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>82.2</i>	<i>34 - 129</i>	<i>4.181</i>	<i>1.73</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.108</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>82.2</i>	<i>40 - 125</i>	<i>3.463</i>	<i>17</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.476</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.5</i>	<i>20 - 120</i>	<i>3.1</i>	<i>11.4</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>5.339</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>107</i>	<i>40 - 135</i>	<i>5.541</i>	<i>3.71</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.75</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>75.0</i>	<i>41 - 120</i>	<i>3.186</i>	<i>16.3</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>3.663</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.3</i>	<i>20 - 120</i>	<i>3.21</i>	<i>13.2</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010958-02 HS20010958-03 HS20010958-04 HS20010958-05  
 HS20010958-06 HS20010958-07

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

<b>Batch ID:</b> R355061 ( 0 )		<b>Instrument:</b> VOA4		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200125</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 11:52</b>				
Client ID:	Run ID: <b>VOA4_355061</b>	SeqNo: <b>5447903</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.99</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.8</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>48.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200125</b>	Units: <b>ug/L</b>			Analysis Date: <b>25-Jan-2020 11:02</b>				
Client ID:	Run ID: <b>VOA4_355061</b>	SeqNo: <b>5447902</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18.78	1.0	20	0	93.9	70 - 124			
Benzene	19.4	1.0	20	0	97.0	74 - 120			
Chlorobenzene	19.61	1.0	20	0	98.0	76 - 113			
Ethylbenzene	19.57	1.0	20	0	97.8	77 - 117			
Methylene chloride	20.36	2.0	20	0	102	70 - 127			
Toluene	19.83	1.0	20	0	99.1	77 - 118			
Vinyl chloride	19.46	1.0	20	0	97.3	70 - 130			
Xylenes, Total	58.33	1.0	60	0	97.2	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.1</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

**Batch ID:** R355061 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20010937-01MS			Units: ug/L		Analysis Date: 25-Jan-2020 15:09			
Client ID:		Run ID: VOA4_355061			SeqNo: 5447911		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.24	1.0	20	0	81.2	70 - 127				
Benzene	17.71	1.0	20	0	88.5	70 - 127				
Chlorobenzene	17.17	1.0	20	0	85.8	70 - 114				
Ethylbenzene	17.53	1.0	20	0	87.6	70 - 124				
Methylene chloride	18.41	2.0	20	0	92.1	70 - 128				
Toluene	18.12	1.0	20	0	90.6	70 - 123				
Vinyl chloride	17.88	1.0	20	0	89.4	70 - 130				
Xylenes, Total	51.84	1.0	60	0	86.4	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.64</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>51.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20010937-01MSD			Units: ug/L		Analysis Date: 25-Jan-2020 15:34			
Client ID:		Run ID: VOA4_355061			SeqNo: 5447912		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.9	1.0	20	0	79.5	70 - 127	16.24	2.11	20	
Benzene	17.13	1.0	20	0	85.6	70 - 127	17.71	3.33	20	
Chlorobenzene	17.02	1.0	20	0	85.1	70 - 114	17.17	0.844	20	
Ethylbenzene	16.92	1.0	20	0	84.6	70 - 124	17.53	3.5	20	
Methylene chloride	17.25	2.0	20	0	86.3	70 - 128	18.41	6.49	20	
Toluene	16.89	1.0	20	0	84.4	70 - 123	18.12	7.08	20	
Vinyl chloride	16.93	1.0	20	0	84.6	70 - 130	17.88	5.47	20	
Xylenes, Total	50.49	1.0	60	0	84.1	70 - 130	51.84	2.64	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.37</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>	<i>51.17</i>	<i>0.402</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.63</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>49.86</i>	<i>3.48</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>51.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>77 - 123</i>	<i>50.64</i>	<i>2.29</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.8</i>	<i>82 - 127</i>	<i>51.21</i>	<i>3.59</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010958-03      HS20010958-04      HS20010958-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

**Batch ID:** R355150 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200127</b>			Units: <b>ug/L</b>		Analysis Date: <b>28-Jan-2020 00:04</b>			
Client ID:		Run ID: <b>VOA4_355150</b>			SeqNo: <b>5449458</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>53.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>107</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>51.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>51.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200127</b>			Units: <b>ug/L</b>		Analysis Date: <b>27-Jan-2020 23:14</b>			
Client ID:		Run ID: <b>VOA4_355150</b>			SeqNo: <b>5449457</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.77	1.0	20	0	88.9	70 - 124				
Benzene	18.82	1.0	20	0	94.1	74 - 120				
Chlorobenzene	18.67	1.0	20	0	93.4	76 - 113				
Ethylbenzene	18.47	1.0	20	0	92.3	77 - 117				
Methylene chloride	20	2.0	20	0	100.0	70 - 127				
Toluene	18.6	1.0	20	0	93.0	77 - 118				
Xylenes, Total	57.11	1.0	60	0	95.2	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>51.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>49.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.7</i>	<i>81 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QC BATCH REPORT**

**Batch ID:** R355150 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20010958-02MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>28-Jan-2020 02:07</b>			
Client ID: <b>WG-1620-MW82B-20200121</b>		Run ID: <b>VOA4_355150</b>			SeqNo: <b>5449463</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.91	1.0	20	0	84.5	70 - 127				
Benzene	18.82	1.0	20	0	94.1	70 - 127				
Chlorobenzene	17.88	1.0	20	0	89.4	70 - 114				
Ethylbenzene	17.99	1.0	20	0	89.9	70 - 124				
Methylene chloride	18.93	2.0	20	0	94.6	70 - 128				
Toluene	18.76	1.0	20	0	93.8	70 - 123				
Xylenes, Total	54.55	1.0	60	0	90.9	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.92</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>51.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.57</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20010958-02MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>28-Jan-2020 02:32</b>			
Client ID: <b>WG-1620-MW82B-20200121</b>		Run ID: <b>VOA4_355150</b>			SeqNo: <b>5449464</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.29	1.0	20	0	81.4	70 - 127	16.91	3.76	20	
Benzene	17.94	1.0	20	0	89.7	70 - 127	18.82	4.77	20	
Chlorobenzene	17.53	1.0	20	0	87.7	70 - 114	17.88	1.94	20	
Ethylbenzene	17.71	1.0	20	0	88.5	70 - 124	17.99	1.57	20	
Methylene chloride	18.74	2.0	20	0	93.7	70 - 128	18.93	0.98	20	
Toluene	17.72	1.0	20	0	88.6	70 - 123	18.76	5.69	20	
Xylenes, Total	53.22	1.0	60	0	88.7	70 - 130	54.55	2.47	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>	<i>51.55</i>	<i>0.229</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.04</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>	<i>50.92</i>	<i>0.238</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>51.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>77 - 123</i>	<i>51.17</i>	<i>0.604</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>	<i>50.57</i>	<i>0.637</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20010958-01    HS20010958-02    HS20010958-06    HS20010958-07

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010958

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010958

Date/Time Received: 22-Jan-2020 11:30
Received by: AC

Checklist completed by: Paresh M. Giga
eSignature
Date: 22-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 27-Jan-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [ ] No [checked]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:206376

Temperature(s)/Thermometer(s): 0.6C; 1.7C; 1.4C u/c IR25
Cooler(s)/Kit(s): 44257/4197/42818
Date/Time sample(s) sent to storage: 1/22/2020 13:25

- Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]

pH adjusted by:

Login Notes: ID Differs. Samples identified using collection times: COC=WG-1620-MW38C-20200121 LBL=WG-1620-MW38B-20200121.
Logged ID/date/time per Label, per client direction.

Client Contacted: Golder Associates Date Contacted: 27-Jan-2020 Person Contacted: Michelle Hermiston
Contacted By: 369 Regarding: anlayte list

Comments: No sample in this WO requires reporting of vinyl chloride.

Corrective Action: Logged per client instructions to report site list VOCs only. Log in sample ID as MW38B per sample labels.



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

## HS20010958

Golder Associates Inc.  
Houston TX - Wood Preserving Works

Page \_\_\_\_ of \_\_\_\_

COC ID: 206376



ALS Project Manager:

Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0.SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive	Address	1400 Douglas Street
	Suite 4004		Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TB07-20200121	1-21-20	-	Water	1	2		X									
2	WG-1620-MW-82B-20200121	1-21-20	9:10		1	6	X	X	X	X							
3	WG-1620-MW-82BMS-20200121	1-21-20	9:10		1	6	X	X	X	X							
4	WG-1620-MW-82BMSD-20200121	1-21-20	9:10		1	6	X	X	X	X							
5	WG-1620-MW-38A-20200121	1-21-20	11:10		1	6	X	X	X	X							
6	WG-1620-MW-38C-20200121	1-21-20	13:45		1	6	X	X	X	X							
7	WG-1620-MW-03-20200121	1-21-20	15:00		1	6	X	X	X	X							
8	WG-1620-MW-04-20200121	1-21-20	16:00		1	6	X	X	X	X							
9	WG-1620-FB-11-20200121	1-21-20	15:30		1	6	X	X	X	X							
10																	

Sampler(s) Please Print & Sign <i>TIM M Spalden</i>		Shipment Method	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Deys <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24-hour		Results Due Date:
Relinquished by: <i>T. M. Spalden</i>	Date: 1-22-2020 Time: 11:30	Received by:	Notes: UPRR Houston MWPW		
Relinquished by:	Date: 1/22/2020 Time: 11:30	Received by (Laboratory): AC	Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)
Logged by (Laboratory):	Date:	Checked by (Laboratory):	44257	0.6	<input type="checkbox"/> Level II Std OC
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035			4197	1.7	<input type="checkbox"/> Level III Std OC/Raw Data
			42818	1.4	<input type="checkbox"/> Level IV SV643/CLP
					<input checked="" type="checkbox"/> TRRP Check st
					<input type="checkbox"/> TRRP Level IV

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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January 31, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20011116**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 5 sample(s) on Jan 23, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "Dane J. Wacasey".

Generated By: DAYNA.FISHER  
Dane J. Wacasey

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey



Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/31/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20011116			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 150059, 150108, R355244, R355245			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?		X			2
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			3
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				4
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/31/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20011116			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 150059, 150108, R355244, R355245			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?		X			5
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 01/31/2020
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS20011116
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 150059, 150108, R355244, R355245
ER# <sup>5</sup>	Description	
1	Batch 150108, Semivolatile Organics Method SW8270, samples WG-1620-MW41B-20200123, WG-1620-MW12B-20200123, WG-1620-MW68B-20200123: surrogate recoveries could not be determined due to dilution below the calibration range.	
2	Batch 150108, Semivolatiles by Method SW8270, LCSD RPD recovered above upper limits for 2-Methylnaphthalene due to possible matrix effect.	
3	Batch 150059, Metals by method SW6020, Sample WG-1620-MW41B-20200123, MSD recovered below lower limits for Arsenic; however, the result in the parent sample is greater than 4x the spike amount.	
4	Batch R355245, Volatiles by Method SW8260, Samples WG-1620-MW68B-20200123, WG-1620-MW12B-20200123: Lowest practical dilution due to high concentration of non-target analyte(s). Batch 150108, Semivolatiles by Method SW8270, Sample WG-1620-MW12B-20200123, The GCMS semi-volatile extract of this sample was run at a dilution due to a high level of matrix interference.	
5	Batch 150059, Metals by method SW6020, Sample WG-1620-MW41B-20200123, The percent difference between the results of the sample and the serial dilution were greater than 10% for Arsenic due to possible matrix effect.	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.</p> <p>O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);                  NA = Not Applicable;                  NR = Not Reviewed;                  R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20011116

**SAMPLE SUMMARY**

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Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20011116-01	WQ-1620-TB08-20200123	Water		23-Jan-2020 16:00	23-Jan-2020 17:50	<input type="checkbox"/>
HS20011116-02	WG-1620-MW41B-20200123	Water		23-Jan-2020 12:40	23-Jan-2020 17:50	<input type="checkbox"/>
HS20011116-03	WG-1620-MW12B-20200123	Water		23-Jan-2020 14:00	23-Jan-2020 17:50	<input type="checkbox"/>
HS20011116-04	WG-1620-MW68B-20200123	Water		23-Jan-2020 15:25	23-Jan-2020 17:50	<input type="checkbox"/>
HS20011116-05	WG-1620-FB12-20200123	Water		23-Jan-2020 13:00	23-Jan-2020 17:50	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB08-20200123  
 Collection Date: 23-Jan-2020 16:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:27
Benzene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:27
Chlorobenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 14:27
Ethylbenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 14:27
Methylene chloride	U		0.0010	0.0020	mg/L	1	28-Jan-2020 14:27
Toluene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:27
Vinyl chloride	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:27
Xylenes, Total	U		0.00030	0.0010	mg/L	1	28-Jan-2020 14:27
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:27</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:27</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:27</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:27</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW41B-20200123  
 Collection Date: 23-Jan-2020 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.00020	0.0010	mg/L	1	28-Jan-2020 15:14
<b>Benzene</b>	<b>0.012</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	28-Jan-2020 15:14
Chlorobenzene		U	0.00030	0.0010	mg/L	1	28-Jan-2020 15:14
<b>Ethylbenzene</b>	<b>0.066</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	28-Jan-2020 15:14
Methylene chloride		U	0.0010	0.0020	mg/L	1	28-Jan-2020 15:14
<b>Toluene</b>	<b>0.087</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	28-Jan-2020 15:14
Vinyl chloride		U	0.00020	0.0010	mg/L	1	28-Jan-2020 15:14
<b>Xylenes, Total</b>	<b>0.16</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	28-Jan-2020 15:14
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.2</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:14</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>102</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:14</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:14</i>
<i>Surr: Toluene-d8</i>	<i>99.5</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 15:14</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW41B-20200123  
 Collection Date: 23-Jan-2020 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	30-Jan-2020 18:04
<b>2,4-Dimethylphenol</b>	<b>0.043</b>		<b>0.00080</b>	<b>0.0040</b>	<b>mg/L</b>	20	31-Jan-2020 11:30
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	30-Jan-2020 18:04
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	30-Jan-2020 18:04
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	30-Jan-2020 18:04
<b>2-Methylnaphthalene</b>	<b>0.67</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 12:27
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	30-Jan-2020 18:04
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	30-Jan-2020 18:04
<b>Acenaphthene</b>	<b>0.49</b>		<b>0.0054</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 12:27
<b>Acenaphthylene</b>	<b>0.0089</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 18:04
<b>Anthracene</b>	<b>0.045</b>		<b>0.00028</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 11:30
<b>Benz(a)anthracene</b>	<b>0.0040</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 18:04
<b>Benzo(a)pyrene</b>	<b>0.00090</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 18:04
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	30-Jan-2020 18:04
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00032</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	30-Jan-2020 18:04
<b>Chrysene</b>	<b>0.0033</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 18:04
<b>Dibenzofuran</b>	<b>0.39</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 12:27
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	30-Jan-2020 18:04
<b>Fluoranthene</b>	<b>0.052</b>		<b>0.00020</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 11:30
<b>Fluorene</b>	<b>0.35</b>		<b>0.0060</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 12:27
<b>Naphthalene</b>	<b>10</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	31-Jan-2020 13:24
Nitrobenzene	U		0.000024	0.00020	mg/L	1	30-Jan-2020 18:04
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	30-Jan-2020 18:04
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	30-Jan-2020 18:04
<b>Phenanthrene</b>	<b>0.41</b>		<b>0.0042</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 12:27
Phenol	U		0.000035	0.00020	mg/L	1	30-Jan-2020 18:04
<b>Pyrene</b>	<b>0.029</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 11:30
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 13:24</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>86.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 18:04</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>114</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>31-Jan-2020 11:30</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>31-Jan-2020 12:27</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>104</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>31-Jan-2020 11:30</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>112</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 18:04</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>31-Jan-2020 12:27</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 13:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>31-Jan-2020 12:27</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 13:24</i>
<i>Surr: 2-Fluorophenol</i>	<i>104</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 18:04</i>
<i>Surr: 2-Fluorophenol</i>	<i>110</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>31-Jan-2020 11:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW41B-20200123  
 Collection Date: 23-Jan-2020 12:40

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	111			40-135	%REC	20	31-Jan-2020 11:30
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	31-Jan-2020 12:27
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	31-Jan-2020 13:24
Surr: 4-Terphenyl-d14	108			40-135	%REC	1	30-Jan-2020 18:04
Surr: Nitrobenzene-d5	109			41-120	%REC	20	31-Jan-2020 11:30
Surr: Nitrobenzene-d5	84.4			41-120	%REC	1	30-Jan-2020 18:04
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	31-Jan-2020 12:27
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	31-Jan-2020 13:24
Surr: Phenol-d6	0	JS		20-120	%REC	200	31-Jan-2020 12:27
Surr: Phenol-d6	0	JS		20-120	%REC	2000	31-Jan-2020 13:24
Surr: Phenol-d6	92.3			20-120	%REC	1	30-Jan-2020 18:04
Surr: Phenol-d6	116			20-120	%REC	20	31-Jan-2020 11:30
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 28-Jan-2020		Analyst: JHD	
Arsenic	0.0842		0.000400	0.00200	mg/L	1	29-Jan-2020 15:59

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12B-20200123  
 Collection Date: 23-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.0050	0.025	mg/L	25	29-Jan-2020 07:52
<b>Benzene</b>	<b>0.030</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	29-Jan-2020 07:52
Chlorobenzene	U		0.0075	0.025	mg/L	25	29-Jan-2020 07:52
<b>Ethylbenzene</b>	<b>0.014</b>	J	<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	29-Jan-2020 07:52
Methylene chloride	U		0.025	0.050	mg/L	25	29-Jan-2020 07:52
Toluene	U		0.0050	0.025	mg/L	25	29-Jan-2020 07:52
Vinyl chloride	U		0.0050	0.025	mg/L	25	29-Jan-2020 07:52
Xylenes, Total	U		0.0075	0.025	mg/L	25	29-Jan-2020 07:52
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>96.7</i>			<i>70-126</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:52</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.4</i>			<i>81-113</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:52</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.6</i>			<i>77-123</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:52</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:52</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12B-20200123  
 Collection Date: 23-Jan-2020 14:00

**ANALYTICAL REPORT**

WorkOrder:HS20011116  
 Lab ID:HS20011116-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.0021	0.020	mg/L	10	30-Jan-2020 18:23
2,4-Dimethylphenol	U		0.0040	0.020	mg/L	10	30-Jan-2020 18:23
2,4-Dinitrotoluene	U		0.0058	0.020	mg/L	10	30-Jan-2020 18:23
2,6-Dinitrotoluene	U		0.0042	0.020	mg/L	10	30-Jan-2020 18:23
2-Chloronaphthalene	U		0.0021	0.020	mg/L	10	30-Jan-2020 18:23
<b>2-Methylnaphthalene</b>	<b>320</b>		<b>3.8</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
4,6-Dinitro-2-methylphenol	U		0.0020	0.020	mg/L	10	30-Jan-2020 18:23
4-Nitrophenol	U		0.0047	0.10	mg/L	10	30-Jan-2020 18:23
<b>Acenaphthene</b>	<b>350</b>		<b>5.4</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
<b>Acenaphthylene</b>	<b>5.7</b>		<b>0.015</b>	<b>0.10</b>	<b>mg/L</b>	100	30-Jan-2020 18:42
<b>Anthracene</b>	<b>160</b>		<b>0.28</b>	<b>2.0</b>	<b>mg/L</b>	2000	31-Jan-2020 11:49
<b>Benz(a)anthracene</b>	<b>79</b>		<b>1.0</b>	<b>2.0</b>	<b>mg/L</b>	2000	31-Jan-2020 11:49
<b>Benzo(a)pyrene</b>	<b>22</b>		<b>0.40</b>	<b>2.0</b>	<b>mg/L</b>	2000	31-Jan-2020 11:49
Bis(2-chloroethoxy)methane	U		0.0030	0.020	mg/L	10	30-Jan-2020 18:23
Bis(2-ethylhexyl)phthalate	U		0.0037	0.020	mg/L	10	30-Jan-2020 18:23
<b>Chrysene</b>	<b>64</b>		<b>0.42</b>	<b>2.0</b>	<b>mg/L</b>	2000	31-Jan-2020 11:49
<b>Dibenzofuran</b>	<b>320</b>		<b>4.0</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
Di-n-butyl phthalate	U		0.0020	0.020	mg/L	10	30-Jan-2020 18:23
<b>Fluoranthene</b>	<b>520</b>		<b>2.0</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
<b>Fluorene</b>	<b>420</b>		<b>6.0</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
<b>Naphthalene</b>	<b>760</b>		<b>4.0</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
Nitrobenzene	U		0.0024	0.020	mg/L	10	30-Jan-2020 18:23
N-Nitrosodiphenylamine	U		0.0025	0.020	mg/L	10	30-Jan-2020 18:23
Pentachlorophenol	U		0.0079	0.020	mg/L	10	30-Jan-2020 18:23
<b>Phenanthrene</b>	<b>1,200</b>		<b>4.2</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
Phenol	U		0.0035	0.020	mg/L	10	30-Jan-2020 18:23
<b>Pyrene</b>	<b>330</b>		<b>3.8</b>	<b>20</b>	<b>mg/L</b>	20000	31-Jan-2020 12:46
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>20000</i>	<i>31-Jan-2020 12:46</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 11:49</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>40.5</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>30-Jan-2020 18:23</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>100</i>	<i>30-Jan-2020 18:42</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>71.2</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>30-Jan-2020 18:23</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>100</i>	<i>30-Jan-2020 18:42</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 11:49</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>20000</i>	<i>31-Jan-2020 12:46</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>20000</i>	<i>31-Jan-2020 12:46</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 11:49</i>
<i>Surr: 2-Fluorophenol</i>	<i>101</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>30-Jan-2020 18:23</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>100</i>	<i>30-Jan-2020 18:42</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW12B-20200123  
 Collection Date: 23-Jan-2020 14:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	99.5	J		40-135	%REC	10	30-Jan-2020 18:23
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	31-Jan-2020 11:49
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	20000	31-Jan-2020 12:46
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	30-Jan-2020 18:42
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	20000	31-Jan-2020 12:46
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	30-Jan-2020 18:42
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	31-Jan-2020 11:49
Surr: Nitrobenzene-d5	111	J		41-120	%REC	10	30-Jan-2020 18:23
Surr: Phenol-d6	105	J		20-120	%REC	10	30-Jan-2020 18:23
Surr: Phenol-d6	0	JS		20-120	%REC	100	30-Jan-2020 18:42
Surr: Phenol-d6	0	JS		20-120	%REC	20000	31-Jan-2020 12:46
Surr: Phenol-d6	0	JS		20-120	%REC	2000	31-Jan-2020 11:49
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 28-Jan-2020		Analyst: JC	
Arsenic	0.0491		0.000400	0.00200	mg/L	1	29-Jan-2020 16:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68B-20200123  
 Collection Date: 23-Jan-2020 15:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.0050	0.025	mg/L	25	29-Jan-2020 07:26
<b>Benzene</b>	<b>1.4</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	29-Jan-2020 07:26
Chlorobenzene		U	0.0075	0.025	mg/L	25	29-Jan-2020 07:26
<b>Ethylbenzene</b>	<b>0.39</b>		<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	29-Jan-2020 07:26
Methylene chloride		U	0.025	0.050	mg/L	25	29-Jan-2020 07:26
<b>Toluene</b>	<b>0.23</b>		<b>0.0050</b>	<b>0.025</b>	<b>mg/L</b>	25	29-Jan-2020 07:26
Vinyl chloride		U	0.0050	0.025	mg/L	25	29-Jan-2020 07:26
<b>Xylenes, Total</b>	<b>1.1</b>		<b>0.0075</b>	<b>0.025</b>	<b>mg/L</b>	25	29-Jan-2020 07:26
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>96.7</i>			<i>70-126</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:26</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.4</i>			<i>81-113</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:26</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.3</i>			<i>77-123</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:26</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	25	<i>29-Jan-2020 07:26</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68B-20200123  
 Collection Date: 23-Jan-2020 15:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	30-Jan-2020 19:01
<b>2,4-Dimethylphenol</b>	<b>0.075</b>		<b>0.00080</b>	<b>0.0040</b>	<b>mg/L</b>	20	31-Jan-2020 12:08
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	30-Jan-2020 19:01
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	30-Jan-2020 19:01
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	30-Jan-2020 19:01
<b>2-Methylnaphthalene</b>	<b>1.1</b>		<b>0.0038</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 13:05
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	30-Jan-2020 19:01
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	30-Jan-2020 19:01
<b>Acenaphthene</b>	<b>0.44</b>		<b>0.0054</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 13:05
<b>Acenaphthylene</b>	<b>0.0037</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 19:01
<b>Anthracene</b>	<b>0.15</b>		<b>0.00028</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 12:08
<b>Benz(a)anthracene</b>	<b>0.047</b>		<b>0.0010</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 12:08
<b>Benzo(a)pyrene</b>	<b>0.0089</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	30-Jan-2020 19:01
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	30-Jan-2020 19:01
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.0011</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	30-Jan-2020 19:01
<b>Chrysene</b>	<b>0.040</b>		<b>0.00042</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 12:08
<b>Dibenzofuran</b>	<b>0.50</b>		<b>0.0040</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 13:05
<b>Di-n-butyl phthalate</b>	<b>0.00018</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	30-Jan-2020 19:01
<b>Fluoranthene</b>	<b>0.31</b>		<b>0.0020</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 13:05
<b>Fluorene</b>	<b>0.32</b>		<b>0.0060</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 13:05
<b>Naphthalene</b>	<b>12</b>		<b>0.040</b>	<b>0.20</b>	<b>mg/L</b>	2000	31-Jan-2020 13:43
Nitrobenzene	U		0.000024	0.00020	mg/L	1	30-Jan-2020 19:01
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	30-Jan-2020 19:01
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	30-Jan-2020 19:01
<b>Phenanthrene</b>	<b>0.95</b>		<b>0.0042</b>	<b>0.020</b>	<b>mg/L</b>	200	31-Jan-2020 13:05
Phenol	U		0.000035	0.00020	mg/L	1	30-Jan-2020 19:01
<b>Pyrene</b>	<b>0.19</b>		<b>0.00038</b>	<b>0.0020</b>	<b>mg/L</b>	20	31-Jan-2020 12:08
<i>Surr: 2,4,6-Tribromophenol</i>	<i>112</i>			<i>34-129</i>	<i>%REC</i>	<i>20</i>	<i>31-Jan-2020 12:08</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>31-Jan-2020 13:05</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>53.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 19:01</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 13:43</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 13:43</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>56.6</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 19:01</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>115</i>			<i>40-125</i>	<i>%REC</i>	<i>20</i>	<i>31-Jan-2020 12:08</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>31-Jan-2020 13:05</i>
<i>Surr: 2-Fluorophenol</i>	<i>113</i>			<i>20-120</i>	<i>%REC</i>	<i>20</i>	<i>31-Jan-2020 12:08</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>31-Jan-2020 13:05</i>
<i>Surr: 2-Fluorophenol</i>	<i>54.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 19:01</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>31-Jan-2020 13:43</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW68B-20200123  
 Collection Date: 23-Jan-2020 15:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	31-Jan-2020 13:43
Surr: 4-Terphenyl-d14	89.9			40-135	%REC	1	30-Jan-2020 19:01
Surr: 4-Terphenyl-d14	106			40-135	%REC	20	31-Jan-2020 12:08
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	31-Jan-2020 13:05
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	31-Jan-2020 13:43
Surr: Nitrobenzene-d5	110			41-120	%REC	1	30-Jan-2020 19:01
Surr: Nitrobenzene-d5	93.0			41-120	%REC	20	31-Jan-2020 12:08
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	31-Jan-2020 13:05
Surr: Phenol-d6	99.0			20-120	%REC	20	31-Jan-2020 12:08
Surr: Phenol-d6	0	JS		20-120	%REC	200	31-Jan-2020 13:05
Surr: Phenol-d6	48.7			20-120	%REC	1	30-Jan-2020 19:01
Surr: Phenol-d6	0	JS		20-120	%REC	2000	31-Jan-2020 13:43
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 28-Jan-2020		Analyst: JC	
Arsenic	0.00944		0.000400	0.00200	mg/L	1	29-Jan-2020 16:39

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB12-20200123  
 Collection Date: 23-Jan-2020 13:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:50
Benzene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:50
Chlorobenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 14:50
Ethylbenzene	U		0.00030	0.0010	mg/L	1	28-Jan-2020 14:50
Methylene chloride	U		0.0010	0.0020	mg/L	1	28-Jan-2020 14:50
Toluene	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:50
Vinyl chloride	U		0.00020	0.0010	mg/L	1	28-Jan-2020 14:50
Xylenes, Total	U		0.00030	0.0010	mg/L	1	28-Jan-2020 14:50
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>96.0</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:50</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.7</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:50</i>
<i>Surr: Dibromofluoromethane</i>		<i>101</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:50</i>
<i>Surr: Toluene-d8</i>		<i>99.9</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>28-Jan-2020 14:50</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB12-20200123  
 Collection Date: 23-Jan-2020 13:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011116  
 Lab ID:HS20011116-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 29-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	31-Jan-2020 10:33
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	31-Jan-2020 10:33
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	31-Jan-2020 10:33
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	31-Jan-2020 10:33
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	31-Jan-2020 10:33
<b>2-Methylnaphthalene</b>	<b>0.000099</b>	<b>J</b>	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>31-Jan-2020 10:33</b>
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	31-Jan-2020 10:33
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	31-Jan-2020 10:33
Acenaphthene	U		0.000027	0.00010	mg/L	1	31-Jan-2020 10:33
Acenaphthylene	U		0.000015	0.00010	mg/L	1	31-Jan-2020 10:33
Anthracene	U		0.000014	0.00010	mg/L	1	31-Jan-2020 10:33
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	31-Jan-2020 10:33
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	31-Jan-2020 10:33
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	31-Jan-2020 10:33
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	31-Jan-2020 10:33
Chrysene	U		0.000021	0.00010	mg/L	1	31-Jan-2020 10:33
Dibenzofuran	U		0.000020	0.00010	mg/L	1	31-Jan-2020 10:33
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	31-Jan-2020 10:33
Fluoranthene	U		0.000010	0.00010	mg/L	1	31-Jan-2020 10:33
Fluorene	U		0.000030	0.00010	mg/L	1	31-Jan-2020 10:33
<b>Naphthalene</b>	<b>0.00083</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>31-Jan-2020 10:33</b>
Nitrobenzene	U		0.000024	0.00020	mg/L	1	31-Jan-2020 10:33
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	31-Jan-2020 10:33
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	31-Jan-2020 10:33
<b>Phenanthrene</b>	<b>0.000073</b>	<b>J</b>	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>31-Jan-2020 10:33</b>
Phenol	U		0.000035	0.00020	mg/L	1	31-Jan-2020 10:33
Pyrene	U		0.000019	0.00010	mg/L	1	31-Jan-2020 10:33
<i>Surr: 2,4,6-Tribromophenol</i>	<i>75.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>31-Jan-2020 10:33</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>107</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>31-Jan-2020 10:33</i>
<i>Surr: 2-Fluorophenol</i>	<i>86.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>31-Jan-2020 10:33</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>107</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>31-Jan-2020 10:33</i>
<i>Surr: Nitrobenzene-d5</i>	<i>94.4</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>31-Jan-2020 10:33</i>
<i>Surr: Phenol-d6</i>	<i>91.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>31-Jan-2020 10:33</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 28-Jan-2020		Analyst: JC	
Arsenic	U		0.000400	0.00200	mg/L	1	29-Jan-2020 16:52

Note: See Qualifiers Page for a list of qualifiers and their explanation.



## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**Batch ID:** 150059      **Start Date:** 28 Jan 2020 11:00      **End Date:** 28 Jan 2020 15:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20011116-02		10 (mL)	10 (mL)	1
HS20011116-03		10 (mL)	10 (mL)	1
HS20011116-04		10 (mL)	10 (mL)	1
HS20011116-05		10 (mL)	10 (mL)	1

**Batch ID:** 150108      **Start Date:** 29 Jan 2020 08:59      **End Date:** 29 Jan 2020 14:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20011116-02	1	1000 (mL)	1 (mL)	0.001
HS20011116-03	1	1000 (mL)	10 (mL)	0.01
HS20011116-04	1	1000 (mL)	1 (mL)	0.001
HS20011116-05	1	1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID:</b> 150059 ( 0 )		<b>Test Name :</b> ICP-MS METALS BY SW6020A			<b>Matrix:</b> Water	
HS20011116-02	WG-1620-MW41B-20200123	23 Jan 2020 12:40		28 Jan 2020 11:00	29 Jan 2020 15:59	1
HS20011116-03	WG-1620-MW12B-20200123	23 Jan 2020 14:00		28 Jan 2020 11:00	29 Jan 2020 16:37	1
HS20011116-04	WG-1620-MW68B-20200123	23 Jan 2020 15:25		28 Jan 2020 11:00	29 Jan 2020 16:39	1
HS20011116-05	WG-1620-FB12-20200123	23 Jan 2020 13:00		28 Jan 2020 11:00	29 Jan 2020 16:52	1
<b>Batch ID:</b> 150108 ( 0 )		<b>Test Name :</b> LOW-LEVEL SEMIVOLATILES BY 8270D			<b>Matrix:</b> Water	
HS20011116-02	WG-1620-MW41B-20200123	23 Jan 2020 12:40		29 Jan 2020 08:59	31 Jan 2020 12:27	200
HS20011116-02	WG-1620-MW41B-20200123	23 Jan 2020 12:40		29 Jan 2020 08:59	31 Jan 2020 13:24	2000
HS20011116-02	WG-1620-MW41B-20200123	23 Jan 2020 12:40		29 Jan 2020 08:59	31 Jan 2020 11:30	20
HS20011116-02	WG-1620-MW41B-20200123	23 Jan 2020 12:40		29 Jan 2020 08:59	30 Jan 2020 18:04	1
HS20011116-03	WG-1620-MW12B-20200123	23 Jan 2020 14:00		29 Jan 2020 08:59	31 Jan 2020 12:46	2000 0
HS20011116-03	WG-1620-MW12B-20200123	23 Jan 2020 14:00		29 Jan 2020 08:59	31 Jan 2020 11:49	2000
HS20011116-03	WG-1620-MW12B-20200123	23 Jan 2020 14:00		29 Jan 2020 08:59	30 Jan 2020 18:42	100
HS20011116-03	WG-1620-MW12B-20200123	23 Jan 2020 14:00		29 Jan 2020 08:59	30 Jan 2020 18:23	10
HS20011116-04	WG-1620-MW68B-20200123	23 Jan 2020 15:25		29 Jan 2020 08:59	31 Jan 2020 13:05	200
HS20011116-04	WG-1620-MW68B-20200123	23 Jan 2020 15:25		29 Jan 2020 08:59	31 Jan 2020 13:43	2000
HS20011116-04	WG-1620-MW68B-20200123	23 Jan 2020 15:25		29 Jan 2020 08:59	31 Jan 2020 12:08	20
HS20011116-04	WG-1620-MW68B-20200123	23 Jan 2020 15:25		29 Jan 2020 08:59	30 Jan 2020 19:01	1
HS20011116-05	WG-1620-FB12-20200123	23 Jan 2020 13:00		29 Jan 2020 08:59	31 Jan 2020 10:33	1
<b>Batch ID:</b> R355244 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Water	
HS20011116-01	WQ-1620-TB08-20200123	23 Jan 2020 16:00			28 Jan 2020 14:27	1
HS20011116-02	WG-1620-MW41B-20200123	23 Jan 2020 12:40			28 Jan 2020 15:14	1
HS20011116-05	WG-1620-FB12-20200123	23 Jan 2020 13:00			28 Jan 2020 14:50	1
<b>Batch ID:</b> R355245 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Water	
HS20011116-03	WG-1620-MW12B-20200123	23 Jan 2020 14:00			29 Jan 2020 07:52	25
HS20011116-04	WG-1620-MW68B-20200123	23 Jan 2020 15:25			29 Jan 2020 07:26	25

WorkOrder: HS20011116  
 InstrumentID: ICPMS06  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000479	0.000400	0.00200

WorkOrder: HS20011116  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20011116  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

Matrix: Aqueous

Units: mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20011116  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

Batch ID: 150059 ( 0 )		Instrument: ICPMS06		Method: ICP-MS METALS BY SW6020A						
<b>MBLK</b>	Sample ID: <b>MBLK-150059</b>	Units: <b>mg/L</b>		Analysis Date: <b>29-Jan-2020 15:56</b>						
Client ID:		Run ID: <b>ICPMS06_355299</b>	SeqNo: <b>5452423</b>	PrepDate: <b>28-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-150059</b>	Units: <b>mg/L</b>		Analysis Date: <b>29-Jan-2020 15:58</b>						
Client ID:		Run ID: <b>ICPMS06_355299</b>	SeqNo: <b>5452424</b>	PrepDate: <b>28-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.04514	0.00200	0.05	0	90.3	80 - 120				
<b>MS</b>	Sample ID: <b>HS20011116-02MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>29-Jan-2020 16:03</b>						
Client ID: <b>WG-1620-MW41B-20200123</b>		Run ID: <b>ICPMS06_355299</b>	SeqNo: <b>5452427</b>	PrepDate: <b>28-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.1253	0.00200	0.05	0.08425	82.0	80 - 120				
<b>MSD</b>	Sample ID: <b>HS20011116-02MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>29-Jan-2020 16:04</b>						
Client ID: <b>WG-1620-MW41B-20200123</b>		Run ID: <b>ICPMS06_355299</b>	SeqNo: <b>5452428</b>	PrepDate: <b>28-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.1164	0.00200	0.05	0.08425	64.3	80 - 120	0.1253	7.35	20	S
<b>PDS</b>	Sample ID: <b>HS20011116-02PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>29-Jan-2020 16:06</b>						
Client ID: <b>WG-1620-MW41B-20200123</b>		Run ID: <b>ICPMS06_355299</b>	SeqNo: <b>5452429</b>	PrepDate: <b>28-Jan-2020</b>	DF: <b>1</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit	Qual
Arsenic	0.1758	0.00200	0.1	0.08425	91.6	75 - 125				
<b>SD</b>	Sample ID: <b>HS20011116-02SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>29-Jan-2020 16:01</b>						
Client ID: <b>WG-1620-MW41B-20200123</b>		Run ID: <b>ICPMS06_355299</b>	SeqNo: <b>5452426</b>	PrepDate: <b>28-Jan-2020</b>	DF: <b>5</b>					
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual
Arsenic	0.07491	0.0100					0.08425	11.1	10	R

The following samples were analyzed in this batch: HS20011116-02 HS20011116-03 HS20011116-04 HS20011116-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

Batch ID: 150108 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-150108	Units: ug/L			Analysis Date: 29-Jan-2020 20:03					
Client ID:	Run ID: SV-6_355351	SeqNo: 5453256	PrepDate: 29-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.224</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>84.5</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.201</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>104</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.679</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.6</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.066</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.664</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.624</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>92.5</i>	<i>20 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

Batch ID: 150108 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-150108	Units: ug/L			Analysis Date: 29-Jan-2020 20:23					
Client ID:	Run ID: SV-6_355351	SeqNo: 5453257	PrepDate: 29-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	5.107	0.20	5	0	102	39 - 127				
2,4-Dimethylphenol	4.701	0.20	5	0	94.0	35 - 120				
2,4-Dinitrotoluene	5.219	0.20	5	0	104	50 - 122				
2,6-Dinitrotoluene	5.149	0.20	5	0	103	50 - 120				
2-Chloronaphthalene	4.991	0.20	5	0	99.8	50 - 120				
2-Methylnaphthalene	4.769	0.10	5	0	95.4	50 - 120				
4,6-Dinitro-2-methylphenol	5.839	0.20	5	0	117	25 - 121				
4-Nitrophenol	5.028	1.0	5	0	101	30 - 130				
Acenaphthene	5.038	0.10	5	0	101	45 - 120				
Acenaphthylene	4.847	0.10	5	0	96.9	47 - 120				
Anthracene	5.231	0.10	5	0	105	45 - 120				
Benz(a)anthracene	4.934	0.10	5	0	98.7	40 - 120				
Benzo(a)pyrene	4.902	0.10	5	0	98.0	45 - 120				
Bis(2-chloroethoxy)methane	4.639	0.20	5	0	92.8	45 - 120				
Bis(2-ethylhexyl)phthalate	5.985	0.20	5	0	120	40 - 139				
Chrysene	4.767	0.10	5	0	95.3	43 - 120				
Dibenzofuran	4.752	0.10	5	0	95.0	50 - 120				
Di-n-butyl phthalate	5.839	0.20	5	0	117	45 - 123				
Fluoranthene	5.012	0.10	5	0	100	45 - 125				
Fluorene	4.867	0.10	5	0	97.3	49 - 120				
Naphthalene	4.782	0.10	5	0	95.6	45 - 120				
Nitrobenzene	4.42	0.20	5	0	88.4	44 - 120				
N-Nitrosodiphenylamine	5.618	0.20	5	0	112	40 - 125				
Pentachlorophenol	4.372	0.20	5	0	87.4	19 - 121				
Phenanthrene	4.959	0.10	5	0	99.2	45 - 121				
Phenol	4.747	0.20	5	0	94.9	20 - 124				
Pyrene	5.005	0.10	5	0	100	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.299</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>86.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>5.038</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>101</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.503</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.1</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.95</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>99.0</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.54</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.8</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.729</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.6</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

Batch ID: 150108 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-150108		Units: ug/L		Analysis Date: 29-Jan-2020 20:42				
Client ID:		Run ID: SV-6_355351		SeqNo: 5453258		PrepDate: 29-Jan-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	5.262	0.20	5	0	105	39 - 127	5.107	3	20	
2,4-Dimethylphenol	4.491	0.20	5	0	89.8	35 - 120	4.701	4.56	20	
2,4-Dinitrotoluene	5.375	0.20	5	0	107	50 - 122	5.219	2.94	20	
2,6-Dinitrotoluene	5.408	0.20	5	0	108	50 - 120	5.149	4.9	20	
2-Chloronaphthalene	4.958	0.20	5	0	99.2	50 - 120	4.991	0.651	20	
2-Methylnaphthalene	5.835	0.10	5	0	117	50 - 120	4.769	20.1	20 R	
4,6-Dinitro-2-methylphenol	5.428	0.20	5	0	109	25 - 121	5.839	7.3	30	
4-Nitrophenol	5.067	1.0	5	0	101	30 - 130	5.028	0.77	20	
Acenaphthene	5.053	0.10	5	0	101	45 - 120	5.038	0.302	20	
Acenaphthylene	4.875	0.10	5	0	97.5	47 - 120	4.847	0.584	20	
Anthracene	5.218	0.10	5	0	104	45 - 120	5.231	0.257	20	
Benz(a)anthracene	5.073	0.10	5	0	101	40 - 120	4.934	2.78	20	
Benzo(a)pyrene	4.915	0.10	5	0	98.3	45 - 120	4.902	0.266	20	
Bis(2-chloroethoxy)methane	4.715	0.20	5	0	94.3	45 - 120	4.639	1.63	20	
Bis(2-ethylhexyl)phthalate	5.942	0.20	5	0	119	40 - 139	5.985	0.725	20	
Chrysene	4.855	0.10	5	0	97.1	43 - 120	4.767	1.83	20	
Dibenzofuran	4.808	0.10	5	0	96.2	50 - 120	4.752	1.19	20	
Di-n-butyl phthalate	5.854	0.20	5	0	117	45 - 123	5.839	0.253	20	
Fluoranthene	4.9	0.10	5	0	98.0	45 - 125	5.012	2.27	20	
Fluorene	4.841	0.10	5	0	96.8	49 - 120	4.867	0.542	20	
Naphthalene	4.712	0.10	5	0	94.2	45 - 120	4.782	1.48	20	
Nitrobenzene	4.576	0.20	5	0	91.5	44 - 120	4.42	3.47	20	
N-Nitrosodiphenylamine	5.763	0.20	5	0	115	40 - 125	5.618	2.55	20	
Pentachlorophenol	4.251	0.20	5	0	85.0	19 - 121	4.372	2.79	20	
Phenanthrene	4.876	0.10	5	0	97.5	45 - 121	4.959	1.69	20	
Phenol	4.406	0.20	5	0	88.1	20 - 124	4.747	7.46	20	
Pyrene	4.965	0.10	5	0	99.3	40 - 130	5.005	0.811	20	
Surr: 2,4,6-Tribromophenol	4.039	0.20	5	0	80.8	34 - 129	4.299	6.25	20	
Surr: 2-Fluorobiphenyl	4.979	0.20	5	0	99.6	40 - 125	5.038	1.19	20	
Surr: 2-Fluorophenol	4.378	0.20	5	0	87.6	20 - 120	4.503	2.81	20	
Surr: 4-Terphenyl-d14	4.766	0.20	5	0	95.3	40 - 135	4.95	3.77	20	
Surr: Nitrobenzene-d5	4.575	0.20	5	0	91.5	41 - 120	4.54	0.784	20	
Surr: Phenol-d6	4.363	0.20	5	0	87.3	20 - 120	4.729	8.06	20	

The following samples were analyzed in this batch: HS20011116-02 HS20011116-03 HS20011116-04 HS20011116-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

<b>Batch ID:</b> R355244 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200128</b>	Units: <b>ug/L</b>			Analysis Date: <b>28-Jan-2020 13:40</b>				
Client ID:	Run ID: <b>VOA2_355244</b>	SeqNo: <b>5451226</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>47.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.8</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.9</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200128</b>	Units: <b>ug/L</b>			Analysis Date: <b>28-Jan-2020 12:53</b>				
Client ID:	Run ID: <b>VOA2_355244</b>	SeqNo: <b>5451225</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	18.02	1.0	20	0	90.1	70 - 124			
Benzene	17.64	1.0	20	0	88.2	74 - 120			
Chlorobenzene	17.96	1.0	20	0	89.8	76 - 113			
Ethylbenzene	17.99	1.0	20	0	90.0	77 - 117			
Methylene chloride	19.04	2.0	20	0	95.2	70 - 127			
Toluene	19.88	1.0	20	0	99.4	77 - 118			
Vinyl chloride	17.17	1.0	20	0	85.9	70 - 130			
Xylenes, Total	55.27	1.0	60	0	92.1	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.41</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.8</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

Batch ID: R355244 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>		Sample ID: <b>HS20011116-02MS</b>		Units: <b>ug/L</b>		Analysis Date: <b>28-Jan-2020 15:37</b>				
Client ID: <b>WG-1620-MW41B-20200123</b>		Run ID: <b>VOA2_355244</b>		SeqNo: <b>5451231</b>		PrepDate:		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	16.84	1.0	20	0	84.2	70 - 127				
Benzene	29.47	1.0	20	12.37	85.5	70 - 127				
Chlorobenzene	17.11	1.0	20	0	85.5	70 - 114				
Ethylbenzene	84.99	1.0	20	66.49	92.5	70 - 124				
Methylene chloride	17.37	2.0	20	0	86.9	70 - 128				
Toluene	110.3	1.0	20	86.56	119	70 - 123			O	
Vinyl chloride	20.48	1.0	20	0	102	70 - 130				
Xylenes, Total	219.7	1.0	60	164.8	91.4	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>				
<b>MSD</b>		Sample ID: <b>HS20011116-02MSD</b>		Units: <b>ug/L</b>		Analysis Date: <b>28-Jan-2020 16:01</b>				
Client ID: <b>WG-1620-MW41B-20200123</b>		Run ID: <b>VOA2_355244</b>		SeqNo: <b>5451232</b>		PrepDate:		DF: <b>1</b>		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	16.56	1.0	20	0	82.8	70 - 127	16.84	1.64	20	
Benzene	28.93	1.0	20	12.37	82.8	70 - 127	29.47	1.84	20	
Chlorobenzene	16.61	1.0	20	0	83.0	70 - 114	17.11	2.97	20	
Ethylbenzene	82.23	1.0	20	66.49	78.7	70 - 124	84.99	3.3	20	
Methylene chloride	17.54	2.0	20	0	87.7	70 - 128	17.37	0.981	20	
Toluene	106.3	1.0	20	86.56	98.7	70 - 123	110.3	3.69	20 O	
Vinyl chloride	19.88	1.0	20	0	99.4	70 - 130	20.48	2.99	20	
Xylenes, Total	212.9	1.0	60	164.8	80.1	70 - 130	219.7	3.15	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.45</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 126</i>	<i>50.84</i>	<i>1.18</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>51.97</i>	<i>0.464</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>77 - 123</i>	<i>49.5</i>	<i>0.578</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.75</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>82 - 127</i>	<i>50.03</i>	<i>0.56</i>	<i>20</i>	
The following samples were analyzed in this batch: HS20011116-01 HS20011116-02 HS20011116-05										

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

<b>Batch ID:</b> R355245 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200128</b>	Units: <b>ug/L</b>			Analysis Date: <b>29-Jan-2020 01:03</b>				
Client ID:	Run ID: <b>VOA2_355245</b>	SeqNo: <b>5451281</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.86</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.5</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.02</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200128</b>	Units: <b>ug/L</b>			Analysis Date: <b>29-Jan-2020 00:16</b>				
Client ID:	Run ID: <b>VOA2_355245</b>	SeqNo: <b>5451280</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	19.02	1.0	20	0	95.1	70 - 124			
Benzene	18.45	1.0	20	0	92.3	74 - 120			
Chlorobenzene	18.55	1.0	20	0	92.7	76 - 113			
Ethylbenzene	18.58	1.0	20	0	92.9	77 - 117			
Methylene chloride	19.05	2.0	20	0	95.2	70 - 127			
Toluene	20.61	1.0	20	0	103	77 - 118			
Vinyl chloride	17.95	1.0	20	0	89.8	70 - 130			
Xylenes, Total	57.33	1.0	60	0	95.6	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.22</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.6</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.7</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QC BATCH REPORT**

**Batch ID:** R355245 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20011181-02MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>29-Jan-2020 03:23</b>			
Client ID:		Run ID: <b>VOA2_355245</b>			SeqNo: <b>5451287</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.01	1.0	20	0	90.1	70 - 127				
Benzene	18.01	1.0	20	0	90.0	70 - 127				
Chlorobenzene	18.25	1.0	20	0	91.2	70 - 114				
Ethylbenzene	18.69	1.0	20	0	93.4	70 - 124				
Methylene chloride	18.42	2.0	20	0	92.1	70 - 128				
Toluene	20.55	1.0	20	0	103	70 - 123				
Vinyl chloride	17.79	1.0	20	0	89.0	70 - 130				
Xylenes, Total	57.19	1.0	60	0	95.3	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.9</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.79</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.38</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20011181-02MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>29-Jan-2020 03:47</b>			
Client ID:		Run ID: <b>VOA2_355245</b>			SeqNo: <b>5451288</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.35	1.0	20	0	86.7	70 - 127	18.01	3.74	20	
Benzene	17.3	1.0	20	0	86.5	70 - 127	18.01	3.99	20	
Chlorobenzene	17.41	1.0	20	0	87.0	70 - 114	18.25	4.73	20	
Ethylbenzene	17.98	1.0	20	0	89.9	70 - 124	18.69	3.85	20	
Methylene chloride	17.49	2.0	20	0	87.5	70 - 128	18.42	5.17	20	
Toluene	19.27	1.0	20	0	96.3	70 - 123	20.55	6.44	20	
Vinyl chloride	16.9	1.0	20	0	84.5	70 - 130	17.79	5.17	20	
Xylenes, Total	54.86	1.0	60	0	91.4	70 - 130	57.19	4.16	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>70 - 126</i>	<i>51.14</i>	<i>2.81</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>81 - 113</i>	<i>49.9</i>	<i>0.117</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.63</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.3</i>	<i>77 - 123</i>	<i>49.79</i>	<i>0.325</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.46</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>50.38</i>	<i>0.144</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20011116-03      HS20011116-04

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011116

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

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**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20011116

**SAMPLE TRACKING**

Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS20011116-01	WQ-1620-TB08-20200123	Login	1/24/2020 2:47:57 PM	JRM	VOA160
HS20011116-02	WG-1620-MW41B-20200123	Login	1/24/2020 2:47:57 PM	JRM	MET011
HS20011116-02	WG-1620-MW41B-20200123	Login	1/24/2020 2:47:57 PM	JRM	EXT043
HS20011116-02	WG-1620-MW41B-20200123	Login	1/24/2020 2:47:57 PM	JRM	VOA160
HS20011116-03	WG-1620-MW12B-20200123	Login	1/24/2020 2:47:57 PM	JRM	MET011
HS20011116-03	WG-1620-MW12B-20200123	Login	1/24/2020 2:47:57 PM	JRM	EXT043
HS20011116-03	WG-1620-MW12B-20200123	Login	1/24/2020 2:47:57 PM	JRM	VOA160
HS20011116-04	WG-1620-MW68B-20200123	Login	1/24/2020 2:47:57 PM	JRM	MET011
HS20011116-04	WG-1620-MW68B-20200123	Login	1/24/2020 2:47:57 PM	JRM	EXT043
HS20011116-04	WG-1620-MW68B-20200123	Login	1/24/2020 2:47:57 PM	JRM	VOA160
HS20011116-05	WG-1620-FB12-20200123	Login	1/24/2020 2:47:57 PM	JRM	MET011
HS20011116-05	WG-1620-FB12-20200123	Login	1/24/2020 2:47:57 PM	JRM	EXT043
HS20011116-05	WG-1620-FB12-20200123	Login	1/24/2020 2:47:57 PM	JRM	VOA160

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20011116

Date/Time Received: 23-Jan-2020 17:50
Received by: DDG

Checklist completed by: Jared R. Makan
eSignature
Date: 24-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 27-Jan-2020

Matrices: Water

Carrier name: ALS Courier

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:214254

Temperature(s)/Thermometer(s): 2.0°C/2.0°C UC/C IR25
Cooler(s)/Kit(s): 44479
Date/Time sample(s) sent to storage: 01/24/2020 14:55
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by: [ ]

Login Notes: Samples refrigerated prior to login.

Client Contacted: Golder Associates Date Contacted: 27-Jan-2020 Person Contacted: Michelle Hermiston
Contacted By: 369 Regarding: anlayte list

Comments: No sample in this WO requires reporting of vinyl chloride.

Corrective Action: Logged per client instructions to report site list VOCs only.



Cincinnati, OH  
+1 513 733 5336  
Everett, WA  
+1 425 356 2600

Fort Collins, CO  
+1 970 490 1511  
Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page 1 of 1

COC ID: 214254

HS20011116

Golder Associates Inc.  
Houston TX-Wood Preserving Works

WV



Customer Information		Project Information		ALS Project Manager:											
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632528 Volatile Organics Site Specific)										
Work Order		Project Number	1620-07-Rev0 SR 92688	B	8260_LL_W (5632528 VOC Site Specific + V.C.)										
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)										
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 5652646 Metals - As)										
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E											
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	F											
Phone	(512) 671-3434	Phone		G											
Fax	(512) 671-3446	Fax		H											
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		I											
				J											

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TB08-20200123	1-23-20	16:00	Water	1	2		X									
2	WG-1620-MW41B-20200123	1-23-20	12:40	H <sub>2</sub> O	1	6	X	X	X	X							
3	WG-1620-MW12B-20200123	1-23-20	14:00	H <sub>2</sub> O	1	6	X	X	X	X							
4	WG-1620-MW68B-20200123	1-23-20	15:25	H <sub>2</sub> O	1	6	X	X	X	X							
5	WG-1620-FB12-20200123	1-23-20	13:00	H <sub>2</sub> O	1	6	X	X	X	X							
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>Tim McSpadden</i> <i>T. McSpadden</i>		Shipment Method	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 1 Wk Days <input type="checkbox"/> 24 Hour				Results Due Date:	
Relinquished by: <i>T. McSpadden</i>	Date: 1-23-20	Time: 16:30	Received by: <i>D.S.</i>		Notes: UPRR Houston MWPW			
Relinquished by: <i>D.S.</i>	Date: 1-23-20	Time: 17:50	Received by (Laboratory): <i>D.S.</i>		Cooler ID 444-701	Cooler Temp. 4K 2.0	QC Package: (Check One Box Below)	
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):		<input type="checkbox"/> Level II Std QC	<input type="checkbox"/> Level III Std QC/Raw Data	<input checked="" type="checkbox"/> TTRP Checklist	<input type="checkbox"/> TTRP Level IV
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035					<input type="checkbox"/> Level IV SIMMS/CLP			

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
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February 04, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20011249**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 5 sample(s) on Jan 27, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

---

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 02/04/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20011249			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 150153,150187,R355340,R355399			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?		X			2
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?		X			3
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				4
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 02/04/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20011249			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 150153,150187,R355340,R355399			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?		X			5
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).



**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 02/04/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20011249
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 150153,150187,R355340,R355399

ER# <sup>5</sup>	Description
1	Semivolatile Organics Method SW8270, sample WG-1620-MW49B-20200127, the surrogate recoveries could not be determined due to dilution below the calibration range.
2	Batch R355399, Volatile Organics Method SW8260, LCS recovery was above the control limits for Vinyl chloride. The analyte was not detected in the associated samples.
3	Batch 150187, Semivolatile Organics Method SW8270, sample HS20011317-01, MS and MSD were performed on unrelated sample.
4	Batch 150187, Semivolatile Organics Method SW8270, sample WG-1620-MW49B-20200127: The GCMS semi-volatile extract of this sample was run at a dilution due to a high level of matrix interference.
5	Batch R355399, Volatile Organics Method SW8260, Vinyl chloride exceeded % recovery limits on CCV. The analyte was not detected in the associated samples.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
 O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
 NA = Not Applicable;  
 NR = Not Reviewed;  
 R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20011249

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20011249-01	WQ-1620-TB09-20200127	Water		27-Jan-2020 15:00	27-Jan-2020 17:15	<input type="checkbox"/>
HS20011249-02	WG-1620-MW62B-20200127	Water		27-Jan-2020 09:25	27-Jan-2020 17:15	<input type="checkbox"/>
HS20011249-03	WG-1620-MW64A-20200127	Water		27-Jan-2020 11:15	27-Jan-2020 17:15	<input type="checkbox"/>
HS20011249-04	WG-1620-MW49B-20200127	Water		27-Jan-2020 13:30	27-Jan-2020 17:15	<input type="checkbox"/>
HS20011249-05	WG-1620-FB13-20200127	Water		27-Jan-2020 12:00	27-Jan-2020 17:15	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB09-20200127  
 Collection Date: 27-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	29-Jan-2020 13:55
Benzene	U		0.00020	0.0010	mg/L	1	29-Jan-2020 13:55
Chlorobenzene	U		0.00030	0.0010	mg/L	1	29-Jan-2020 13:55
Ethylbenzene	U		0.00030	0.0010	mg/L	1	29-Jan-2020 13:55
Methylene chloride	U		0.0010	0.0020	mg/L	1	29-Jan-2020 13:55
Toluene	U		0.00020	0.0010	mg/L	1	29-Jan-2020 13:55
Vinyl chloride	U		0.00020	0.0010	mg/L	1	29-Jan-2020 13:55
Xylenes, Total	U		0.00030	0.0010	mg/L	1	29-Jan-2020 13:55
<i>Surr: 1,2-Dichloroethane-d4</i>		<i>97.2</i>		<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 13:55</i>
<i>Surr: 4-Bromofluorobenzene</i>		<i>94.5</i>		<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 13:55</i>
<i>Surr: Dibromofluoromethane</i>		<i>101</i>		<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 13:55</i>
<i>Surr: Toluene-d8</i>		<i>102</i>		<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 13:55</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW62B-20200127  
 Collection Date: 27-Jan-2020 09:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	30-Jan-2020 17:20
Benzene	U		0.00020	0.0010	mg/L	1	30-Jan-2020 17:20
Chlorobenzene	U		0.00030	0.0010	mg/L	1	30-Jan-2020 17:20
Ethylbenzene	U		0.00030	0.0010	mg/L	1	30-Jan-2020 17:20
Methylene chloride	U		0.0010	0.0020	mg/L	1	30-Jan-2020 17:20
Toluene	U		0.00020	0.0010	mg/L	1	30-Jan-2020 17:20
Xylenes, Total	U		0.00030	0.0010	mg/L	1	30-Jan-2020 17:20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>97.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.4</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:20</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:20</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:20</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW62B-20200127  
 Collection Date: 27-Jan-2020 09:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 31-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	03-Feb-2020 12:44
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	03-Feb-2020 12:44
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	03-Feb-2020 12:44
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	03-Feb-2020 12:44
<b>2-Chloronaphthalene</b>	<b>0.00016</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>2-Methylnaphthalene</b>	<b>0.00011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	03-Feb-2020 12:44
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	03-Feb-2020 12:44
<b>Acenaphthene</b>	<b>0.040</b>		<b>0.00011</b>	<b>0.00040</b>	<b>mg/L</b>	4	04-Feb-2020 12:03
<b>Acenaphthylene</b>	<b>0.00029</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>Anthracene</b>	<b>0.00096</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	03-Feb-2020 12:44
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	03-Feb-2020 12:44
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	03-Feb-2020 12:44
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000095</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>Chrysene</b>	<b>0.000027</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>Dibenzofuran</b>	<b>0.0042</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>Di-n-butyl phthalate</b>	<b>0.00014</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>Fluoranthene</b>	<b>0.0020</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<b>Fluorene</b>	<b>0.011</b>		<b>0.00012</b>	<b>0.00040</b>	<b>mg/L</b>	4	04-Feb-2020 12:03
<b>Naphthalene</b>	<b>0.0010</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
Nitrobenzene	U		0.000024	0.00020	mg/L	1	03-Feb-2020 12:44
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	03-Feb-2020 12:44
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	03-Feb-2020 12:44
<b>Phenanthrene</b>	<b>0.00087</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
Phenol	U		0.000035	0.00020	mg/L	1	03-Feb-2020 12:44
<b>Pyrene</b>	<b>0.0013</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 12:44
<i>Surr: 2,4,6-Tribromophenol</i>	<i>73.7</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 12:44</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>65.6</i>			<i>34-129</i>	<i>%REC</i>	<i>4</i>	<i>04-Feb-2020 12:03</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>83.1</i>			<i>40-125</i>	<i>%REC</i>	<i>4</i>	<i>04-Feb-2020 12:03</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>79.2</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 12:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 12:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>52.6</i>			<i>20-120</i>	<i>%REC</i>	<i>4</i>	<i>04-Feb-2020 12:03</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.6</i>			<i>40-135</i>	<i>%REC</i>	<i>4</i>	<i>04-Feb-2020 12:03</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>94.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 12:44</i>
<i>Surr: Nitrobenzene-d5</i>	<i>71.9</i>			<i>41-120</i>	<i>%REC</i>	<i>4</i>	<i>04-Feb-2020 12:03</i>
<i>Surr: Nitrobenzene-d5</i>	<i>72.8</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 12:44</i>
<i>Surr: Phenol-d6</i>	<i>71.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 12:44</i>
<i>Surr: Phenol-d6</i>	<i>69.0</i>			<i>20-120</i>	<i>%REC</i>	<i>4</i>	<i>04-Feb-2020 12:03</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW62B-20200127  
 Collection Date: 27-Jan-2020 09:25

**ANALYTICAL REPORT**

WorkOrder:HS20011249  
 Lab ID:HS20011249-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 30-Jan-2020		Analyst: JHD
Arsenic	0.0161		0.000400	0.00200	mg/L	1	31-Jan-2020 16:08

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW64A-20200127  
 Collection Date: 27-Jan-2020 11:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	30-Jan-2020 17:44
Benzene	U		0.00020	0.0010	mg/L	1	30-Jan-2020 17:44
Chlorobenzene	U		0.00030	0.0010	mg/L	1	30-Jan-2020 17:44
Ethylbenzene	U		0.00030	0.0010	mg/L	1	30-Jan-2020 17:44
Methylene chloride	U		0.0010	0.0020	mg/L	1	30-Jan-2020 17:44
Toluene	U		0.00020	0.0010	mg/L	1	30-Jan-2020 17:44
Xylenes, Total	U		0.00030	0.0010	mg/L	1	30-Jan-2020 17:44
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.9</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:44</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:44</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:44</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>30-Jan-2020 17:44</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW64A-20200127  
 Collection Date: 27-Jan-2020 11:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 31-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	03-Feb-2020 13:04
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	03-Feb-2020 13:04
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	03-Feb-2020 13:04
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	03-Feb-2020 13:04
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	03-Feb-2020 13:04
<b>2-Methylnaphthalene</b>	<b>0.00037</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	03-Feb-2020 13:04
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	03-Feb-2020 13:04
<b>Acenaphthene</b>	<b>0.00016</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
Acenaphthylene	U		0.000015	0.00010	mg/L	1	03-Feb-2020 13:04
<b>Anthracene</b>	<b>0.000096</b>	J	<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	03-Feb-2020 13:04
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	03-Feb-2020 13:04
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	03-Feb-2020 13:04
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00031</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
Chrysene	U		0.000021	0.00010	mg/L	1	03-Feb-2020 13:04
<b>Dibenzofuran</b>	<b>0.000100</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
<b>Di-n-butyl phthalate</b>	<b>0.000079</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
<b>Fluoranthene</b>	<b>0.00028</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
<b>Fluorene</b>	<b>0.000082</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
<b>Naphthalene</b>	<b>0.0020</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
Nitrobenzene	U		0.000024	0.00020	mg/L	1	03-Feb-2020 13:04
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	03-Feb-2020 13:04
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	03-Feb-2020 13:04
Phenanthrene	U		0.000021	0.00010	mg/L	1	03-Feb-2020 13:04
Phenol	U		0.000035	0.00020	mg/L	1	03-Feb-2020 13:04
<b>Pyrene</b>	<b>0.00015</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	03-Feb-2020 13:04
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.5</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 13:04</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>75.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 13:04</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 13:04</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 13:04</i>
<i>Surr: Nitrobenzene-d5</i>	<i>71.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 13:04</i>
<i>Surr: Phenol-d6</i>	<i>69.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>03-Feb-2020 13:04</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jan-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00126</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	31-Jan-2020 16:27

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49B-20200127  
 Collection Date: 27-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane		U	0.010	0.050	mg/L	50	30-Jan-2020 18:07
<b>Benzene</b>	<b>0.60</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	50	30-Jan-2020 18:07
Chlorobenzene		U	0.015	0.050	mg/L	50	30-Jan-2020 18:07
<b>Ethylbenzene</b>	<b>0.26</b>		<b>0.015</b>	<b>0.050</b>	<b>mg/L</b>	50	30-Jan-2020 18:07
Methylene chloride		U	0.050	0.10	mg/L	50	30-Jan-2020 18:07
<b>Toluene</b>	<b>0.72</b>		<b>0.010</b>	<b>0.050</b>	<b>mg/L</b>	50	30-Jan-2020 18:07
Vinyl chloride		U	0.010	0.050	mg/L	50	30-Jan-2020 18:07
<b>Xylenes, Total</b>	<b>0.71</b>		<b>0.015</b>	<b>0.050</b>	<b>mg/L</b>	50	30-Jan-2020 18:07
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.9</i>			<i>70-126</i>	<i>%REC</i>	<i>50</i>	<i>30-Jan-2020 18:07</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.8</i>			<i>81-113</i>	<i>%REC</i>	<i>50</i>	<i>30-Jan-2020 18:07</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.0</i>			<i>77-123</i>	<i>%REC</i>	<i>50</i>	<i>30-Jan-2020 18:07</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>50</i>	<i>30-Jan-2020 18:07</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49B-20200127  
 Collection Date: 27-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 31-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.0013	0.012	mg/L	10	03-Feb-2020 19:05
<b>2,4-Dimethylphenol</b>	<b>25</b>		<b>0.48</b>	<b>2.4</b>	<b>mg/L</b>	2000	04-Feb-2020 12:22
2,4-Dinitrotoluene	U		0.0035	0.012	mg/L	10	03-Feb-2020 19:05
2,6-Dinitrotoluene	U		0.0025	0.012	mg/L	10	03-Feb-2020 19:05
2-Chloronaphthalene	U		0.0013	0.012	mg/L	10	03-Feb-2020 19:05
<b>2-Methylnaphthalene</b>	<b>250</b>		<b>2.3</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
4,6-Dinitro-2-methylphenol	U		0.0012	0.012	mg/L	10	03-Feb-2020 19:05
4-Nitrophenol	U		0.0028	0.060	mg/L	10	03-Feb-2020 19:05
<b>Acenaphthene</b>	<b>190</b>		<b>3.2</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
<b>Acenaphthylene</b>	<b>2.1</b>		<b>0.018</b>	<b>0.12</b>	<b>mg/L</b>	200	04-Feb-2020 11:44
<b>Anthracene</b>	<b>87</b>		<b>0.17</b>	<b>1.2</b>	<b>mg/L</b>	2000	04-Feb-2020 12:22
<b>Benz(a)anthracene</b>	<b>23</b>		<b>0.60</b>	<b>1.2</b>	<b>mg/L</b>	2000	04-Feb-2020 12:22
<b>Benzo(a)pyrene</b>	<b>7.5</b>		<b>0.024</b>	<b>0.12</b>	<b>mg/L</b>	200	04-Feb-2020 11:44
Bis(2-chloroethoxy)methane	U		0.0018	0.012	mg/L	10	03-Feb-2020 19:05
Bis(2-ethylhexyl)phthalate	U		0.0022	0.012	mg/L	10	03-Feb-2020 19:05
<b>Chrysene</b>	<b>23</b>		<b>0.25</b>	<b>1.2</b>	<b>mg/L</b>	2000	04-Feb-2020 12:22
<b>Dibenzofuran</b>	<b>160</b>		<b>2.4</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
Di-n-butyl phthalate	U		0.0012	0.012	mg/L	10	03-Feb-2020 19:05
<b>Fluoranthene</b>	<b>170</b>		<b>1.2</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
<b>Fluorene</b>	<b>170</b>		<b>3.6</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
<b>Naphthalene</b>	<b>1,200</b>		<b>2.4</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
Nitrobenzene	U		0.0014	0.012	mg/L	10	03-Feb-2020 19:05
N-Nitrosodiphenylamine	U		0.0015	0.012	mg/L	10	03-Feb-2020 19:05
Pentachlorophenol	U		0.0047	0.012	mg/L	10	03-Feb-2020 19:05
<b>Phenanthrene</b>	<b>500</b>		<b>2.5</b>	<b>12</b>	<b>mg/L</b>	20000	04-Feb-2020 12:40
Phenol	U		0.0021	0.012	mg/L	10	03-Feb-2020 19:05
<b>Pyrene</b>	<b>96</b>		<b>0.23</b>	<b>1.2</b>	<b>mg/L</b>	2000	04-Feb-2020 12:22
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>200</i>	<i>04-Feb-2020 11:44</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>125</i>	<i>J</i>		<i>34-129</i>	<i>%REC</i>	<i>10</i>	<i>03-Feb-2020 19:05</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>2000</i>	<i>04-Feb-2020 12:22</i>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>0</i>	<i>JS</i>		<i>34-129</i>	<i>%REC</i>	<i>20000</i>	<i>04-Feb-2020 12:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>2000</i>	<i>04-Feb-2020 12:22</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>20000</i>	<i>04-Feb-2020 12:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>95.1</i>	<i>J</i>		<i>40-125</i>	<i>%REC</i>	<i>10</i>	<i>03-Feb-2020 19:05</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>0</i>	<i>JS</i>		<i>40-125</i>	<i>%REC</i>	<i>200</i>	<i>04-Feb-2020 11:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>200</i>	<i>04-Feb-2020 11:44</i>
<i>Surr: 2-Fluorophenol</i>	<i>94.0</i>	<i>J</i>		<i>20-120</i>	<i>%REC</i>	<i>10</i>	<i>03-Feb-2020 19:05</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>2000</i>	<i>04-Feb-2020 12:22</i>
<i>Surr: 2-Fluorophenol</i>	<i>0</i>	<i>JS</i>		<i>20-120</i>	<i>%REC</i>	<i>20000</i>	<i>04-Feb-2020 12:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW49B-20200127  
 Collection Date: 27-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 31-Jan-2020		Analyst: LG	
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	2000	04-Feb-2020 12:22
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	20000	04-Feb-2020 12:40
Surr: 4-Terphenyl-d14	67.0	J		40-135	%REC	10	03-Feb-2020 19:05
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	200	04-Feb-2020 11:44
Surr: Nitrobenzene-d5	86.3	J		41-120	%REC	10	03-Feb-2020 19:05
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	200	04-Feb-2020 11:44
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	2000	04-Feb-2020 12:22
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	20000	04-Feb-2020 12:40
Surr: Phenol-d6	0	JS		20-120	%REC	200	04-Feb-2020 11:44
Surr: Phenol-d6	0	JS		20-120	%REC	2000	04-Feb-2020 12:22
Surr: Phenol-d6	0	JS		20-120	%REC	20000	04-Feb-2020 12:40
Surr: Phenol-d6	51.0	J		20-120	%REC	10	03-Feb-2020 19:05
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jan-2020		Analyst: JHD	
Arsenic	0.0107		0.000400	0.00200	mg/L	1	31-Jan-2020 16:29

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB13-20200127  
 Collection Date: 27-Jan-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	29-Jan-2020 14:18
Benzene	U		0.00020	0.0010	mg/L	1	29-Jan-2020 14:18
Chlorobenzene	U		0.00030	0.0010	mg/L	1	29-Jan-2020 14:18
Ethylbenzene	U		0.00030	0.0010	mg/L	1	29-Jan-2020 14:18
Methylene chloride	U		0.0010	0.0020	mg/L	1	29-Jan-2020 14:18
Toluene	U		0.00020	0.0010	mg/L	1	29-Jan-2020 14:18
Vinyl chloride	U		0.00020	0.0010	mg/L	1	29-Jan-2020 14:18
Xylenes, Total	U		0.00030	0.0010	mg/L	1	29-Jan-2020 14:18
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>95.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 14:18</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 14:18</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 14:18</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>29-Jan-2020 14:18</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB13-20200127  
 Collection Date: 27-Jan-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20011249  
 Lab ID:HS20011249-05  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 31-Jan-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	03-Feb-2020 13:23
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	03-Feb-2020 13:23
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	03-Feb-2020 13:23
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	03-Feb-2020 13:23
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	03-Feb-2020 13:23
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	03-Feb-2020 13:23
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	03-Feb-2020 13:23
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	03-Feb-2020 13:23
Acenaphthene	U		0.000027	0.00010	mg/L	1	03-Feb-2020 13:23
Acenaphthylene	U		0.000015	0.00010	mg/L	1	03-Feb-2020 13:23
Anthracene	U		0.000014	0.00010	mg/L	1	03-Feb-2020 13:23
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	03-Feb-2020 13:23
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	03-Feb-2020 13:23
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	03-Feb-2020 13:23
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	03-Feb-2020 13:23
Chrysene	U		0.000021	0.00010	mg/L	1	03-Feb-2020 13:23
Dibenzofuran	U		0.000020	0.00010	mg/L	1	03-Feb-2020 13:23
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	03-Feb-2020 13:23
Fluoranthene	U		0.000010	0.00010	mg/L	1	03-Feb-2020 13:23
Fluorene	U		0.000030	0.00010	mg/L	1	03-Feb-2020 13:23
Naphthalene	U		0.000020	0.00010	mg/L	1	03-Feb-2020 13:23
Nitrobenzene	U		0.000024	0.00020	mg/L	1	03-Feb-2020 13:23
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	03-Feb-2020 13:23
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	03-Feb-2020 13:23
Phenanthrene	U		0.000021	0.00010	mg/L	1	03-Feb-2020 13:23
Phenol	U		0.000035	0.00020	mg/L	1	03-Feb-2020 13:23
Pyrene	U		0.000019	0.00010	mg/L	1	03-Feb-2020 13:23
<i>Surr: 2,4,6-Tribromophenol</i>		55.7		34-129	%REC	1	03-Feb-2020 13:23
<i>Surr: 2-Fluorobiphenyl</i>		83.6		40-125	%REC	1	03-Feb-2020 13:23
<i>Surr: 2-Fluorophenol</i>		68.1		20-120	%REC	1	03-Feb-2020 13:23
<i>Surr: 4-Terphenyl-d14</i>		92.5		40-135	%REC	1	03-Feb-2020 13:23
<i>Surr: Nitrobenzene-d5</i>		83.5		41-120	%REC	1	03-Feb-2020 13:23
<i>Surr: Phenol-d6</i>		69.4		20-120	%REC	1	03-Feb-2020 13:23
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 30-Jan-2020		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	31-Jan-2020 16:31

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**Batch ID:** 150153      **Start Date:** 30 Jan 2020 09:00      **End Date:** 30 Jan 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20011249-02		10 (mL)	10 (mL)	1
HS20011249-03		10 (mL)	10 (mL)	1
HS20011249-04		10 (mL)	10 (mL)	1
HS20011249-05		10 (mL)	10 (mL)	1

**Batch ID:** 150187      **Start Date:** 31 Jan 2020 08:41      **End Date:** 31 Jan 2020 14:00  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20011249-02	1	1000 (mL)	1 (mL)	0.001
HS20011249-03	1	1000 (mL)	1 (mL)	0.001
HS20011249-04	1	1000 (mL)	6 (mL)	0.006
HS20011249-05	1	1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 150153 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20011249-02	WG-1620-MW62B-20200127	27 Jan 2020 09:25		30 Jan 2020 13:00	31 Jan 2020 16:08	1
HS20011249-03	WG-1620-MW64A-20200127	27 Jan 2020 11:15		30 Jan 2020 13:00	31 Jan 2020 16:27	1
HS20011249-04	WG-1620-MW49B-20200127	27 Jan 2020 13:30		30 Jan 2020 13:00	31 Jan 2020 16:29	1
HS20011249-05	WG-1620-FB13-20200127	27 Jan 2020 12:00		30 Jan 2020 13:00	31 Jan 2020 16:31	1
<b>Batch ID: 150187 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20011249-02	WG-1620-MW62B-20200127	27 Jan 2020 09:25		31 Jan 2020 08:41	04 Feb 2020 12:03	4
HS20011249-02	WG-1620-MW62B-20200127	27 Jan 2020 09:25		31 Jan 2020 08:41	03 Feb 2020 12:44	1
HS20011249-03	WG-1620-MW64A-20200127	27 Jan 2020 11:15		31 Jan 2020 08:41	03 Feb 2020 13:04	1
HS20011249-04	WG-1620-MW49B-20200127	27 Jan 2020 13:30		31 Jan 2020 08:41	04 Feb 2020 12:40	2000 0
HS20011249-04	WG-1620-MW49B-20200127	27 Jan 2020 13:30		31 Jan 2020 08:41	04 Feb 2020 12:22	2000
HS20011249-04	WG-1620-MW49B-20200127	27 Jan 2020 13:30		31 Jan 2020 08:41	04 Feb 2020 11:44	200
HS20011249-04	WG-1620-MW49B-20200127	27 Jan 2020 13:30		31 Jan 2020 08:41	03 Feb 2020 19:05	10
HS20011249-05	WG-1620-FB13-20200127	27 Jan 2020 12:00		31 Jan 2020 08:41	03 Feb 2020 13:23	1
<b>Batch ID: R355340 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20011249-01	WQ-1620-TB09-20200127	27 Jan 2020 15:00			29 Jan 2020 13:55	1
HS20011249-05	WG-1620-FB13-20200127	27 Jan 2020 12:00			29 Jan 2020 14:18	1
<b>Batch ID: R355399 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20011249-02	WG-1620-MW62B-20200127	27 Jan 2020 09:25			30 Jan 2020 17:20	1
HS20011249-03	WG-1620-MW64A-20200127	27 Jan 2020 11:15			30 Jan 2020 17:44	1
HS20011249-04	WG-1620-MW49B-20200127	27 Jan 2020 13:30			30 Jan 2020 18:07	50

WorkOrder: HS20011249  
InstrumentID: ICPMS05  
Test Code: ICP\_TW  
Test Number: SW6020  
Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200



WorkOrder: HS20011249  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20011249  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Vinyl chloride	75-01-4	0.00050	0.00044	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

**Batch ID:** 150153 ( 0 )      **Instrument:** ICPMS05      **Method:** ICP-MS METALS BY SW6020A

**MBLK**      Sample ID: **MBLK-150153**      Units: **mg/L**      Analysis Date: **31-Jan-2020 15:23**  
 Client ID:      Run ID: **ICPMS05\_355457**      SeqNo: **5456059**      PrepDate: **30-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      U      0.00200

**LCS**      Sample ID: **LCS-150153**      Units: **mg/L**      Analysis Date: **31-Jan-2020 15:25**  
 Client ID:      Run ID: **ICPMS05\_355457**      SeqNo: **5456060**      PrepDate: **30-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.04913      0.00200      0.05      0      98.3      80 - 120

**MS**      Sample ID: **HS20011246-19MS**      Units: **mg/L**      Analysis Date: **31-Jan-2020 15:32**  
 Client ID:      Run ID: **ICPMS05\_355457**      SeqNo: **5456063**      PrepDate: **30-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.05845      0.00200      0.05      0.01101      94.9      80 - 120

**MSD**      Sample ID: **HS20011246-19MSD**      Units: **mg/L**      Analysis Date: **31-Jan-2020 15:34**  
 Client ID:      Run ID: **ICPMS05\_355457**      SeqNo: **5456064**      PrepDate: **30-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.06129      0.00200      0.05      0.01101      101      80 - 120      0.05845      4.73      20

**PDS**      Sample ID: **HS20011246-19PDS**      Units: **mg/L**      Analysis Date: **31-Jan-2020 15:36**  
 Client ID:      Run ID: **ICPMS05\_355457**      SeqNo: **5456065**      PrepDate: **30-Jan-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

Arsenic      0.1101      0.00200      0.1      0.01101      99.1      75 - 125

**SD**      Sample ID: **HS20011246-19SD**      Units: **mg/L**      Analysis Date: **31-Jan-2020 15:30**  
 Client ID:      Run ID: **ICPMS05\_355457**      SeqNo: **5456062**      PrepDate: **30-Jan-2020**      DF: **5**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %D      %D Limit Qual

Arsenic      0.01196      0.0100      0.01101      8.65      10

The following samples were analyzed in this batch: HS20011249-02      HS20011249-03      HS20011249-04      HS20011249-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

Batch ID: 150187 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-150187	Units: ug/L			Analysis Date: 03-Feb-2020 10:47					
Client ID:	Run ID: SV-6_355515	SeqNo: 5458500	PrepDate: 31-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	3.568	0.20	5	0	71.4	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.784	0.20	5	0	95.7	40 - 125				
<i>Surr: 2-Fluorophenol</i>	3.407	0.20	5	0	68.1	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.819	0.20	5	0	96.4	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	4.429	0.20	5	0	88.6	41 - 120				
<i>Surr: Phenol-d6</i>	4.155	0.20	5	0	83.1	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

Batch ID: 150187 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-150187	Units: ug/L			Analysis Date: 03-Feb-2020 11:06					
Client ID:	Run ID: SV-6_355515	SeqNo: 5458501	PrepDate: 31-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.974	0.20	5	0	99.5	39 - 127				
2,4-Dimethylphenol	4.102	0.20	5	0	82.0	35 - 120				
2,4-Dinitrotoluene	5.156	0.20	5	0	103	50 - 122				
2,6-Dinitrotoluene	5.308	0.20	5	0	106	50 - 120				
2-Chloronaphthalene	4.849	0.20	5	0	97.0	50 - 120				
2-Methylnaphthalene	4.446	0.10	5	0	88.9	50 - 120				
4,6-Dinitro-2-methylphenol	5.753	0.20	5	0	115	25 - 121				
4-Nitrophenol	5.112	1.0	5	0	102	30 - 130				
Acenaphthene	4.544	0.10	5	0	90.9	45 - 120				
Acenaphthylene	4.692	0.10	5	0	93.8	47 - 120				
Anthracene	4.942	0.10	5	0	98.8	45 - 120				
Benz(a)anthracene	4.879	0.10	5	0	97.6	40 - 120				
Benzo(a)pyrene	4.857	0.10	5	0	97.1	45 - 120				
Bis(2-chloroethoxy)methane	4.461	0.20	5	0	89.2	45 - 120				
Bis(2-ethylhexyl)phthalate	5.993	0.20	5	0	120	40 - 139				
Chrysene	4.777	0.10	5	0	95.5	43 - 120				
Dibenzofuran	4.611	0.10	5	0	92.2	50 - 120				
Di-n-butyl phthalate	5.627	0.20	5	0	113	45 - 123				
Fluoranthene	4.764	0.10	5	0	95.3	45 - 125				
Fluorene	4.707	0.10	5	0	94.1	49 - 120				
Naphthalene	4.784	0.10	5	0	95.7	45 - 120				
Nitrobenzene	4.389	0.20	5	0	87.8	44 - 120				
N-Nitrosodiphenylamine	5.269	0.20	5	0	105	40 - 125				
Pentachlorophenol	3.426	0.20	5	0	68.5	19 - 121				
Phenanthrene	4.766	0.10	5	0	95.3	45 - 121				
Phenol	4.401	0.20	5	0	88.0	20 - 124				
Pyrene	4.891	0.10	5	0	97.8	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.879</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.6</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.775</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>95.5</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.464</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>89.3</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.729</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.364</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.354</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.1</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

Batch ID: 150187 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS	Sample ID: HS20011317-01MS	Units: ug/L			Analysis Date: 03-Feb-2020 11:44					
Client ID:	Run ID: SV-6_355515	SeqNo: 5458503	PrepDate: 31-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.376	0.20	5	0	87.5	39 - 127				
2,4-Dimethylphenol	2.967	0.20	5	0	59.3	35 - 120				
2,4-Dinitrotoluene	5.103	0.20	5	0	102	50 - 122				
2,6-Dinitrotoluene	4.756	0.20	5	0	95.1	50 - 120				
2-Chloronaphthalene	3.688	0.20	5	0	73.8	50 - 120				
2-Methylnaphthalene	3.335	0.10	5	0	66.7	50 - 120				
4,6-Dinitro-2-methylphenol	6.191	0.20	5	0	124	25 - 121				S
4-Nitrophenol	5.628	1.0	5	0	113	30 - 130				
Acenaphthene	3.612	0.10	5	0	72.2	45 - 120				
Acenaphthylene	3.752	0.10	5	0	75.0	47 - 120				
Anthracene	4.894	0.10	5	0	97.9	45 - 120				
Benz(a)anthracene	5.344	0.10	5	0	107	40 - 120				
Benzo(a)pyrene	5.125	0.10	5	0	102	45 - 120				
Bis(2-chloroethoxy)methane	3.215	0.20	5	0	64.3	45 - 120				
Bis(2-ethylhexyl)phthalate	6.461	0.20	5	0	129	40 - 139				
Chrysene	4.879	0.10	5	0	97.6	43 - 120				
Dibenzofuran	3.829	0.10	5	0	76.6	50 - 120				
Di-n-butyl phthalate	5.862	0.20	5	0	117	45 - 123				
Fluoranthene	4.836	0.10	5	0	96.7	45 - 125				
Fluorene	4.126	0.10	5	0	82.5	49 - 120				
Naphthalene	3.422	0.10	5	0	68.4	45 - 120				
Nitrobenzene	3.178	0.20	5	0	63.6	44 - 120				
N-Nitrosodiphenylamine	5.08	0.20	5	0	102	40 - 125				
Pentachlorophenol	4.285	0.20	5	0	85.7	19 - 121				
Phenanthrene	4.616	0.10	5	0	92.3	45 - 121				
Phenol	3.402	0.20	5	0	68.0	20 - 124				
Pyrene	5.192	0.10	5	0	104	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.916</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.634</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.7</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.269</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>65.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.243</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.151</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.0</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.227</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>64.5</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

Batch ID: 150187 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD	Sample ID: HS20011317-01MSD	Units: ug/L			Analysis Date: 03-Feb-2020 12:03					
Client ID:	Run ID: SV-6_355515	SeqNo: 5458504	PrepDate: 31-Jan-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.949	0.20	5	0	99.0	39 - 127	4.376	12.3	20	
2,4-Dimethylphenol	3.31	0.20	5	0	66.2	35 - 120	2.967	10.9	20	
2,4-Dinitrotoluene	5.611	0.20	5	0	112	50 - 122	5.103	9.48	20	
2,6-Dinitrotoluene	5.203	0.20	5	0	104	50 - 120	4.756	8.98	20	
2-Chloronaphthalene	4.007	0.20	5	0	80.1	50 - 120	3.688	8.28	20	
2-Methylnaphthalene	3.668	0.10	5	0	73.4	50 - 120	3.335	9.52	20	
4,6-Dinitro-2-methylphenol	6.078	0.20	5	0	122	25 - 121	6.191	1.83	30	S
4-Nitrophenol	5.579	1.0	5	0	112	30 - 130	5.628	0.874	20	
Acenaphthene	4.043	0.10	5	0	80.9	45 - 120	3.612	11.3	20	
Acenaphthylene	4.189	0.10	5	0	83.8	47 - 120	3.752	11	20	
Anthracene	5.189	0.10	5	0	104	45 - 120	4.894	5.85	20	
Benz(a)anthracene	5.317	0.10	5	0	106	40 - 120	5.344	0.514	20	
Benzo(a)pyrene	5.152	0.10	5	0	103	45 - 120	5.125	0.537	20	
Bis(2-chloroethoxy)methane	3.514	0.20	5	0	70.3	45 - 120	3.215	8.88	20	
Bis(2-ethylhexyl)phthalate	6.589	0.20	5	0	132	40 - 139	6.461	1.97	20	
Chrysene	4.918	0.10	5	0	98.4	43 - 120	4.879	0.794	20	
Dibenzofuran	4.316	0.10	5	0	86.3	50 - 120	3.829	12	20	
Di-n-butyl phthalate	6.118	0.20	5	0	122	45 - 123	5.862	4.28	20	
Fluoranthene	5.049	0.10	5	0	101	45 - 125	4.836	4.3	20	
Fluorene	4.612	0.10	5	0	92.2	49 - 120	4.126	11.1	20	
Naphthalene	3.753	0.10	5	0	75.1	45 - 120	3.422	9.23	20	
Nitrobenzene	3.535	0.20	5	0	70.7	44 - 120	3.178	10.6	20	
N-Nitrosodiphenylamine	5.664	0.20	5	0	113	40 - 125	5.08	10.9	20	
Pentachlorophenol	4.308	0.20	5	0	86.2	19 - 121	4.285	0.544	20	
Phenanthrene	5.046	0.10	5	0	101	45 - 121	4.616	8.9	20	
Phenol	3.764	0.20	5	0	75.3	20 - 124	3.402	10.1	20	
Pyrene	5.302	0.10	5	0	106	40 - 130	5.192	2.08	20	
Surr: 2,4,6-Tribromophenol	4.079	0.20	5	0	81.6	34 - 129	3.916	4.06	20	
Surr: 2-Fluorobiphenyl	3.951	0.20	5	0	79.0	40 - 125	3.634	8.37	20	
Surr: 2-Fluorophenol	3.335	0.20	5	0	66.7	20 - 120	3.269	1.99	20	
Surr: 4-Terphenyl-d14	5.006	0.20	5	0	100	40 - 135	5.243	4.61	20	
Surr: Nitrobenzene-d5	3.487	0.20	5	0	69.7	41 - 120	3.151	10.1	20	
Surr: Phenol-d6	3.341	0.20	5	0	66.8	20 - 120	3.227	3.47	20	

The following samples were analyzed in this batch: HS20011249-02 HS20011249-03 HS20011249-04 HS20011249-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

<b>Batch ID:</b> R355340 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200129</b>	Units: <b>ug/L</b>			Analysis Date: <b>29-Jan-2020 13:32</b>				
Client ID:	Run ID: <b>VOA2_355340</b>	SeqNo: <b>5452958</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.0</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.81</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.6</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>50.34</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>50.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200129</b>	Units: <b>ug/L</b>			Analysis Date: <b>29-Jan-2020 12:45</b>				
Client ID:	Run ID: <b>VOA2_355340</b>	SeqNo: <b>5452957</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	19.14	1.0	20	0	95.7	70 - 124			
Benzene	18.94	1.0	20	0	94.7	74 - 120			
Chlorobenzene	19.34	1.0	20	0	96.7	76 - 113			
Ethylbenzene	19.56	1.0	20	0	97.8	77 - 117			
Methylene chloride	19.55	2.0	20	0	97.8	70 - 127			
Toluene	21.49	1.0	20	0	107	77 - 118			
Vinyl chloride	25.95	1.0	20	0	130	70 - 130			
Xylenes, Total	59.72	1.0	60	0	99.5	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.15</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.47</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.73</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>81 - 120</i>			



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

**Batch ID:** R355340 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20011265-01MS			Units: ug/L		Analysis Date: 29-Jan-2020 15:29			
Client ID:		Run ID: VOA2_355340			SeqNo: 5452963		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	15.88	1.0	20	0	79.4	70 - 127				
Benzene	16.38	1.0	20	0	81.9	70 - 127				
Chlorobenzene	16.56	1.0	20	0	82.8	70 - 114				
Ethylbenzene	16.93	1.0	20	0	84.7	70 - 124				
Methylene chloride	16.47	2.0	20	0	82.4	70 - 128				
Toluene	18.24	1.0	20	0	91.2	70 - 123				
Vinyl chloride	16.8	1.0	20	0	84.0	70 - 130				
Xylenes, Total	51.65	1.0	60	0	86.1	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.12</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.5</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.3</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.6</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.51</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20011265-01MSD			Units: ug/L		Analysis Date: 29-Jan-2020 15:52			
Client ID:		Run ID: VOA2_355340			SeqNo: 5452964		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.17	1.0	20	0	80.9	70 - 127	15.88	1.84	20	
Benzene	16.4	1.0	20	0	82.0	70 - 127	16.38	0.119	20	
Chlorobenzene	16.88	1.0	20	0	84.4	70 - 114	16.56	1.96	20	
Ethylbenzene	17.43	1.0	20	0	87.2	70 - 124	16.93	2.9	20	
Methylene chloride	16.38	2.0	20	0	81.9	70 - 128	16.47	0.562	20	
Toluene	18.87	1.0	20	0	94.4	70 - 123	18.24	3.44	20	
Vinyl chloride	16.92	1.0	20	0	84.6	70 - 130	16.8	0.697	20	
Xylenes, Total	52.68	1.0	60	0	87.8	70 - 130	51.65	1.98	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>70 - 126</i>	<i>50.12</i>	<i>0.362</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.3</i>	<i>81 - 113</i>	<i>49.26</i>	<i>0.814</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.1</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.2</i>	<i>77 - 123</i>	<i>49.3</i>	<i>0.406</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.69</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>49.51</i>	<i>2.35</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20011249-01      HS20011249-05

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

<b>Batch ID:</b> R355399 ( 0 )		<b>Instrument:</b> VOA2		<b>Method:</b> LOW LEVEL VOLATILES BY SW8260C					
<b>MBLK</b>	Sample ID: <b>VBLKW-200130</b>	Units: <b>ug/L</b>			Analysis Date: <b>30-Jan-2020 12:03</b>				
Client ID:	Run ID: <b>VOA2_355399</b>	SeqNo: <b>5454133</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual

1,2-Dichloroethane	U	1.0							
Benzene	U	1.0							
Chlorobenzene	U	1.0							
Ethylbenzene	U	1.0							
Methylene chloride	U	2.0							
Toluene	U	1.0							
Vinyl chloride	U	1.0							
Xylenes, Total	U	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.1</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.69</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.4</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.74</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>			

<b>LCS</b>	Sample ID: <b>VLCSW-200130</b>	Units: <b>ug/L</b>			Analysis Date: <b>30-Jan-2020 10:53</b>				
Client ID:	Run ID: <b>VOA2_355399</b>	SeqNo: <b>5454132</b>		PrepDate:		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	18.56	1.0	20	0	92.8	70 - 124			
Benzene	18.33	1.0	20	0	91.7	74 - 120			
Chlorobenzene	18.39	1.0	20	0	91.9	76 - 113			
Ethylbenzene	18.57	1.0	20	0	92.8	77 - 117			
Methylene chloride	19.02	2.0	20	0	95.1	70 - 127			
Toluene	20.59	1.0	20	0	103	77 - 118			
Vinyl chloride	26.92	1.0	20	0	135	70 - 130			S
Xylenes, Total	57.19	1.0	60	0	95.3	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.23</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.05</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.85</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.7</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.6</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QC BATCH REPORT**

**Batch ID:** R355399 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

MS		Sample ID: HS20011350-01MS			Units: ug/L		Analysis Date: 30-Jan-2020 19:20			
Client ID:		Run ID: VOA2_355399			SeqNo: 5456624		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.07	1.0	20	0	85.4	70 - 127				
Benzene	18.93	1.0	20	0	94.6	70 - 127				
Chlorobenzene	17.61	1.0	20	0	88.0	70 - 114				
Ethylbenzene	18.17	1.0	20	0	90.9	70 - 124				
Methylene chloride	17.9	2.0	20	0	89.5	70 - 128				
Toluene	20.39	1.0	20	0	102	70 - 123				
Vinyl chloride	17.48	1.0	20	0	87.4	70 - 130				
Xylenes, Total	54.48	1.0	60	0	90.8	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.06</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.3</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.26</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

MSD		Sample ID: HS20011350-01MSD			Units: ug/L		Analysis Date: 30-Jan-2020 19:43			
Client ID:		Run ID: VOA2_355399			SeqNo: 5456625		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	16.78	1.0	20	0	83.9	70 - 127	17.07	1.7	20	
Benzene	17.77	1.0	20	0	88.9	70 - 127	18.93	6.29	20	
Chlorobenzene	17.08	1.0	20	0	85.4	70 - 114	17.61	3.01	20	
Ethylbenzene	17.53	1.0	20	0	87.6	70 - 124	18.17	3.61	20	
Methylene chloride	17.33	2.0	20	0	86.6	70 - 128	17.9	3.26	20	
Toluene	19.29	1.0	20	0	96.5	70 - 123	20.39	5.51	20	
Vinyl chloride	16.52	1.0	20	0	82.6	70 - 130	17.48	5.61	20	
Xylenes, Total	52.79	1.0	60	0	88.0	70 - 130	54.48	3.16	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.87</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>	<i>51.06</i>	<i>0.375</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>81 - 113</i>	<i>49.88</i>	<i>1.86</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>49.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>77 - 123</i>	<i>49.67</i>	<i>0.362</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.5</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>50.26</i>	<i>0.473</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20011249-02      HS20011249-03      HS20011249-04

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20011249

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

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**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231 V009	22-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20011249

**SAMPLE TRACKING**

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Lab Samp ID	Client Sample ID	Action	Date	Person	New Location
HS20011249-01	WQ-1620-TB09-20200127	Login	1/27/2020 8:50:44 PM	NDR	VOA186
HS20011249-02	WG-1620-MW62B-20200127	Login	1/27/2020 8:50:44 PM	NDR	MET040
HS20011249-02	WG-1620-MW62B-20200127	Login	1/27/2020 8:50:44 PM	NDR	EXT053
HS20011249-02	WG-1620-MW62B-20200127	Login	1/27/2020 8:50:44 PM	NDR	VOA186

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20011249

Date/Time Received: 27-Jan-2020 17:15
Received by: TW

Checklist completed by: Nilesh D. Ranchod
eSignature Date 27-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature Date 29-Jan-2020

Matrices: Water

Carrier name: ALS Courier

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:206379

Temperature(s)/Thermometer(s): 0.6°C UC/C IR25
Cooler(s)/Kit(s): 45495
Date/Time sample(s) sent to storage: 01/27/2020 9:20PM
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page      of     

COC ID: 206379

HS20011249

Golder Associates Inc.  
Houston TX-Wood Preserving Works

n, wv  
3



Customer Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-07-Rev0 SR 92688
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive	Address	1400 Douglas Street
	Suite 4004		Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address	

A	8260_LL_W (5632528 Volatile Organics Site Specific)
B	8260_LL_W (5632528 VOC Site Specific + V.C.)
C	8270_LOW_W (5632532 SemiVolatiles Site specific)
D	ICP_TW (5636002 5652646 Metals - As)
E	Vinyl Chloride
F	
G	
H	
I	
J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TB09-20200127	1-27-20	15:00	Water	1	2		X									
2	WG-1620 MW 62B 20200127	1-27-20	9:25	H <sub>2</sub> O	1	6	X	X	X	X							
3	WG-1620 MW 64A 20200127	1-27-20	11:15	H <sub>2</sub> O	1	6	X	X	X	X							
4	WG-1620 MW 49B 20200127	1-27-20	13:30	H <sub>2</sub> O	1	6	X	X	X	X	X						
5	WG-1620 FB 13 20200127	1-27-20	12:00	H <sub>2</sub> O	1	6	X	X	X	X							
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>Tim McSpadden</i>		Shipment Method	Required Turnaround Time: (Check Box)		Results Due Date:
<i>Tim McSpadden</i>			<input checked="" type="checkbox"/> STD 10 Wk Days	<input type="checkbox"/> 5 Wk Days	<input type="checkbox"/> 2 Wk Days
Relinquished by:	Date: 1-27-20	Time: 4:20	Received by:		Notes: UPRR Houston MWPW
<i>Tim McSpadden</i>	Date: 1/27/20	Time: 17:15	Received by (Laboratory):		Cooler ID: 45495
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):		Cooler Temp: 0.6
					QC Package: (Check One Box Below)
					<input type="checkbox"/> Level II Std QC
					<input type="checkbox"/> Level III Std QC/Raw Data
					<input checked="" type="checkbox"/> TRRP Check st
					<input type="checkbox"/> TRRP Level IV
					<input type="checkbox"/> Level IV SW846/CLP
					<input type="checkbox"/> Other

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

ote: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.





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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

February 18, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20020456**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 4 sample(s) on Feb 11, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 02/18/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20020456			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 150549,150594,R356238			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?	X				
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			1
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 02/18/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20020456			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 150549,150594,R356238			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section</b>					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods (NELAC Chap 5 or ISO/IEC 17025 Section 5)</b>					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 02/18/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20020456
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 150549,150594,R356238

<b>ER#<sup>5</sup></b>	<b>Description</b>
1	Batch 150549, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20020456

**SAMPLE SUMMARY**

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Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20020456-01	WQ-1620-TB11-20200211	Water	CG-121719 -91	11-Feb-2020 12:00	11-Feb-2020 14:40	<input type="checkbox"/>
HS20020456-02	WG-1620-MW22AR-20200210	Water		10-Feb-2020 10:00	11-Feb-2020 14:40	<input type="checkbox"/>
HS20020456-03	WG-1620-MW22BR-20200210	Water		10-Feb-2020 11:00	11-Feb-2020 14:40	<input type="checkbox"/>
HS20020456-04	WG-1620-FB14-20200210	Water		10-Feb-2020 11:30	11-Feb-2020 14:40	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB11-20200211  
 Collection Date: 11-Feb-2020 12:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	13-Feb-2020 12:57
Benzene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 12:57
Chlorobenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 12:57
Ethylbenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 12:57
Methylene chloride	U		0.0010	0.0020	mg/L	1	13-Feb-2020 12:57
Toluene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 12:57
Xylenes, Total	U		0.00030	0.0010	mg/L	1	13-Feb-2020 12:57
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 12:57</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>101</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 12:57</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 12:57</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 12:57</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22AR-20200210  
 Collection Date: 10-Feb-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	13-Feb-2020 17:37
Benzene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 17:37
Chlorobenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 17:37
Ethylbenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 17:37
Methylene chloride	U		0.0010	0.0020	mg/L	1	13-Feb-2020 17:37
Toluene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 17:37
Xylenes, Total	U		0.00030	0.0010	mg/L	1	13-Feb-2020 17:37
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>98.7</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 17:37</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>99.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 17:37</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 17:37</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>13-Feb-2020 17:37</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22AR-20200210  
 Collection Date: 10-Feb-2020 10:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 12-Feb-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	12-Feb-2020 16:19
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	12-Feb-2020 16:19
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	12-Feb-2020 16:19
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	12-Feb-2020 16:19
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	12-Feb-2020 16:19
<b>2-Methylnaphthalene</b>	<b>0.00012</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	12-Feb-2020 16:19
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	12-Feb-2020 16:19
<b>Acenaphthene</b>	<b>0.00015</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
Acenaphthylene	U		0.000015	0.00010	mg/L	1	12-Feb-2020 16:19
<b>Anthracene</b>	<b>0.00020</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Benz(a)anthracene</b>	<b>0.00029</b>		<b>0.000050</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Benzo(a)pyrene</b>	<b>0.00010</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	12-Feb-2020 16:19
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Chrysene</b>	<b>0.00032</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Dibenzofuran</b>	<b>0.00015</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	12-Feb-2020 16:19
<b>Fluoranthene</b>	<b>0.0015</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Fluorene</b>	<b>0.00019</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Naphthalene</b>	<b>0.00044</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
Nitrobenzene	U		0.000024	0.00020	mg/L	1	12-Feb-2020 16:19
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	12-Feb-2020 16:19
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	12-Feb-2020 16:19
<b>Phenanthrene</b>	<b>0.0015</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Phenol</b>	<b>0.000041</b>	J	<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<b>Pyrene</b>	<b>0.0010</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 16:19
<i>Surr: 2,4,6-Tribromophenol</i>	75.2			34-129	%REC	1	12-Feb-2020 16:19
<i>Surr: 2-Fluorobiphenyl</i>	62.6			40-125	%REC	1	12-Feb-2020 16:19
<i>Surr: 2-Fluorophenol</i>	46.1			20-120	%REC	1	12-Feb-2020 16:19
<i>Surr: 4-Terphenyl-d14</i>	92.4			40-135	%REC	1	12-Feb-2020 16:19
<i>Surr: Nitrobenzene-d5</i>	51.8			41-120	%REC	1	12-Feb-2020 16:19
<i>Surr: Phenol-d6</i>	51.8			20-120	%REC	1	12-Feb-2020 16:19
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Feb-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00352</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	14-Feb-2020 14:31

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22BR-20200210  
 Collection Date: 10-Feb-2020 11:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	13-Feb-2020 17:59
Benzene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 17:59
Chlorobenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 17:59
Ethylbenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 17:59
Methylene chloride	U		0.0010	0.0020	mg/L	1	13-Feb-2020 17:59
Toluene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 17:59
Xylenes, Total	U		0.00030	0.0010	mg/L	1	13-Feb-2020 17:59
<i>Surr: 1,2-Dichloroethane-d4</i>		96.8		70-126	%REC	1	13-Feb-2020 17:59
<i>Surr: 4-Bromofluorobenzene</i>		101		81-113	%REC	1	13-Feb-2020 17:59
<i>Surr: Dibromofluoromethane</i>		99.7		77-123	%REC	1	13-Feb-2020 17:59
<i>Surr: Toluene-d8</i>		99.1		82-127	%REC	1	13-Feb-2020 17:59

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22BR-20200210  
 Collection Date: 10-Feb-2020 11:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 12-Feb-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	12-Feb-2020 14:25
2,4-Dimethylphenol	U		0.000040	0.00020	mg/L	1	12-Feb-2020 14:25
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	12-Feb-2020 14:25
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	12-Feb-2020 14:25
<b>2-Chloronaphthalene</b>	<b>0.00011</b>	J	<b>0.000021</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
<b>2-Methylnaphthalene</b>	<b>0.000043</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	12-Feb-2020 14:25
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	12-Feb-2020 14:25
<b>Acenaphthene</b>	<b>0.018</b>		<b>0.00014</b>	<b>0.00050</b>	<b>mg/L</b>	5	12-Feb-2020 16:38
<b>Acenaphthylene</b>	<b>0.00018</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
<b>Anthracene</b>	<b>0.00042</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	12-Feb-2020 14:25
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	12-Feb-2020 14:25
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	12-Feb-2020 14:25
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000075</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
Chrysene	U		0.000021	0.00010	mg/L	1	12-Feb-2020 14:25
<b>Dibenzofuran</b>	<b>0.00029</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	12-Feb-2020 14:25
<b>Fluoranthene</b>	<b>0.0017</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
<b>Fluorene</b>	<b>0.0026</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
<b>Naphthalene</b>	<b>0.00013</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
Nitrobenzene	U		0.000024	0.00020	mg/L	1	12-Feb-2020 14:25
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	12-Feb-2020 14:25
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	12-Feb-2020 14:25
<b>Phenanthrene</b>	<b>0.00013</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
Phenol	U		0.000035	0.00020	mg/L	1	12-Feb-2020 14:25
<b>Pyrene</b>	<b>0.00082</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:25
<i>Surr: 2,4,6-Tribromophenol</i>	79.8			34-129	%REC	1	12-Feb-2020 14:25
<i>Surr: 2,4,6-Tribromophenol</i>	72.7			34-129	%REC	5	12-Feb-2020 16:38
<i>Surr: 2-Fluorobiphenyl</i>	68.4			40-125	%REC	1	12-Feb-2020 14:25
<i>Surr: 2-Fluorobiphenyl</i>	71.5			40-125	%REC	5	12-Feb-2020 16:38
<i>Surr: 2-Fluorophenol</i>	60.6			20-120	%REC	1	12-Feb-2020 14:25
<i>Surr: 2-Fluorophenol</i>	59.5			20-120	%REC	5	12-Feb-2020 16:38
<i>Surr: 4-Terphenyl-d14</i>	104			40-135	%REC	1	12-Feb-2020 14:25
<i>Surr: 4-Terphenyl-d14</i>	102			40-135	%REC	5	12-Feb-2020 16:38
<i>Surr: Nitrobenzene-d5</i>	59.6			41-120	%REC	1	12-Feb-2020 14:25
<i>Surr: Nitrobenzene-d5</i>	58.8			41-120	%REC	5	12-Feb-2020 16:38
<i>Surr: Phenol-d6</i>	59.4			20-120	%REC	1	12-Feb-2020 14:25
<i>Surr: Phenol-d6</i>	56.5			20-120	%REC	5	12-Feb-2020 16:38

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW22BR-20200210  
 Collection Date: 10-Feb-2020 11:00

**ANALYTICAL REPORT**

WorkOrder:HS20020456  
 Lab ID:HS20020456-03  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Feb-2020		Analyst: JHD	
Arsenic	0.0160		0.000400	0.00200	mg/L	1	14-Feb-2020 14:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB14-20200210  
 Collection Date: 10-Feb-2020 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	U		0.00020	0.0010	mg/L	1	13-Feb-2020 13:20
Benzene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 13:20
Chlorobenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 13:20
Ethylbenzene	U		0.00030	0.0010	mg/L	1	13-Feb-2020 13:20
Methylene chloride	U		0.0010	0.0020	mg/L	1	13-Feb-2020 13:20
Toluene	U		0.00020	0.0010	mg/L	1	13-Feb-2020 13:20
Xylenes, Total	U		0.00030	0.0010	mg/L	1	13-Feb-2020 13:20
<i>Surr: 1,2-Dichloroethane-d4</i>	96.9			70-126	%REC	1	13-Feb-2020 13:20
<i>Surr: 4-Bromofluorobenzene</i>	99.6			81-113	%REC	1	13-Feb-2020 13:20
<i>Surr: Dibromofluoromethane</i>	99.5			77-123	%REC	1	13-Feb-2020 13:20
<i>Surr: Toluene-d8</i>	101			82-127	%REC	1	13-Feb-2020 13:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB14-20200210  
 Collection Date: 10-Feb-2020 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20020456  
 Lab ID:HS20020456-04  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 12-Feb-2020		Analyst: LG	
1,2-Diphenylhydrazine	U		0.000021	0.00020	mg/L	1	12-Feb-2020 14:44
<b>2,4-Dimethylphenol</b>	<b>0.00095</b>		<b>0.000040</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
2,4-Dinitrotoluene	U		0.000058	0.00020	mg/L	1	12-Feb-2020 14:44
2,6-Dinitrotoluene	U		0.000042	0.00020	mg/L	1	12-Feb-2020 14:44
2-Chloronaphthalene	U		0.000021	0.00020	mg/L	1	12-Feb-2020 14:44
<b>2-Methylnaphthalene</b>	<b>0.00039</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
4,6-Dinitro-2-methylphenol	U		0.000020	0.00020	mg/L	1	12-Feb-2020 14:44
4-Nitrophenol	U		0.000047	0.0010	mg/L	1	12-Feb-2020 14:44
<b>Acenaphthene</b>	<b>0.00020</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
Acenaphthylene	U		0.000015	0.00010	mg/L	1	12-Feb-2020 14:44
Anthracene	U		0.000014	0.00010	mg/L	1	12-Feb-2020 14:44
Benz(a)anthracene	U		0.000050	0.00010	mg/L	1	12-Feb-2020 14:44
Benzo(a)pyrene	U		0.000020	0.00010	mg/L	1	12-Feb-2020 14:44
Bis(2-chloroethoxy)methane	U		0.000030	0.00020	mg/L	1	12-Feb-2020 14:44
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	12-Feb-2020 14:44
Chrysene	U		0.000021	0.00010	mg/L	1	12-Feb-2020 14:44
<b>Dibenzofuran</b>	<b>0.00019</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
<b>Di-n-butyl phthalate</b>	<b>0.000025</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
Fluoranthene	U		0.000010	0.00010	mg/L	1	12-Feb-2020 14:44
<b>Fluorene</b>	<b>0.000081</b>	J	<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
<b>Naphthalene</b>	<b>0.0034</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
Nitrobenzene	U		0.000024	0.00020	mg/L	1	12-Feb-2020 14:44
N-Nitrosodiphenylamine	U		0.000025	0.00020	mg/L	1	12-Feb-2020 14:44
Pentachlorophenol	U		0.000079	0.00020	mg/L	1	12-Feb-2020 14:44
<b>Phenanthrene</b>	<b>0.000029</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
<b>Phenol</b>	<b>0.00053</b>		<b>0.000035</b>	<b>0.00020</b>	<b>mg/L</b>	1	12-Feb-2020 14:44
Pyrene	U		0.000019	0.00010	mg/L	1	12-Feb-2020 14:44
<i>Surr: 2,4,6-Tribromophenol</i>	69.2			34-129	%REC	1	12-Feb-2020 14:44
<i>Surr: 2-Fluorobiphenyl</i>	96.0			40-125	%REC	1	12-Feb-2020 14:44
<i>Surr: 2-Fluorophenol</i>	81.2			20-120	%REC	1	12-Feb-2020 14:44
<i>Surr: 4-Terphenyl-d14</i>	92.6			40-135	%REC	1	12-Feb-2020 14:44
<i>Surr: Nitrobenzene-d5</i>	85.1			41-120	%REC	1	12-Feb-2020 14:44
<i>Surr: Phenol-d6</i>	79.4			20-120	%REC	1	12-Feb-2020 14:44
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 13-Feb-2020		Analyst: JHD	
Arsenic	U		0.000400	0.00200	mg/L	1	14-Feb-2020 15:15

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**Batch ID:** 150549      **Start Date:** 12 Feb 2020 08:10      **End Date:** 12 Feb 2020 13:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20020456-02	1	1000 (mL)	1 (mL)	0.001
HS20020456-03	1	1000 (mL)	1 (mL)	0.001
HS20020456-04	1	1000 (mL)	1 (mL)	0.001

**Batch ID:** 150594      **Start Date:** 13 Feb 2020 09:00      **End Date:** 13 Feb 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20020456-02		10 (mL)	10 (mL)	1
HS20020456-03		10 (mL)	10 (mL)	1
HS20020456-04		10 (mL)	10 (mL)	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 150549 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20020456-02	WG-1620-MW22AR-20200210	10 Feb 2020 10:00		12 Feb 2020 08:10	12 Feb 2020 16:19	1
HS20020456-03	WG-1620-MW22BR-20200210	10 Feb 2020 11:00		12 Feb 2020 08:10	12 Feb 2020 16:38	5
HS20020456-03	WG-1620-MW22BR-20200210	10 Feb 2020 11:00		12 Feb 2020 08:10	12 Feb 2020 14:25	1
HS20020456-04	WG-1620-FB14-20200210	10 Feb 2020 11:30		12 Feb 2020 08:10	12 Feb 2020 14:44	1
<b>Batch ID: 150594 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Water</b>	
HS20020456-02	WG-1620-MW22AR-20200210	10 Feb 2020 10:00		13 Feb 2020 13:00	14 Feb 2020 14:31	1
HS20020456-03	WG-1620-MW22BR-20200210	10 Feb 2020 11:00		13 Feb 2020 13:00	14 Feb 2020 14:34	1
HS20020456-04	WG-1620-FB14-20200210	10 Feb 2020 11:30		13 Feb 2020 13:00	14 Feb 2020 15:15	1
<b>Batch ID: R356238 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20020456-01	WQ-1620-TB11-20200211	11 Feb 2020 12:00			13 Feb 2020 12:57	1
HS20020456-02	WG-1620-MW22AR-20200210	10 Feb 2020 10:00			13 Feb 2020 17:37	1
HS20020456-03	WG-1620-MW22BR-20200210	10 Feb 2020 11:00			13 Feb 2020 17:59	1
HS20020456-04	WG-1620-FB14-20200210	10 Feb 2020 11:30			13 Feb 2020 13:20	1

WorkOrder: HS20020456  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20020456  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.000098	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000051	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.000094	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.000083	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.000093	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000031	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000037	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000059	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000050	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.000081	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000060	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00012	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00011	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.0000094	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20020456  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QC BATCH REPORT**

Batch ID: 150594 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A						
<b>MBLK</b>	Sample ID: <b>MBLK-150594</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Feb-2020 13:14</b>						
Client ID:	Run ID: <b>ICPMS05_356303</b>	SeqNo: <b>5474006</b>		PrepDate: <b>13-Feb-2020</b>		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	U	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-150594</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Feb-2020 13:17</b>						
Client ID:	Run ID: <b>ICPMS05_356303</b>	SeqNo: <b>5474007</b>		PrepDate: <b>13-Feb-2020</b>		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.04604	0.00200	0.05	0	92.1	80 - 120				
<b>MS</b>	Sample ID: <b>HS20020453-01MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Feb-2020 13:23</b>						
Client ID:	Run ID: <b>ICPMS05_356303</b>	SeqNo: <b>5474010</b>		PrepDate: <b>13-Feb-2020</b>		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.0889	0.00200	0.05	0.04209	93.6	80 - 120				
<b>MSD</b>	Sample ID: <b>HS20020453-01MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Feb-2020 13:26</b>						
Client ID:	Run ID: <b>ICPMS05_356303</b>	SeqNo: <b>5474011</b>		PrepDate: <b>13-Feb-2020</b>		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.09146	0.00200	0.05	0.04209	98.7	80 - 120	0.0889	2.83	20	
<b>PDS</b>	Sample ID: <b>HS20020453-01PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Feb-2020 13:28</b>						
Client ID:	Run ID: <b>ICPMS05_356303</b>	SeqNo: <b>5474012</b>		PrepDate: <b>13-Feb-2020</b>		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
Arsenic	0.1563	0.00200	0.1	0.04209	114	75 - 125				
<b>SD</b>	Sample ID: <b>HS20020453-01SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>14-Feb-2020 13:21</b>						
Client ID:	Run ID: <b>ICPMS05_356303</b>	SeqNo: <b>5474009</b>		PrepDate: <b>13-Feb-2020</b>		DF: <b>5</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit Qual	
Arsenic	0.04136	0.0100					0.04209	1.73	10	

The following samples were analyzed in this batch: HS20020456-02 HS20020456-03 HS20020456-04

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QC BATCH REPORT**

Batch ID: 150549 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-150549	Units: ug/L			Analysis Date: 12-Feb-2020 09:58					
Client ID:	Run ID: SV-6_356146	SeqNo: 5470142	PrepDate: 12-Feb-2020	DF: 1						
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	U	0.20								
2,4-Dimethylphenol	U	0.20								
2,4-Dinitrotoluene	U	0.20								
2,6-Dinitrotoluene	U	0.20								
2-Chloronaphthalene	U	0.20								
2-Methylnaphthalene	U	0.10								
4,6-Dinitro-2-methylphenol	U	0.20								
4-Nitrophenol	U	1.0								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Benz(a)anthracene	U	0.10								
Benzo(a)pyrene	U	0.10								
Bis(2-chloroethoxy)methane	U	0.20								
Bis(2-ethylhexyl)phthalate	U	0.20								
Chrysene	U	0.10								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Nitrobenzene	U	0.20								
N-Nitrosodiphenylamine	U	0.20								
Pentachlorophenol	U	0.20								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	3.329	0.20	5	0	66.6	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	4.669	0.20	5	0	93.4	40 - 125				
<i>Surr: 2-Fluorophenol</i>	4.4	0.20	5	0	88.0	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.725	0.20	5	0	94.5	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	4.118	0.20	5	0	82.4	41 - 120				
<i>Surr: Phenol-d6</i>	3.995	0.20	5	0	79.9	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QC BATCH REPORT**

Batch ID: 150549 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-150549	Units: ug/L			Analysis Date: 12-Feb-2020 10:17					
Client ID:	Run ID: SV-6_356146	SeqNo: 5470143		PrepDate: 12-Feb-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.832	0.20	5	0	96.6	39 - 127				
2,4-Dimethylphenol	4.124	0.20	5	0	82.5	35 - 120				
2,4-Dinitrotoluene	5.313	0.20	5	0	106	50 - 122				
2,6-Dinitrotoluene	5.337	0.20	5	0	107	50 - 120				
2-Chloronaphthalene	4.964	0.20	5	0	99.3	50 - 120				
2-Methylnaphthalene	4.367	0.10	5	0	87.3	50 - 120				
4,6-Dinitro-2-methylphenol	5.848	0.20	5	0	117	25 - 121				
4-Nitrophenol	4.061	1.0	5	0	81.2	30 - 130				
Acenaphthene	4.707	0.10	5	0	94.1	45 - 120				
Acenaphthylene	4.744	0.10	5	0	94.9	47 - 120				
Anthracene	5.024	0.10	5	0	100	45 - 120				
Benz(a)anthracene	4.981	0.10	5	0	99.6	40 - 120				
Benzo(a)pyrene	4.733	0.10	5	0	94.7	45 - 120				
Bis(2-chloroethoxy)methane	4.378	0.20	5	0	87.6	45 - 120				
Bis(2-ethylhexyl)phthalate	6.092	0.20	5	0	122	40 - 139				
Chrysene	4.751	0.10	5	0	95.0	43 - 120				
Dibenzofuran	4.607	0.10	5	0	92.1	50 - 120				
Di-n-butyl phthalate	5.595	0.20	5	0	112	45 - 123				
Fluoranthene	4.741	0.10	5	0	94.8	45 - 125				
Fluorene	4.695	0.10	5	0	93.9	49 - 120				
Naphthalene	4.404	0.10	5	0	88.1	45 - 120				
Nitrobenzene	4.15	0.20	5	0	83.0	44 - 120				
N-Nitrosodiphenylamine	5.326	0.20	5	0	107	40 - 125				
Pentachlorophenol	3.971	0.20	5	0	79.4	19 - 121				
Phenanthrene	4.698	0.10	5	0	94.0	45 - 121				
Phenol	4.283	0.20	5	0	85.7	20 - 124				
Pyrene	5.051	0.10	5	0	101	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.83</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.6</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.712</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>94.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>4.272</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>85.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.806</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>96.1</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>4.018</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>80.4</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>4.152</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.0</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QC BATCH REPORT**

**Batch ID:** 150549 ( 0 )      **Instrument:** SV-6      **Method:** LOW-LEVEL SEMIVOLATILES BY 8270D

LCSD	Sample ID: LCSD-150549	Units: ug/L			Analysis Date: 12-Feb-2020 10:36					
Client ID:	Run ID: SV-6_356146	SeqNo: 5470144	PrepDate: 12-Feb-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.646	0.20	5	0	92.9	39 - 127	4.832	3.91	20	
2,4-Dimethylphenol	4.085	0.20	5	0	81.7	35 - 120	4.124	0.951	20	
2,4-Dinitrotoluene	4.973	0.20	5	0	99.5	50 - 122	5.313	6.61	20	
2,6-Dinitrotoluene	4.989	0.20	5	0	99.8	50 - 120	5.337	6.73	20	
2-Chloronaphthalene	4.655	0.20	5	0	93.1	50 - 120	4.964	6.43	20	
2-Methylnaphthalene	4.232	0.10	5	0	84.6	50 - 120	4.367	3.13	20	
4,6-Dinitro-2-methylphenol	5.8	0.20	5	0	116	25 - 121	5.848	0.825	30	
4-Nitrophenol	3.845	1.0	5	0	76.9	30 - 130	4.061	5.46	20	
Acenaphthene	4.442	0.10	5	0	88.8	45 - 120	4.707	5.8	20	
Acenaphthylene	4.414	0.10	5	0	88.3	47 - 120	4.744	7.22	20	
Anthracene	4.909	0.10	5	0	98.2	45 - 120	5.024	2.31	20	
Benz(a)anthracene	4.706	0.10	5	0	94.1	40 - 120	4.981	5.69	20	
Benzo(a)pyrene	4.536	0.10	5	0	90.7	45 - 120	4.733	4.26	20	
Bis(2-chloroethoxy)methane	4.276	0.20	5	0	85.5	45 - 120	4.378	2.36	20	
Bis(2-ethylhexyl)phthalate	5.715	0.20	5	0	114	40 - 139	6.092	6.39	20	
Chrysene	4.467	0.10	5	0	89.3	43 - 120	4.751	6.16	20	
Dibenzofuran	4.373	0.10	5	0	87.5	50 - 120	4.607	5.22	20	
Di-n-butyl phthalate	5.419	0.20	5	0	108	45 - 123	5.595	3.2	20	
Fluoranthene	4.584	0.10	5	0	91.7	45 - 125	4.741	3.36	20	
Fluorene	4.429	0.10	5	0	88.6	49 - 120	4.695	5.83	20	
Naphthalene	4.286	0.10	5	0	85.7	45 - 120	4.404	2.72	20	
Nitrobenzene	4.078	0.20	5	0	81.6	44 - 120	4.15	1.75	20	
N-Nitrosodiphenylamine	5.247	0.20	5	0	105	40 - 125	5.326	1.5	20	
Pentachlorophenol	4.014	0.20	5	0	80.3	19 - 121	3.971	1.09	20	
Phenanthrene	4.567	0.10	5	0	91.3	45 - 121	4.698	2.82	20	
Phenol	3.965	0.20	5	0	79.3	20 - 124	4.283	7.71	20	
Pyrene	4.755	0.10	5	0	95.1	40 - 130	5.051	6.03	20	
Surr: 2,4,6-Tribromophenol	3.716	0.20	5	0	74.3	34 - 129	3.83	3.02	20	
Surr: 2-Fluorobiphenyl	4.416	0.20	5	0	88.3	40 - 125	4.712	6.49	20	
Surr: 2-Fluorophenol	3.948	0.20	5	0	79.0	20 - 120	4.272	7.88	20	
Surr: 4-Terphenyl-d14	4.493	0.20	5	0	89.9	40 - 135	4.806	6.74	20	
Surr: Nitrobenzene-d5	3.853	0.20	5	0	77.1	41 - 120	4.018	4.21	20	
Surr: Phenol-d6	3.79	0.20	5	0	75.8	20 - 120	4.152	9.1	20	

The following samples were analyzed in this batch: HS20020456-02      HS20020456-03      HS20020456-04



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QC BATCH REPORT**

**Batch ID:** R356238 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200213</b>			Units: <b>ug/L</b>		Analysis Date: <b>13-Feb-2020 12:34</b>			
Client ID:		Run ID: <b>VOA2_356238</b>			SeqNo: <b>5472049</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	U	1.0								
Benzene	U	1.0								
Chlorobenzene	U	1.0								
Ethylbenzene	U	1.0								
Methylene chloride	U	2.0								
Toluene	U	1.0								
Xylenes, Total	U	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.97</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.9</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.41</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.56</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200213</b>			Units: <b>ug/L</b>		Analysis Date: <b>13-Feb-2020 11:48</b>			
Client ID:		Run ID: <b>VOA2_356238</b>			SeqNo: <b>5472048</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.97	1.0	20	0	94.8	70 - 124				
Benzene	18.57	1.0	20	0	92.8	74 - 120				
Chlorobenzene	18.83	1.0	20	0	94.2	76 - 113				
Ethylbenzene	18.94	1.0	20	0	94.7	77 - 117				
Methylene chloride	19.11	2.0	20	0	95.5	70 - 127				
Toluene	20.9	1.0	20	0	104	77 - 118				
Xylenes, Total	57.59	1.0	60	0	96.0	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>52.12</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.03</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QC BATCH REPORT**

Batch ID: R356238 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>		Sample ID: HS20020304-21MS		Units: ug/L		Analysis Date: 13-Feb-2020 14:56				
Client ID:		Run ID: VOA2_356238		SeqNo: 5472137		PrepDate:		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	18.92	1.0	20	0.7205	91.0	70 - 127				
Benzene	18.21	1.0	20	0	91.0	70 - 127				
Chlorobenzene	17.96	1.0	20	0	89.8	70 - 114				
Ethylbenzene	18.68	1.0	20	0	93.4	70 - 124				
Methylene chloride	18.3	2.0	20	0	91.5	70 - 128				
Toluene	20.23	1.0	20	0	101	70 - 123				
Xylenes, Total	56.88	1.0	60	0	94.8	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.96</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>51.8</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>104</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.16</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.0</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: HS20020304-21MSD		Units: ug/L		Analysis Date: 13-Feb-2020 15:19			
Client ID:		Run ID: VOA2_356238		SeqNo: 5472138		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	18.04	1.0	20	0.7205	86.6	70 - 127	18.92	4.78	20
Benzene	17.41	1.0	20	0	87.0	70 - 127	18.21	4.5	20
Chlorobenzene	17.6	1.0	20	0	88.0	70 - 114	17.96	2.02	20
Ethylbenzene	18.38	1.0	20	0	91.9	70 - 124	18.68	1.63	20
Methylene chloride	17.76	2.0	20	0	88.8	70 - 128	18.3	2.98	20
Toluene	19.85	1.0	20	0	99.3	70 - 123	20.23	1.88	20
Xylenes, Total	55.31	1.0	60	0	92.2	70 - 130	56.88	2.81	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>50.93</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>70 - 126</i>	<i>50.96</i>	<i>0.0592</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.77</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 113</i>	<i>51.8</i>	<i>2.01</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>49.13</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>77 - 123</i>	<i>49.16</i>	<i>0.0527</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>50.25</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>	<i>49.49</i>	<i>1.51</i>	<i>20</i>

The following samples were analyzed in this batch: HS20020456-01 HS20020456-02 HS20020456-03 HS20020456-04

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20020456

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

---

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231 V009	22-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20020456

Date/Time Received: 11-Feb-2020 14:40
Received by: PMG

Checklist completed by: Paresh M. Giga
eSignature
Date: 11-Feb-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 12-Feb-2020

Matrices: Water

Carrier name: Client

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:206375

Temperature(s)/Thermometer(s): 0.2°C U/C IR25
Cooler(s)/Kit(s): 45623
Date/Time sample(s) sent to storage: 2/11/2020 18:30

- Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]

pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Fort Collins, CO  
+1 970 490 1511

Everett, WA  
+1 425 356 2600

Holland, MI  
+1 616 399 6070

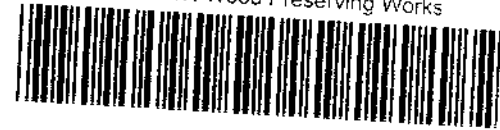
# Chain of Custody Form

Page 1 of 1

COC ID: 206375

HS20020456

Golder Associates Inc.  
Houston TX-Wood Preserving Works



ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A
Work Order		Project Number	1620-07-Rev0 SR 92688	B
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750	E
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	F
Phone	(512) 671-3434	Phone		G
Fax	(512) 671-3446	Fax		H
e-Mail Address	eric.matzner@pbwllc.com	e-Mail Address		I
				J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TBO/LL-20200211	2-11-2020	12:00	Water	1	2		X									
2	WG-1620-MW 2-HAR 20200210	2-10-2020	10:00	Water	1	6	X		X	X							
3	WG-1620-MW 2-BR 20200210	2-10-2020	11:00	Water	1	6	X		X	X							
4	WG-1620-FB 14 20200210	2-10-2020	11:30	Water	1	6	X		X	X							
5																	
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign: *Tim Spalding*

Relinquished by: *Tim Spalding* Date: 2-11-2020 Time: 12:25

Received by (Laboratory): *Blake Sobora* Date: 2-11-20 Time: 1440

Shipment Method: \_\_\_\_\_ Required Turnaround Time: (Check Box)  STD 10 Wk Days  5 Wk Days  2 Wk Days  24-hour

Results Due Date: \_\_\_\_\_

Notes: UPRR Houston MWPV

Cooler ID: 43623 Cooler Temp: 0.2

QC Package: (Check One Box Below)  Level II Std QC  TRRP Check st  TRRP Level IV  Level III Std QC/Raw Data  Level IV SV483/CLP  Other

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>3</sub> 7-Other 8-4°C 9-5035

ote: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.



# Memorandum

March 10, 2020

To: Eric Matzner Ref. No.: 11183954-1620

From: <sup>CK</sup> Chris G. Knight/eew/559-NF Tel: 512-506-8803

CC: Jesse Orth, Jon Lang; Julie Lidstone

**Subject: Data Usability Summary  
Semiannual Groundwater Monitoring Event  
Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works  
Houston, Texas  
January 2020**

## 1. Scope of Data Usability Study

This document details a Data Usability Summary (DUS) of analytical results for groundwater samples collected in support of the Semiannual Groundwater Monitoring Event at the Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works site during January 2020. Samples were submitted to ALS Environmental (ALS), located in Houston, Texas and are reported in data package HS20010618. The intended use of the data is to support the Semiannual Groundwater Monitoring Event at the site by providing current concentration of chemicals of concern.

Data were reviewed and validated by Chris G. Knight of GHD, in accordance with Title 30 of the Texas Administrative Code Section 350.54 (30 TAC 350.54) as described in the Texas Commission on Environmental Quality (TCEQ) Regulatory Guidance document entitled "Review and Reporting of COC Concentration Data under TRRP", (RG-366/TRRP-13), revised May 2010, herein referred to as "TRRP-13 Guidance". Evaluation of the data was based on information obtained from the chain of custody forms, the finished report forms, method blank data, recovery data from surrogate spikes/laboratory control samples (LCS)/matrix spikes (MS), duplicate data, field quality assurance/quality control (QA/QC) samples, the laboratory review checklists (LRC), and the laboratory exception report (ER).

A sample collection and analysis summary is presented in Table 1. This summary provides a cross-reference of field sample identification numbers and location identification. Each sample is assigned a unique field identification number.

The validated sample results are presented in Table 2. A summary of the analytical methodology is presented in Table 3.



## 2. Laboratory Qualifications

The Laboratory's quality assurance program is consistent with the quality standards outlined in the National Environmental Laboratory Accreditation Program (NELAP). This laboratory was accredited under Texas Certification number # TX104704231 at the time the analysis was performed and the certificate is included in Attachment A.

## 3. Project Objectives

### 3.1 Sampling/Analytical QA/QC Objectives

The QA/QC program was designed to identify contamination resulting from the sampling, sample transport and analytical process through the analysis of field blank samples, a field duplicate sample set, and method blanks. The QA/QC program was designed to evaluate the quality of the resulting data with respect to bias and precision through analysis of LCS and MS.

## 4. Data Review/Validation Results

### 4.1 Sample Holding Time and Preservation

Samples were shipped with a chain of custody and the paper work was filled out properly. All samples were properly preserved, delivered on ice, and stored by the laboratory at the required temperature (0-6°C).

The sample chain of custody documents and the analytical report were used to determine sample holding times. All samples were prepared and analyzed within the required holding times.

### 4.2 Sample Containers

Sample containers used were certified pre-cleaned glass containers provided by the laboratory. These containers meet or exceed analyte specifications established in the United States Environmental Protection Agency (USEPA) *Specifications and Guidance for Contaminant-free Sample Containers*.

### 4.3 Calibrations

According to the LRC, initial calibration and continuing calibration data met the criteria for the selected method.

### 4.4 Laboratory Method Blank Analyses

Method blanks are prepared from a purified matrix and analyzed with investigative samples to determine the existence and magnitude of sample contamination introduced during the analytical procedures. As these were not discrete samples handled in the field, these blanks are not listed on the sample identification cross-reference list found in the data package.





For this study, laboratory method blanks were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch and results are reported in the laboratory data package.

The method blank results were non-detect or below the method quantitation limit (MQL), indicating that laboratory contamination was not a factor for this investigation.

#### **4.5 Internal Standard and Surrogate Spike Recoveries**

Recoveries of internal standards are addressed in the LRC of the data package. All internal standard recoveries associated with the compounds of interest were acceptable per the LRC.

In accordance with the methods employed, all samples, blanks, and QC samples analyzed for semi-volatile organic compounds (SVOCs) are spiked with surrogate compounds prior to sample analysis. Surrogate recoveries provide a means to evaluate the effects of laboratory performance on individual sample matrices. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project. Each individual surrogate compound is expected to meet the laboratory control limits. According to the TRRP-13 Guidelines, one outlying surrogate is acceptable for methods with multiple surrogate spike compounds as long as the recovery is at least ten percent. Sample analyzed at elevated sample dilutions (five times or greater) were not assessed.

Surrogate recoveries were assessed against laboratory control limits and/or the guidance in TRRP-13. All surrogate recoveries met the above criteria.

#### **4.6 Laboratory Control Sample Analysis**

LCS are prepared and analyzed as samples to assess the analytical efficiencies of the methods employed, independent of sample matrix effects. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project.

For this study, LCS were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch.

The LCS contained all compounds specified in the method. All LCS recoveries were within the laboratory control limits, demonstrating acceptable analytical accuracy.

#### **4.7 Matrix Spike Analysis**

To evaluate the effects of sample matrices on the preparation process, measurement procedures, and accuracy of a particular analysis, samples are spiked with known concentrations of the analytes of interest and analyzed as MS/matrix spike duplicate (MSD) samples. The RPD between the MS and MSD is used to assess analytical precision.

An MS/MSD analysis was performed as specified in Table 1. The recovery ranges established by the laboratory is adopted as the acceptance criteria for the project.



The MS/MSD samples were spiked with all compounds specified in the method. All percent recoveries and the RPD value were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision.

#### **4.8 Field QA/QC Samples**

The field QA/QC consisted of two field blank samples and one field duplicate sample set.

##### ***Field Blank Sample Analysis***

To assess ambient conditions at the site, two field blank samples were submitted for analysis, as identified in Table 1. All results were non-detect for the compounds of interest.

##### ***Field Duplicate Sample Analysis***

To assess the analytical and sampling protocol precision, one field duplicate sample set was collected and submitted "blind" to the laboratory, as specified in Table 1. The RPDs associated with these duplicate samples must be less than thirty percent for water samples. The RPDs are only used when sample concentrations are above the estimated regions of detection.

Field duplicate summary data are presented in Table 2. All field duplicate results were within acceptable agreement, demonstrating acceptable sampling and analytical precision.

#### **4.9 Field Procedures**

Golder Associates, Inc. collected groundwater samples in accordance with their Standard Operating Procedures (SOP) for sample collection.

#### **4.10 Analyte Reporting**

The laboratory reported detected results for each analyte down to the sample detection limit (SDL), which is defined as the method detection limit (MDL) with sample-specific adjustments for dilutions, aliquot size, volumes, etc. Positive analyte detections less than the MQL but greater than the SDL were qualified as estimated (J) in Table 2.

The detectability check standard (DCS) results supported the laboratory MDLs.

## **5. Conclusion**

Based on the assessment detailed in the foregoing, the data summarized in Table 2 are usable for the purpose of supporting the Semiannual Groundwater Monitoring Event at the site by providing current concentration of chemicals of concern without qualification.

Table 1

**Sample Collection and Analysis Summary**  
**Semiannual Groundwater Monitoring Event**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	<u>Analysis/Parameters</u>	
					SVOCs	Comments
WG-1620-P12-20200113	P-12	Water	01/13/2020	10:30	X	MS/MSD
WG-1620-MW08-20200113	MW-08	Water	01/13/2020	11:30	X	
WG-1620-FB01-20200113	-	Water	01/13/2020	11:30	X	Field Blank
WG-1620-MW07-20200113	MW-07	Water	01/13/2020	13:15	X	
WG-1620-P10-20200113	P-10	Water	01/13/2020	14:20	X	
WG-1620-MW11B-20200114	MW-11B	Water	01/14/2020	09:25	X	
WG-1620-MW11A-20200114	MW-11A	Water	01/14/2020	10:35	X	
WG-1620-MW10B-20200114	MW-10B	Water	01/14/2020	11:35	X	
WG-1620-MW10A-20200114	MW-10A	Water	01/14/2020	12:25	X	
WG-1620-MW02-20200114	MW-02	Water	01/14/2020	13:30	X	
WG-1620-MW01A-20200114	MW-01A	Water	01/14/2020	14:35	X	
WG-1620-FD01-20200114	MW-01A	Water	01/14/2020	14:35	X	Field duplicate of MW-01A
WG-1620-FB02-20200114	-	Water	01/14/2020	15:00	X	Field Blank

## Notes:

- SVOCs - Semi-volatile Organic Compounds  
MS/MSD - Matrix Spike/ Matrix Spike Duplicate  
"- " - Not Applicable

**Table 2**  
**Analytical Results Summary**  
**Semiannual Groundwater Monitoring Event**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January 2020**

Location ID:	MW-01A	MW-01A	MW-02	MW-07	
Sample Name:	WG-1620-MW01A-20200114	WG-1620-FD01-20200114	WG-1620-MW02-20200114	WG-1620-MW07-20200113	
Sample Date:	01/14/2020	01/14/2020 Duplicate	01/14/2020	01/13/2020	
Parameters	Unit				
<b>Semi-volatile Organic Compounds</b>					
2-Methylnaphthalene	mg/L	0.00019	<0.000019	<0.000019	0.000066 J
Acenaphthene	mg/L	0.024	0.018	0.0030	<0.000027
Acenaphthylene	mg/L	0.00084	0.00066	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	0.00011	<0.000014
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000037	0.000074 J	<0.000037	<0.000037
Di-n-butylphthalate (DBP)	mg/L	--	--	--	--
Dibenzofuran	mg/L	0.0036	0.0021	0.00039	0.000057 J
Fluoranthene	mg/L	0.0011	0.0012	0.00024	0.00010
Fluorene	mg/L	0.0064	0.0038	0.0017	<0.000030
Naphthalene	mg/L	0.00052	<0.000020	<0.000020	0.00017
Phenanthrene	mg/L	<0.000021	<0.000021	0.00011	0.00014
Phenol	mg/L	--	--	--	--
Pyrene	mg/L	0.00052	0.00059	0.00011	<0.000019

**Table 2**  
**Analytical Results Summary**  
**Semiannual Groundwater Monitoring Event**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January 2020**

<b>Location ID:</b>	<b>MW-08</b>	<b>MW-10A</b>	<b>MW-10B</b>	<b>MW-11A</b>
<b>Sample Name:</b>	<b>WG-1620-MW08-20200113</b>	<b>WG-1620-MW10A-20200114</b>	<b>WG-1620-MW10B-20200114</b>	<b>WG-1620-MW11A-20200114</b>
<b>Sample Date:</b>	<b>01/13/2020</b>	<b>01/14/2020</b>	<b>01/14/2020</b>	<b>01/14/2020</b>

<b>Parameters</b>	<b>Unit</b>				
<b>Semi-volatile Organic Compounds</b>					
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	--	<0.000019
Acenaphthene	mg/L	<0.000027	0.00011	0.069	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	0.00066	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	0.0028	<0.000014
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00021	<0.000037	0.00020	<0.000037
Di-n-butylphthalate (DBP)	mg/L	--	--	<0.000020	--
Dibenzofuran	mg/L	<0.000020	<0.000020	0.022	<0.000020
Fluoranthene	mg/L	<0.000010	<0.000010	0.0029	<0.000010
Fluorene	mg/L	<0.000030	<0.000030	0.036	<0.000030
Naphthalene	mg/L	<0.000020	<0.000020	0.0021	<0.000020
Phenanthrene	mg/L	0.000046 J	<0.000021	--	<0.000021
Phenol	mg/L	--	--	<0.000035	--
Pyrene	mg/L	<0.000019	<0.000019	0.0013	<0.000019

Table 2

**Analytical Results Summary**  
**Semiannual Groundwater Monitoring Event**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January 2020**

Location ID:	MW-11B	P-10	P-12
Sample Name:	WG-1620-MW11B-20200114	WG-1620-P10-20200113	WG-1620-P12-20200113
Sample Date:	01/14/2020	01/13/2020	01/13/2020
Parameters	Unit		
<b>Semi-volatile Organic Compounds</b>			
2-Methylnaphthalene	mg/L	--	--
Acenaphthene	mg/L	0.033	<0.000027
Acenaphthylene	mg/L	0.0016	<0.000015
Anthracene	mg/L	<0.000014	<0.000014
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000095 J	<0.000037
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020
Fluoranthene	mg/L	0.0024	<0.000010
Fluorene	mg/L	0.00035	<0.000030
Naphthalene	mg/L	<0.000020	0.00017
Phenanthrene	mg/L	--	--
Phenol	mg/L	<0.000035	<0.000035
Pyrene	mg/L	0.0023	<0.000019

## Notes:

- < - Not detected at the associated reporting limit
- J - Estimated concentration
- " - Not applicable

Table 3

**Analytical Methods**  
**Semiannual Groundwater Monitoring Event**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**January 2020**

Parameter	Method	Matrix	Holding Time	
			Collection to Extraction (Days)	Extraction to Analysis (Days)
SVOCs	SW-846 8270D	Water	7	40

## Notes:

SVOCs - Semi-volatile Organic Compounds

## Method References:

SW-846 - "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition, 1986, with subsequent revisions

# Attachment A

## Laboratory NELAP Certificate





# Texas Commission on Environmental Quality

## NELAP - Recognized Laboratory Fields of Accreditation



ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: *Drinking Water***

**Method** EPA 1613

Analyte	AB	Analyte ID	Method ID
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408

**Method** EPA 200.8

Analyte	AB	Analyte ID	Method ID
Copper	TX	1055	10014605
Lead	TX	1075	10014605



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 120.1			
Analyte Conductivity	TX	1610	10006403
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 160.4			
Analyte Residue-volatile	TX	1970	10010409
Method EPA 1613			
Analyte 1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10120408
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10120408
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10120408
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10120408
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10120408
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10120408
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10120408
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10120408
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,6,7,8-HxCDD)	TX	9456	10120408
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10120408
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10120408
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10120408
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10120408
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10120408



# Texas Commission on Environmental Quality



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Issue Date: 5/1/2019

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**Matrix: Non-Potable Water**

2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10120408
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10120408
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10120408
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10120408
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10120408
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10120408
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10120408
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10120408
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10120408
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10120408
<b>Method EPA 1664</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10127807
<b>Method EPA 180.1</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Turbidity	TX	2055	10011606
<b>Method EPA 200.8</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Aluminum	TX	1000	10014605
Antimony	TX	1005	10014605
Arsenic	TX	1010	10014605
Barium	TX	1015	10014605
Beryllium	TX	1020	10014605
Boron	TX	1025	10014605
Cadmium	TX	1030	10014605
Calcium	TX	1035	10014605
Chromium	TX	1040	10014605
Cobalt	TX	1050	10014605
Copper	TX	1055	10014605
Iron	TX	1070	10014605



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**Matrix: Non-Potable Water**

Lead	TX	1075	10014605
Magnesium	TX	1085	10014605
Manganese	TX	1090	10014605
Molybdenum	TX	1100	10014605
Nickel	TX	1105	10014605
Potassium	TX	1125	10014605
Selenium	TX	1140	10014605
Silver	TX	1150	10014605
Sodium	TX	1155	10014605
Strontium	TX	1160	10014605
Thallium	TX	1165	10014605
Tin	TX	1175	10014605
Titanium	TX	1180	10014605
Uranium	TX	3035	10014605
Vanadium	TX	1185	10014605
Zinc	TX	1190	10014605

**Method EPA 245.1**

Analyte	AB	Analyte ID	Method ID
Mercury	TX	1095	10036609

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 325.1**

Analyte	AB	Analyte ID	Method ID
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**Matrix: Non-Potable Water**

Chloride	TX	1575	10056801
<b>Method</b> EPA 335.1			
<b>Analyte</b> Amenable cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1510	<b>Method ID</b> 10060001
<b>Method</b> EPA 335.2			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10278203
<b>Method</b> EPA 335.4			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10061402
<b>Method</b> EPA 350.3			
<b>Analyte</b> Ammonia as N	<b>AB</b> TX	<b>Analyte ID</b> 1515	<b>Method ID</b> 10064401
<b>Method</b> EPA 365.3			
<b>Analyte</b> Orthophosphate as P Phosphorus	<b>AB</b> TX TX	<b>Analyte ID</b> 1870 1910	<b>Method ID</b> 10070801 10070801
<b>Method</b> EPA 375.4			
<b>Analyte</b> Sulfate	<b>AB</b> TX	<b>Analyte ID</b> 2000	<b>Method ID</b> 10073800
<b>Method</b> EPA 376.1			
<b>Analyte</b> Sulfide	<b>AB</b> TX	<b>Analyte ID</b> 2005	<b>Method ID</b> 10074201
<b>Method</b> EPA 410.4			
<b>Analyte</b> Chemical oxygen demand (COD)	<b>AB</b> TX	<b>Analyte ID</b> 1565	<b>Method ID</b> 10077404
<b>Method</b> EPA 415.1			
<b>Analyte</b> Total Organic Carbon (TOC)	<b>AB</b> TX	<b>Analyte ID</b> 2040	<b>Method ID</b> 10078407
<b>Method</b> EPA 420.1			
<b>Analyte</b> Total phenolics	<b>AB</b> TX	<b>Analyte ID</b> 1905	<b>Method ID</b> 10079400



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**Matrix: Non-Potable Water**

**Method EPA 420.4**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10080203

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156419
Antimony	TX	1005	10156419
Arsenic	TX	1010	10156419
Barium	TX	1015	10156419
Beryllium	TX	1020	10156419
Boron	TX	1025	10156419
Cadmium	TX	1030	10156419
Calcium	TX	1035	10156419
Chromium	TX	1040	10156419
Cobalt	TX	1050	10156419
Copper	TX	1055	10156419
Iron	TX	1070	10156419
Lead	TX	1075	10156419
Lithium	TX	1080	10156419
Magnesium	TX	1085	10156419
Manganese	TX	1090	10156419
Molybdenum	TX	1100	10156419
Nickel	TX	1105	10156419
Potassium	TX	1125	10156419
Selenium	TX	1140	10156419
Silver	TX	1150	10156419
Sodium	TX	1155	10156419
Strontium	TX	1160	10156419
Thallium	TX	1165	10156419
Tin	TX	1175	10156419
Titanium	TX	1180	10156419



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**Matrix: Non-Potable Water**

Vanadium	TX	1185	10156419
Zinc	TX	1190	10156419
<b>Method EPA 608</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
4,4'-DDD	TX	7355	10103603
4,4'-DDE	TX	7360	10103603
4,4'-DDT	TX	7365	10103603
Aldrin	TX	7025	10103603
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10103603
alpha-Chlordane	TX	7240	10103603
Aroclor-1016 (PCB-1016)	TX	8880	10103603
Aroclor-1221 (PCB-1221)	TX	8885	10103603
Aroclor-1232 (PCB-1232)	TX	8890	10103603
Aroclor-1242 (PCB-1242)	TX	8895	10103603
Aroclor-1248 (PCB-1248)	TX	8900	10103603
Aroclor-1254 (PCB-1254)	TX	8905	10103603
Aroclor-1260 (PCB-1260)	TX	8910	10103603
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10103603
Chlordane (tech.)	TX	7250	10103603
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10103603
Dieldrin	TX	7470	10103603
Endosulfan I	TX	7510	10103603
Endosulfan II	TX	7515	10103603
Endosulfan sulfate	TX	7520	10103603
Endrin	TX	7540	10103603
Endrin aldehyde	TX	7530	10103603
Endrin ketone	TX	7535	10103603
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10103603
gamma-Chlordane	TX	7245	10103603
Heptachlor	TX	7685	10103603





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**Matrix: Non-Potable Water**

Heptachlor epoxide	TX	7690	10103603
Methoxychlor	TX	7810	10103603
Toxaphene (Chlorinated camphene)	TX	8250	10103603

**Method EPA 624**

Analyte	AB	Analyte ID	Method ID
1,1,1-Trichloroethane	TX	5160	10107207
1,1,2,2-Tetrachloroethane	TX	5110	10107207
1,1,2-Trichloroethane	TX	5165	10107207
1,1-Dichloroethane	TX	4630	10107207
1,1-Dichloroethylene	TX	4640	10107207
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10107207
1,2-Dichlorobenzene	TX	4610	10107207
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10107207
1,2-Dichloropropane	TX	4655	10107207
1,3-Dichlorobenzene	TX	4615	10107207
1,4-Dichlorobenzene	TX	4620	10107207
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10107207
2-Chloroethyl vinyl ether	TX	4500	10107207
Acetone (2-Propanone)	TX	4315	10107207
Acrolein (Propenal)	TX	4325	10107207
Acrylonitrile	TX	4340	10107207
Benzene	TX	4375	10107207
Bromodichloromethane	TX	4395	10107207
Bromoform	TX	4400	10107207
Carbon tetrachloride	TX	4455	10107207
Chlorobenzene	TX	4475	10107207
Chlorodibromomethane	TX	4575	10107207
Chloroethane (Ethyl chloride)	TX	4485	10107207
Chloroform	TX	4505	10107207
cis-1,2-Dichloroethylene	TX	4645	10107207





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**Matrix: Non-Potable Water**

cis-1,3-Dichloropropene	TX	4680	10107207
Ethylbenzene	TX	4765	10107207
m+p-xylene	TX	5240	10107207
Methyl bromide (Bromomethane)	TX	4950	10107207
Methyl chloride (Chloromethane)	TX	4960	10107207
Methyl tert-butyl ether (MTBE)	TX	5000	10107207
Methylene chloride (Dichloromethane)	TX	4975	10107207
Naphthalene	TX	5005	10107207
o-Xylene	TX	5250	10107207
Tetrachloroethylene (Perchloroethylene)	TX	5115	10107207
Toluene	TX	5140	10107207
trans-1,2-Dichloroethylene	TX	4700	10107207
trans-1,3-Dichloropropylene	TX	4685	10107207
Trichloroethene (Trichloroethylene)	TX	5170	10107207
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10107207
Vinyl chloride	TX	5235	10107207
Xylene (total)	TX	5260	10107207

**Method EPA 625**

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10107401
1,2,4-Trichlorobenzene	TX	5155	10107401
1,2-Dichlorobenzene	TX	4610	10107401
1,2-Diphenylhydrazine	TX	6220	10107401
1,3-Dichlorobenzene	TX	4615	10107401
1,4-Dichlorobenzene	TX	4620	10107401
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10107401
2,4,5-Trichlorophenol	TX	6835	10107401
2,4,6-Trichlorophenol	TX	6840	10107401
2,4-Dichlorophenol	TX	6000	10107401
2,4-Dimethylphenol	TX	6130	10107401



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**Matrix: Non-Potable Water**

2,4-Dinitrophenol	TX	6175	10107401
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10107401
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10107401
2-Chloronaphthalene	TX	5795	10107401
2-Chlorophenol	TX	5800	10107401
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10107401
2-Methylphenol (o-Cresol)	TX	6400	10107401
2-Nitrophenol	TX	6490	10107401
3,3'-Dichlorobenzidine	TX	5945	10107401
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10107401
4-Chloro-3-methylphenol	TX	5700	10107401
4-Chlorophenyl phenylether	TX	5825	10107401
4-Methylphenol (p-Cresol)	TX	6410	10107401
4-Nitrophenol	TX	6500	10107401
Acenaphthene	TX	5500	10107401
Acenaphthylene	TX	5505	10107401
Anthracene	TX	5555	10107401
Benzidine	TX	5595	10107401
Benzo(a)anthracene	TX	5575	10107401
Benzo(a)pyrene	TX	5580	10107401
Benzo(b)fluoranthene	TX	5585	10107401
Benzo(g,h,i)perylene	TX	5590	10107401
Benzo(k)fluoranthene	TX	5600	10107401
bis(2-Chloroethoxy)methane	TX	5760	10107401
bis(2-Chloroethyl) ether	TX	5765	10107401
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10107401
Butyl benzyl phthalate	TX	5670	10107401
Chrysene	TX	5855	10107401
Dibenz(a,h) anthracene	TX	5895	10107401
Diethyl phthalate	TX	6070	10107401



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**Matrix: Non-Potable Water**

Dimethyl phthalate	TX	6135	10107401
Di-n-butyl phthalate	TX	5925	10107401
Di-n-octyl phthalate	TX	6200	10107401
Fluoranthene	TX	6265	10107401
Fluorene	TX	6270	10107401
Hexachlorobenzene	TX	6275	10107401
Hexachlorobutadiene	TX	4835	10107401
Hexachlorocyclopentadiene	TX	6285	10107401
Hexachloroethane	TX	4840	10107401
Indeno(1,2,3-cd) pyrene	TX	6315	10107401
Isophorone	TX	6320	10107401
Naphthalene	TX	5005	10107401
Nitrobenzene	TX	5015	10107401
n-Nitrosodiethylamine	TX	6525	10107401
n-Nitrosodimethylamine	TX	6530	10107401
n-Nitrosodi-n-butylamine	TX	5025	10107401
n-Nitrosodi-n-propylamine	TX	6545	10107401
n-Nitrosodiphenylamine	TX	6535	10107401
Pentachlorobenzene	TX	6590	10107401
Pentachlorophenol	TX	6605	10107401
Phenanthrene	TX	6615	10107401
Phenol	TX	6625	10107401
Pyrene	TX	6665	10107401
Pyridine	TX	5095	10107401
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603



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**Matrix: Non-Potable Water**

**Method EPA 8011**

Analyte	AB	Analyte ID	Method ID
1,2,3-Trichloropropane	TX	5180	10173009
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10173009
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10173009

**Method EPA 8015**

Analyte	AB	Analyte ID	Method ID
Diesel range organics (DRO)	TX	9369	10173203
Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402



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**Matrix: Non-Potable Water**

4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402
Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Hexachlorobenzene	TX	6275	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201



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**Matrix: Non-Potable Water**

Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8151**

Analyte	AB	Analyte ID	Method ID
2,4,5-T	TX	8655	10183003
2,4-D	TX	8545	10183003
2,4-DB	TX	8560	10183003
Dalapon	TX	8555	10183003
Dicamba	TX	8595	10183003
Dichloroprop (Dichloroprop, Weedone)	TX	8605	10183003
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10183003
MCPA	TX	7775	10183003
MCPP	TX	7780	10183003
Silvex (2,4,5-TP)	TX	8650	10183003

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404
1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404





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**Matrix: Non-Potable Water**

1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
2-Pentanone	TX	5045	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404
Allyl alcohol	TX	4350	10184404
Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404



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**Matrix: Non-Potable Water**

Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Di-isopropylether (DIPE)	TX	9375	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Ethyl-t-butylether (ETBE) (2-Ethoxy-2-methylpropane)	TX	4770	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404
Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404





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**Matrix: Non-Potable Water**

Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
T-amylmethylether (TAME)	TX	4370	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404
Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404



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**Matrix: Non-Potable Water**

Method EPA 8270

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203
2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203



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**Matrix: Non-Potable Water**

2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Dimethyl aminoazobenzene	TX	6105	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrobiphenyl	TX	6480	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Chloro-2-methylaniline	TX	5695	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203



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**Matrix: Non-Potable Water**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Captan	TX	7190	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203



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**Matrix: Non-Potable Water**

Chrysene	TX	5855	10185203
Coumaphos	TX	7315	10185203
Demeton	TX	7390	10185203
Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Dioxathion	TX	7495	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethion	TX	7565	10185203
Ethyl methanesulfonate	TX	6260	10185203
Famphur	TX	7580	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203



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**Matrix: Non-Potable Water**

Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203
Kepone	TX	7740	10185203
Maleic anhydride	TX	6335	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naled	TX	7905	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203





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**Matrix: Non-Potable Water**

Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Phosmet (Imidan)	TX	8000	10185203
Phthalic anhydride	TX	6640	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203
Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Resorcinol	TX	6680	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203
Trifluralin (Treflan)	TX	8295	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209



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**Matrix: Non-Potable Water**

1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209

**Method EPA 8316**

Analyte	AB	Analyte ID	Method ID
Acrylamide	TX	4330	10188202

**Method EPA 8330**

Analyte	AB	Analyte ID	Method ID
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807





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**Matrix: Non-Potable Water**

Methyl-2,4,6-trinitrophenylamine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807
<b>Method EPA 9014</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803
<b>Method EPA 9038</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Sulfate	TX	2000	10196608
<b>Method EPA 9040</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
pH	TX	1900	10196802
<b>Method EPA 9050</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	10198604
<b>Method EPA 9056</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209
<b>Method EPA 9060</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	10200201



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**Matrix: Non-Potable Water**

Method	Analyte	AB	Analyte ID	Method ID
EPA 9065	Total phenolics	TX	1905	10200405
EPA 9066	Total phenolics	TX	1905	10200609
EPA 9250	Chloride	TX	1575	10207202
EPA RSK 175	2-methylpropane (Isobutane)	TX	4942	10212905
	Ethane	TX	4747	10212905
	Ethene	TX	4752	10212905
	Methane	TX	4926	10212905
	n-Butane	TX	5007	10212905
	n-Propane	TX	5029	10212905
HACH 8000	Chemical oxygen demand (COD)	TX	1565	60003001
SM 2120 B	Color	TX	1605	20223807
SM 2310 B (4a)	Acidity, as CaCO3	TX	1500	20002806
SM 2320 B	Alkalinity as CaCO3	TX	1505	20045005
SM 2340 B		AB	Analyte ID	Method ID



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**Matrix: Non-Potable Water**

Total hardness as CaCO <sub>3</sub>	TX	1755	20046008
<b>Method</b> SM 2510 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	20048004
<b>Method</b> SM 2540 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-total (total solids)	TX	1950	20004608
<b>Method</b> SM 2540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-filterable (TDS)	TX	1955	20049803
<b>Method</b> SM 2540 D			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-nonfilterable (TSS)	TX	1960	20004802
<b>Method</b> SM 3500-Cr B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	20065809
<b>Method</b> SM 4500-Cl F			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total residual chlorine	TX	1940	20080482
<b>Method</b> SM 4500-Cl <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chloride	TX	1575	20019209
<b>Method</b> SM 4500-CN <sup>-</sup> C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20020808
<b>Method</b> SM 4500-CN <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20021209
<b>Method</b> SM 4500-CN <sup>-</sup> G			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	20021607



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**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method SM 4500-H+ B			
Analyte	AB	Analyte ID	Method ID
pH	TX	1900	20104603
Method SM 4500-NH3 D			
Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	20108809
Kjeldahl Nitrogen (Total Kjeldahl Nitrogen-TKN)	TX	1790	20108809
Method SM 4500-NH3 F			
Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	20023001
Method SM 4500-O G			
Analyte	AB	Analyte ID	Method ID
Oxygen, dissolved	TX	1880	20025405
Method SM 4500-P E			
Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	20025803
Phosphorus	TX	1910	20025803
Method SM 4500-S2 <sup>-</sup> F			
Analyte	AB	Analyte ID	Method ID
Sulfide	TX	2005	20126209
Method SM 4500-SiO2 D			
Analyte	AB	Analyte ID	Method ID
Silica as SiO2	TX	1990	20127202
Method SM 4500-SO3 <sup>-</sup> B			
Analyte	AB	Analyte ID	Method ID
Sulfite	TX	2015	20026806
Method SM 5210 B			
Analyte	AB	Analyte ID	Method ID
Biochemical oxygen demand (BOD)	TX	1530	20027401
Carbonaceous BOD, CBOD	TX	1555	20027401
Method SM 5310 B			
Analyte	AB	Analyte ID	Method ID



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**Matrix: Non-Potable Water**

Total Organic Carbon (TOC)	TX	2040	20137206
<b>Method</b> SM 5310 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	20138209
<b>Method</b> SM 5540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Surfactants - MBAS	TX	2025	20144405
<b>Method</b> TCEQ 1005			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
Method ASTM D2216			
Analyte Moisture	TX	10337	ASTM D2216-05
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 1030			
Analyte Ignitability	TX	1780	10117201
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 1668			
Analyte Decachlorobiphenyls	TX	10332	10262007
Dichlorobiphenyls	TX	464	10262007
Heptachlorobiphenyls	TX	486	10262007
Hexachlorobiphenyls	TX	487	10262007
Monochlorobiphenyls	TX	501	10262007
Nonachlorobiphenyls	TX	507	10262007
Octachlorobiphenyls	TX	508	10262007
Pentachlorobiphenyls	TX	515	10262007
Tetrachlorobiphenyls	TX	528	10262007
Trichlorobiphenyls	TX	541	10262007
Method EPA 200.8			
Analyte Uranium	TX	3035	10014605



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Matrix: **Solid & Chemical Materials**

**Method** EPA 300.0

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method** EPA 310.1

Analyte	AB	Analyte ID	Method ID
Alkalinity as CaCO3	TX	1505	10054805

**Method** EPA 350.3

Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	10064401

**Method** EPA 365.3

Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	10070801
Phosphorus	TX	1910	10070801

**Method** EPA 6020

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156204
Antimony	TX	1005	10156204
Arsenic	TX	1010	10156204
Barium	TX	1015	10156204
Beryllium	TX	1020	10156204
Boron	TX	1025	10156204
Cadmium	TX	1030	10156204
Calcium	TX	1035	10156204
Chromium	TX	1040	10156204





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**Matrix: Solid & Chemical Materials**

Cobalt	TX	1050	10156204
Copper	TX	1055	10156204
Iron	TX	1070	10156204
Lead	TX	1075	10156204
Lithium	TX	1080	10156204
Magnesium	TX	1085	10156204
Manganese	TX	1090	10156204
Molybdenum	TX	1100	10156204
Nickel	TX	1105	10156204
Potassium	TX	1125	10156204
Selenium	TX	1140	10156204
Silver	TX	1150	10156204
Sodium	TX	1155	10156204
Strontium	TX	1160	10156204
Thallium	TX	1165	10156204
Tin	TX	1175	10156204
Titanium	TX	1180	10156204
Vanadium	TX	1185	10156204
Zinc	TX	1190	10156204
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603
<b>Method EPA 7471</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10166004
<b>Method EPA 8015</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Diesel range organics (DRO)	TX	9369	10173203





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**Matrix: Solid & Chemical Materials**

Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402
4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402



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**Matrix: Solid & Chemical Materials**

Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201
Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404



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**Matrix: Solid & Chemical Materials**

1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404



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**Matrix: Solid & Chemical Materials**

Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404
Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404



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**Matrix: Solid & Chemical Materials**

Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404
Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: Solid & Chemical Materials**

Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404
<b>Method EPA 8270</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203





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**Matrix: Solid & Chemical Materials**

2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203
2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203



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**Matrix: Solid & Chemical Materials**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203
Chrysene	TX	5855	10185203





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**Matrix: Solid & Chemical Materials**

Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzo(a,e) pyrene	TX	5890	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethyl methanesulfonate	TX	6260	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203
Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203



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**Matrix: Solid & Chemical Materials**

Kepona	TX	7740	10185203
Malathion	TX	7770	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203
Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203



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**Matrix: Solid & Chemical Materials**

Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209



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**Matrix: Solid & Chemical Materials**

Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209

**Method EPA 8316**

Analyte	AB	Analyte ID	Method ID
Acrylamide	TX	4330	10188202

**Method EPA 8330**

Analyte	AB	Analyte ID	Method ID
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807
Methyl-2,4,6-trinitrophenylnitramine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807

**Method EPA 9014**

Analyte	AB	Analyte ID	Method ID
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803

**Method EPA 9038**

Analyte	AB	Analyte ID	Method ID
Sulfate	TX	2000	10196608



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**Matrix: Solid & Chemical Materials**

**Method EPA 9040**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197203
pH	TX	1900	10196802

**Method EPA 9045**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197805
pH	TX	1900	10197805

**Method EPA 9050**

Analyte	AB	Analyte ID	Method ID
Conductivity	TX	1610	10198604

**Method EPA 9056**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209

**Method EPA 9060**

Analyte	AB	Analyte ID	Method ID
Total Organic Carbon (TOC)	TX	2040	10200201

**Method EPA 9065**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10200405

**Method EPA 9071**

Analyte	AB	Analyte ID	Method ID
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10201204



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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
EPA 9095			
<b>Analyte</b> Paint Filter Liquids Test	TX	10312	10204009
EPA 9250			
<b>Analyte</b> Chloride	TX	1575	10207202
SM 2320 B			
<b>Analyte</b> Alkalinity as CaCO3	TX	1505	20045005
SM 2510 B			
<b>Analyte</b> Conductivity	TX	1610	20048004
SM 2540 G			
<b>Analyte</b> Residue-total (total solids)	TX	1950	20005203
SSA/ASA Part 3:34			
<b>Analyte</b> Carbon, organic (Walkley-Black)	TX	10340	SSA/ASA Pt 3:34
TCEQ 1005			
<b>Analyte</b> Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

January 23, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20010618**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 13 sample(s) on Jan 15, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey



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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.



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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/23/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010618			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149789			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X			1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?	X				
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?		X			2
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

<b>Laboratory Review Checklist: Supporting Data</b>							
Laboratory Name: ALS Laboratory Group				LRC Date: 01/23/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20010618			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 149789			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?			X		
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?			X		
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?			X		
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group		LRC Date: 01/23/2020
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS20010618
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 149789
ER# <sup>5</sup>	Description	
1	Semivolatile Organics Method SW3510/8270, sample WG-1620-FD01-20200114, surrogate 4-Terphenyl-d14 recovered above the control limit due to possible matrix interference.	
1,2	Batch 149789, Semivolatile Organics Method 8270, sample WG-1620-P12-20200113, MS/MSD RPD recovered above the RPD limits for surrogates 2,4,6-Tribromophenol and 2-Fluorobiphenyl. The individual recoveries met acceptance criteria.	
<p>Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.</p> <p>O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);          NA = Not Applicable;          NR = Not Reviewed;          R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).</p>		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20010618

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20010618-01	WG-1620-P12-20200113	Groundwater		13-Jan-2020 10:30	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-02	WG-1620-FB01-20200113	Water		13-Jan-2020 11:30	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-03	WG-1620-MW08-20200113	Groundwater		13-Jan-2020 11:30	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-04	WG-1620-MW07-20200113	Groundwater		13-Jan-2020 13:15	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-05	WG-1620-P10-20200113	Groundwater		13-Jan-2020 14:20	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-06	WG-1620-MW11B-20200114	Groundwater		14-Jan-2020 09:25	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-07	WG-1620-MW11A-20200114	Groundwater		14-Jan-2020 10:35	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-08	WG-1620-MW10B-20200114	Groundwater		14-Jan-2020 11:35	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-09	WG-1620-MW10A-20200114	Groundwater		14-Jan-2020 12:25	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-10	WG-1620-MW02-20200114	Groundwater		14-Jan-2020 13:30	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-11	WG-1620-MW01A-20200114	Groundwater		14-Jan-2020 14:35	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-12	WG-1620-FD01-20200114	Groundwater		14-Jan-2020 14:35	15-Jan-2020 11:50	<input type="checkbox"/>
HS20010618-13	WG-1620-FB02-20200114	Water		14-Jan-2020 15:00	15-Jan-2020 11:50	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-P12-20200113  
 Collection Date: 13-Jan-2020 10:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-01  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
Acenaphthene	U		0.000027	0.00010	mg/L	1	20-Jan-2020 18:20
Acenaphthylene	U		0.000015	0.00010	mg/L	1	20-Jan-2020 18:20
<b>Anthracene</b>	<b>0.00010</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:20
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	20-Jan-2020 18:20
Dibenzofuran	U		0.000020	0.00010	mg/L	1	20-Jan-2020 18:20
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	20-Jan-2020 18:20
Fluoranthene	U		0.000010	0.00010	mg/L	1	20-Jan-2020 18:20
Fluorene	U		0.000030	0.00010	mg/L	1	20-Jan-2020 18:20
<b>Naphthalene</b>	<b>0.00016</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:20
Phenol	U		0.000035	0.00020	mg/L	1	20-Jan-2020 18:20
<b>Pyrene</b>	<b>0.00063</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	20-Jan-2020 18:20
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 18:20</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>84.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 18:20</i>
<i>Surr: 2-Fluorophenol</i>	<i>68.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 18:20</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>92.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 18:20</i>
<i>Surr: Nitrobenzene-d5</i>	<i>59.2</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 18:20</i>
<i>Surr: Phenol-d6</i>	<i>66.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>20-Jan-2020 18:20</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB01-20200113  
 Collection Date: 13-Jan-2020 11:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-02  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>	<b>Method:SW8270</b>				Prep:SW3510 / 19-Jan-2020		Analyst: LG
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 16:43
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 16:43
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 16:43
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 16:43
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 16:43
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 16:43
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 16:43
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 16:43
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 16:43
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 16:43
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 16:43
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 16:43
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 16:43
<i>Surr: 2,4,6-Tribromophenol</i>		73.7		34-129	%REC	1	21-Jan-2020 16:43
<i>Surr: 2-Fluorobiphenyl</i>		91.9		40-125	%REC	1	21-Jan-2020 16:43
<i>Surr: 2-Fluorophenol</i>		70.8		20-120	%REC	1	21-Jan-2020 16:43
<i>Surr: 4-Terphenyl-d14</i>		97.3		40-135	%REC	1	21-Jan-2020 16:43
<i>Surr: Nitrobenzene-d5</i>		80.7		41-120	%REC	1	21-Jan-2020 16:43
<i>Surr: Phenol-d6</i>		78.3		20-120	%REC	1	21-Jan-2020 16:43

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW08-20200113  
 Collection Date: 13-Jan-2020 11:30

**ANALYTICAL REPORT**

WorkOrder:HS20010618  
 Lab ID:HS20010618-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 17:02
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 17:02
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 17:02
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 17:02
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00021</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 17:02
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 17:02
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 17:02
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 17:02
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 17:02
<b>Phenanthrene</b>	<b>0.000046</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:02
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 17:02
<i>Surr: 2,4,6-Tribromophenol</i>	63.9			34-129	%REC	1	21-Jan-2020 17:02
<i>Surr: 2-Fluorobiphenyl</i>	70.0			40-125	%REC	1	21-Jan-2020 17:02
<i>Surr: 2-Fluorophenol</i>	55.9			20-120	%REC	1	21-Jan-2020 17:02
<i>Surr: 4-Terphenyl-d14</i>	96.0			40-135	%REC	1	21-Jan-2020 17:02
<i>Surr: Nitrobenzene-d5</i>	62.1			41-120	%REC	1	21-Jan-2020 17:02
<i>Surr: Phenol-d6</i>	60.3			20-120	%REC	1	21-Jan-2020 17:02

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW07-20200113  
 Collection Date: 13-Jan-2020 13:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-04  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
<b>2-Methylnaphthalene</b>	<b>0.000066</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:21
Acenaphthene		U	0.000027	0.00010	mg/L	1	21-Jan-2020 17:21
Acenaphthylene		U	0.000015	0.00010	mg/L	1	21-Jan-2020 17:21
Anthracene		U	0.000014	0.00010	mg/L	1	21-Jan-2020 17:21
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	21-Jan-2020 17:21
<b>Dibenzofuran</b>	<b>0.000057</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:21
<b>Fluoranthene</b>	<b>0.00010</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:21
Fluorene		U	0.000030	0.00010	mg/L	1	21-Jan-2020 17:21
<b>Naphthalene</b>	<b>0.00017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:21
<b>Phenanthrene</b>	<b>0.00014</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:21
Pyrene		U	0.000019	0.00010	mg/L	1	21-Jan-2020 17:21
<i>Surr: 2,4,6-Tribromophenol</i>	<i>74.6</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:21</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>80.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:21</i>
<i>Surr: 2-Fluorophenol</i>	<i>59.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:21</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>85.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:21</i>
<i>Surr: Nitrobenzene-d5</i>	<i>66.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:21</i>
<i>Surr: Phenol-d6</i>	<i>71.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:21</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-P10-20200113  
 Collection Date: 13-Jan-2020 14:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-05  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 17:40
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 17:40
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 17:40
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 17:40
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 17:40
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 17:40
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 17:40
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 17:40
<b>Naphthalene</b>	<b>0.00017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 17:40
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 17:40
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 17:40
<i>Surr: 2,4,6-Tribromophenol</i>	<i>79.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:40</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>93.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:40</i>
<i>Surr: 2-Fluorophenol</i>	<i>78.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:40</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>99.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:40</i>
<i>Surr: Nitrobenzene-d5</i>	<i>82.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:40</i>
<i>Surr: Phenol-d6</i>	<i>78.2</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 17:40</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW11B-20200114  
 Collection Date: 14-Jan-2020 09:25

**ANALYTICAL REPORT**

WorkOrder:HS20010618  
 Lab ID:HS20010618-06  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
<b>Acenaphthene</b>	<b>0.033</b>		<b>0.00014</b>	<b>0.00050</b>	<b>mg/L</b>	5	22-Jan-2020 12:03
<b>Acenaphthylene</b>	<b>0.0016</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 18:00
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 18:00
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000095</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 18:00
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 18:00
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 18:00
<b>Fluoranthene</b>	<b>0.0024</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 18:00
<b>Fluorene</b>	<b>0.00035</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 18:00
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 18:00
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 18:00
<b>Pyrene</b>	<b>0.0023</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 18:00
Surr: 2,4,6-Tribromophenol	81.4			34-129	%REC	1	21-Jan-2020 18:00
Surr: 2,4,6-Tribromophenol	101			34-129	%REC	5	22-Jan-2020 12:03
Surr: 2-Fluorobiphenyl	90.1			40-125	%REC	5	22-Jan-2020 12:03
Surr: 2-Fluorobiphenyl	72.9			40-125	%REC	1	21-Jan-2020 18:00
Surr: 2-Fluorophenol	74.1			20-120	%REC	5	22-Jan-2020 12:03
Surr: 2-Fluorophenol	56.2			20-120	%REC	1	21-Jan-2020 18:00
Surr: 4-Terphenyl-d14	92.8			40-135	%REC	1	21-Jan-2020 18:00
Surr: 4-Terphenyl-d14	123			40-135	%REC	5	22-Jan-2020 12:03
Surr: Nitrobenzene-d5	80.4			41-120	%REC	5	22-Jan-2020 12:03
Surr: Nitrobenzene-d5	61.3			41-120	%REC	1	21-Jan-2020 18:00
Surr: Phenol-d6	60.8			20-120	%REC	1	21-Jan-2020 18:00
Surr: Phenol-d6	77.3			20-120	%REC	5	22-Jan-2020 12:03

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW11A-20200114  
 Collection Date: 14-Jan-2020 10:35

**ANALYTICAL REPORT**

WorkOrder:HS20010618  
 Lab ID:HS20010618-07  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 18:19
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 18:19
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 18:19
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 18:19
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 18:19
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 18:19
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 18:19
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 18:19
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 18:19
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 18:19
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 18:19
<i>Surr: 2,4,6-Tribromophenol</i>		<i>78.8</i>		<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:19</i>
<i>Surr: 2-Fluorobiphenyl</i>		<i>80.3</i>		<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:19</i>
<i>Surr: 2-Fluorophenol</i>		<i>56.8</i>		<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:19</i>
<i>Surr: 4-Terphenyl-d14</i>		<i>92.1</i>		<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:19</i>
<i>Surr: Nitrobenzene-d5</i>		<i>63.9</i>		<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:19</i>
<i>Surr: Phenol-d6</i>		<i>65.0</i>		<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:19</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW10B-20200114  
 Collection Date: 14-Jan-2020 11:35

**ANALYTICAL REPORT**

WorkOrder:HS20010618  
 Lab ID:HS20010618-08  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
Acenaphthene	0.069		0.00027	0.0010	mg/L	10	22-Jan-2020 12:22
Acenaphthylene	0.00066		0.000015	0.00010	mg/L	1	21-Jan-2020 18:38
Anthracene	0.0028		0.000014	0.00010	mg/L	1	21-Jan-2020 18:38
Bis(2-ethylhexyl)phthalate	0.00020		0.000037	0.00020	mg/L	1	21-Jan-2020 18:38
Dibenzofuran	0.022		0.00020	0.0010	mg/L	10	22-Jan-2020 12:22
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 18:38
Fluoranthene	0.0029		0.000010	0.00010	mg/L	1	21-Jan-2020 18:38
Fluorene	0.036		0.00030	0.0010	mg/L	10	22-Jan-2020 12:22
Naphthalene	0.0021		0.000020	0.00010	mg/L	1	21-Jan-2020 18:38
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 18:38
Pyrene	0.0013		0.000019	0.00010	mg/L	1	21-Jan-2020 18:38
Surr: 2,4,6-Tribromophenol	70.8			34-129	%REC	10	22-Jan-2020 12:22
Surr: 2,4,6-Tribromophenol	80.7			34-129	%REC	1	21-Jan-2020 18:38
Surr: 2-Fluorobiphenyl	77.4			40-125	%REC	1	21-Jan-2020 18:38
Surr: 2-Fluorobiphenyl	85.2			40-125	%REC	10	22-Jan-2020 12:22
Surr: 2-Fluorophenol	60.1			20-120	%REC	10	22-Jan-2020 12:22
Surr: 2-Fluorophenol	62.8			20-120	%REC	1	21-Jan-2020 18:38
Surr: 4-Terphenyl-d14	103			40-135	%REC	10	22-Jan-2020 12:22
Surr: 4-Terphenyl-d14	89.0			40-135	%REC	1	21-Jan-2020 18:38
Surr: Nitrobenzene-d5	75.6			41-120	%REC	10	22-Jan-2020 12:22
Surr: Nitrobenzene-d5	68.2			41-120	%REC	1	21-Jan-2020 18:38
Surr: Phenol-d6	69.2			20-120	%REC	1	21-Jan-2020 18:38
Surr: Phenol-d6	74.4			20-120	%REC	10	22-Jan-2020 12:22

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW10A-20200114  
 Collection Date: 14-Jan-2020 12:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-09  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 18:57
<b>Acenaphthene</b>	<b>0.00011</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 18:57
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 18:57
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 18:57
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 18:57
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 18:57
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 18:57
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 18:57
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 18:57
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 18:57
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 18:57
<i>Surr: 2,4,6-Tribromophenol</i>	<i>60.3</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:57</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>77.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:57</i>
<i>Surr: 2-Fluorophenol</i>	<i>64.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:57</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>110</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:57</i>
<i>Surr: Nitrobenzene-d5</i>	<i>68.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:57</i>
<i>Surr: Phenol-d6</i>	<i>64.7</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>21-Jan-2020 18:57</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW02-20200114  
 Collection Date: 14-Jan-2020 13:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-10  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
2-Methylnaphthalene		U	0.000019	0.00010	mg/L	1	21-Jan-2020 19:16
<b>Acenaphthene</b>	<b>0.0030</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
Acenaphthylene		U	0.000015	0.00010	mg/L	1	21-Jan-2020 19:16
<b>Anthracene</b>	<b>0.00011</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
Bis(2-ethylhexyl)phthalate		U	0.000037	0.00020	mg/L	1	21-Jan-2020 19:16
<b>Dibenzofuran</b>	<b>0.00039</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
<b>Fluoranthene</b>	<b>0.00024</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
<b>Fluorene</b>	<b>0.0017</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
Naphthalene		U	0.000020	0.00010	mg/L	1	21-Jan-2020 19:16
<b>Phenanthrene</b>	<b>0.00011</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
<b>Pyrene</b>	<b>0.00011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:16
Surr: 2,4,6-Tribromophenol	64.5			34-129	%REC	1	21-Jan-2020 19:16
Surr: 2-Fluorobiphenyl	83.6			40-125	%REC	1	21-Jan-2020 19:16
Surr: 2-Fluorophenol	62.6			20-120	%REC	1	21-Jan-2020 19:16
Surr: 4-Terphenyl-d14	99.3			40-135	%REC	1	21-Jan-2020 19:16
Surr: Nitrobenzene-d5	70.1			41-120	%REC	1	21-Jan-2020 19:16
Surr: Phenol-d6	66.9			20-120	%REC	1	21-Jan-2020 19:16

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW01A-20200114  
 Collection Date: 14-Jan-2020 14:35

**ANALYTICAL REPORT**

WorkOrder:HS20010618  
 Lab ID:HS20010618-11  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MLL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
<b>2-Methylnaphthalene</b>	<b>0.00019</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
<b>Acenaphthene</b>	<b>0.024</b>		<b>0.00014</b>	<b>0.00050</b>	<b>mg/L</b>	5	22-Jan-2020 12:41
<b>Acenaphthylene</b>	<b>0.00084</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 19:35
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 19:35
<b>Dibenzofuran</b>	<b>0.0036</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
<b>Fluoranthene</b>	<b>0.0011</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
<b>Fluorene</b>	<b>0.0064</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
<b>Naphthalene</b>	<b>0.00052</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 19:35
<b>Pyrene</b>	<b>0.00052</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:35
Surr: 2,4,6-Tribromophenol	81.3			34-129	%REC	1	21-Jan-2020 19:35
Surr: 2,4,6-Tribromophenol	82.2			34-129	%REC	5	22-Jan-2020 12:41
Surr: 2-Fluorobiphenyl	95.4			40-125	%REC	5	22-Jan-2020 12:41
Surr: 2-Fluorobiphenyl	90.6			40-125	%REC	1	21-Jan-2020 19:35
Surr: 2-Fluorophenol	68.6			20-120	%REC	1	21-Jan-2020 19:35
Surr: 2-Fluorophenol	77.7			20-120	%REC	5	22-Jan-2020 12:41
Surr: 4-Terphenyl-d14	103			40-135	%REC	5	22-Jan-2020 12:41
Surr: 4-Terphenyl-d14	94.4			40-135	%REC	1	21-Jan-2020 19:35
Surr: Nitrobenzene-d5	88.0			41-120	%REC	5	22-Jan-2020 12:41
Surr: Nitrobenzene-d5	80.0			41-120	%REC	1	21-Jan-2020 19:35
Surr: Phenol-d6	75.8			20-120	%REC	1	21-Jan-2020 19:35
Surr: Phenol-d6	77.6			20-120	%REC	5	22-Jan-2020 12:41

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200114  
 Collection Date: 14-Jan-2020 14:35

**ANALYTICAL REPORT**

WorkOrder:HS20010618  
 Lab ID:HS20010618-12  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 19:54
<b>Acenaphthene</b>	<b>0.018</b>		<b>0.00011</b>	<b>0.00040</b>	<b>mg/L</b>	4	22-Jan-2020 13:00
<b>Acenaphthylene</b>	<b>0.00066</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:54
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 19:54
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000074</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	21-Jan-2020 19:54
<b>Dibenzofuran</b>	<b>0.0021</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:54
<b>Fluoranthene</b>	<b>0.0012</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:54
<b>Fluorene</b>	<b>0.0038</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:54
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 19:54
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 19:54
<b>Pyrene</b>	<b>0.00059</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	21-Jan-2020 19:54
Surr: 2,4,6-Tribromophenol	107			34-129	%REC	1	21-Jan-2020 19:54
Surr: 2,4,6-Tribromophenol	117			34-129	%REC	4	22-Jan-2020 13:00
Surr: 2-Fluorobiphenyl	98.7			40-125	%REC	4	22-Jan-2020 13:00
Surr: 2-Fluorobiphenyl	88.3			40-125	%REC	1	21-Jan-2020 19:54
Surr: 2-Fluorophenol	59.0			20-120	%REC	1	21-Jan-2020 19:54
Surr: 2-Fluorophenol	66.0			20-120	%REC	4	22-Jan-2020 13:00
Surr: 4-Terphenyl-d14	142	S		40-135	%REC	4	22-Jan-2020 13:00
Surr: 4-Terphenyl-d14	129			40-135	%REC	1	21-Jan-2020 19:54
Surr: Nitrobenzene-d5	63.6			41-120	%REC	1	21-Jan-2020 19:54
Surr: Nitrobenzene-d5	71.5			41-120	%REC	4	22-Jan-2020 13:00
Surr: Phenol-d6	80.1			20-120	%REC	4	22-Jan-2020 13:00
Surr: Phenol-d6	67.3			20-120	%REC	1	21-Jan-2020 19:54

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FB02-20200114  
 Collection Date: 14-Jan-2020 15:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20010618  
 Lab ID:HS20010618-13  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Jan-2020		Analyst: LG	
2-Methylnaphthalene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 20:13
Acenaphthene	U		0.000027	0.00010	mg/L	1	21-Jan-2020 20:13
Acenaphthylene	U		0.000015	0.00010	mg/L	1	21-Jan-2020 20:13
Anthracene	U		0.000014	0.00010	mg/L	1	21-Jan-2020 20:13
Bis(2-ethylhexyl)phthalate	U		0.000037	0.00020	mg/L	1	21-Jan-2020 20:13
Dibenzofuran	U		0.000020	0.00010	mg/L	1	21-Jan-2020 20:13
Di-n-butyl phthalate	U		0.000020	0.00020	mg/L	1	21-Jan-2020 20:13
Fluoranthene	U		0.000010	0.00010	mg/L	1	21-Jan-2020 20:13
Fluorene	U		0.000030	0.00010	mg/L	1	21-Jan-2020 20:13
Naphthalene	U		0.000020	0.00010	mg/L	1	21-Jan-2020 20:13
Phenanthrene	U		0.000021	0.00010	mg/L	1	21-Jan-2020 20:13
Phenol	U		0.000035	0.00020	mg/L	1	21-Jan-2020 20:13
Pyrene	U		0.000019	0.00010	mg/L	1	21-Jan-2020 20:13
<i>Surr: 2,4,6-Tribromophenol</i>		61.9		34-129	%REC	1	21-Jan-2020 20:13
<i>Surr: 2-Fluorobiphenyl</i>		109		40-125	%REC	1	21-Jan-2020 20:13
<i>Surr: 2-Fluorophenol</i>		79.0		20-120	%REC	1	21-Jan-2020 20:13
<i>Surr: 4-Terphenyl-d14</i>		100		40-135	%REC	1	21-Jan-2020 20:13
<i>Surr: Nitrobenzene-d5</i>		91.2		41-120	%REC	1	21-Jan-2020 20:13
<i>Surr: Phenol-d6</i>		88.1		20-120	%REC	1	21-Jan-2020 20:13

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**Batch ID:** 149789      **Start Date:** 19 Jan 2020 07:30      **End Date:** 19 Jan 2020 16:00  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20010618-01	1	1000 (mL)	1 (mL)	0.001
HS20010618-02	1	1000 (mL)	1 (mL)	0.001
HS20010618-03	1	1000 (mL)	1 (mL)	0.001
HS20010618-04	1	1000 (mL)	1 (mL)	0.001
HS20010618-05	1	1000 (mL)	1 (mL)	0.001
HS20010618-06	1	1000 (mL)	1 (mL)	0.001
HS20010618-07	1	1000 (mL)	1 (mL)	0.001
HS20010618-08	1	1000 (mL)	1 (mL)	0.001
HS20010618-09	1	1000 (mL)	1 (mL)	0.001
HS20010618-10	1	1000 (mL)	1 (mL)	0.001
HS20010618-11	1	1000 (mL)	1 (mL)	0.001
HS20010618-12	1	1000 (mL)	1 (mL)	0.001
HS20010618-13	1	1000 (mL)	1 (mL)	0.001

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 149789 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Water</b>	
HS20010618-02	WG-1620-FB01-20200113	13 Jan 2020 11:30		19 Jan 2020 15:02	21 Jan 2020 16:43	1
HS20010618-13	WG-1620-FB02-20200114	14 Jan 2020 15:00		19 Jan 2020 15:02	21 Jan 2020 20:13	1
<b>Batch ID: 149789 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Groundwater</b>	
HS20010618-01	WG-1620-P12-20200113	13 Jan 2020 10:30		19 Jan 2020 15:02	20 Jan 2020 18:20	1
HS20010618-03	WG-1620-MW08-20200113	13 Jan 2020 11:30		19 Jan 2020 15:02	21 Jan 2020 17:02	1
HS20010618-04	WG-1620-MW07-20200113	13 Jan 2020 13:15		19 Jan 2020 15:02	21 Jan 2020 17:21	1
HS20010618-05	WG-1620-P10-20200113	13 Jan 2020 14:20		19 Jan 2020 15:02	21 Jan 2020 17:40	1
HS20010618-06	WG-1620-MW11B-20200114	14 Jan 2020 09:25		19 Jan 2020 15:02	22 Jan 2020 12:03	5
HS20010618-06	WG-1620-MW11B-20200114	14 Jan 2020 09:25		19 Jan 2020 15:02	21 Jan 2020 18:00	1
HS20010618-07	WG-1620-MW11A-20200114	14 Jan 2020 10:35		19 Jan 2020 15:02	21 Jan 2020 18:19	1
HS20010618-08	WG-1620-MW10B-20200114	14 Jan 2020 11:35		19 Jan 2020 15:02	22 Jan 2020 12:22	10
HS20010618-08	WG-1620-MW10B-20200114	14 Jan 2020 11:35		19 Jan 2020 15:02	21 Jan 2020 18:38	1
HS20010618-09	WG-1620-MW10A-20200114	14 Jan 2020 12:25		19 Jan 2020 15:02	21 Jan 2020 18:57	1
HS20010618-10	WG-1620-MW02-20200114	14 Jan 2020 13:30		19 Jan 2020 15:02	21 Jan 2020 19:16	1
HS20010618-11	WG-1620-MW01A-20200114	14 Jan 2020 14:35		19 Jan 2020 15:02	22 Jan 2020 12:41	5
HS20010618-11	WG-1620-MW01A-20200114	14 Jan 2020 14:35		19 Jan 2020 15:02	21 Jan 2020 19:35	1
HS20010618-12	WG-1620-FD01-20200114	14 Jan 2020 14:35		19 Jan 2020 15:02	22 Jan 2020 13:00	4
HS20010618-12	WG-1620-FD01-20200114	14 Jan 2020 14:35		19 Jan 2020 15:02	21 Jan 2020 19:54	1

WorkOrder: HS20010618  
 InstrumentID: SV-6  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	2-Methylnaphthalene	91-57-6	0.000050	0.000060	0.000019	0.00010
A	Acenaphthene	83-32-9	0.000050	0.000059	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000050	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000047	0.000014	0.00010
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00012	0.000037	0.00020
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00011	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000049	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000053	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000049	0.000020	0.00010
A	Phenanthrene	85-01-8	0.000050	0.000049	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.000096	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000047	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**QC BATCH REPORT**

Batch ID: 149789 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
<b>MBLK</b>	Sample ID: <b>MBLK-149789</b>	Units: <b>ug/L</b>			Analysis Date: <b>20-Jan-2020 14:13</b>					
Client ID:	Run ID: <b>SV-6_354667</b>	SeqNo: <b>5441804</b>		PrepDate: <b>19-Jan-2020</b>		DF: <b>1</b>				
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
2-Methylnaphthalene	U	0.10								
Acenaphthene	U	0.10								
Acenaphthylene	U	0.10								
Anthracene	U	0.10								
Bis(2-ethylhexyl)phthalate	U	0.20								
Dibenzofuran	U	0.10								
Di-n-butyl phthalate	U	0.20								
Fluoranthene	U	0.10								
Fluorene	U	0.10								
Naphthalene	U	0.10								
Phenanthrene	U	0.10								
Phenol	U	0.20								
Pyrene	U	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.413</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>68.3</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.677</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>93.5</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.815</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.3</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.529</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>90.6</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.693</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.9</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.844</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.9</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**QC BATCH REPORT**

Batch ID: 149789 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
<b>LCS</b>	Sample ID: <b>LCS-149789</b>	Units: <b>ug/L</b>			Analysis Date: <b>20-Jan-2020 14:32</b>					
Client ID:	Run ID: <b>SV-6_354667</b>	SeqNo: <b>5441805</b>		PrepDate: <b>19-Jan-2020</b>		DF: <b>1</b>				
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD	RPD Limit Qual	
2-Methylnaphthalene	4.12	0.10	5	0	82.4	50 - 120				
Acenaphthene	4.616	0.10	5	0	92.3	45 - 120				
Acenaphthylene	4.438	0.10	5	0	88.8	47 - 120				
Anthracene	4.793	0.10	5	0	95.9	45 - 120				
Bis(2-ethylhexyl)phthalate	5.208	0.20	5	0	104	40 - 139				
Dibenzofuran	4.382	0.10	5	0	87.6	50 - 120				
Di-n-butyl phthalate	5.327	0.20	5	0	107	45 - 123				
Fluoranthene	4.608	0.10	5	0	92.2	45 - 125				
Fluorene	4.509	0.10	5	0	90.2	49 - 120				
Naphthalene	4.352	0.10	5	0	87.0	45 - 120				
Phenanthrene	4.579	0.10	5	0	91.6	45 - 121				
Phenol	4.045	0.20	5	0	80.9	20 - 124				
Pyrene	4.543	0.10	5	0	90.9	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.728</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.6</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.553</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.1</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.92</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.4</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.553</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.1</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.663</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.3</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.85</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>77.0</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**QC BATCH REPORT**

Batch ID: 149789 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MS		Sample ID: HS20010618-01MS		Units: ug/L		Analysis Date: 20-Jan-2020 18:40				
Client ID: WG-1620-P12-20200113		Run ID: SV-6_354667		SeqNo: 5441807		PrepDate: 19-Jan-2020		DF: 1		
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
2-Methylnaphthalene	4.077	0.10	5	0	81.5	50 - 120				
Acenaphthene	4.554	0.10	5	0	91.1	45 - 120				
Acenaphthylene	4.432	0.10	5	0	88.6	47 - 120				
Anthracene	5.422	0.10	5	0.1005	106	45 - 120				
Bis(2-ethylhexyl)phthalate	5.953	0.20	5	0	119	40 - 139				
Dibenzofuran	4.513	0.10	5	0	90.3	50 - 120				
Di-n-butyl phthalate	5.9	0.20	5	0	118	45 - 123				
Fluoranthene	5.303	0.10	5	0	106	45 - 125				
Fluorene	4.678	0.10	5	0	93.6	49 - 120				
Naphthalene	4.398	0.10	5	0.1605	84.8	45 - 120				
Phenanthrene	5.217	0.10	5	0.0662	103	45 - 121				
Phenol	3.858	0.20	5	0	77.2	20 - 124				
Pyrene	5.985	0.10	5	0.6322	107	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.348</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>87.0</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>4.593</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.9</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.499</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>70.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>5.237</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>105</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.192</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>63.8</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.6</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.0</i>	<i>20 - 120</i>				



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**QC BATCH REPORT**

Batch ID: 149789 ( 0 )		Instrument: SV-6		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MSD		Sample ID: HS20010618-01MSD		Units: ug/L		Analysis Date: 20-Jan-2020 18:59				
Client ID: WG-1620-P12-20200113		Run ID: SV-6_354667		SeqNo: 5441808		PrepDate: 19-Jan-2020		DF: 1		
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
2-Methylnaphthalene	3.617	0.10	5	0	72.3	50 - 120	4.077	12	20	
Acenaphthene	3.858	0.10	5	0	77.2	45 - 120	4.554	16.6	20	
Acenaphthylene	3.761	0.10	5	0	75.2	47 - 120	4.432	16.4	20	
Anthracene	4.577	0.10	5	0.1005	89.5	45 - 120	5.422	16.9	20	
Bis(2-ethylhexyl)phthalate	5.317	0.20	5	0	106	40 - 139	5.953	11.3	20	
Dibenzofuran	3.783	0.10	5	0	75.7	50 - 120	4.513	17.6	20	
Di-n-butyl phthalate	5.105	0.20	5	0	102	45 - 123	5.9	14.5	20	
Fluoranthene	4.616	0.10	5	0	92.3	45 - 125	5.303	13.8	20	
Fluorene	4	0.10	5	0	80.0	49 - 120	4.678	15.6	20	
Naphthalene	3.832	0.10	5	0.1605	73.4	45 - 120	4.398	13.8	20	
Phenanthrene	4.366	0.10	5	0.0662	86.0	45 - 121	5.217	17.8	20	
Phenol	3.371	0.20	5	0	67.4	20 - 124	3.858	13.5	20	
Pyrene	5.245	0.10	5	0.6322	92.2	40 - 130	5.985	13.2	20	
<i>Surr: 2,4,6-Tribromophenol</i>	3.502	0.20	5	0	70.0	34 - 129	4.348	21.5	20 R	
<i>Surr: 2-Fluorobiphenyl</i>	3.715	0.20	5	0	74.3	40 - 125	4.593	21.1	20 R	
<i>Surr: 2-Fluorophenol</i>	2.97	0.20	5	0	59.4	20 - 120	3.499	16.3	20	
<i>Surr: 4-Terphenyl-d14</i>	4.522	0.20	5	0	90.4	40 - 135	5.237	14.7	20	
<i>Surr: Nitrobenzene-d5</i>	2.709	0.20	5	0	54.2	41 - 120	3.192	16.4	20	
<i>Surr: Phenol-d6</i>	3.121	0.20	5	0	62.4	20 - 120	3.6	14.3	20	

The following samples were analyzed in this batch:

HS20010618-01	HS20010618-02	HS20010618-03	HS20010618-04
HS20010618-05	HS20010618-06	HS20010618-07	HS20010618-08
HS20010618-09	HS20010618-10	HS20010618-11	HS20010618-12
HS20010618-13			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20010618

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231	20-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20010618

Date/Time Received: 15-Jan-2020 11:50
Received by: AC

Checklist completed by: Paris Frazier
eSignature
Date: 15-Jan-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 15-Jan-2020

Matrices: WATER

Carrier name: Client

- Shipping container/cooler in good condition?
Custody seals intact on shipping container/cooler?
Custody seals intact on sample bottles?
VOA/TX1005/TX1006 Solids in hermetically sealed vials?
Chain of custody present?
Chain of custody signed when relinquished and received?
Samplers name present on COC?
Chain of custody agrees with sample labels?
Samples in proper container/bottle?
Sample containers intact?
Sufficient sample volume for indicated test?
All samples received within holding time?
Container/Temp Blank temperature in compliance?

- Yes/No/Not Present checkboxes for each item in the list above.

2 Page(s)
COC IDs:215992/215990

Temperature(s)/Thermometer(s): 2.3C/2.3C,2.1C/2.1C. 1.2C/1.2C UC/C IR25
Cooler(s)/Kit(s): 44307/45460/BLUE
Date/Time sample(s) sent to storage: 01.15.2020 17:00

- Water - VOA vials have zero headspace?
Water - pH acceptable upon receipt?
pH adjusted?
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



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Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page 1 of 2

COC ID: 215996

## HS20010618

Golder Associates Inc.  
Houston TX-Wood Preserving Works

Houston, WV  
1168

1280

Customer Information		Project Information	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works
Work Order		Project Number	1620-08-Rev0 SR 92688 SWMU1
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable
Address	2201 Double Creek Drive Suite 4004	Address	1400 Douglas Street Stop 0750
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750
Phone	(512) 671-3434	Phone	
Fax	(512) 671-3446	Fax	
e-Mail Address	eric_matzner@golder.com	e-Mail Address	



A 8270\_LOW\_W(5632532 ATZ SemiVolatiles)  
 B 8270\_LOW\_W(5632532 BTZ SemiVolatiles)  
 C 8270\_LOW\_W(5632532 ATZ & BTZ SemiVolatiles)  
 D  
 E  
 F  
 G  
 H  
 I  
 J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-P12-20200113	1-13-20	1030	Groundwa	6	2		X									
2	WG-1620-P12MS-20200113		1030	W		2		X									
3	WG-1620-P12MSD-20200113		1030	W		2		X									
4	WG-1620-FB01-20200113		1130	W		2		X									
5	WG-1620-MW08-20200113		1130	W		2	X										
6	WG-1620-MW07-20200113		1315	W		2	X										
7	WG-1620-P10-20200113		1420	W		2		X									
8	WG-1620-MW11B-20200114	1-14-20	0925	W		2		X									
9	WG-1620-MW11A-20200114		1035	W		2	X										
10	WG-1620-MW10B-20200114		1135	W		2		X									

Sampler(s) Please Print & Sign <i>JOHN BRAYTON jbr</i>		Shipment Method HAND DELIVERED		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour				Results Due Date:			
Relinquished by: <i>John</i>	Date: 1-15-20	Time: 11:50	Received by:	Notes: UPRR Houston MWPW							
Relinquished by:	Date: 1-15-2020	Time: 11:50	Received by (Laboratory): AC	Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)					
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	44307	2.3	<input type="checkbox"/> Level II Std GC	<input checked="" type="checkbox"/> TRRP Checklist				
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				45460	2.1	<input type="checkbox"/> Level III Std GC/Faw Date	<input type="checkbox"/> TRRP Level IV				
				BLUE	1.2	<input type="checkbox"/> Level IV SW846/CLP	<input type="checkbox"/> Other				

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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# Chain of Custody Form

Page 2 of 2

COC ID: 215990

## HS20010618

Golder Associates Inc.  
Houston TX-Wood Preserving Works

n, WV  
B  
D



Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8270_LOW_W (5632532 ATZ SemiVolatiles)
Work Order		Project Number	1620-08-Rev0 SR 92688 SWMU1	B	8270_LOW_W (5632532 BTZ SemiVolatiles)
Company Name	Golder Associates	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 ATZ & BTZ SemiVolatiles)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	eric_matzner@golder.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-MW10A-20200114	1-14-20	1225	Groundwa	8	2	X										
2	WG-1620-MW02-20200114		1330	W		2	X										
3	UG-1620-MW01A-20200114		1435	W		2	X										
4	WG-1620-FD01-20200114		1435	W		2	X										
5	WG-1620-FB02-20200114		1500	W		2	X										
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <b>JOHN BRAYTON</b>		Shipment Method <b>HAND DELIVERED</b>		Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 10 Wk Days <input type="checkbox"/> 5 Wk Days <input type="checkbox"/> 2 Wk Days <input type="checkbox"/> 24 Hour			Results Due Date:	
Relinquished by: <b>[Signature]</b>	Date: <b>1-15-20</b>	Time: <b>1:50</b>	Received by:	Notes: <b>UPRR Houston MWPW</b>				
Relinquished by: <b>[Signature]</b>	Date: <b>1-15-2020</b>	Time: <b>11:50</b>	Received by (Laboratory): <b>AC</b>	Cooler ID		Cooler Temp.		QC Package: (Check One Box Below)
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	<input type="checkbox"/> Level II Std QC		<input checked="" type="checkbox"/> TRRP Checklist		
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				<input type="checkbox"/> Level III Std QC/Raw Date		<input type="checkbox"/> TRRP Level IV		
				<input type="checkbox"/> Level IV SW846/CLP		<input type="checkbox"/> Other		

- Note:
- Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.
  - Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.
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# Memorandum

April 6, 2020

To: Eric Matzner Ref. No.: 11183954-1620

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From: <sup>CK</sup> Chris G. Knight/eew/585-NF Tel: 512-506-8803

---

cc: Jesse Orth, Jon Lang; Julie Lidstone

---

**Subject: Data Usability Summary  
HWPW - Groundwater Monitoring (New Wells)  
Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works  
Houston, Texas  
March 2020**

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## 1. Scope of Data Usability Study

This document details a Data Usability Summary (DUS) of analytical results for groundwater samples collected in support of the HWPW - Groundwater Monitoring (New Wells) at the Union Pacific Railroad (UPRR) / Houston TX-Wood Preserving Works site during March 2020. Samples were submitted to ALS Environmental (ALS), located in Houston, Texas and are reported in data packages HS20030619, HS20030819, and HS20030961. The intended use of the data is to support the HWPW - Groundwater Monitoring (New Wells) at the site by providing current concentration of chemicals of concern.

Data were reviewed and validated by Chris G. Knight of GHD, in accordance with Title 30 of the Texas Administrative Code Section 350.54 (30 TAC 350.54) as described in the Texas Commission on Environmental Quality (TCEQ) Regulatory Guidance document entitled "Review and Reporting of COC Concentration Data under TRRP", (RG-366/TRRP-13), revised May 2010, herein referred to as "TRRP-13 Guidance". Evaluation of the data was based on information obtained from the chain of custody forms, the finished report forms, method blank data, recovery data from surrogate spikes/laboratory control samples (LCS)/matrix spikes (MS), duplicate data, field quality assurance/quality control (QA/QC) samples, the laboratory review checklists (LRC), and the laboratory exception reports (ER).

A sample collection and analysis summary is presented in Table 1. This summary provides a cross-reference of field sample identification numbers and location identification. Each sample is assigned a unique field identification number.

The validated sample results are presented in Table 2. A summary of the analytical methodology is presented in Table 3.



## 2. Laboratory Qualifications

The Laboratory's quality assurance program is consistent with the quality standards outlined in the National Environmental Laboratory Accreditation Program (NELAP). This laboratory was accredited under Texas Certification number # TX104704231 at the time the analysis was performed and the certificate is included in Attachment A.

## 3. Project Objectives

### 3.1 Sampling/Analytical QA/QC Objectives

The QA/QC program was designed to identify contamination resulting from the sampling, sample transport and analytical process through the analysis of trip blank samples, field duplicate sample sets, and method blanks. The QA/QC program was designed to evaluate the quality of the resulting data with respect to bias and precision through analysis of LCS, MS, and duplicate analyses.

## 4. Data Review/Validation Results

### 4.1 Sample Holding Time and Preservation

Samples were shipped with a chain of custody and the paper work was filled out properly. All samples were properly preserved, delivered on ice, and stored by the laboratory at the required temperature (0-6°C).

The sample chain of custody documents and the analytical reports were used to determine sample holding times. All samples were prepared and analyzed within the required holding times.

### 4.2 Sample Containers

Sample containers used were certified pre-cleaned glass and plastic containers provided by the laboratory. These containers meet or exceed analyte specifications established in the United States Environmental Protection Agency (USEPA) *Specifications and Guidance for Contaminant-free Sample Containers*.

### 4.3 Calibrations

According to the LRC, initial calibration and continuing calibration data met the criteria for the selected method.

### 4.4 Laboratory Method Blank Analyses

Method blanks are prepared from a purified matrix and analyzed with investigative samples to determine the existence and magnitude of sample contamination introduced during the analytical procedures. As these were not discrete samples handled in the field, these blanks are not listed on the sample identification cross-reference list found in the data packages.





For this study, laboratory method blanks were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch and results are reported in the laboratory data packages.

The method blank results were non-detect or below the method quantitation limit (MQL), indicating that laboratory contamination was not a factor for this investigation.

#### **4.5 Internal Standard and Surrogate Spike Recoveries**

Recoveries of internal standards are addressed in the LRC of the data packages. All internal standard recoveries associated with the compounds of interest were acceptable per the LRC.

In accordance with the methods employed, all samples, blanks, and QC samples analyzed for volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), and total petroleum hydrocarbons (TPH) are spiked with surrogate compounds prior to sample analysis. Surrogate recoveries provide a means to evaluate the effects of laboratory performance on individual sample matrices. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project. Each individual surrogate compound is expected to meet the laboratory control limits. According to the TRRP-13 Guidelines, one outlying surrogate is acceptable for methods with multiple surrogate spike compounds as long as the recovery is at least ten percent. Samples analyzed at elevated sample dilutions (five times or greater) were not assessed.

Surrogate recoveries were assessed against laboratory control limits and/or the guidance in TRRP-13. All surrogate recoveries met the above criteria.

#### **4.6 Laboratory Control Sample Analysis**

LCS or LCS/laboratory control sample duplicate (LCSD) are prepared and analyzed as samples to assess the analytical efficiencies of the methods employed, independent of sample matrix effects. The relative percent difference (RPD) of the LCS/LCSD recoveries is used to evaluate analytical precision. The recovery ranges established by the laboratory are adopted as the acceptance criteria for the project.

For this study, LCS or LCS/LCSD were analyzed at a minimum frequency of one per twenty investigative samples and/or one per analytical batch.

The LCS or LCS/LCSD contained all analytes specified in the methods. All LCS recoveries and/or RPDs were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision (where applicable) with the following exception:

- i) One LCS/LCSD was reported with an elevated RPD for TPH (C6-C12). The associated sample result was non-detect and would not be affected by the indicated variability. No further action was required.

#### **4.7 Matrix Spike Analysis**

To evaluate the effects of sample matrices on the preparation process, measurement procedures, and accuracy of a particular analysis, samples are spiked with known concentrations of the analytes of interest



and analyzed as MS/matrix spike duplicate (MSD) samples. The RPD between the MS and MSD is used to assess analytical precision.

MS/MSD analyses were performed as specified in Table 1. The recovery ranges established by the laboratory is adopted as the acceptance criteria for the project.

The MS/MSD samples were spiked with all analytes specified in the methods. All percent recoveries and the RPD value were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision.

The laboratory also performed additional MS/MSD analyses on non-site samples. These cannot be used to assess accuracy and precision for the site samples.

#### **4.8 Duplicate Sample Analyses**

Analytical precision is evaluated based on the analysis of laboratory duplicate samples. For this study, duplicate samples were prepared and analyzed by the laboratory for arsenic analysis. The RPD established by the laboratory are adopted as the acceptance criteria for the project.

The laboratory performed duplicate analyses on non-site samples. These cannot be used to assess precision for the site samples.

#### **4.9 Field QA/QC Samples**

The field QA/QC consisted of three trip blank sample and two field duplicate sample sets.

##### ***Trip Blank Sample Analysis***

To evaluate contamination from sample collection, transportation, storage, and analytical activities, one trip blank sample was submitted to the laboratory for VOCs analysis. All results were non-detect for the compounds of interest.

##### ***Field Duplicate Sample Analysis***

To assess the analytical and sampling protocol precision, two field duplicate sample sets were collected and submitted "blind" to the laboratory, as specified in Table 1. The RPDs associated with these duplicate samples must be less than thirty percent for water samples. The RPDs are only used when sample concentrations are above the estimated regions of detection.

Field duplicate summary data are presented in Table 2. All field duplicate results were within acceptable agreement, demonstrating acceptable sampling and analytical precision with the following exceptions (see Table 4):

- i) WG-1620-MW70C-20200312 and WG-1620-FD01-20200312 did show some variability in SVOCs results and were qualified as estimated.



#### **4.10 Field Procedures**

Golder Associates, Inc. collected groundwater samples in accordance with their Standard Operating Procedures (SOP) for sample collection.

#### **4.11 Analyte Reporting**

The laboratory reported detected results for each analyte down to the sample detection limit (SDL), which is defined as the method detection limit (MDL) with sample-specific adjustments for dilutions, aliquot size, volumes, etc. Positive analyte detections less than the MQL but greater than the SDL were qualified as estimated (J) in Table 2 unless qualified otherwise in this memorandum.

The detectability check standard (DCS) results supported the laboratory MDLs.

### **5. Conclusion**

Based on the assessment detailed in the foregoing, the data summarized in Table 2 are usable for the purpose of supporting the HWPW - Groundwater Monitoring (New Wells) at the site by providing current concentration of chemicals of concern with the specific qualifications noted herein.

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-MW54B-20200311	MW-54B	Water	03/11/2020	14:20	X	X		X	
WQ-1620-MW99C-20200311	MW-99C	Water	03/11/2020	15:35	X	X		X	MS/MSD-P
WG-1620-MW91A-20200312	MW-91A	Water	03/12/2020	09:20	X	X		X	
WG-1620-MW70C-20200312	MW-70C	Water	03/12/2020	10:35	X	X		X	
WG-1620-FD01-20200312	MW-70C	Water	03/12/2020	10:35	X	X		X	Field duplicate of MW70C
WG-1620-MW76B-20200312	MW-76B	Water	03/12/2020	10:35	X	X	X	X	
WG-1620-MW97A-20200312	MW-97A	Water	03/12/2020	11:47	X	X	X	X	MS/MSD-P
WG-1620-MW84A-20200312	MW-84A	Water	03/12/2020	12:15	X	X		X	MS/MSD-P
WG-1620-MW98A-20200312	MW-98A	Water	03/12/2020	12:54	X	X	X	X	
WG-1620-MW92B-20200312	MW-92B	Water	03/12/2020	13:20	X	X		X	
WG-1620-MW98B-20200312	MW-98B	Water	03/12/2020	13:53	X	X	X	X	
WG-1620-MW93B-20200312	MW-93B	Water	03/12/2020	14:35	X	X		X	
WG-1620-MW50B-20200312	MW-50B	Water	03/12/2020	15:20	X	X	X	X	
WG-1620-MW94A-20200312	MW-94A	Water	03/12/2020	15:45	X	X		X	
WQ-1620-TB01-20200312	-	Water	03/12/2020	-	X				Trip Blank
WQ-1620-TB02-20200312	-	Water	03/12/2020	-	X				Trip Blank
WG-1620-MW95A-20200317	MW-95A	Water	03/17/2020	09:10	X	X		X	
WG-1620-MW96B-20200317	MW-96B	Water	03/17/2020	10:25	X	X		X	

Table 1

**Sample Collection and Analysis Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	Analysis/Parameters				Comments
					VOCs	SVOCs	TPH	Arsenic	
WG-1620-FD02-20200317	MW-96B	Water	03/17/2020	10:25	X	X		X	Field duplicate of MW96B
WG-1620-MW61B-20200317	MW61B	Water	03/17/2020	11:45	X	X		X	
WG-1620-MW88A-20200317	MW-88A	Water	03/17/2020	12:55	X	X		X	MS/MSD-P
WG-1620-MW88B-20200317	MW-88B	Water	03/17/2020	13:50	X	X		X	
WG-1620-MW60B-20200317	MW60B	Water	03/17/2020	15:15	X	X		X	
WQ-1620-TB01-20200317	-	Water	03/17/2020	-	X				Trip Blank
WG-1620-MW60AR-20200320	MW-60AR	Water	03/20/2020	08:55	X	X		X	
WG-1620-MW47A-20200320	MW-47A	Water	03/20/2020	10:30	X	X		X	
WQ-1620-TB01-20200320	-	Water	03/20/2020	-	X				Trip Blank

## Notes:

- VOCs - Volatile Organic Compounds
- SVOCs - Semi-volatile Organic Compounds
- TPH - Total Petroleum Hydrocarbons
- MS/MSD-P - Matrix Spike/ Matrix Spike Duplicate (partial parameters)
- "-" - Not Applicable

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-47A</b>	<b>MW-50B</b>	<b>MW-54B</b>	<b>MW-60AR</b>
	<b>Sample Name:</b>	<b>WG-1620-MW47A-20200320</b>	<b>WG-1620-MW50B-20200312</b>	<b>WG-1620-MW54B-20200311</b>	<b>WG-1620-MW60AR-20200320</b>
	<b>Sample Date:</b>	<b>03/20/2020</b>	<b>03/12/2020</b>	<b>03/11/2020</b>	<b>03/20/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	0.00043 J	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	0.0068	0.0011	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	0.0033	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	0.0056	0.0018	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000022	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000042	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000060	<0.000059	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000044	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000022	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	0.016	0.00011	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000021	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000049	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000027	0.020	0.0093	<0.000027
Acenaphthylene	mg/L	<0.000015	0.00028	0.000087 J	<0.000015
Anthracene	mg/L	<0.000014	0.0016	0.00017	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000052	<0.000051	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000021	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000031	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00011 J	<0.000039	0.000089 J	0.000063 J
Chrysene	mg/L	<0.000021	<0.000022	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.000076 J	0.000038 J	0.000033 J	<0.000020
Dibenzofuran	mg/L	<0.000020	0.015	0.000076 J	<0.000020

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-47A</b>	<b>MW-50B</b>	<b>MW-54B</b>	<b>MW-60AR</b>
	<b>Sample Name:</b>	<b>WG-1620-MW47A-20200320</b>	<b>WG-1620-MW50B-20200312</b>	<b>WG-1620-MW54B-20200311</b>	<b>WG-1620-MW60AR-20200320</b>
	<b>Sample Date:</b>	<b>03/20/2020</b>	<b>03/12/2020</b>	<b>03/11/2020</b>	<b>03/20/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Fluoranthene	mg/L	<0.000010	0.0011	0.000022 J	<0.000010
Fluorene	mg/L	<0.000030	0.011	0.00015	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000026	<0.000025	<0.000025
Naphthalene	mg/L	<0.000020	0.31	0.0021	<0.000020
Nitrobenzene	mg/L	<0.000024	<0.000025	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000082	<0.000080	<0.000079
Phenanthrene	mg/L	<0.000021	0.016	0.0028	<0.000021
Phenol	mg/L	<0.000035	<0.000036	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	0.00059	<0.000019	<0.000019
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	--	2.10	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	1.1	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	<0.19	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	1.0	--	--
<b>Metals</b>					
Arsenic	mg/L	0.000566 J	0.00372	0.00117 J	0.00189 J

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-70C</b>	<b>MW-70C</b>	<b>MW-76B</b>	<b>MW-84A</b>
	<b>Sample Name:</b>	<b>WG-1620-MW70C-20200312</b>	<b>WG-1620-FD01-20200312</b>	<b>WG-1620-MW76B-20200312</b>	<b>WG-1620-MW84A-20200312</b>
	<b>Sample Date:</b>	<b>03/12/2020</b>	<b>03/12/2020</b> <b>Duplicate</b>	<b>03/12/2020</b>	<b>03/12/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	0.010	0.0089	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	0.057	0.052	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.034	0.032	<0.00020	<0.00020
Xylenes (total)	mg/L	0.064	0.061	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000022	<0.000022	<0.000022	<0.000021
2,4-Dimethylphenol	mg/L	0.00041	<0.000041	<0.000042	<0.000041
2,4-Dinitrotoluene	mg/L	<0.000060	<0.000060	<0.000060	<0.000059
2,6-Dinitrotoluene	mg/L	<0.000044	<0.000043	<0.000044	<0.000043
2-Chloronaphthalene	mg/L	<0.000022	<0.000022	<0.000022	<0.000021
2-Methylnaphthalene	mg/L	0.067 J	0.098 J	<0.000020	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000021	<0.000021	<0.000021	<0.000020
4-Nitrophenol	mg/L	<0.000049	<0.000048	<0.000049	<0.000048
Acenaphthene	mg/L	0.059 J	0.087 J	<0.000028	<0.000028
Acenaphthylene	mg/L	<0.000016	0.00074	<0.000016	<0.000015
Anthracene	mg/L	0.0049 J	0.0082 J	<0.000015	<0.000014
Benzo(a)anthracene	mg/L	<0.000052	0.000079 J	<0.000052	<0.000051
Benzo(a)pyrene	mg/L	<0.000021	<0.000021	<0.000021	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000031	<0.000031	<0.000031	<0.000031
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000039	0.000052 J	<0.000039	<0.000038
Chrysene	mg/L	<0.000022	0.000061 J	<0.000022	<0.000021
Di-n-butylphthalate (DBP)	mg/L	0.000061 J	0.00010 J	0.000025 J	<0.000020
Dibenzofuran	mg/L	0.055 J	0.078 J	<0.000021	<0.000020



Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-70C</b>	<b>MW-70C</b>	<b>MW-76B</b>	<b>MW-84A</b>
	<b>Sample Name:</b>	<b>WG-1620-MW70C-20200312</b>	<b>WG-1620-FD01-20200312</b>	<b>WG-1620-MW76B-20200312</b>	<b>WG-1620-MW84A-20200312</b>
	<b>Sample Date:</b>	<b>03/12/2020</b>	<b>03/12/2020</b> <b>Duplicate</b>	<b>03/12/2020</b>	<b>03/12/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Fluoranthene	mg/L	0.0030 J	0.0053 J	<0.000010	<0.000010
Fluorene	mg/L	0.028 J	0.038 J	<0.000031	<0.000031
N-Nitrosodiphenylamine	mg/L	<0.000026	<0.000026	<0.000026	<0.000026
Naphthalene	mg/L	1.6 J	2.5 J	0.000096 J	0.000023 J
Nitrobenzene	mg/L	<0.000025	<0.000025	<0.000025	<0.000024
Pentachlorophenol	mg/L	<0.000082	<0.000081	<0.000082	<0.000081
Phenanthrene	mg/L	0.040	0.054	<0.000022	<0.000021
Phenol	mg/L	<0.000036	<0.000036	<0.000036	<0.000036
Pyrene	mg/L	0.0017 J	0.0029 J	<0.000020	<0.000019
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	--	--	<0.19	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	<0.19	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	<0.19	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	<0.19	--
<b>Metals</b>					
Arsenic	mg/L	0.00598	0.00470	0.00115 J	0.00464

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-88A</b>	<b>MW-88B</b>	<b>MW-91A</b>	<b>MW-92B</b>
	<b>Sample Name:</b>	<b>WG-1620-MW88A-20200317</b>	<b>WG-1620-MW88B-20200317</b>	<b>WG-1620-MW91A-20200312</b>	<b>WG-1620-MW92B-20200312</b>
	<b>Sample Date:</b>	<b>03/17/2020</b>	<b>03/17/2020</b>	<b>03/12/2020</b>	<b>03/12/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	<0.00020	0.00049 J
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000022	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	<0.000042	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000058	<0.000061	<0.000059
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000044	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000022	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	<0.000020	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000021	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000049	<0.000047
Acenaphthene	mg/L	0.0015	<0.000027	<0.000028	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000016	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000015	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000050	<0.000053	<0.000051
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000021	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000032	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00010 J	<0.000037	<0.000039	<0.000037
Chrysene	mg/L	<0.000021	<0.000021	<0.000022	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	0.000031 J	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	<0.000021	<0.000020

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-88A</b>	<b>MW-88B</b>	<b>MW-91A</b>	<b>MW-92B</b>
	<b>Sample Name:</b>	<b>WG-1620-MW88A-20200317</b>	<b>WG-1620-MW88B-20200317</b>	<b>WG-1620-MW91A-20200312</b>	<b>WG-1620-MW92B-20200312</b>
	<b>Sample Date:</b>	<b>03/17/2020</b>	<b>03/17/2020</b>	<b>03/12/2020</b>	<b>03/12/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Fluoranthene	mg/L	0.00025	<0.000010	<0.000011	<0.000010
Fluorene	mg/L	0.00026	<0.000030	<0.000032	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000026	<0.000025
Naphthalene	mg/L	<0.000020	<0.000020	0.000083 J	<0.000020
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000025	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000079	<0.000083	<0.000080
Phenanthrene	mg/L	<0.000021	<0.000021	<0.000022	<0.000021
Phenol	mg/L	<0.000035	<0.000035	<0.000037	<0.000035
Pyrene	mg/L	0.00053	<0.000019	<0.000020	<0.000019
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--
<b>Metals</b>					
Arsenic	mg/L	0.00355	0.00166 J	0.00989	0.00201

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	Location ID:	MW-93B	MW-94A	MW-95A	MW-96B
	Sample Name:	WG-1620-MW93B-20200312	WG-1620-MW94A-20200312	WG-1620-MW95A-20200317	WG-1620-MW96B-20200317
	Sample Date:	03/12/2020	03/12/2020	03/17/2020	03/17/2020
Parameters	Unit				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.00092 J	<0.00020	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000022	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000041	<0.000041	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000059	<0.000060	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000043	<0.000043	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000022	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000020	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000021	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000048	<0.000048	<0.000047	<0.000047
Acenaphthene	mg/L	0.000039 J	<0.000028	<0.000027	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000051	<0.000052	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000021	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000031	<0.000031	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	<0.000038	0.000056 J	<0.000037	0.00035
Chrysene	mg/L	<0.000021	<0.000022	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000021	<0.000020	<0.000020
Dibenzofuran	mg/L	0.000071 J	<0.000021	<0.000020	<0.000020

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-93B</b>	<b>MW-94A</b>	<b>MW-95A</b>	<b>MW-96B</b>
	<b>Sample Name:</b>	<b>WG-1620-MW93B-20200312</b>	<b>WG-1620-MW94A-20200312</b>	<b>WG-1620-MW95A-20200317</b>	<b>WG-1620-MW96B-20200317</b>
	<b>Sample Date:</b>	<b>03/12/2020</b>	<b>03/12/2020</b>	<b>03/17/2020</b>	<b>03/17/2020</b>
<b>Parameters</b>	<b>Unit</b>				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Fluoranthene	mg/L	<0.000010	<0.000010	<0.000010	<0.000010
Fluorene	mg/L	0.000058 J	<0.000031	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000026	<0.000026	<0.000025	<0.000025
Naphthalene	mg/L	<0.000020	<0.000021	<0.000020	<0.000020
Nitrobenzene	mg/L	<0.000024	<0.000025	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000081	<0.000081	<0.000079	<0.000079
Phenanthrene	mg/L	0.00016	<0.000022	<0.000021	<0.000021
Phenol	mg/L	<0.000036	<0.000036	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000020	<0.000019	<0.000019
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	--	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--	--
<b>Metals</b>					
Arsenic	mg/L	0.00455	0.00540	0.000977 J	0.00312

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-96B</b>	<b>MW-97A</b>	<b>MW-98A</b>	<b>MW-98B</b>
	<b>Sample Name:</b>	<b>WG-1620-FD02-20200317</b>	<b>WG-1620-MW97A-20200312</b>	<b>WG-1620-MW98A-20200312</b>	<b>WG-1620-MW98B-20200312</b>
	<b>Sample Date:</b>	<b>03/17/2020</b>	<b>03/12/2020</b>	<b>03/12/2020</b>	<b>03/12/2020</b>
		<b>Duplicate</b>			
<b>Parameters</b>	<b>Unit</b>				
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010	<0.0010
Toluene	mg/L	<0.00020	<0.00020	0.00065 J	0.0020
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>					
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000040	<0.000040	0.000078 J	<0.000041
2,4-Dinitrotoluene	mg/L	<0.000058	<0.000059	<0.000059	<0.000059
2,6-Dinitrotoluene	mg/L	<0.000042	<0.000042	<0.000043	<0.000043
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	<0.000019	<0.000019	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000047	<0.000047	<0.000048	<0.000048
Acenaphthene	mg/L	<0.000027	<0.000027	0.000040 J	<0.000028
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000050	<0.000051	<0.000051	<0.000051
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000030	<0.000030	<0.000031	<0.000031
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.00013 J	<0.000037	<0.000038	<0.000038
Chrysene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	<0.000020	<0.000020	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	<0.000020	<0.000020

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-96B</b>	<b>MW-97A</b>	<b>MW-98A</b>	<b>MW-98B</b>
	<b>Sample Name:</b>	<b>WG-1620-FD02-20200317</b>	<b>WG-1620-MW97A-20200312</b>	<b>WG-1620-MW98A-20200312</b>	<b>WG-1620-MW98B-20200312</b>
	<b>Sample Date:</b>	<b>03/17/2020</b>	<b>03/12/2020</b>	<b>03/12/2020</b>	<b>03/12/2020</b>
		<b>Duplicate</b>			
<b>Parameters</b>	<b>Unit</b>				
<b>Semi-volatile Organic Compounds (Continued)</b>					
Fluoranthene	mg/L	<0.000010	<0.000010	0.000018 J	<0.000010
Fluorene	mg/L	<0.000030	<0.000030	<0.000031	<0.000031
N-Nitrosodiphenylamine	mg/L	<0.000025	<0.000025	<0.000026	<0.000026
Naphthalene	mg/L	<0.000020	<0.000020	0.000042 J	0.000036 J
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000079	<0.000080	<0.000081	<0.000081
Phenanthrene	mg/L	<0.000021	<0.000021	<0.000021	<0.000021
Phenol	mg/L	<0.000035	<0.000035	<0.000036	<0.000036
Pyrene	mg/L	<0.000019	<0.000019	<0.000019	<0.000019
<b>Total Petroleum Hydrocarbons</b>					
Total Petroleum Hydrocarbons	mg/L	--	<0.19	1.00	0.380 J
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	<0.19	1.0	0.38 J
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	<0.19	<0.19	<0.19
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	<0.19	<0.19	<0.19
<b>Metals</b>					
Arsenic	mg/L	0.00311	0.00121 J	0.00963	0.00165 J

Table 2

**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	Location ID:	MW-99C	MW60B	MW61B
	Sample Name:	WQ-1620-MW99C-20200311	WG-1620-MW60B-20200317	WG-1620-MW61B-20200317
	Sample Date:	03/11/2020	03/17/2020	03/17/2020
Parameters	Unit			
<b>Volatile Organic Compounds</b>				
1,2-Dichloroethane	mg/L	<0.00020	<0.00020	<0.00020
Benzene	mg/L	<0.00020	<0.00020	<0.00020
Chlorobenzene	mg/L	<0.00030	<0.00030	<0.00030
Ethylbenzene	mg/L	<0.00030	<0.00030	<0.00030
Methylene chloride	mg/L	<0.0010	<0.0010	<0.0010
Toluene	mg/L	0.00078 J	<0.00020	<0.00020
Xylenes (total)	mg/L	<0.00030	<0.00030	<0.00030
<b>Semi-volatile Organic Compounds</b>				
1,2-Diphenylhydrazine	mg/L	<0.000021	<0.000021	<0.000021
2,4-Dimethylphenol	mg/L	<0.000041	<0.000040	<0.000040
2,4-Dinitrotoluene	mg/L	<0.000059	<0.000058	<0.000058
2,6-Dinitrotoluene	mg/L	<0.000043	<0.000042	<0.000042
2-Chloronaphthalene	mg/L	<0.000021	<0.000021	<0.000021
2-Methylnaphthalene	mg/L	0.000086 J	<0.000019	<0.000019
4,6-Dinitro-2-methylphenol	mg/L	<0.000020	<0.000020	<0.000020
4-Nitrophenol	mg/L	<0.000048	<0.000047	<0.000047
Acenaphthene	mg/L	<0.000028	<0.000027	<0.000027
Acenaphthylene	mg/L	<0.000015	<0.000015	<0.000015
Anthracene	mg/L	<0.000014	<0.000014	<0.000014
Benzo(a)anthracene	mg/L	<0.000051	<0.000050	<0.000050
Benzo(a)pyrene	mg/L	<0.000020	<0.000020	<0.000020
bis(2-Chloroethoxy)methane	mg/L	<0.000031	<0.000030	<0.000030
bis(2-Ethylhexyl)phthalate (DEHP)	mg/L	0.000070 J	0.000092	0.00015 J
Chrysene	mg/L	<0.000021	<0.000021	<0.000021
Di-n-butylphthalate (DBP)	mg/L	<0.000020	0.000047 J	<0.000020
Dibenzofuran	mg/L	<0.000020	<0.000020	<0.000020



**Table 2**  
**Analytical Results Summary**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

	<b>Location ID:</b>	<b>MW-99C</b>	<b>MW60B</b>	<b>MW61B</b>
	<b>Sample Name:</b>	<b>WQ-1620-MW99C-20200311</b>	<b>WG-1620-MW60B-20200317</b>	<b>WG-1620-MW61B-20200317</b>
	<b>Sample Date:</b>	<b>03/11/2020</b>	<b>03/17/2020</b>	<b>03/17/2020</b>
<b>Parameters</b>	<b>Unit</b>			
<b>Semi-volatile Organic Compounds (Continued)</b>				
Fluoranthene	mg/L	<0.000010	0.000015 J	<0.000010
Fluorene	mg/L	<0.000031	<0.000030	<0.000030
N-Nitrosodiphenylamine	mg/L	<0.000026	<0.000025	<0.000025
Naphthalene	mg/L	0.00018	<0.000020	<0.000020
Nitrobenzene	mg/L	<0.000024	<0.000024	<0.000024
Pentachlorophenol	mg/L	<0.000081	<0.000079	<0.000079
Phenanthrene	mg/L	0.000053 J	<0.000021	<0.000021
Phenol	mg/L	<0.000036	<0.000035	<0.000035
Pyrene	mg/L	<0.000019	<0.000019	<0.000019
<b>Total Petroleum Hydrocarbons</b>				
Total Petroleum Hydrocarbons	mg/L	--	--	--
Total Petroleum Hydrocarbons (>C12-C28)	mg/L	--	--	--
Total Petroleum Hydrocarbons (>C28-C35)	mg/L	--	--	--
Total Petroleum Hydrocarbons (C6-C12)	mg/L	--	--	--
<b>Metals</b>				
Arsenic	mg/L	0.000866 J	0.00234	0.00461

## Notes:

&lt; - Not detected at the associated reporting limit

J - Estimated concentration

"--" - Not analyzed

Table 3

**Analytical Methods**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

Parameter	Method	Matrix	Holding Time	
			Collection to Extraction (Days)	Extraction to Analysis (Days)
VOCs	SW-846 8260C	Water	-	14
SVOCs	SW-846 8270D	Water	7	40
TPH	TX1005	Water	14	40
Arsenic	SW-846 6020A	Water	-	180

## Notes:

- VOCs - Volatile Organic Compounds  
SVOCs - Semi-volatile Organic Compounds  
TPH - Total Petroleum Hydrocarbons  
"-" - Not Applicable

## Method References:

- SW-846 - "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition, 1986, with subsequent revisions

Table 4

**Qualified Sample Data Due to Variability in Field Duplicate Results**  
**HWPW - Groundwater Monitoring (New Wells)**  
**Union Pacific Railroad (UPRR)/Houston, TX-Wood Preserving Works**  
**Houston, Texas**  
**March 2020**

Parameter	Analyte	RPD	Sample ID	Qualified Result	Field Duplicate Sample ID	Qualified Result	Units
SVOCs	2-Methylnaphthalene	37.6	WG-1620-MW70C-20200312	0.067 J	WG-1620-FD01-20200312	0.098 J	mg/L
	Acenaphthene	38.4		0.059 J		0.087 J	mg/L
	Anthracene	50.4		0.0049 J		0.0082 J	mg/L
	Dibenzofuran	34.6		0.055 J		0.078 J	mg/L
	Di-n-butylphthalate (DBP)	48.4		0.000061 J		0.00010 J	mg/L
	Fluoranthene	55.4		0.0030 J		0.0053 J	mg/L
	Fluorene	30.3		0.028 J		0.038 J	mg/L
	Naphthalene	43.9		1.6 J		2.5 J	mg/L
	Pyrene	52.2		0.0017 J		0.0029 J	mg/L

## Notes:

- RPD - Relative Percent Difference  
SVOCs - Semi-volatile Organic Compounds  
J - Estimated concentration

# Attachment A

## Laboratory NELAP Certificate



# Texas Commission on Environmental Quality

## NELAP - Recognized Laboratory Fields of Accreditation



ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
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Issue Date: 5/1/2019

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**Matrix: *Drinking Water***

**Method** EPA 1613

Analyte	AB	Analyte ID	Method ID
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408

**Method** EPA 200.8

Analyte	AB	Analyte ID	Method ID
Copper	TX	1055	10014605
Lead	TX	1075	10014605



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**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 120.1			
Analyte Conductivity	TX	1610	10006403
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 160.4			
Analyte Residue-volatile	TX	1970	10010409
Method EPA 1613			
Analyte 1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10120408
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10120408
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10120408
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10120408
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10120408
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10120408
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10120408
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10120408
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,6,7,8-HxCDD)	TX	9456	10120408
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10120408
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10120408
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10120408
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10120408
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10120408



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**Matrix: Non-Potable Water**

2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10120408
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10120408
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10120408
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10120408
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10120408
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10120408
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10120408
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10120408
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10120408
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10120408
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10120408
<b>Method EPA 1664</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10127807
<b>Method EPA 180.1</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Turbidity	TX	2055	10011606
<b>Method EPA 200.8</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Aluminum	TX	1000	10014605
Antimony	TX	1005	10014605
Arsenic	TX	1010	10014605
Barium	TX	1015	10014605
Beryllium	TX	1020	10014605
Boron	TX	1025	10014605
Cadmium	TX	1030	10014605
Calcium	TX	1035	10014605
Chromium	TX	1040	10014605
Cobalt	TX	1050	10014605
Copper	TX	1055	10014605
Iron	TX	1070	10014605



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**Matrix: Non-Potable Water**

Lead	TX	1075	10014605
Magnesium	TX	1085	10014605
Manganese	TX	1090	10014605
Molybdenum	TX	1100	10014605
Nickel	TX	1105	10014605
Potassium	TX	1125	10014605
Selenium	TX	1140	10014605
Silver	TX	1150	10014605
Sodium	TX	1155	10014605
Strontium	TX	1160	10014605
Thallium	TX	1165	10014605
Tin	TX	1175	10014605
Titanium	TX	1180	10014605
Uranium	TX	3035	10014605
Vanadium	TX	1185	10014605
Zinc	TX	1190	10014605

**Method EPA 245.1**

Analyte	AB	Analyte ID	Method ID
Mercury	TX	1095	10036609

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 325.1**

Analyte	AB	Analyte ID	Method ID
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**Matrix: Non-Potable Water**

Chloride	TX	1575	10056801
<b>Method</b> EPA 335.1			
<b>Analyte</b> Amenable cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1510	<b>Method ID</b> 10060001
<b>Method</b> EPA 335.2			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10278203
<b>Method</b> EPA 335.4			
<b>Analyte</b> Total cyanide	<b>AB</b> TX	<b>Analyte ID</b> 1645	<b>Method ID</b> 10061402
<b>Method</b> EPA 350.3			
<b>Analyte</b> Ammonia as N	<b>AB</b> TX	<b>Analyte ID</b> 1515	<b>Method ID</b> 10064401
<b>Method</b> EPA 365.3			
<b>Analyte</b> Orthophosphate as P Phosphorus	<b>AB</b> TX TX	<b>Analyte ID</b> 1870 1910	<b>Method ID</b> 10070801 10070801
<b>Method</b> EPA 375.4			
<b>Analyte</b> Sulfate	<b>AB</b> TX	<b>Analyte ID</b> 2000	<b>Method ID</b> 10073800
<b>Method</b> EPA 376.1			
<b>Analyte</b> Sulfide	<b>AB</b> TX	<b>Analyte ID</b> 2005	<b>Method ID</b> 10074201
<b>Method</b> EPA 410.4			
<b>Analyte</b> Chemical oxygen demand (COD)	<b>AB</b> TX	<b>Analyte ID</b> 1565	<b>Method ID</b> 10077404
<b>Method</b> EPA 415.1			
<b>Analyte</b> Total Organic Carbon (TOC)	<b>AB</b> TX	<b>Analyte ID</b> 2040	<b>Method ID</b> 10078407
<b>Method</b> EPA 420.1			
<b>Analyte</b> Total phenolics	<b>AB</b> TX	<b>Analyte ID</b> 1905	<b>Method ID</b> 10079400



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**Matrix: Non-Potable Water**

**Method EPA 420.4**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10080203

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156419
Antimony	TX	1005	10156419
Arsenic	TX	1010	10156419
Barium	TX	1015	10156419
Beryllium	TX	1020	10156419
Boron	TX	1025	10156419
Cadmium	TX	1030	10156419
Calcium	TX	1035	10156419
Chromium	TX	1040	10156419
Cobalt	TX	1050	10156419
Copper	TX	1055	10156419
Iron	TX	1070	10156419
Lead	TX	1075	10156419
Lithium	TX	1080	10156419
Magnesium	TX	1085	10156419
Manganese	TX	1090	10156419
Molybdenum	TX	1100	10156419
Nickel	TX	1105	10156419
Potassium	TX	1125	10156419
Selenium	TX	1140	10156419
Silver	TX	1150	10156419
Sodium	TX	1155	10156419
Strontium	TX	1160	10156419
Thallium	TX	1165	10156419
Tin	TX	1175	10156419
Titanium	TX	1180	10156419



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**Matrix: Non-Potable Water**

Vanadium	TX	1185	10156419
Zinc	TX	1190	10156419
<b>Method EPA 608</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
4,4'-DDD	TX	7355	10103603
4,4'-DDE	TX	7360	10103603
4,4'-DDT	TX	7365	10103603
Aldrin	TX	7025	10103603
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10103603
alpha-Chlordane	TX	7240	10103603
Aroclor-1016 (PCB-1016)	TX	8880	10103603
Aroclor-1221 (PCB-1221)	TX	8885	10103603
Aroclor-1232 (PCB-1232)	TX	8890	10103603
Aroclor-1242 (PCB-1242)	TX	8895	10103603
Aroclor-1248 (PCB-1248)	TX	8900	10103603
Aroclor-1254 (PCB-1254)	TX	8905	10103603
Aroclor-1260 (PCB-1260)	TX	8910	10103603
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10103603
Chlordane (tech.)	TX	7250	10103603
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10103603
Dieldrin	TX	7470	10103603
Endosulfan I	TX	7510	10103603
Endosulfan II	TX	7515	10103603
Endosulfan sulfate	TX	7520	10103603
Endrin	TX	7540	10103603
Endrin aldehyde	TX	7530	10103603
Endrin ketone	TX	7535	10103603
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10103603
gamma-Chlordane	TX	7245	10103603
Heptachlor	TX	7685	10103603



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**Matrix: Non-Potable Water**

Heptachlor epoxide	TX	7690	10103603
Methoxychlor	TX	7810	10103603
Toxaphene (Chlorinated camphene)	TX	8250	10103603

**Method EPA 624**

Analyte	AB	Analyte ID	Method ID
1,1,1-Trichloroethane	TX	5160	10107207
1,1,2,2-Tetrachloroethane	TX	5110	10107207
1,1,2-Trichloroethane	TX	5165	10107207
1,1-Dichloroethane	TX	4630	10107207
1,1-Dichloroethylene	TX	4640	10107207
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10107207
1,2-Dichlorobenzene	TX	4610	10107207
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10107207
1,2-Dichloropropane	TX	4655	10107207
1,3-Dichlorobenzene	TX	4615	10107207
1,4-Dichlorobenzene	TX	4620	10107207
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10107207
2-Chloroethyl vinyl ether	TX	4500	10107207
Acetone (2-Propanone)	TX	4315	10107207
Acrolein (Propenal)	TX	4325	10107207
Acrylonitrile	TX	4340	10107207
Benzene	TX	4375	10107207
Bromodichloromethane	TX	4395	10107207
Bromoform	TX	4400	10107207
Carbon tetrachloride	TX	4455	10107207
Chlorobenzene	TX	4475	10107207
Chlorodibromomethane	TX	4575	10107207
Chloroethane (Ethyl chloride)	TX	4485	10107207
Chloroform	TX	4505	10107207
cis-1,2-Dichloroethylene	TX	4645	10107207



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: Non-Potable Water**

cis-1,3-Dichloropropene	TX	4680	10107207
Ethylbenzene	TX	4765	10107207
m+p-xylene	TX	5240	10107207
Methyl bromide (Bromomethane)	TX	4950	10107207
Methyl chloride (Chloromethane)	TX	4960	10107207
Methyl tert-butyl ether (MTBE)	TX	5000	10107207
Methylene chloride (Dichloromethane)	TX	4975	10107207
Naphthalene	TX	5005	10107207
o-Xylene	TX	5250	10107207
Tetrachloroethylene (Perchloroethylene)	TX	5115	10107207
Toluene	TX	5140	10107207
trans-1,2-Dichloroethylene	TX	4700	10107207
trans-1,3-Dichloropropylene	TX	4685	10107207
Trichloroethene (Trichloroethylene)	TX	5170	10107207
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10107207
Vinyl chloride	TX	5235	10107207
Xylene (total)	TX	5260	10107207

**Method EPA 625**

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10107401
1,2,4-Trichlorobenzene	TX	5155	10107401
1,2-Dichlorobenzene	TX	4610	10107401
1,2-Diphenylhydrazine	TX	6220	10107401
1,3-Dichlorobenzene	TX	4615	10107401
1,4-Dichlorobenzene	TX	4620	10107401
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10107401
2,4,5-Trichlorophenol	TX	6835	10107401
2,4,6-Trichlorophenol	TX	6840	10107401
2,4-Dichlorophenol	TX	6000	10107401
2,4-Dimethylphenol	TX	6130	10107401



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**Matrix: Non-Potable Water**

2,4-Dinitrophenol	TX	6175	10107401
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10107401
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10107401
2-Chloronaphthalene	TX	5795	10107401
2-Chlorophenol	TX	5800	10107401
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10107401
2-Methylphenol (o-Cresol)	TX	6400	10107401
2-Nitrophenol	TX	6490	10107401
3,3'-Dichlorobenzidine	TX	5945	10107401
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10107401
4-Chloro-3-methylphenol	TX	5700	10107401
4-Chlorophenyl phenylether	TX	5825	10107401
4-Methylphenol (p-Cresol)	TX	6410	10107401
4-Nitrophenol	TX	6500	10107401
Acenaphthene	TX	5500	10107401
Acenaphthylene	TX	5505	10107401
Anthracene	TX	5555	10107401
Benzidine	TX	5595	10107401
Benzo(a)anthracene	TX	5575	10107401
Benzo(a)pyrene	TX	5580	10107401
Benzo(b)fluoranthene	TX	5585	10107401
Benzo(g,h,i)perylene	TX	5590	10107401
Benzo(k)fluoranthene	TX	5600	10107401
bis(2-Chloroethoxy)methane	TX	5760	10107401
bis(2-Chloroethyl) ether	TX	5765	10107401
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10107401
Butyl benzyl phthalate	TX	5670	10107401
Chrysene	TX	5855	10107401
Dibenz(a,h) anthracene	TX	5895	10107401
Diethyl phthalate	TX	6070	10107401





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**Matrix: Non-Potable Water**

Dimethyl phthalate	TX	6135	10107401
Di-n-butyl phthalate	TX	5925	10107401
Di-n-octyl phthalate	TX	6200	10107401
Fluoranthene	TX	6265	10107401
Fluorene	TX	6270	10107401
Hexachlorobenzene	TX	6275	10107401
Hexachlorobutadiene	TX	4835	10107401
Hexachlorocyclopentadiene	TX	6285	10107401
Hexachloroethane	TX	4840	10107401
Indeno(1,2,3-cd) pyrene	TX	6315	10107401
Isophorone	TX	6320	10107401
Naphthalene	TX	5005	10107401
Nitrobenzene	TX	5015	10107401
n-Nitrosodiethylamine	TX	6525	10107401
n-Nitrosodimethylamine	TX	6530	10107401
n-Nitrosodi-n-butylamine	TX	5025	10107401
n-Nitrosodi-n-propylamine	TX	6545	10107401
n-Nitrosodiphenylamine	TX	6535	10107401
Pentachlorobenzene	TX	6590	10107401
Pentachlorophenol	TX	6605	10107401
Phenanthrene	TX	6615	10107401
Phenol	TX	6625	10107401
Pyrene	TX	6665	10107401
Pyridine	TX	5095	10107401
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603



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**Matrix: Non-Potable Water**

**Method EPA 8011**

Analyte	AB	Analyte ID	Method ID
1,2,3-Trichloropropane	TX	5180	10173009
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10173009
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10173009

**Method EPA 8015**

Analyte	AB	Analyte ID	Method ID
Diesel range organics (DRO)	TX	9369	10173203
Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402





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**Matrix: Non-Potable Water**

4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402
Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Hexachlorobenzene	TX	6275	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201



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**Matrix: Non-Potable Water**

Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8151**

Analyte	AB	Analyte ID	Method ID
2,4,5-T	TX	8655	10183003
2,4-D	TX	8545	10183003
2,4-DB	TX	8560	10183003
Dalapon	TX	8555	10183003
Dicamba	TX	8595	10183003
Dichloroprop (Dichloroprop, Weedone)	TX	8605	10183003
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10183003
MCPA	TX	7775	10183003
MCPP	TX	7780	10183003
Silvex (2,4,5-TP)	TX	8650	10183003

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404
1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404



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**Matrix: Non-Potable Water**

1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
2-Pentanone	TX	5045	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404
Allyl alcohol	TX	4350	10184404
Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404



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**Matrix: Non-Potable Water**

Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Di-isopropylether (DIPE)	TX	9375	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Ethyl-t-butylether (ETBE) (2-Ethoxy-2-methylpropane)	TX	4770	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404
Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404



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**Matrix: Non-Potable Water**

Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
T-amylmethylether (TAME)	TX	4370	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404
Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404



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**Matrix: Non-Potable Water**

Method EPA 8270

Analyte	AB	Analyte ID	Method ID
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203
2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203





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**Matrix: Non-Potable Water**

2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Dimethyl aminoazobenzene	TX	6105	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrobiphenyl	TX	6480	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Chloro-2-methylaniline	TX	5695	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203



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**Matrix: Non-Potable Water**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Captan	TX	7190	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203





# Texas Commission on Environmental Quality



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**Matrix: Non-Potable Water**

Chrysene	TX	5855	10185203
Coumaphos	TX	7315	10185203
Demeton	TX	7390	10185203
Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Dioxathion	TX	7495	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethion	TX	7565	10185203
Ethyl methanesulfonate	TX	6260	10185203
Famphur	TX	7580	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203



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**Matrix: Non-Potable Water**

Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203
Kepone	TX	7740	10185203
Maleic anhydride	TX	6335	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naled	TX	7905	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203



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**Matrix: Non-Potable Water**

Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Phosmet (Imidan)	TX	8000	10185203
Phthalic anhydride	TX	6640	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203
Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Resorcinol	TX	6680	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203
Trifluralin (Treflan)	TX	8295	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209



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**Matrix: Non-Potable Water**

1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209
Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209

**Method EPA 8316**

Analyte	AB	Analyte ID	Method ID
Acrylamide	TX	4330	10188202

**Method EPA 8330**

Analyte	AB	Analyte ID	Method ID
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807



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**Matrix: Non-Potable Water**

Methyl-2,4,6-trinitrophenylamine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807
<b>Method EPA 9014</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803
<b>Method EPA 9038</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Sulfate	TX	2000	10196608
<b>Method EPA 9040</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
pH	TX	1900	10196802
<b>Method EPA 9050</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	10198604
<b>Method EPA 9056</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209
<b>Method EPA 9060</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	10200201



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**Matrix: Non-Potable Water**

Method	Analyte	AB	Analyte ID	Method ID
EPA 9065	Total phenolics	TX	1905	10200405
EPA 9066	Total phenolics	TX	1905	10200609
EPA 9250	Chloride	TX	1575	10207202
EPA RSK 175	2-methylpropane (Isobutane)	TX	4942	10212905
	Ethane	TX	4747	10212905
	Ethene	TX	4752	10212905
	Methane	TX	4926	10212905
	n-Butane	TX	5007	10212905
	n-Propane	TX	5029	10212905
HACH 8000	Chemical oxygen demand (COD)	TX	1565	60003001
SM 2120 B	Color	TX	1605	20223807
SM 2310 B (4a)	Acidity, as CaCO <sub>3</sub>	TX	1500	20002806
SM 2320 B	Alkalinity as CaCO <sub>3</sub>	TX	1505	20045005
SM 2340 B		AB	Analyte ID	Method ID





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**Matrix: Non-Potable Water**

Total hardness as CaCO <sub>3</sub>	TX	1755	20046008
<b>Method</b> SM 2510 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Conductivity	TX	1610	20048004
<b>Method</b> SM 2540 B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-total (total solids)	TX	1950	20004608
<b>Method</b> SM 2540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-filterable (TDS)	TX	1955	20049803
<b>Method</b> SM 2540 D			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Residue-nonfilterable (TSS)	TX	1960	20004802
<b>Method</b> SM 3500-Cr B			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	20065809
<b>Method</b> SM 4500-Cl F			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total residual chlorine	TX	1940	20080482
<b>Method</b> SM 4500-Cl <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chloride	TX	1575	20019209
<b>Method</b> SM 4500-CN <sup>-</sup> C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20020808
<b>Method</b> SM 4500-CN <sup>-</sup> E			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total cyanide	TX	1645	20021209
<b>Method</b> SM 4500-CN <sup>-</sup> G			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Amenable cyanide	TX	1510	20021607



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**Matrix: Non-Potable Water**

Method	AB	Analyte ID	Method ID
Method SM 4500-H+ B			
Analyte	AB	Analyte ID	Method ID
pH	TX	1900	20104603
Method SM 4500-NH3 D			
Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	20108809
Kjeldahl Nitrogen (Total Kjeldahl Nitrogen-TKN)	TX	1790	20108809
Method SM 4500-NH3 F			
Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	20023001
Method SM 4500-O G			
Analyte	AB	Analyte ID	Method ID
Oxygen, dissolved	TX	1880	20025405
Method SM 4500-P E			
Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	20025803
Phosphorus	TX	1910	20025803
Method SM 4500-S2 <sup>-</sup> F			
Analyte	AB	Analyte ID	Method ID
Sulfide	TX	2005	20126209
Method SM 4500-SiO2 D			
Analyte	AB	Analyte ID	Method ID
Silica as SiO2	TX	1990	20127202
Method SM 4500-SO3 <sup>-</sup> B			
Analyte	AB	Analyte ID	Method ID
Sulfite	TX	2015	20026806
Method SM 5210 B			
Analyte	AB	Analyte ID	Method ID
Biochemical oxygen demand (BOD)	TX	1530	20027401
Carbonaceous BOD, CBOD	TX	1555	20027401
Method SM 5310 B			
Analyte	AB	Analyte ID	Method ID





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**Matrix: Non-Potable Water**

Total Organic Carbon (TOC)	TX	2040	20137206
<b>Method</b> SM 5310 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Organic Carbon (TOC)	TX	2040	20138209
<b>Method</b> SM 5540 C			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Surfactants - MBAS	TX	2025	20144405
<b>Method</b> TCEQ 1005			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
Method ASTM D2216			
Analyte Moisture	TX	10337	ASTM D2216-05
Method EPA 1010			
Analyte Ignitability	TX	1780	10116606
Method EPA 1030			
Analyte Ignitability	TX	1780	10117201
Method EPA 1311			
Analyte TCLP	TX	849	10118806
Method EPA 1312			
Analyte SPLP	TX	850	10119003
Method EPA 1668			
Analyte Decachlorobiphenyls	TX	10332	10262007
Dichlorobiphenyls	TX	464	10262007
Heptachlorobiphenyls	TX	486	10262007
Hexachlorobiphenyls	TX	487	10262007
Monochlorobiphenyls	TX	501	10262007
Nonachlorobiphenyls	TX	507	10262007
Octachlorobiphenyls	TX	508	10262007
Pentachlorobiphenyls	TX	515	10262007
Tetrachlorobiphenyls	TX	528	10262007
Trichlorobiphenyls	TX	541	10262007
Method EPA 200.8			
Analyte Uranium	TX	3035	10014605



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**Matrix: Solid & Chemical Materials**

**Method EPA 300.0**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10053200
Chloride	TX	1575	10053200
Fluoride	TX	1730	10053200
Nitrate as N	TX	1810	10053200
Nitrate-nitrite	TX	1820	10053200
Nitrite as N	TX	1840	10053200
Orthophosphate as P	TX	1870	10053200
Sulfate	TX	2000	10053200

**Method EPA 310.1**

Analyte	AB	Analyte ID	Method ID
Alkalinity as CaCO3	TX	1505	10054805

**Method EPA 350.3**

Analyte	AB	Analyte ID	Method ID
Ammonia as N	TX	1515	10064401

**Method EPA 365.3**

Analyte	AB	Analyte ID	Method ID
Orthophosphate as P	TX	1870	10070801
Phosphorus	TX	1910	10070801

**Method EPA 6020**

Analyte	AB	Analyte ID	Method ID
Aluminum	TX	1000	10156204
Antimony	TX	1005	10156204
Arsenic	TX	1010	10156204
Barium	TX	1015	10156204
Beryllium	TX	1020	10156204
Boron	TX	1025	10156204
Cadmium	TX	1030	10156204
Calcium	TX	1035	10156204
Chromium	TX	1040	10156204



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**Matrix: Solid & Chemical Materials**

Cobalt	TX	1050	10156204
Copper	TX	1055	10156204
Iron	TX	1070	10156204
Lead	TX	1075	10156204
Lithium	TX	1080	10156204
Magnesium	TX	1085	10156204
Manganese	TX	1090	10156204
Molybdenum	TX	1100	10156204
Nickel	TX	1105	10156204
Potassium	TX	1125	10156204
Selenium	TX	1140	10156204
Silver	TX	1150	10156204
Sodium	TX	1155	10156204
Strontium	TX	1160	10156204
Thallium	TX	1165	10156204
Tin	TX	1175	10156204
Titanium	TX	1180	10156204
Vanadium	TX	1185	10156204
Zinc	TX	1190	10156204
<b>Method EPA 7196</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Chromium (VI)	TX	1045	10162206
<b>Method EPA 7470</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10165603
<b>Method EPA 7471</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Mercury	TX	1095	10166004
<b>Method EPA 8015</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
Diesel range organics (DRO)	TX	9369	10173203



# Texas Commission on Environmental Quality



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Issue Date: 5/1/2019

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**Matrix: Solid & Chemical Materials**

Ethanol	TX	4750	10173203
Ethylene glycol	TX	4785	10173203
Gasoline range organics (GRO)	TX	9408	10173203
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10173203
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10173203
Methanol	TX	4930	10173203
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10173203
n-Propanol (1-Propanol)	TX	5055	10173203
Propylene Glycol	TX	6657	10173203
tert-Butyl alcohol	TX	4420	10173203

**Method EPA 8021**

Analyte	AB	Analyte ID	Method ID
Benzene	TX	4375	10174400
Ethylbenzene	TX	4765	10174400
m+p-xylene	TX	5240	10174400
Methyl tert-butyl ether (MTBE)	TX	5000	10174400
o-Xylene	TX	5250	10174400
Toluene	TX	5140	10174400
Xylene (total)	TX	5260	10174400

**Method EPA 8081**

Analyte	AB	Analyte ID	Method ID
4,4'-DDD	TX	7355	10178402
4,4'-DDE	TX	7360	10178402
4,4'-DDT	TX	7365	10178402
Aldrin	TX	7025	10178402
alpha-BHC (alpha-Hexachlorocyclohexane)	TX	7110	10178402
alpha-Chlordane	TX	7240	10178402
beta-BHC (beta-Hexachlorocyclohexane)	TX	7115	10178402
Chlordane (tech.)	TX	7250	10178402
delta-BHC (delta-Hexachlorocyclohexane)	TX	7105	10178402
Dieldrin	TX	7470	10178402



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**Matrix: Solid & Chemical Materials**

Endosulfan I	TX	7510	10178402
Endosulfan II	TX	7515	10178402
Endosulfan sulfate	TX	7520	10178402
Endrin	TX	7540	10178402
Endrin aldehyde	TX	7530	10178402
Endrin ketone	TX	7535	10178402
gamma-BHC (Lindane, gamma-Hexachlorocyclohexane)	TX	7120	10178402
gamma-Chlordane	TX	7245	10178402
Heptachlor	TX	7685	10178402
Heptachlor epoxide	TX	7690	10178402
Methoxychlor	TX	7810	10178402
Mirex	TX	7870	10178402
Toxaphene (Chlorinated camphene)	TX	8250	10178402

**Method EPA 8082**

Analyte	AB	Analyte ID	Method ID
Aroclor-1016 (PCB-1016)	TX	8880	10179201
Aroclor-1221 (PCB-1221)	TX	8885	10179201
Aroclor-1232 (PCB-1232)	TX	8890	10179201
Aroclor-1242 (PCB-1242)	TX	8895	10179201
Aroclor-1248 (PCB-1248)	TX	8900	10179201
Aroclor-1254 (PCB-1254)	TX	8905	10179201
Aroclor-1260 (PCB-1260)	TX	8910	10179201
PCBs (total)	TX	8870	10179201

**Method EPA 8260**

Analyte	AB	Analyte ID	Method ID
1,1,1,2-Tetrachloroethane	TX	5105	10184404
1,1,1-Trichloroethane	TX	5160	10184404
1,1,2,2-Tetrachloroethane	TX	5110	10184404
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	TX	5195	10184404
1,1,2-Trichloroethane	TX	5165	10184404
1,1-Dichloroethane	TX	4630	10184404



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**Matrix: Solid & Chemical Materials**

1,1-Dichloroethylene	TX	4640	10184404
1,1-Dichloropropene	TX	4670	10184404
1,2,3-Trichlorobenzene	TX	5150	10184404
1,2,3-Trichloropropane	TX	5180	10184404
1,2,4-Trichlorobenzene	TX	5155	10184404
1,2,4-Trimethylbenzene	TX	5210	10184404
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10184404
1,2-Dibromoethane (EDB, Ethylene dibromide)	TX	4585	10184404
1,2-Dichlorobenzene	TX	4610	10184404
1,2-Dichloroethane (Ethylene dichloride)	TX	4635	10184404
1,2-Dichloropropane	TX	4655	10184404
1,3,5-Trimethylbenzene	TX	5215	10184404
1,3-Dichlorobenzene	TX	4615	10184404
1,3-Dichloropropane	TX	4660	10184404
1,4-Dichlorobenzene	TX	4620	10184404
1,4-Dioxane (1,4-Diethyleneoxide)	TX	4735	10184404
1-Chlorohexane	TX	4510	10184404
1-Propanol	TX	5060	10184404
2,2-Dichloropropane	TX	4665	10184404
2-Butanone (Methyl ethyl ketone, MEK)	TX	4410	10184404
2-Chloroethyl vinyl ether	TX	4500	10184404
2-Chlorotoluene	TX	4535	10184404
2-Hexanone (MBK)	TX	4860	10184404
4-Chlorotoluene	TX	4540	10184404
4-Isopropyltoluene (p-Cymene)	TX	4915	10184404
4-Methyl-2-pentanone (MIBK)	TX	4995	10184404
Acetone (2-Propanone)	TX	4315	10184404
Acetonitrile	TX	4320	10184404
Acrolein (Propenal)	TX	4325	10184404
Acrylonitrile	TX	4340	10184404





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**Matrix: Solid & Chemical Materials**

Allyl chloride (3-Chloropropene)	TX	4355	10184404
Benzene	TX	4375	10184404
Benzyl chloride	TX	5635	10184404
Bromobenzene	TX	4385	10184404
Bromochloromethane	TX	4390	10184404
Bromodichloromethane	TX	4395	10184404
Bromoform	TX	4400	10184404
Carbon disulfide	TX	4450	10184404
Carbon tetrachloride	TX	4455	10184404
Chlorobenzene	TX	4475	10184404
Chlorodibromomethane	TX	4575	10184404
Chloroethane (Ethyl chloride)	TX	4485	10184404
Chloroform	TX	4505	10184404
Chloroprene (2-Chloro-1,3-butadiene)	TX	4525	10184404
cis-1,2-Dichloroethylene	TX	4645	10184404
cis-1,3-Dichloropropene	TX	4680	10184404
Dibromofluoromethane	TX	4590	10184404
Dibromomethane (Methylene bromide)	TX	4595	10184404
Dichlorodifluoromethane (Freon-12)	TX	4625	10184404
Diethyl ether	TX	4725	10184404
Epichlorohydrin (1-Chloro-2,3-epoxypropane)	TX	4745	10184404
Ethanol	TX	4750	10184404
Ethyl acetate	TX	4755	10184404
Ethyl methacrylate	TX	4810	10184404
Ethylbenzene	TX	4765	10184404
Ethylene oxide	TX	4795	10184404
Hexachlorobutadiene	TX	4835	10184404
Iodomethane (Methyl iodide)	TX	4870	10184404
Isobutyl alcohol (2-Methyl-1-propanol)	TX	4875	10184404
Isopropyl alcohol (2-Propanol, Isopropanol)	TX	4895	10184404





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**Matrix: Solid & Chemical Materials**

Isopropylbenzene (Cumene)	TX	4900	10184404
m+p-xylene	TX	5240	10184404
Methacrylonitrile	TX	4925	10184404
Methyl acetate	TX	4940	10184404
Methyl acrylate	TX	4945	10184404
Methyl bromide (Bromomethane)	TX	4950	10184404
Methyl chloride (Chloromethane)	TX	4960	10184404
Methyl methacrylate	TX	4990	10184404
Methyl tert-butyl ether (MTBE)	TX	5000	10184404
Methylcyclohexane	TX	4965	10184404
Methylene chloride (Dichloromethane)	TX	4975	10184404
Naphthalene	TX	5005	10184404
n-Butyl alcohol (1-Butanol, n-Butanol)	TX	4425	10184404
n-Butylbenzene	TX	4435	10184404
n-Propylbenzene	TX	5090	10184404
o-Xylene	TX	5250	10184404
Pentachloroethane	TX	5035	10184404
Propionitrile (Ethyl cyanide)	TX	5080	10184404
Pyridine	TX	5095	10184404
sec-Butylbenzene	TX	4440	10184404
Styrene	TX	5100	10184404
tert-Butyl alcohol	TX	4420	10184404
tert-Butylbenzene	TX	4445	10184404
Tetrachloroethylene (Perchloroethylene)	TX	5115	10184404
Toluene	TX	5140	10184404
trans-1,2-Dichloroethylene	TX	4700	10184404
trans-1,3-Dichloropropylene	TX	4685	10184404
trans-1,4-Dichloro-2-butene	TX	4605	10184404
Trichloroethene (Trichloroethylene)	TX	5170	10184404
Trichlorofluoromethane (Fluorotrichloromethane, Freon 11)	TX	5175	10184404



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**Matrix: Solid & Chemical Materials**

Vinyl acetate	TX	5225	10184404
Vinyl chloride	TX	5235	10184404
Xylene (total)	TX	5260	10184404
<b>Method EPA 8270</b>			
<b>Analyte</b>	<b>AB</b>	<b>Analyte ID</b>	<b>Method ID</b>
1,2,4,5-Tetrachlorobenzene	TX	6715	10185203
1,2,4-Trichlorobenzene	TX	5155	10185203
1,2-Dibromo-3-chloropropane (DBCP)	TX	4570	10185203
1,2-Dichlorobenzene	TX	4610	10185203
1,2-Dinitrobenzene	TX	6155	10185203
1,2-Diphenylhydrazine	TX	6220	10185203
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10185203
1,3-Dichlorobenzene	TX	4615	10185203
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10185203
1,4-Dichlorobenzene	TX	4620	10185203
1,4-Dinitrobenzene	TX	6165	10185203
1,4-Naphthoquinone	TX	6420	10185203
1,4-Phenylenediamine	TX	6630	10185203
1-Chloronaphthalene	TX	5790	10185203
1-Naphthylamine	TX	6425	10185203
2,2'-Oxybis(1-chloropropane) (bis(2-Chloro-1-methylethyl)ether)	TX	4659	10185203
2,3,4,6-Tetrachlorophenol	TX	6735	10185203
2,4,5-Trichlorophenol	TX	6835	10185203
2,4,5-Trimethylaniline	TX	6880	10185203
2,4,6-Trichlorophenol	TX	6840	10185203
2,4-Diaminotoluene	TX	5880	10185203
2,4-Dichlorophenol	TX	6000	10185203
2,4-Dimethylphenol	TX	6130	10185203
2,4-Dinitrophenol	TX	6175	10185203
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10185203



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**Matrix: Solid & Chemical Materials**

2,6-Dichlorophenol	TX	6005	10185203
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10185203
2-Acetylaminofluorene	TX	5515	10185203
2-Chloronaphthalene	TX	5795	10185203
2-Chlorophenol	TX	5800	10185203
2-Methyl-4,6-dinitrophenol (4,6-Dinitro-2-methylphenol)	TX	6360	10185203
2-Methylaniline (o-Toluidine)	TX	5145	10185203
2-Methylnaphthalene	TX	6385	10185203
2-Methylphenol (o-Cresol)	TX	6400	10185203
2-Naphthylamine	TX	6430	10185203
2-Nitroaniline	TX	6460	10185203
2-Nitrophenol	TX	6490	10185203
2-Picoline (2-Methylpyridine)	TX	5050	10185203
3,3'-Dichlorobenzidine	TX	5945	10185203
3,3'-Dimethylbenzidine	TX	6120	10185203
3-Methylcholanthrene	TX	6355	10185203
3-Methylphenol (m-Cresol)	TX	6405	10185203
3-Nitroaniline	TX	6465	10185203
4-Aminobiphenyl	TX	5540	10185203
4-Bromophenyl phenyl ether (BDE-3)	TX	5660	10185203
4-Chloro-3-methylphenol	TX	5700	10185203
4-Chloroaniline	TX	5745	10185203
4-Chlorophenyl phenylether	TX	5825	10185203
4-Methylphenol (p-Cresol)	TX	6410	10185203
4-Nitroaniline	TX	6470	10185203
4-Nitrophenol	TX	6500	10185203
4-Nitroquinoline-1-oxide	TX	6510	10185203
5-Nitro-o-toluidine	TX	6570	10185203
7,12-Dimethylbenz(a) anthracene	TX	6115	10185203
a-a-Dimethylphenethylamine	TX	6125	10185203



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**Matrix: Solid & Chemical Materials**

Acenaphthene	TX	5500	10185203
Acenaphthylene	TX	5505	10185203
Acetophenone	TX	5510	10185203
Aniline	TX	5545	10185203
Anthracene	TX	5555	10185203
Aramite	TX	5560	10185203
Atrazine	TX	7065	10185203
Azinphos-methyl (Guthion)	TX	7075	10185203
Azobenzene	TX	5562	10185203
Benzenethiol (Thiophenol)	TX	6750	10185203
Benzidine	TX	5595	10185203
Benzo(a)anthracene	TX	5575	10185203
Benzo(a)pyrene	TX	5580	10185203
Benzo(b)fluoranthene	TX	5585	10185203
Benzo(e)pyrene	TX	5605	10185203
Benzo(g,h,i)perylene	TX	5590	10185203
Benzo(k)fluoranthene	TX	5600	10185203
Benzoic acid	TX	5610	10185203
Benzyl alcohol	TX	5630	10185203
Biphenyl	TX	5640	10185203
bis(2-Chloroethoxy)methane	TX	5760	10185203
bis(2-Chloroethyl) ether	TX	5765	10185203
bis(2-Ethylhexyl) phthalate (Di(2-Ethylhexyl) phthalate, DEHP)	TX	6065	10185203
Butyl benzyl phthalate	TX	5670	10185203
Caprolactam	TX	7180	10185203
Carbaryl (Sevin)	TX	7195	10185203
Carbazole	TX	5680	10185203
Carbophenothion	TX	7220	10185203
Chlorobenzilate	TX	7260	10185203
Chrysene	TX	5855	10185203



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**Matrix: Solid & Chemical Materials**

Demeton	TX	7390	10185203
Demeton-o	TX	7395	10185203
Demeton-s	TX	7385	10185203
Diallate	TX	7405	10185203
Dibenz(a,h) anthracene	TX	5895	10185203
Dibenz(a,j) acridine	TX	5900	10185203
Dibenzo(a,e) pyrene	TX	5890	10185203
Dibenzofuran	TX	5905	10185203
Dichlorovos (DDVP, Dichlorvos)	TX	8610	10185203
Diethyl phthalate	TX	6070	10185203
Dimethoate	TX	7475	10185203
Dimethyl phthalate	TX	6135	10185203
Di-n-butyl phthalate	TX	5925	10185203
Di-n-octyl phthalate	TX	6200	10185203
Dinoseb (2-sec-butyl-4,6-dinitrophenol, DNBP)	TX	8620	10185203
Diphenylamine	TX	6205	10185203
Disulfoton	TX	8625	10185203
Ethyl methanesulfonate	TX	6260	10185203
Fluoranthene	TX	6265	10185203
Fluorene	TX	6270	10185203
Hexachlorobenzene	TX	6275	10185203
Hexachlorobutadiene	TX	4835	10185203
Hexachlorocyclopentadiene	TX	6285	10185203
Hexachloroethane	TX	4840	10185203
Hexachlorophene	TX	6290	10185203
Hexachloropropene	TX	6295	10185203
Indeno(1,2,3-cd) pyrene	TX	6315	10185203
Isodrin	TX	7725	10185203
Isophorone	TX	6320	10185203
Isosafrole	TX	6325	10185203



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**Matrix: Solid & Chemical Materials**

Kepona	TX	7740	10185203
Malathion	TX	7770	10185203
Methapyrilene	TX	6345	10185203
Methyl methanesulfonate	TX	6375	10185203
Methyl parathion (Parathion, methyl)	TX	7825	10185203
Mevinphos	TX	7850	10185203
Naphthalene	TX	5005	10185203
Nitrobenzene	TX	5015	10185203
n-Nitrosodiethylamine	TX	6525	10185203
n-Nitrosodimethylamine	TX	6530	10185203
n-Nitrosodi-n-butylamine	TX	5025	10185203
n-Nitrosodi-n-propylamine	TX	6545	10185203
n-Nitrosodiphenylamine	TX	6535	10185203
n-Nitrosomethylethylamine	TX	6550	10185203
n-Nitrosomorpholine	TX	6555	10185203
n-Nitrosopiperidine	TX	6560	10185203
n-Nitrosopyrrolidine	TX	6565	10185203
o,o,o-Triethyl phosphorothioate	TX	8290	10185203
o-Anisidine	TX	5550	10185203
Parathion, ethyl	TX	7955	10185203
p-Cresidine	TX	5860	10185203
Pentachlorobenzene	TX	6590	10185203
Pentachloronitrobenzene (PCNB)	TX	6600	10185203
Pentachlorophenol	TX	6605	10185203
Phenacetin	TX	6610	10185203
Phenanthrene	TX	6615	10185203
Phenol	TX	6625	10185203
Phorate	TX	7985	10185203
Pronamide (Kerb)	TX	6650	10185203
Pyrene	TX	6665	10185203





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**Matrix: Solid & Chemical Materials**

Pyridine	TX	5095	10185203
Quinoline	TX	6670	10185203
Safrole	TX	6685	10185203
Sulfotepp	TX	8155	10185203
Terbufos	TX	8185	10185203
Tetrachlorvinphos (Stirophos, Gardona)	TX	8197	10185203
Thionazin (Zinophos)	TX	8235	10185203
Toluene diisocyanate	TX	6775	10185203

**Method EPA 8290**

Analyte	AB	Analyte ID	Method ID
1,2,3,4,6,7,8,9-Octachlorodibenzofuran (OCDF)	TX	9516	10187209
1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin (OCDD)	TX	9519	10187209
1,2,3,4,6,7,8-Heptachlorodibenzofuran (1,2,3,4,6,7,8-HpCDF)	TX	9420	10187209
1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (1,2,3,4,6,7,8-HpCDD)	TX	9426	10187209
1,2,3,4,7,8,9-Heptachlorodibenzofuran (1,2,3,4,7,8,9-HpCDF)	TX	9423	10187209
1,2,3,4,7,8-Hexachlorodibenzofuran (1,2,3,4,7,8-HxCDF)	TX	9471	10187209
1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (1,2,3,4,7,8-HxCDD)	TX	9453	10187209
1,2,3,6,7,8-Hexachlorodibenzofuran (1,2,3,6,7,8-HxCDF)	TX	9474	10187209
1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin(1,2,3,6,7,8-HxCDD)	TX	9456	10187209
1,2,3,7,8,9-Hexachlorodibenzofuran (1,2,3,7,8,9-HxCDF)	TX	9477	10187209
1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (1,2,3,7,8,9-HxCDD)	TX	9459	10187209
1,2,3,7,8-Pentachlorodibenzofuran (1,2,3,7,8-PeCDF)	TX	9543	10187209
1,2,3,7,8-Pentachlorodibenzo-p-dioxin (1,2,3,7,8-PeCDD)	TX	9540	10187209
2,3,4,6,7,8-Hexachlorodibenzofuran (2,3,4,6,7,8-HxCDF)	TX	9480	10187209
2,3,4,7,8-Pentachlorodibenzofuran (2,3,4,7,8-PeCDF)	TX	9549	10187209
2,3,7,8-Tetrachlorodibenzofuran (2,3,7,8-TCDF)	TX	9612	10187209
2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD)	TX	9618	10187209
Total Heptachlorodibenzofuran (Total HpCDF)	TX	9444	10187209
Total Heptachlorodibenzo-p-dioxin (Total HpCDD)	TX	9438	10187209
Total Hexachlorodibenzofuran (Total HxCDF)	TX	9483	10187209



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
Houston, TX 77099-4338

Certificate: T104704231-19-23  
Expiration Date: 4/30/2020  
Issue Date: 5/1/2019

These fields of accreditation supercede all previous fields. The Texas Commission on Environmental Quality urges customers to verify the laboratory's current accreditation status for particular methods and analyses.

**Matrix: Solid & Chemical Materials**

Total Hexachlorodibenzo-p-dioxin (Total HxCDD)	TX	9468	10187209
Total Pentachlorodibenzofuran (Total PeCDF)	TX	9552	10187209
Total Pentachlorodibenzo-p-dioxin (Total PeCDD)	TX	9555	10187209
Total Tetrachlorodibenzofuran (Total TCDF)	TX	9615	10187209
Total Tetrachlorodibenzo-p-dioxin (Total TCDD)	TX	9609	10187209

**Method EPA 8316**

Analyte	AB	Analyte ID	Method ID
Acrylamide	TX	4330	10188202

**Method EPA 8330**

Analyte	AB	Analyte ID	Method ID
1,3,5-Trinitrobenzene (1,3,5-TNB)	TX	6885	10189807
1,3-Dinitrobenzene (1,3-DNB)	TX	6160	10189807
2,4,6-Trinitrotoluene (2,4,6-TNT)	TX	9651	10189807
2,4-Dinitrotoluene (2,4-DNT)	TX	6185	10189807
2,6-Dinitrotoluene (2,6-DNT)	TX	6190	10189807
2-Amino-4,6-dinitrotoluene (2-am-dnt)	TX	9303	10189807
2-Nitrotoluene	TX	9507	10189807
3-Nitrotoluene	TX	9510	10189807
4-Amino-2,6-dinitrotoluene (4-am-dnt)	TX	9306	10189807
4-Nitrotoluene	TX	9513	10189807
Methyl-2,4,6-trinitrophenylnitramine (tetryl)	TX	6415	10189807
Nitrobenzene	TX	5015	10189807
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	TX	9522	10189807
RDX (hexahydro-1,3,5-trinitro-1,3,5-triazine)	TX	9432	10189807

**Method EPA 9014**

Analyte	AB	Analyte ID	Method ID
Amenable cyanide	TX	1510	10193803
Total cyanide	TX	1645	10193803

**Method EPA 9038**

Analyte	AB	Analyte ID	Method ID
Sulfate	TX	2000	10196608





# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

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Certificate: T104704231-19-23  
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**Matrix: Solid & Chemical Materials**

**Method EPA 9040**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197203
pH	TX	1900	10196802

**Method EPA 9045**

Analyte	AB	Analyte ID	Method ID
Corrosivity	TX	1615	10197805
pH	TX	1900	10197805

**Method EPA 9050**

Analyte	AB	Analyte ID	Method ID
Conductivity	TX	1610	10198604

**Method EPA 9056**

Analyte	AB	Analyte ID	Method ID
Bromide	TX	1540	10199209
Chloride	TX	1575	10199209
Fluoride	TX	1730	10199209
Nitrate as N	TX	1810	10199209
Nitrate-nitrite	TX	1820	10199209
Nitrite as N	TX	1840	10199209
Orthophosphate as P	TX	1870	10199209
Sulfate	TX	2000	10199209

**Method EPA 9060**

Analyte	AB	Analyte ID	Method ID
Total Organic Carbon (TOC)	TX	2040	10200201

**Method EPA 9065**

Analyte	AB	Analyte ID	Method ID
Total phenolics	TX	1905	10200405

**Method EPA 9071**

Analyte	AB	Analyte ID	Method ID
n-Hexane Extractable Material (HEM) (O&G)	TX	1803	10201204



# Texas Commission on Environmental Quality



## NELAP - Recognized Laboratory Fields of Accreditation

ALS Laboratory Group, Environmental Services Division (Houston, Texas)

10450 Stancliff Road, Suite 210  
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**Matrix: Solid & Chemical Materials**

Method	AB	Analyte ID	Method ID
EPA 9095			
<b>Analyte</b> Paint Filter Liquids Test	TX	10312	10204009
EPA 9250			
<b>Analyte</b> Chloride	TX	1575	10207202
SM 2320 B			
<b>Analyte</b> Alkalinity as CaCO3	TX	1505	20045005
SM 2510 B			
<b>Analyte</b> Conductivity	TX	1610	20048004
SM 2540 G			
<b>Analyte</b> Residue-total (total solids)	TX	1950	20005203
SSA/ASA Part 3:34			
<b>Analyte</b> Carbon, organic (Walkley-Black)	TX	10340	SSA/ASA Pt 3:34
TCEQ 1005			
<b>Analyte</b> Total Petroleum Hydrocarbons (TPH)	TX	2050	90019208



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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

March 23, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20030619**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 16 sample(s) on Mar 13, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data								
Laboratory Name: ALS Laboratory Group				LRC Date: 03/23/2020				
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20030619				
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 151716, 151763, 151785, 151787, 151813, R358243, R358244				
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>	
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>						
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X					
		Were all departures from standard conditions described in an exception report?	X					
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>						
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X					
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X					
<b>R3</b>	OI	<b>Test reports</b>						
		Were all samples prepared and analyzed within holding times?	X					
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X					
		Were calculations checked by a peer or supervisor?	X					
		Were all analyte identifications checked by a peer or supervisor?	X					
		Were sample detection limits reported for all analytes not detected?	X					
		Were all results for soil and sediment samples reported on a dry weight basis?			X			
		Were % moisture (or solids) reported for all soil and sediment samples?			X			
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X			
		If required for the project, TICs reported?			X			
<b>R4</b>	O	<b>Surrogate recovery data</b>						
		Were surrogates added prior to extraction?	X					
		Were surrogate percent recoveries in all samples within the laboratory QC limits?		X				1
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>						
		Were appropriate type(s) of blanks analyzed?	X					
		Were blanks analyzed at the appropriate frequency?	X					
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X					
		Were blank concentrations < MQL?	X					
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>						
		Were all COCs included in the LCS?	X					
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X					
		Were LCSs analyzed at the required frequency?	X					
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X					
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X					
		Was the LCSD RPD within QC limits?		X				2
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>						
		Were the project/method specified analytes included in the MS and MSD?	X					
		Were MS/MSD analyzed at the appropriate frequency?		X				3
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X					
		Were MS/MSD RPDs within laboratory QC limits?	X					
<b>R8</b>	OI	<b>Analytical duplicate data</b>						
		Were appropriate analytical duplicates analyzed for each matrix?	X					
		Were analytical duplicates analyzed at the appropriate frequency?	X					
		Were RPDs or relative standard deviations within the laboratory QC limits?	X					
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>						
		Are the MQLs for each method analyte included in the laboratory data package?	X					
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X					
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X					
<b>R10</b>	OI	<b>Other problems/anomalies</b>						
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X					
		Were all necessary corrective actions performed for the reported data?	X					
		Was applicable and available technology used to lower the SDL and minimize the matrix interference affects on the sample results?	X					
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X					

**Laboratory Review Checklist: Supporting Data**

Laboratory Name: ALS Laboratory Group		LRC Date: 03/23/2020					
Project Name: Houston TX-Wood Preserving Works		Laboratory Job Number: HS20030619					
Reviewer Name: Dane Wacasey		Prep Batch Number(s): 151716, 151763, 151785, 151787, 151813, R358243, R358244					
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>S1</b>	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
<b>S2</b>	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
<b>S3</b>	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
<b>S4</b>	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
<b>S5</b>	OI	<b>Raw data (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section</b>					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
<b>S6</b>	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
<b>S7</b>	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
<b>S8</b>	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
<b>S9</b>	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
<b>S10</b>	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
<b>S11</b>	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
<b>S12</b>	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
<b>S13</b>	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
<b>S14</b>	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
<b>S15</b>	OI	<b>Verification/validation documentation for methods (NELAC Chap 5 or ISO/IEC 17025 Section 5)</b>					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
<b>S16</b>	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 03/23/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20030619
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 151716, 151763, 151785, 151787, 151813, R358243, R358244

ER# <sup>5</sup>	Description
1	Semivolatile Organics Method SW8270, samples WG-1620-MW70C-20200312, WG-1620-FD01-20200312, the surrogates could not be determined due to dilution below the calibration range in the 1000x run.  Semivolatile Organics Method SW8270, sample WG-1620-MW50B-20200312, the surrogates could not be determined due to dilution below the calibration range in the 100x run.
2	Batch 151785, Texas TPH by TX 1005, LCS/LCSD RPD was above the RPD limit for nC6 to nC12. The individual recoveries met acceptance criteria.
3	Batch 151716, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
 O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
 NA = Not Applicable;  
 NR = Not Reviewed;  
 R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20030619

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20030619-01	WQ-1620-TB01-20200312	Water		12-Mar-2020 16:30	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-02	WQ-1620-MW99C-20200311	Groundwater		11-Mar-2020 15:35	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-03	WG-1620-MW54B-20200311	Groundwater		11-Mar-2020 14:20	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-04	WG-1620-MW94A-20200312	Groundwater		12-Mar-2020 15:45	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-05	WG-1620-MW93B-20200312	Groundwater		12-Mar-2020 14:35	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-06	WG-1620-MW92B-20200312	Groundwater		12-Mar-2020 13:20	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-07	WG-1620-MW84A-20200312	Groundwater		12-Mar-2020 12:15	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-08	WG-1620-MW70C-20200312	Groundwater		12-Mar-2020 10:35	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-09	WG-1620-MW91A-20200312	Groundwater		12-Mar-2020 09:20	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-10	WG-1620-FD01-20200312	Groundwater		12-Mar-2020 00:00	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-11	WQ-1620-TB02-20200312	Water		12-Mar-2020 12:36	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-12	WG-1620-MW76B-20200312	Groundwater		12-Mar-2020 10:35	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-13	WG-1620-MW97A-20200312	Groundwater		12-Mar-2020 11:47	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-14	WG-1620-MW98A-20200312	Groundwater		12-Mar-2020 12:54	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-15	WG-1620-MW98B-20200312	Groundwater		12-Mar-2020 13:53	13-Mar-2020 14:00	<input type="checkbox"/>
HS20030619-16	WG-1620-MW50B-20200312	Groundwater		12-Mar-2020 15:20	13-Mar-2020 14:00	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB01-20200312  
 Collection Date: 12-Mar-2020 16:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>					Analyst: AKP
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 06:28
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 06:28
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 06:28
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 06:28
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 06:28
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 06:28
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 06:28
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>104</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 06:28</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 06:28</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 06:28</i>
<i>Surr: Toluene-d8</i>	<i>100.0</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 06:28</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-MW99C-20200311  
 Collection Date: 11-Mar-2020 15:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-02  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:05
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:05
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:05
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:05
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	14-Mar-2020 18:05
<b>Toluene</b>	<b>0.00078</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Mar-2020 18:05
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:05
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>104</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:05</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:05</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:05</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:05</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-MW99C-20200311  
 Collection Date: 11-Mar-2020 15:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-02  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 17:59
2,4-Dimethylphenol	< 0.000041		0.000041	0.00020	mg/L	1	16-Mar-2020 17:59
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	16-Mar-2020 17:59
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00020	mg/L	1	16-Mar-2020 17:59
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 17:59
<b>2-Methylnaphthalene</b>	<b>0.000086</b>	J	<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 17:59
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 17:59
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	16-Mar-2020 17:59
Acenaphthene	< 0.000028		0.000028	0.00010	mg/L	1	16-Mar-2020 17:59
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	16-Mar-2020 17:59
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	16-Mar-2020 17:59
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	16-Mar-2020 17:59
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 17:59
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00020	mg/L	1	16-Mar-2020 17:59
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000070</b>	J	<b>0.000038</b>	<b>0.00020</b>	<b>mg/L</b>	1	16-Mar-2020 17:59
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 17:59
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 17:59
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 17:59
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	16-Mar-2020 17:59
Fluorene	< 0.000031		0.000031	0.00010	mg/L	1	16-Mar-2020 17:59
<b>Naphthalene</b>	<b>0.00018</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 17:59
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	16-Mar-2020 17:59
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00020	mg/L	1	16-Mar-2020 17:59
Pentachlorophenol	< 0.000081		0.000081	0.00020	mg/L	1	16-Mar-2020 17:59
<b>Phenanthrene</b>	<b>0.000053</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 17:59
Phenol	< 0.000036		0.000036	0.00020	mg/L	1	16-Mar-2020 17:59
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	16-Mar-2020 17:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>77.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>16-Mar-2020 17:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>57.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>16-Mar-2020 17:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>57.1</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>16-Mar-2020 17:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>65.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>16-Mar-2020 17:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>47.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>16-Mar-2020 17:59</i>
<i>Surr: Phenol-d6</i>	<i>51.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>16-Mar-2020 17:59</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Mar-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.000866</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Mar-2020 18:20

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54B-20200311  
 Collection Date: 11-Mar-2020 14:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:30
<b>Benzene</b>	<b>0.00043</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Mar-2020 18:30
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:30
<b>Ethylbenzene</b>	<b>0.0011</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Mar-2020 18:30
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	14-Mar-2020 18:30
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:30
<b>Xylenes, Total</b>	<b>0.0018</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Mar-2020 18:30
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>106</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:30</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:30</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:30</i>
<i>Surr: Toluene-d8</i>	<i>99.6</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW54B-20200311  
 Collection Date: 11-Mar-2020 14:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 18:18
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	16-Mar-2020 18:18
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	16-Mar-2020 18:18
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	16-Mar-2020 18:18
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 18:18
<b>2-Methylnaphthalene</b>	<b>0.00011</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 18:18
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	16-Mar-2020 18:18
<b>Acenaphthene</b>	<b>0.0093</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
<b>Acenaphthylene</b>	<b>0.000087</b>	J	<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
<b>Anthracene</b>	<b>0.00017</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	16-Mar-2020 18:18
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 18:18
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	16-Mar-2020 18:18
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000089</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 18:18
<b>Dibenzofuran</b>	<b>0.000076</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
<b>Di-n-butyl phthalate</b>	<b>0.000033</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
<b>Fluoranthene</b>	<b>0.000022</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
<b>Fluorene</b>	<b>0.00015</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
<b>Naphthalene</b>	<b>0.0021</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	16-Mar-2020 18:18
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	16-Mar-2020 18:18
Pentachlorophenol	< 0.000080		0.000080	0.00020	mg/L	1	16-Mar-2020 18:18
<b>Phenanthrene</b>	<b>0.0028</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:18
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	16-Mar-2020 18:18
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	16-Mar-2020 18:18
<i>Surr: 2,4,6-Tribromophenol</i>	86.7			34-129	%REC	1	16-Mar-2020 18:18
<i>Surr: 2-Fluorobiphenyl</i>	52.5			40-125	%REC	1	16-Mar-2020 18:18
<i>Surr: 2-Fluorophenol</i>	54.7			20-120	%REC	1	16-Mar-2020 18:18
<i>Surr: 4-Terphenyl-d14</i>	88.2			40-135	%REC	1	16-Mar-2020 18:18
<i>Surr: Nitrobenzene-d5</i>	49.7			41-120	%REC	1	16-Mar-2020 18:18
<i>Surr: Phenol-d6</i>	54.3			20-120	%REC	1	16-Mar-2020 18:18
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Mar-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00117</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Mar-2020 18:23

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW94A-20200312  
 Collection Date: 12-Mar-2020 15:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-04  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:54
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:54
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:54
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:54
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	14-Mar-2020 18:54
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 18:54
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 18:54
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>105</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:54</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:54</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:54</i>
<i>Surr: Toluene-d8</i>	<i>95.7</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 18:54</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW94A-20200312  
 Collection Date: 12-Mar-2020 15:45

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-04  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000022		0.000022	0.00021	mg/L	1	16-Mar-2020 18:37
2,4-Dimethylphenol	< 0.000041		0.000041	0.00021	mg/L	1	16-Mar-2020 18:37
2,4-Dinitrotoluene	< 0.000060		0.000060	0.00021	mg/L	1	16-Mar-2020 18:37
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00021	mg/L	1	16-Mar-2020 18:37
2-Chloronaphthalene	< 0.000022		0.000022	0.00021	mg/L	1	16-Mar-2020 18:37
2-Methylnaphthalene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 18:37
4,6-Dinitro-2-methylphenol	< 0.000021		0.000021	0.00021	mg/L	1	16-Mar-2020 18:37
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	16-Mar-2020 18:37
Acenaphthene	< 0.000028		0.000028	0.00010	mg/L	1	16-Mar-2020 18:37
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	16-Mar-2020 18:37
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	16-Mar-2020 18:37
Benz(a)anthracene	< 0.000052		0.000052	0.00010	mg/L	1	16-Mar-2020 18:37
Benzo(a)pyrene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 18:37
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00021	mg/L	1	16-Mar-2020 18:37
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000056</b>	J	<b>0.000038</b>	<b>0.00021</b>	<b>mg/L</b>	1	16-Mar-2020 18:37
Chrysene	< 0.000022		0.000022	0.00010	mg/L	1	16-Mar-2020 18:37
Dibenzofuran	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 18:37
Di-n-butyl phthalate	< 0.000021		0.000021	0.00021	mg/L	1	16-Mar-2020 18:37
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	16-Mar-2020 18:37
Fluorene	< 0.000031		0.000031	0.00010	mg/L	1	16-Mar-2020 18:37
Naphthalene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 18:37
Nitrobenzene	< 0.000025		0.000025	0.00021	mg/L	1	16-Mar-2020 18:37
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00021	mg/L	1	16-Mar-2020 18:37
Pentachlorophenol	< 0.000081		0.000081	0.00021	mg/L	1	16-Mar-2020 18:37
Phenanthrene	< 0.000022		0.000022	0.00010	mg/L	1	16-Mar-2020 18:37
Phenol	< 0.000036		0.000036	0.00021	mg/L	1	16-Mar-2020 18:37
Pyrene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 18:37
<i>Surr: 2,4,6-Tribromophenol</i>	94.2			34-129	%REC	1	16-Mar-2020 18:37
<i>Surr: 2-Fluorobiphenyl</i>	62.9			40-125	%REC	1	16-Mar-2020 18:37
<i>Surr: 2-Fluorophenol</i>	65.9			20-120	%REC	1	16-Mar-2020 18:37
<i>Surr: 4-Terphenyl-d14</i>	92.0			40-135	%REC	1	16-Mar-2020 18:37
<i>Surr: Nitrobenzene-d5</i>	63.1			41-120	%REC	1	16-Mar-2020 18:37
<i>Surr: Phenol-d6</i>	76.1			20-120	%REC	1	16-Mar-2020 18:37
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Mar-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00540</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Mar-2020 18:25

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW93B-20200312  
 Collection Date: 12-Mar-2020 14:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-05  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 19:19
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 19:19
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 19:19
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 19:19
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	14-Mar-2020 19:19
<b>Toluene</b>	<b>0.00092</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Mar-2020 19:19
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 19:19
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>104</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:19</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:19</i>
<i>Surr: Dibromofluoromethane</i>	<i>99.5</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:19</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:19</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW93B-20200312  
 Collection Date: 12-Mar-2020 14:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-05  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 18:56
2,4-Dimethylphenol	< 0.000041		0.000041	0.00020	mg/L	1	16-Mar-2020 18:56
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	16-Mar-2020 18:56
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00020	mg/L	1	16-Mar-2020 18:56
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 18:56
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	16-Mar-2020 18:56
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 18:56
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	16-Mar-2020 18:56
<b>Acenaphthene</b>	<b>0.000039</b>	J	<b>0.000028</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:56
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	16-Mar-2020 18:56
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	16-Mar-2020 18:56
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	16-Mar-2020 18:56
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 18:56
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00020	mg/L	1	16-Mar-2020 18:56
Bis(2-ethylhexyl)phthalate	< 0.000038		0.000038	0.00020	mg/L	1	16-Mar-2020 18:56
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 18:56
<b>Dibenzofuran</b>	<b>0.000071</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:56
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 18:56
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	16-Mar-2020 18:56
<b>Fluorene</b>	<b>0.000058</b>	J	<b>0.000031</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:56
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 18:56
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	16-Mar-2020 18:56
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00020	mg/L	1	16-Mar-2020 18:56
Pentachlorophenol	< 0.000081		0.000081	0.00020	mg/L	1	16-Mar-2020 18:56
<b>Phenanthrene</b>	<b>0.00016</b>		<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	16-Mar-2020 18:56
Phenol	< 0.000036		0.000036	0.00020	mg/L	1	16-Mar-2020 18:56
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	16-Mar-2020 18:56
<i>Surr: 2,4,6-Tribromophenol</i>	83.6			34-129	%REC	1	16-Mar-2020 18:56
<i>Surr: 2-Fluorobiphenyl</i>	73.6			40-125	%REC	1	16-Mar-2020 18:56
<i>Surr: 2-Fluorophenol</i>	29.5			20-120	%REC	1	16-Mar-2020 18:56
<i>Surr: 4-Terphenyl-d14</i>	65.6			40-135	%REC	1	16-Mar-2020 18:56
<i>Surr: Nitrobenzene-d5</i>	61.4			41-120	%REC	1	16-Mar-2020 18:56
<i>Surr: Phenol-d6</i>	69.7			20-120	%REC	1	16-Mar-2020 18:56
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Mar-2020		Analyst: JC	
<b>Arsenic</b>	<b>0.00455</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	18-Mar-2020 18:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW92B-20200312  
 Collection Date: 12-Mar-2020 13:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-06  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 19:44
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 19:44
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 19:44
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 19:44
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	14-Mar-2020 19:44
<b>Toluene</b>	<b>0.00049</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	14-Mar-2020 19:44
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 19:44
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>105</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:44</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:44</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:44</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 19:44</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW92B-20200312  
 Collection Date: 12-Mar-2020 13:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-06  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 19:15
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	16-Mar-2020 19:15
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	16-Mar-2020 19:15
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	16-Mar-2020 19:15
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	16-Mar-2020 19:15
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	16-Mar-2020 19:15
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 19:15
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	16-Mar-2020 19:15
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	16-Mar-2020 19:15
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	16-Mar-2020 19:15
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	16-Mar-2020 19:15
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	16-Mar-2020 19:15
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 19:15
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	16-Mar-2020 19:15
Bis(2-ethylhexyl)phthalate	< 0.000037		0.000037	0.00020	mg/L	1	16-Mar-2020 19:15
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 19:15
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 19:15
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	16-Mar-2020 19:15
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	16-Mar-2020 19:15
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	16-Mar-2020 19:15
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	16-Mar-2020 19:15
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	16-Mar-2020 19:15
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	16-Mar-2020 19:15
Pentachlorophenol	< 0.000080		0.000080	0.00020	mg/L	1	16-Mar-2020 19:15
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	16-Mar-2020 19:15
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	16-Mar-2020 19:15
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	16-Mar-2020 19:15
<i>Surr: 2,4,6-Tribromophenol</i>	90.6			34-129	%REC	1	16-Mar-2020 19:15
<i>Surr: 2-Fluorobiphenyl</i>	62.5			40-125	%REC	1	16-Mar-2020 19:15
<i>Surr: 2-Fluorophenol</i>	60.4			20-120	%REC	1	16-Mar-2020 19:15
<i>Surr: 4-Terphenyl-d14</i>	88.7			40-135	%REC	1	16-Mar-2020 19:15
<i>Surr: Nitrobenzene-d5</i>	57.7			41-120	%REC	1	16-Mar-2020 19:15
<i>Surr: Phenol-d6</i>	70.9			20-120	%REC	1	16-Mar-2020 19:15
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 17-Mar-2020		Analyst: JC	
Arsenic	0.00201		0.000400	0.00200	mg/L	1	18-Mar-2020 18:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84A-20200312  
 Collection Date: 12-Mar-2020 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-07  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 00:13
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 00:13
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 00:13
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 00:13
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 00:13
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 00:13
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 00:13
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>104</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 00:13</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 00:13</i>
<i>Surr: Dibromofluoromethane</i>	<i>100.0</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 00:13</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 00:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW84A-20200312  
 Collection Date: 12-Mar-2020 12:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-07  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 17:59
2,4-Dimethylphenol	< 0.000041		0.000041	0.00020	mg/L	1	17-Mar-2020 17:59
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	17-Mar-2020 17:59
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00020	mg/L	1	17-Mar-2020 17:59
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 17:59
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 17:59
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 17:59
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	17-Mar-2020 17:59
Acenaphthene	< 0.000028		0.000028	0.00010	mg/L	1	17-Mar-2020 17:59
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	17-Mar-2020 17:59
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	17-Mar-2020 17:59
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	17-Mar-2020 17:59
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 17:59
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00020	mg/L	1	17-Mar-2020 17:59
Bis(2-ethylhexyl)phthalate	< 0.000038		0.000038	0.00020	mg/L	1	17-Mar-2020 17:59
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 17:59
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 17:59
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 17:59
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	17-Mar-2020 17:59
Fluorene	< 0.000031		0.000031	0.00010	mg/L	1	17-Mar-2020 17:59
<b>Naphthalene</b>	<b>0.000023</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>17-Mar-2020 17:59</b>
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	17-Mar-2020 17:59
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00020	mg/L	1	17-Mar-2020 17:59
Pentachlorophenol	< 0.000081		0.000081	0.00020	mg/L	1	17-Mar-2020 17:59
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 17:59
Phenol	< 0.000036		0.000036	0.00020	mg/L	1	17-Mar-2020 17:59
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 17:59
<i>Surr: 2,4,6-Tribromophenol</i>	<i>88.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 17:59</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>70.1</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 17:59</i>
<i>Surr: 2-Fluorophenol</i>	<i>66.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 17:59</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 17:59</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 17:59</i>
<i>Surr: Phenol-d6</i>	<i>69.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 17:59</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00464</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	<b>1</b>	<b>20-Mar-2020 16:51</b>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW70C-20200312  
 Collection Date: 12-Mar-2020 10:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-08  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 01:50
<b>Benzene</b>	<b>0.010</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 01:50
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 01:50
<b>Ethylbenzene</b>	<b>0.057</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 01:50
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 01:50
<b>Toluene</b>	<b>0.034</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 01:50
<b>Xylenes, Total</b>	<b>0.064</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 01:50
Surr: 1,2-Dichloroethane-d4	102			70-126	%REC	1	15-Mar-2020 01:50
Surr: 4-Bromofluorobenzene	101			81-113	%REC	1	15-Mar-2020 01:50
Surr: Dibromofluoromethane	102			77-123	%REC	1	15-Mar-2020 01:50
Surr: Toluene-d8	100			82-127	%REC	1	15-Mar-2020 01:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW70C-20200312  
 Collection Date: 12-Mar-2020 10:35

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-08  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 18:18
<b>2,4-Dimethylphenol</b>	<b>0.00041</b>		<b>0.000042</b>	<b>0.00021</b>	<b>mg/L</b>	1	17-Mar-2020 18:18
2,4-Dinitrotoluene	< 0.000060		0.000060	0.00021	mg/L	1	17-Mar-2020 18:18
2,6-Dinitrotoluene	< 0.000044		0.000044	0.00021	mg/L	1	17-Mar-2020 18:18
2-Chloronaphthalene	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 18:18
<b>2-Methylnaphthalene</b>	<b>0.067</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 12:22
4,6-Dinitro-2-methylphenol	< 0.000021		0.000021	0.00021	mg/L	1	17-Mar-2020 18:18
4-Nitrophenol	< 0.000049		0.000049	0.0010	mg/L	1	17-Mar-2020 18:18
<b>Acenaphthene</b>	<b>0.059</b>		<b>0.00028</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 12:22
Acenaphthylene	< 0.000016		0.000016	0.00010	mg/L	1	17-Mar-2020 18:18
<b>Anthracene</b>	<b>0.0049</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:18
Benz(a)anthracene	< 0.000052		0.000052	0.00010	mg/L	1	17-Mar-2020 18:18
Benzo(a)pyrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 18:18
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00021	mg/L	1	17-Mar-2020 18:18
Bis(2-ethylhexyl)phthalate	< 0.000039		0.000039	0.00021	mg/L	1	17-Mar-2020 18:18
Chrysene	< 0.000022		0.000022	0.00010	mg/L	1	17-Mar-2020 18:18
<b>Dibenzofuran</b>	<b>0.055</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 12:22
<b>Di-n-butyl phthalate</b>	<b>0.000061</b>	J	<b>0.000021</b>	<b>0.00021</b>	<b>mg/L</b>	1	17-Mar-2020 18:18
<b>Fluoranthene</b>	<b>0.0030</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:18
<b>Fluorene</b>	<b>0.028</b>		<b>0.00031</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 12:22
<b>Naphthalene</b>	<b>1.6</b>		<b>0.021</b>	<b>0.10</b>	<b>mg/L</b>	1000	18-Mar-2020 18:26
Nitrobenzene	< 0.000025		0.000025	0.00021	mg/L	1	17-Mar-2020 18:18
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00021	mg/L	1	17-Mar-2020 18:18
Pentachlorophenol	< 0.000082		0.000082	0.00021	mg/L	1	17-Mar-2020 18:18
<b>Phenanthrene</b>	<b>0.040</b>		<b>0.00022</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 12:22
Phenol	< 0.000036		0.000036	0.00021	mg/L	1	17-Mar-2020 18:18
<b>Pyrene</b>	<b>0.0017</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:18
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	1000	18-Mar-2020 18:26
Surr: 2,4,6-Tribromophenol	86.3			34-129	%REC	10	18-Mar-2020 12:22
Surr: 2,4,6-Tribromophenol	85.5			34-129	%REC	1	17-Mar-2020 18:18
Surr: 2-Fluorobiphenyl	48.0			40-125	%REC	1	17-Mar-2020 18:18
Surr: 2-Fluorobiphenyl	54.8			40-125	%REC	10	18-Mar-2020 12:22
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	1000	18-Mar-2020 18:26
Surr: 2-Fluorophenol	0	JS		20-120	%REC	1000	18-Mar-2020 18:26
Surr: 2-Fluorophenol	52.4			20-120	%REC	10	18-Mar-2020 12:22
Surr: 2-Fluorophenol	62.1			20-120	%REC	1	17-Mar-2020 18:18
Surr: 4-Terphenyl-d14	85.5			40-135	%REC	1	17-Mar-2020 18:18
Surr: 4-Terphenyl-d14	87.7			40-135	%REC	10	18-Mar-2020 12:22
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	1000	18-Mar-2020 18:26

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW70C-20200312  
 Collection Date: 12-Mar-2020 10:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-08  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	18-Mar-2020 18:26
Surr: Nitrobenzene-d5	56.1			41-120	%REC	1	17-Mar-2020 18:18
Surr: Nitrobenzene-d5	49.1			41-120	%REC	10	18-Mar-2020 12:22
Surr: Phenol-d6	57.0			20-120	%REC	10	18-Mar-2020 12:22
Surr: Phenol-d6	0	JS		20-120	%REC	1000	18-Mar-2020 18:26
Surr: Phenol-d6	51.4			20-120	%REC	1	17-Mar-2020 18:18
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
Arsenic	0.00598		0.000400	0.00200	mg/L	1	20-Mar-2020 16:53

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW91A-20200312  
 Collection Date: 12-Mar-2020 09:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-09  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 03:14
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 03:14
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 03:14
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 03:14
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 03:14
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 03:14
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 03:14
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 03:14</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>98.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 03:14</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 03:14</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 03:14</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW91A-20200312  
 Collection Date: 12-Mar-2020 09:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-09  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 18:37
2,4-Dimethylphenol	< 0.000042		0.000042	0.00021	mg/L	1	17-Mar-2020 18:37
2,4-Dinitrotoluene	< 0.000061		0.000061	0.00021	mg/L	1	17-Mar-2020 18:37
2,6-Dinitrotoluene	< 0.000044		0.000044	0.00021	mg/L	1	17-Mar-2020 18:37
2-Chloronaphthalene	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 18:37
2-Methylnaphthalene	< 0.000020		0.000020	0.00011	mg/L	1	17-Mar-2020 18:37
4,6-Dinitro-2-methylphenol	< 0.000021		0.000021	0.00021	mg/L	1	17-Mar-2020 18:37
4-Nitrophenol	< 0.000049		0.000049	0.0011	mg/L	1	17-Mar-2020 18:37
Acenaphthene	< 0.000028		0.000028	0.00011	mg/L	1	17-Mar-2020 18:37
Acenaphthylene	< 0.000016		0.000016	0.00011	mg/L	1	17-Mar-2020 18:37
Anthracene	< 0.000015		0.000015	0.00011	mg/L	1	17-Mar-2020 18:37
Benz(a)anthracene	< 0.000053		0.000053	0.00011	mg/L	1	17-Mar-2020 18:37
Benzo(a)pyrene	< 0.000021		0.000021	0.00011	mg/L	1	17-Mar-2020 18:37
Bis(2-chloroethoxy)methane	< 0.000032		0.000032	0.00021	mg/L	1	17-Mar-2020 18:37
Bis(2-ethylhexyl)phthalate	< 0.000039		0.000039	0.00021	mg/L	1	17-Mar-2020 18:37
Chrysene	< 0.000022		0.000022	0.00011	mg/L	1	17-Mar-2020 18:37
Dibenzofuran	< 0.000021		0.000021	0.00011	mg/L	1	17-Mar-2020 18:37
<b>Di-n-butyl phthalate</b>	<b>0.000031</b>	J	<b>0.000021</b>	<b>0.00021</b>	<b>mg/L</b>	1	17-Mar-2020 18:37
Fluoranthene	< 0.000011		0.000011	0.00011	mg/L	1	17-Mar-2020 18:37
Fluorene	< 0.000032		0.000032	0.00011	mg/L	1	17-Mar-2020 18:37
<b>Naphthalene</b>	<b>0.000083</b>	J	<b>0.000021</b>	<b>0.00011</b>	<b>mg/L</b>	1	17-Mar-2020 18:37
Nitrobenzene	< 0.000025		0.000025	0.00021	mg/L	1	17-Mar-2020 18:37
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00021	mg/L	1	17-Mar-2020 18:37
Pentachlorophenol	< 0.000083		0.000083	0.00021	mg/L	1	17-Mar-2020 18:37
Phenanthrene	< 0.000022		0.000022	0.00011	mg/L	1	17-Mar-2020 18:37
Phenol	< 0.000037		0.000037	0.00021	mg/L	1	17-Mar-2020 18:37
Pyrene	< 0.000020		0.000020	0.00011	mg/L	1	17-Mar-2020 18:37
<i>Surr: 2,4,6-Tribromophenol</i>	<i>99.1</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 18:37</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>69.9</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 18:37</i>
<i>Surr: 2-Fluorophenol</i>	<i>67.9</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 18:37</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>100.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 18:37</i>
<i>Surr: Nitrobenzene-d5</i>	<i>62.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 18:37</i>
<i>Surr: Phenol-d6</i>	<i>68.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 18:37</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00989</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Mar-2020 18:32

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200312  
 Collection Date: 12-Mar-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-10  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 03:39
<b>Benzene</b>	<b>0.0089</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 03:39
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 03:39
<b>Ethylbenzene</b>	<b>0.052</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 03:39
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 03:39
<b>Toluene</b>	<b>0.032</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 03:39
<b>Xylenes, Total</b>	<b>0.061</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 03:39
Surr: 1,2-Dichloroethane-d4	103			70-126	%REC	1	15-Mar-2020 03:39
Surr: 4-Bromofluorobenzene	102			81-113	%REC	1	15-Mar-2020 03:39
Surr: Dibromofluoromethane	98.7			77-123	%REC	1	15-Mar-2020 03:39
Surr: Toluene-d8	101			82-127	%REC	1	15-Mar-2020 03:39

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200312  
 Collection Date: 12-Mar-2020 00:00

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-10  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 18:56
2,4-Dimethylphenol	< 0.000041		0.000041	0.00021	mg/L	1	17-Mar-2020 18:56
2,4-Dinitrotoluene	< 0.000060		0.000060	0.00021	mg/L	1	17-Mar-2020 18:56
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00021	mg/L	1	17-Mar-2020 18:56
2-Chloronaphthalene	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 18:56
<b>2-Methylnaphthalene</b>	<b>0.098</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 13:00
4,6-Dinitro-2-methylphenol	< 0.000021		0.000021	0.00021	mg/L	1	17-Mar-2020 18:56
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	17-Mar-2020 18:56
<b>Acenaphthene</b>	<b>0.087</b>		<b>0.00028</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 13:00
<b>Acenaphthylene</b>	<b>0.00074</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<b>Anthracene</b>	<b>0.0082</b>		<b>0.000014</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<b>Benz(a)anthracene</b>	<b>0.000079</b>	J	<b>0.000052</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
Benzo(a)pyrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 18:56
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00021	mg/L	1	17-Mar-2020 18:56
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000052</b>	J	<b>0.000038</b>	<b>0.00021</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<b>Chrysene</b>	<b>0.000061</b>	J	<b>0.000022</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<b>Dibenzofuran</b>	<b>0.078</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 13:00
<b>Di-n-butyl phthalate</b>	<b>0.00010</b>	J	<b>0.000021</b>	<b>0.00021</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<b>Fluoranthene</b>	<b>0.0053</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<b>Fluorene</b>	<b>0.038</b>		<b>0.00031</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 13:00
<b>Naphthalene</b>	<b>2.5</b>		<b>0.021</b>	<b>0.10</b>	<b>mg/L</b>	1000	18-Mar-2020 18:45
Nitrobenzene	< 0.000025		0.000025	0.00021	mg/L	1	17-Mar-2020 18:56
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00021	mg/L	1	17-Mar-2020 18:56
Pentachlorophenol	< 0.000081		0.000081	0.00021	mg/L	1	17-Mar-2020 18:56
<b>Phenanthrene</b>	<b>0.054</b>		<b>0.00022</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 13:00
Phenol	< 0.000036		0.000036	0.00021	mg/L	1	17-Mar-2020 18:56
<b>Pyrene</b>	<b>0.0029</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 18:56
<i>Surr: 2,4,6-Tribromophenol</i>	79.3			34-129	%REC	10	18-Mar-2020 13:00
<i>Surr: 2,4,6-Tribromophenol</i>	89.9			34-129	%REC	1	17-Mar-2020 18:56
<i>Surr: 2,4,6-Tribromophenol</i>	0	JS		34-129	%REC	1000	18-Mar-2020 18:45
<i>Surr: 2-Fluorobiphenyl</i>	0	JS		40-125	%REC	1000	18-Mar-2020 18:45
<i>Surr: 2-Fluorobiphenyl</i>	57.7			40-125	%REC	1	17-Mar-2020 18:56
<i>Surr: 2-Fluorobiphenyl</i>	54.1			40-125	%REC	10	18-Mar-2020 13:00
<i>Surr: 2-Fluorophenol</i>	62.7			20-120	%REC	10	18-Mar-2020 13:00
<i>Surr: 2-Fluorophenol</i>	68.7			20-120	%REC	1	17-Mar-2020 18:56
<i>Surr: 2-Fluorophenol</i>	0	JS		20-120	%REC	1000	18-Mar-2020 18:45
<i>Surr: 4-Terphenyl-d14</i>	0	JS		40-135	%REC	1000	18-Mar-2020 18:45
<i>Surr: 4-Terphenyl-d14</i>	89.9			40-135	%REC	1	17-Mar-2020 18:56
<i>Surr: 4-Terphenyl-d14</i>	79.6			40-135	%REC	10	18-Mar-2020 13:00

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD01-20200312  
 Collection Date: 12-Mar-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-10  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	55.8			41-120	%REC	10	18-Mar-2020 13:00
Surr: Nitrobenzene-d5	44.9			41-120	%REC	1	17-Mar-2020 18:56
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	1000	18-Mar-2020 18:45
Surr: Phenol-d6	0	JS		20-120	%REC	1000	18-Mar-2020 18:45
Surr: Phenol-d6	62.1			20-120	%REC	1	17-Mar-2020 18:56
Surr: Phenol-d6	57.1			20-120	%REC	10	18-Mar-2020 13:00
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
Arsenic	0.00470		0.000400	0.00200	mg/L	1	20-Mar-2020 18:35

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB02-20200312  
 Collection Date: 12-Mar-2020 12:36

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-11  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 23:48
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 23:48
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 23:48
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 23:48
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	14-Mar-2020 23:48
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	14-Mar-2020 23:48
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	14-Mar-2020 23:48
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>103</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 23:48</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 23:48</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 23:48</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>14-Mar-2020 23:48</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76B-20200312  
 Collection Date: 12-Mar-2020 10:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-12  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:03
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:03
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:03
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:03
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 04:03
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:03
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:03
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:03</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>100</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:03</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.1</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:03</i>
<i>Surr: Toluene-d8</i>	<i>98.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:03</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76B-20200312  
 Collection Date: 12-Mar-2020 10:35

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-12  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 19:16
2,4-Dimethylphenol	< 0.000042		0.000042	0.00021	mg/L	1	17-Mar-2020 19:16
2,4-Dinitrotoluene	< 0.000060		0.000060	0.00021	mg/L	1	17-Mar-2020 19:16
2,6-Dinitrotoluene	< 0.000044		0.000044	0.00021	mg/L	1	17-Mar-2020 19:16
2-Chloronaphthalene	< 0.000022		0.000022	0.00021	mg/L	1	17-Mar-2020 19:16
2-Methylnaphthalene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:16
4,6-Dinitro-2-methylphenol	< 0.000021		0.000021	0.00021	mg/L	1	17-Mar-2020 19:16
4-Nitrophenol	< 0.000049		0.000049	0.0010	mg/L	1	17-Mar-2020 19:16
Acenaphthene	< 0.000028		0.000028	0.00010	mg/L	1	17-Mar-2020 19:16
Acenaphthylene	< 0.000016		0.000016	0.00010	mg/L	1	17-Mar-2020 19:16
Anthracene	< 0.000015		0.000015	0.00010	mg/L	1	17-Mar-2020 19:16
Benz(a)anthracene	< 0.000052		0.000052	0.00010	mg/L	1	17-Mar-2020 19:16
Benzo(a)pyrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 19:16
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00021	mg/L	1	17-Mar-2020 19:16
Bis(2-ethylhexyl)phthalate	< 0.000039		0.000039	0.00021	mg/L	1	17-Mar-2020 19:16
Chrysene	< 0.000022		0.000022	0.00010	mg/L	1	17-Mar-2020 19:16
Dibenzofuran	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 19:16
<b>Di-n-butyl phthalate</b>	<b>0.000025</b>	J	<b>0.000021</b>	<b>0.00021</b>	<b>mg/L</b>	1	17-Mar-2020 19:16
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	17-Mar-2020 19:16
Fluorene	< 0.000031		0.000031	0.00010	mg/L	1	17-Mar-2020 19:16
<b>Naphthalene</b>	<b>0.000096</b>	J	<b>0.000021</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 19:16
Nitrobenzene	< 0.000025		0.000025	0.00021	mg/L	1	17-Mar-2020 19:16
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00021	mg/L	1	17-Mar-2020 19:16
Pentachlorophenol	< 0.000082		0.000082	0.00021	mg/L	1	17-Mar-2020 19:16
Phenanthrene	< 0.000022		0.000022	0.00010	mg/L	1	17-Mar-2020 19:16
Phenol	< 0.000036		0.000036	0.00021	mg/L	1	17-Mar-2020 19:16
Pyrene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:16
<i>Surr: 2,4,6-Tribromophenol</i>	<i>81.4</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:16</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>63.8</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:16</i>
<i>Surr: 2-Fluorophenol</i>	<i>60.0</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:16</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>91.0</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:16</i>
<i>Surr: Nitrobenzene-d5</i>	<i>55.6</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:16</i>
<i>Surr: Phenol-d6</i>	<i>63.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:16</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW76B-20200312  
 Collection Date: 12-Mar-2020 10:35

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-12  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>			Prep:TX1005PR / 17-Mar-2020		Analyst: MBG
nC6 to nC12	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 09:42
>nC12 to nC28	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 09:42
>nC28 to nC35	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 09:42
Total Petroleum Hydrocarbon	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 09:42
Surr: 2-Fluorobiphenyl	82.2			70-130	%REC	1	18-Mar-2020 09:42
Surr: Trifluoromethyl benzene	96.2			70-130	%REC	1	18-Mar-2020 09:42
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 18-Mar-2020		Analyst: JHD
Arsenic	0.00115	J	0.000400	0.00200	mg/L	1	20-Mar-2020 18:37

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW97A-20200312  
 Collection Date: 12-Mar-2020 11:47

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-13  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:27
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:27
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:27
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:27
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 04:27
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:27
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:27
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:27</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>94.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:27</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.7</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:27</i>
<i>Surr: Toluene-d8</i>	<i>99.4</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:27</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW97A-20200312  
 Collection Date: 12-Mar-2020 11:47

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-13  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 19:35
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	17-Mar-2020 19:35
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	17-Mar-2020 19:35
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	17-Mar-2020 19:35
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 19:35
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 19:35
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 19:35
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	17-Mar-2020 19:35
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	17-Mar-2020 19:35
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	17-Mar-2020 19:35
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	17-Mar-2020 19:35
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	17-Mar-2020 19:35
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:35
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	17-Mar-2020 19:35
Bis(2-ethylhexyl)phthalate	< 0.000037		0.000037	0.00020	mg/L	1	17-Mar-2020 19:35
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 19:35
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:35
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 19:35
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	17-Mar-2020 19:35
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	17-Mar-2020 19:35
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:35
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	17-Mar-2020 19:35
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	17-Mar-2020 19:35
Pentachlorophenol	< 0.000080		0.000080	0.00020	mg/L	1	17-Mar-2020 19:35
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 19:35
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	17-Mar-2020 19:35
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 19:35
<i>Surr: 2,4,6-Tribromophenol</i>	<i>87.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:35</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>64.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:35</i>
<i>Surr: 2-Fluorophenol</i>	<i>60.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:35</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.3</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:35</i>
<i>Surr: Nitrobenzene-d5</i>	<i>54.5</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:35</i>
<i>Surr: Phenol-d6</i>	<i>59.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:35</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW97A-20200312  
 Collection Date: 12-Mar-2020 11:47

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-13  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>			Prep:TX1005PR / 17-Mar-2020		Analyst: PVL
nC6 to nC12	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 18:40
>nC12 to nC28	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 18:40
>nC28 to nC35	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 18:40
Total Petroleum Hydrocarbon	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 18:40
Surr: 2-Fluorobiphenyl	80.6			70-130	%REC	1	18-Mar-2020 18:40
Surr: Trifluoromethyl benzene	92.4			70-130	%REC	1	18-Mar-2020 18:40
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>			Prep:SW3010A / 18-Mar-2020		Analyst: JHD
Arsenic	0.00121	J	0.000400	0.00200	mg/L	1	20-Mar-2020 18:39

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW98A-20200312  
 Collection Date: 12-Mar-2020 12:54

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-14  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:51
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 04:51
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:51
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:51
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 04:51
<b>Toluene</b>	<b>0.00065</b>	J	<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 04:51
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 04:51
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>100.0</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:51</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:51</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:51</i>
<i>Surr: Toluene-d8</i>	<i>98.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 04:51</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW98A-20200312  
 Collection Date: 12-Mar-2020 12:54

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-14  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	ML	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 19:54
<b>2,4-Dimethylphenol</b>	<b>0.000078</b>	J	<b>0.000041</b>	<b>0.00020</b>	<b>mg/L</b>	1	17-Mar-2020 19:54
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	17-Mar-2020 19:54
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00020	mg/L	1	17-Mar-2020 19:54
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 19:54
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 19:54
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 19:54
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	17-Mar-2020 19:54
<b>Acenaphthene</b>	<b>0.000040</b>	J	<b>0.000028</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 19:54
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	17-Mar-2020 19:54
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	17-Mar-2020 19:54
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	17-Mar-2020 19:54
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:54
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00020	mg/L	1	17-Mar-2020 19:54
Bis(2-ethylhexyl)phthalate	< 0.000038		0.000038	0.00020	mg/L	1	17-Mar-2020 19:54
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 19:54
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 19:54
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 19:54
<b>Fluoranthene</b>	<b>0.000018</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 19:54
Fluorene	< 0.000031		0.000031	0.00010	mg/L	1	17-Mar-2020 19:54
<b>Naphthalene</b>	<b>0.000042</b>	J	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	17-Mar-2020 19:54
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	17-Mar-2020 19:54
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00020	mg/L	1	17-Mar-2020 19:54
Pentachlorophenol	< 0.000081		0.000081	0.00020	mg/L	1	17-Mar-2020 19:54
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 19:54
Phenol	< 0.000036		0.000036	0.00020	mg/L	1	17-Mar-2020 19:54
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 19:54
<i>Surr: 2,4,6-Tribromophenol</i>	<i>98.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:54</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>51.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:54</i>
<i>Surr: 2-Fluorophenol</i>	<i>54.6</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:54</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>99.4</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:54</i>
<i>Surr: Nitrobenzene-d5</i>	<i>52.1</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:54</i>
<i>Surr: Phenol-d6</i>	<i>56.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 19:54</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW98A-20200312  
 Collection Date: 12-Mar-2020 12:54

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-14  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 17-Mar-2020		Analyst: PVL	
nC6 to nC12	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 20:10
<b>&gt;nC12 to nC28</b>	<b>1.0</b>		<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	18-Mar-2020 20:10
>nC28 to nC35	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 20:10
<b>Total Petroleum Hydrocarbon</b>	<b>1.00</b>		<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	18-Mar-2020 20:10
Surr: 2-Fluorobiphenyl	93.3			70-130	%REC	1	18-Mar-2020 20:10
Surr: Trifluoromethyl benzene	92.4			70-130	%REC	1	18-Mar-2020 20:10
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00963</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Mar-2020 18:50

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW98B-20200312  
 Collection Date: 12-Mar-2020 13:53

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-15  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 05:15
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 05:15
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 05:15
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 05:15
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 05:15
<b>Toluene</b>	<b>0.0020</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 05:15
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 05:15
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:15</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:15</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:15</i>
<i>Surr: Toluene-d8</i>	<i>99.2</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:15</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW98B-20200312  
 Collection Date: 12-Mar-2020 13:53

**ANALYTICAL REPORT**

WorkOrder:HS20030619  
 Lab ID:HS20030619-15  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 20:13
2,4-Dimethylphenol	< 0.000041		0.000041	0.00020	mg/L	1	17-Mar-2020 20:13
2,4-Dinitrotoluene	< 0.000059		0.000059	0.00020	mg/L	1	17-Mar-2020 20:13
2,6-Dinitrotoluene	< 0.000043		0.000043	0.00020	mg/L	1	17-Mar-2020 20:13
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	17-Mar-2020 20:13
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 20:13
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 20:13
4-Nitrophenol	< 0.000048		0.000048	0.0010	mg/L	1	17-Mar-2020 20:13
Acenaphthene	< 0.000028		0.000028	0.00010	mg/L	1	17-Mar-2020 20:13
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	17-Mar-2020 20:13
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	17-Mar-2020 20:13
Benz(a)anthracene	< 0.000051		0.000051	0.00010	mg/L	1	17-Mar-2020 20:13
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 20:13
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00020	mg/L	1	17-Mar-2020 20:13
Bis(2-ethylhexyl)phthalate	< 0.000038		0.000038	0.00020	mg/L	1	17-Mar-2020 20:13
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 20:13
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	17-Mar-2020 20:13
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	17-Mar-2020 20:13
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	17-Mar-2020 20:13
Fluorene	< 0.000031		0.000031	0.00010	mg/L	1	17-Mar-2020 20:13
<b>Naphthalene</b>	<b>0.000036</b>	<b>J</b>	<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	<b>1</b>	<b>17-Mar-2020 20:13</b>
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	17-Mar-2020 20:13
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00020	mg/L	1	17-Mar-2020 20:13
Pentachlorophenol	< 0.000081		0.000081	0.00020	mg/L	1	17-Mar-2020 20:13
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	17-Mar-2020 20:13
Phenol	< 0.000036		0.000036	0.00020	mg/L	1	17-Mar-2020 20:13
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	17-Mar-2020 20:13
<i>Surr: 2,4,6-Tribromophenol</i>	<i>67.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 20:13</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>50.4</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 20:13</i>
<i>Surr: 2-Fluorophenol</i>	<i>50.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 20:13</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>87.1</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 20:13</i>
<i>Surr: Nitrobenzene-d5</i>	<i>46.7</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 20:13</i>
<i>Surr: Phenol-d6</i>	<i>52.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>17-Mar-2020 20:13</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW98B-20200312  
 Collection Date: 12-Mar-2020 13:53

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-15  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 17-Mar-2020		Analyst: PVL	
nC6 to nC12	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 20:39
<b>&gt;nC12 to nC28</b>	<b>0.38</b>	J	<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	18-Mar-2020 20:39
>nC28 to nC35	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 20:39
<b>Total Petroleum Hydrocarbon</b>	<b>0.380</b>	J	<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	18-Mar-2020 20:39
Surr: 2-Fluorobiphenyl	86.1			70-130	%REC	1	18-Mar-2020 20:39
Surr: Trifluoromethyl benzene	95.2			70-130	%REC	1	18-Mar-2020 20:39
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00165</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	20-Mar-2020 18:52

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50B-20200312  
 Collection Date: 12-Mar-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-16  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 05:39
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	15-Mar-2020 05:39
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	15-Mar-2020 05:39
<b>Ethylbenzene</b>	<b>0.0068</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 05:39
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	15-Mar-2020 05:39
<b>Toluene</b>	<b>0.0033</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 05:39
<b>Xylenes, Total</b>	<b>0.0056</b>		<b>0.00030</b>	<b>0.0010</b>	<b>mg/L</b>	1	15-Mar-2020 05:39
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:39</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:39</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:39</i>
<i>Surr: Toluene-d8</i>	<i>98.9</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>15-Mar-2020 05:39</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50B-20200312  
 Collection Date: 12-Mar-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-16  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000022		0.000022	0.00021	mg/L	1	18-Mar-2020 17:09
2,4-Dimethylphenol	< 0.000042		0.000042	0.00021	mg/L	1	18-Mar-2020 17:09
2,4-Dinitrotoluene	< 0.000060		0.000060	0.00021	mg/L	1	18-Mar-2020 17:09
2,6-Dinitrotoluene	< 0.000044		0.000044	0.00021	mg/L	1	18-Mar-2020 17:09
2-Chloronaphthalene	< 0.000022		0.000022	0.00021	mg/L	1	18-Mar-2020 17:09
<b>2-Methylnaphthalene</b>	<b>0.016</b>		<b>0.00020</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 17:29
4,6-Dinitro-2-methylphenol	< 0.000021		0.000021	0.00021	mg/L	1	18-Mar-2020 17:09
4-Nitrophenol	< 0.000049		0.000049	0.0010	mg/L	1	18-Mar-2020 17:09
<b>Acenaphthene</b>	<b>0.020</b>		<b>0.00028</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 17:29
<b>Acenaphthylene</b>	<b>0.00028</b>		<b>0.000016</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Mar-2020 17:09
<b>Anthracene</b>	<b>0.0016</b>		<b>0.000015</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Mar-2020 17:09
Benz(a)anthracene	< 0.000052		0.000052	0.00010	mg/L	1	18-Mar-2020 17:09
Benzo(a)pyrene	< 0.000021		0.000021	0.00010	mg/L	1	18-Mar-2020 17:09
Bis(2-chloroethoxy)methane	< 0.000031		0.000031	0.00021	mg/L	1	18-Mar-2020 17:09
Bis(2-ethylhexyl)phthalate	< 0.000039		0.000039	0.00021	mg/L	1	18-Mar-2020 17:09
Chrysene	< 0.000022		0.000022	0.00010	mg/L	1	18-Mar-2020 17:09
<b>Dibenzofuran</b>	<b>0.015</b>		<b>0.00021</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 17:29
<b>Di-n-butyl phthalate</b>	<b>0.000038</b>	J	<b>0.000021</b>	<b>0.00021</b>	<b>mg/L</b>	1	18-Mar-2020 17:09
<b>Fluoranthene</b>	<b>0.0011</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Mar-2020 17:09
<b>Fluorene</b>	<b>0.011</b>		<b>0.00031</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 17:29
<b>Naphthalene</b>	<b>0.31</b>		<b>0.0021</b>	<b>0.010</b>	<b>mg/L</b>	100	18-Mar-2020 17:48
Nitrobenzene	< 0.000025		0.000025	0.00021	mg/L	1	18-Mar-2020 17:09
N-Nitrosodiphenylamine	< 0.000026		0.000026	0.00021	mg/L	1	18-Mar-2020 17:09
Pentachlorophenol	< 0.000082		0.000082	0.00021	mg/L	1	18-Mar-2020 17:09
<b>Phenanthrene</b>	<b>0.016</b>		<b>0.00022</b>	<b>0.0010</b>	<b>mg/L</b>	10	18-Mar-2020 17:29
Phenol	< 0.000036		0.000036	0.00021	mg/L	1	18-Mar-2020 17:09
<b>Pyrene</b>	<b>0.00059</b>		<b>0.000020</b>	<b>0.00010</b>	<b>mg/L</b>	1	18-Mar-2020 17:09
Surr: 2,4,6-Tribromophenol	84.4			34-129	%REC	1	18-Mar-2020 17:09
Surr: 2,4,6-Tribromophenol	71.2			34-129	%REC	10	18-Mar-2020 17:29
Surr: 2,4,6-Tribromophenol	0	JS		34-129	%REC	100	18-Mar-2020 17:48
Surr: 2-Fluorobiphenyl	44.0			40-125	%REC	1	18-Mar-2020 17:09
Surr: 2-Fluorobiphenyl	47.0			40-125	%REC	10	18-Mar-2020 17:29
Surr: 2-Fluorobiphenyl	0	JS		40-125	%REC	100	18-Mar-2020 17:48
Surr: 2-Fluorophenol	47.4			20-120	%REC	1	18-Mar-2020 17:09
Surr: 2-Fluorophenol	49.0			20-120	%REC	10	18-Mar-2020 17:29
Surr: 2-Fluorophenol	0	JS		20-120	%REC	100	18-Mar-2020 17:48
Surr: 4-Terphenyl-d14	80.2			40-135	%REC	1	18-Mar-2020 17:09
Surr: 4-Terphenyl-d14	73.5			40-135	%REC	10	18-Mar-2020 17:29
Surr: 4-Terphenyl-d14	0	JS		40-135	%REC	100	18-Mar-2020 17:48

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW50B-20200312  
 Collection Date: 12-Mar-2020 15:20

**ANALYTICAL REPORT**  
 WorkOrder:HS20030619  
 Lab ID:HS20030619-16  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 16-Mar-2020		Analyst: GEY	
Surr: Nitrobenzene-d5	0	JS		41-120	%REC	100	18-Mar-2020 17:48
Surr: Nitrobenzene-d5	45.0			41-120	%REC	1	18-Mar-2020 17:09
Surr: Nitrobenzene-d5	43.1			41-120	%REC	10	18-Mar-2020 17:29
Surr: Phenol-d6	48.9			20-120	%REC	1	18-Mar-2020 17:09
Surr: Phenol-d6	52.5			20-120	%REC	10	18-Mar-2020 17:29
Surr: Phenol-d6	0	JS		20-120	%REC	100	18-Mar-2020 17:48
<b>LOW-LEVEL TEXAS TPH BY TX1005</b>		<b>Method:TX1005</b>		Prep:TX1005PR / 17-Mar-2020		Analyst: PVL	
nC6 to nC12	1.0		0.19	0.48	mg/L	1	18-Mar-2020 21:09
>nC12 to nC28	1.1		0.19	0.48	mg/L	1	18-Mar-2020 21:09
>nC28 to nC35	< 0.19		0.19	0.48	mg/L	1	18-Mar-2020 21:09
<b>Total Petroleum Hydrocarbon</b>	<b>2.10</b>		<b>0.19</b>	<b>0.48</b>	<b>mg/L</b>	1	18-Mar-2020 21:09
Surr: 2-Fluorobiphenyl	101			70-130	%REC	1	18-Mar-2020 21:09
Surr: Trifluoromethyl benzene	95.2			70-130	%REC	1	18-Mar-2020 21:09
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 18-Mar-2020		Analyst: JHD	
Arsenic	0.00372		0.000400	0.00200	mg/L	1	20-Mar-2020 18:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**Batch ID:** 151716      **Start Date:** 16 Mar 2020 07:00      **End Date:** 16 Mar 2020 12:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030619-02	1	980 (mL)	1 (mL)	0.00102
HS20030619-03	1	990 (mL)	1 (mL)	0.00101
HS20030619-04	1	970 (mL)	1 (mL)	0.001031
HS20030619-05	1	980 (mL)	1 (mL)	0.00102
HS20030619-06	1	990 (mL)	1 (mL)	0.00101
HS20030619-07	1	980 (mL)	1 (mL)	0.00102
HS20030619-08	1	960 (mL)	1 (mL)	0.001042
HS20030619-09	1	950 (mL)	1 (mL)	0.001053
HS20030619-10	1	970 (mL)	1 (mL)	0.001031
HS20030619-12	1	960 (mL)	1 (mL)	0.001042
HS20030619-13	1	990 (mL)	1 (mL)	0.00101
HS20030619-14	1	980 (mL)	1 (mL)	0.00102
HS20030619-15	1	980 (mL)	1 (mL)	0.00102
HS20030619-16	1	960 (mL)	1 (mL)	0.001042

**Batch ID:** 151763      **Start Date:** 17 Mar 2020 10:00      **End Date:** 17 Mar 2020 14:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030619-02		10 (mL)	10 (mL)	1
HS20030619-03		10 (mL)	10 (mL)	1
HS20030619-04		10 (mL)	10 (mL)	1
HS20030619-05		10 (mL)	10 (mL)	1
HS20030619-06		10 (mL)	10 (mL)	1

**Batch ID:** 151785      **Start Date:** 17 Mar 2020 11:00      **End Date:** 17 Mar 2020 14:30  
**Method:** TX 1005 PREP      **Prep Code:** TX 1005\_W PR

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030619-12	1	31.11 (g)	3 (mL)	0.09643

**Batch ID:** 151787      **Start Date:** 17 Mar 2020 12:00      **End Date:** 17 Mar 2020 15:30  
**Method:** TX 1005 PREP      **Prep Code:** TX 1005\_W PR

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030619-13	1	31.15 (g)	3 (mL)	0.09631
HS20030619-14	1	31.19 (g)	3 (mL)	0.09618
HS20030619-15	1	31.11 (g)	3 (mL)	0.09643
HS20030619-16	1	31.04 (g)	3 (mL)	0.09665

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**Batch ID:** 151813      **Start Date:** 18 Mar 2020 10:30      **End Date:** 18 Mar 2020 14:30  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030619-07		10 (mL)	10 (mL)	1
HS20030619-08		10 (mL)	10 (mL)	1
HS20030619-09		10 (mL)	10 (mL)	1
HS20030619-10		10 (mL)	10 (mL)	1
HS20030619-12		10 (mL)	10 (mL)	1
HS20030619-13		10 (mL)	10 (mL)	1
HS20030619-14		10 (mL)	10 (mL)	1
HS20030619-15		10 (mL)	10 (mL)	1
HS20030619-16		10 (mL)	10 (mL)	1



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 151716 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Groundwater</b>	
HS20030619-02	WQ-1620-MW99C-20200311	11 Mar 2020 15:35		16 Mar 2020 11:48	16 Mar 2020 17:59	1
HS20030619-03	WG-1620-MW54B-20200311	11 Mar 2020 14:20		16 Mar 2020 11:48	16 Mar 2020 18:18	1
HS20030619-04	WG-1620-MW94A-20200312	12 Mar 2020 15:45		16 Mar 2020 11:48	16 Mar 2020 18:37	1
HS20030619-05	WG-1620-MW93B-20200312	12 Mar 2020 14:35		16 Mar 2020 11:48	16 Mar 2020 18:56	1
HS20030619-06	WG-1620-MW92B-20200312	12 Mar 2020 13:20		16 Mar 2020 11:48	16 Mar 2020 19:15	1
HS20030619-07	WG-1620-MW84A-20200312	12 Mar 2020 12:15		16 Mar 2020 11:48	17 Mar 2020 17:59	1
HS20030619-08	WG-1620-MW70C-20200312	12 Mar 2020 10:35		16 Mar 2020 11:48	18 Mar 2020 18:26	1000
HS20030619-08	WG-1620-MW70C-20200312	12 Mar 2020 10:35		16 Mar 2020 11:48	18 Mar 2020 12:22	10
HS20030619-08	WG-1620-MW70C-20200312	12 Mar 2020 10:35		16 Mar 2020 11:48	17 Mar 2020 18:18	1
HS20030619-09	WG-1620-MW91A-20200312	12 Mar 2020 09:20		16 Mar 2020 11:48	17 Mar 2020 18:37	1
HS20030619-10	WG-1620-FD01-20200312	12 Mar 2020 00:00		16 Mar 2020 11:48	18 Mar 2020 18:45	1000
HS20030619-10	WG-1620-FD01-20200312	12 Mar 2020 00:00		16 Mar 2020 11:48	18 Mar 2020 13:00	10
HS20030619-10	WG-1620-FD01-20200312	12 Mar 2020 00:00		16 Mar 2020 11:48	17 Mar 2020 18:56	1
HS20030619-12	WG-1620-MW76B-20200312	12 Mar 2020 10:35		16 Mar 2020 11:48	17 Mar 2020 19:16	1
HS20030619-13	WG-1620-MW97A-20200312	12 Mar 2020 11:47		16 Mar 2020 11:48	17 Mar 2020 19:35	1
HS20030619-14	WG-1620-MW98A-20200312	12 Mar 2020 12:54		16 Mar 2020 07:00	17 Mar 2020 19:54	1
HS20030619-15	WG-1620-MW98B-20200312	12 Mar 2020 13:53		16 Mar 2020 11:48	17 Mar 2020 20:13	1
HS20030619-16	WG-1620-MW50B-20200312	12 Mar 2020 15:20		16 Mar 2020 11:48	18 Mar 2020 17:48	100
HS20030619-16	WG-1620-MW50B-20200312	12 Mar 2020 15:20		16 Mar 2020 11:48	18 Mar 2020 17:29	10
HS20030619-16	WG-1620-MW50B-20200312	12 Mar 2020 15:20		16 Mar 2020 11:48	18 Mar 2020 17:09	1
<b>Batch ID: 151763 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Groundwater</b>	
HS20030619-02	WQ-1620-MW99C-20200311	11 Mar 2020 15:35		17 Mar 2020 14:00	18 Mar 2020 18:20	1
HS20030619-03	WG-1620-MW54B-20200311	11 Mar 2020 14:20		17 Mar 2020 14:00	18 Mar 2020 18:23	1
HS20030619-04	WG-1620-MW94A-20200312	12 Mar 2020 15:45		17 Mar 2020 14:00	18 Mar 2020 18:25	1
HS20030619-05	WG-1620-MW93B-20200312	12 Mar 2020 14:35		17 Mar 2020 14:00	18 Mar 2020 18:34	1
HS20030619-06	WG-1620-MW92B-20200312	12 Mar 2020 13:20		17 Mar 2020 14:00	18 Mar 2020 18:37	1
<b>Batch ID: 151785 ( 0 )</b>		<b>Test Name : LOW-LEVEL TEXAS TPH BY TX1005</b>			<b>Matrix: Groundwater</b>	
HS20030619-12	WG-1620-MW76B-20200312	12 Mar 2020 10:35		17 Mar 2020 11:00	18 Mar 2020 09:42	1
<b>Batch ID: 151787 ( 0 )</b>		<b>Test Name : LOW-LEVEL TEXAS TPH BY TX1005</b>			<b>Matrix: Groundwater</b>	
HS20030619-13	WG-1620-MW97A-20200312	12 Mar 2020 11:47		17 Mar 2020 12:00	18 Mar 2020 18:40	1
HS20030619-14	WG-1620-MW98A-20200312	12 Mar 2020 12:54		17 Mar 2020 12:00	18 Mar 2020 20:10	1
HS20030619-15	WG-1620-MW98B-20200312	12 Mar 2020 13:53		17 Mar 2020 12:00	18 Mar 2020 20:39	1
HS20030619-16	WG-1620-MW50B-20200312	12 Mar 2020 15:20		17 Mar 2020 12:00	18 Mar 2020 21:09	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 151813 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Groundwater</b>	
HS20030619-07	WG-1620-MW84A-20200312	12 Mar 2020 12:15		18 Mar 2020 14:30	20 Mar 2020 16:51	1
HS20030619-08	WG-1620-MW70C-20200312	12 Mar 2020 10:35		18 Mar 2020 14:30	20 Mar 2020 16:53	1
HS20030619-09	WG-1620-MW91A-20200312	12 Mar 2020 09:20		18 Mar 2020 14:30	20 Mar 2020 18:32	1
HS20030619-10	WG-1620-FD01-20200312	12 Mar 2020 00:00		18 Mar 2020 14:30	20 Mar 2020 18:35	1
HS20030619-12	WG-1620-MW76B-20200312	12 Mar 2020 10:35		18 Mar 2020 14:30	20 Mar 2020 18:37	1
HS20030619-13	WG-1620-MW97A-20200312	12 Mar 2020 11:47		18 Mar 2020 14:30	20 Mar 2020 18:39	1
HS20030619-14	WG-1620-MW98A-20200312	12 Mar 2020 12:54		18 Mar 2020 14:30	20 Mar 2020 18:50	1
HS20030619-15	WG-1620-MW98B-20200312	12 Mar 2020 13:53		18 Mar 2020 14:30	20 Mar 2020 18:52	1
HS20030619-16	WG-1620-MW50B-20200312	12 Mar 2020 15:20		18 Mar 2020 14:30	20 Mar 2020 18:55	1
<b>Batch ID: R358243 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Groundwater</b>	
HS20030619-02	WQ-1620-MW99C-20200311	11 Mar 2020 15:35			14 Mar 2020 18:05	1
HS20030619-03	WG-1620-MW54B-20200311	11 Mar 2020 14:20			14 Mar 2020 18:30	1
HS20030619-04	WG-1620-MW94A-20200312	12 Mar 2020 15:45			14 Mar 2020 18:54	1
HS20030619-05	WG-1620-MW93B-20200312	12 Mar 2020 14:35			14 Mar 2020 19:19	1
HS20030619-06	WG-1620-MW92B-20200312	12 Mar 2020 13:20			14 Mar 2020 19:44	1
<b>Batch ID: R358244 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Groundwater</b>	
HS20030619-07	WG-1620-MW84A-20200312	12 Mar 2020 12:15			15 Mar 2020 00:13	1
HS20030619-08	WG-1620-MW70C-20200312	12 Mar 2020 10:35			15 Mar 2020 01:50	1
HS20030619-09	WG-1620-MW91A-20200312	12 Mar 2020 09:20			15 Mar 2020 03:14	1
HS20030619-10	WG-1620-FD01-20200312	12 Mar 2020 00:00			15 Mar 2020 03:39	1
HS20030619-12	WG-1620-MW76B-20200312	12 Mar 2020 10:35			15 Mar 2020 04:03	1
HS20030619-13	WG-1620-MW97A-20200312	12 Mar 2020 11:47			15 Mar 2020 04:27	1
HS20030619-14	WG-1620-MW98A-20200312	12 Mar 2020 12:54			15 Mar 2020 04:51	1
HS20030619-15	WG-1620-MW98B-20200312	12 Mar 2020 13:53			15 Mar 2020 05:15	1
HS20030619-16	WG-1620-MW50B-20200312	12 Mar 2020 15:20			15 Mar 2020 05:39	1
<b>Batch ID: R358244 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20030619-01	WQ-1620-TB01-20200312	12 Mar 2020 16:30			15 Mar 2020 06:28	1
HS20030619-11	WQ-1620-TB02-20200312	12 Mar 2020 12:36			14 Mar 2020 23:48	1

WorkOrder: HS20030619  
 InstrumentID: FID-12  
 Test Code: TX1005\_W\_Low  
 Test Number: TX1005  
 Test Name: Low-level Texas TPH by TX1005

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	nC6 to nC12	TPH-1005-1	0.25	0.28	0.20	0.50
A	>nC12 to nC28	TPH-1005-2	0.25	0.29	0.20	0.50
A	>nC28 to nC35	TPH-1005-4	0.25	0.28	0.20	0.50
A	Total Petroleum Hydrocarbon	TPH	0.25	0.28	0.20	0.50
S	2-Fluorobiphenyl	321-60-8	0	0	0	0
S	Trifluoromethyl benzene	98-08-8	0	0	0	0

WorkOrder: HS20030619  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20030619  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20030619  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151785 ( 0 )		Instrument: FID-12		Method: LOW-LEVEL TEXAS TPH BY TX1005						
<b>MBLK</b>	Sample ID: <b>MBLK-151785</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 08:26</b>					
Client ID:		Run ID: <b>FID-12_358444</b>		SeqNo: <b>5519192</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	< 0.20	0.50								
>nC12 to nC28	< 0.20	0.50								
>nC28 to nC35	< 0.20	0.50								
Total Petroleum Hydrocarbon	< 0.20	0.50								
Surr: 2-Fluorobiphenyl	2.33	0	2.5	0	93.2	70 - 130				
Surr: Trifluoromethyl benzene	2.515	0	2.5	0	101	70 - 130				
<b>LCS</b>	Sample ID: <b>LCS-151785</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 09:42</b>					
Client ID:		Run ID: <b>FID-12_358444</b>		SeqNo: <b>5519193</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	25.3	0.50	25	0	101	75 - 125				
>nC12 to nC28	26.67	0.50	25	0	107	75 - 125				
Surr: 2-Fluorobiphenyl	2.381	0	2.5	0	95.3	70 - 130				
Surr: Trifluoromethyl benzene	2.604	0	2.5	0	104	70 - 130				
<b>LCSD</b>	Sample ID: <b>LCSD-151785</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 10:11</b>					
Client ID:		Run ID: <b>FID-12_358444</b>		SeqNo: <b>5519194</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	20.14	0.50	25	0	80.6	75 - 125	25.3	22.7	20	R
>nC12 to nC28	22.75	0.50	25	0	91.0	75 - 125	26.67	15.9	20	
Surr: 2-Fluorobiphenyl	2.285	0	2.5	0	91.4	70 - 130	2.381	4.13	20	
Surr: Trifluoromethyl benzene	2.419	0	2.5	0	96.8	70 - 130	2.604	7.37	20	
<b>MS</b>	Sample ID: <b>HS20030619-02MS</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 11:10</b>					
Client ID: <b>WQ-1620-MW99C-20200311</b>		Run ID: <b>FID-12_358444</b>		SeqNo: <b>5519196</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	24.63	0.48	24.19	0	102	75 - 125				
>nC12 to nC28	27.56	0.48	24.19	0	114	75 - 125				
Surr: 2-Fluorobiphenyl	2.389	0	2.419	0	98.8	70 - 130				
Surr: Trifluoromethyl benzene	2.553	0	2.419	0	106	70 - 130				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

**Batch ID:** 151785 ( 0 )      **Instrument:** FID-12      **Method:** LOW-LEVEL TEXAS TPH BY TX1005

**MSD**      Sample ID: **HS20030619-02MSD**      Units: **mg/L**      Analysis Date: **18-Mar-2020 11:39**  
**Client ID:** **WQ-1620-MW99C-20200311**      **Run ID:** **FID-12\_358444**      **SeqNo:** **5519197**      **PrepDate:** **17-Mar-2020**      **DF:** **1**  
**Analyte**      **Result**      **MQL**      **SPK Val**      **SPK Ref Value**      **%REC**      **Control Limit**      **RPD Ref Value**      **%RPD**      **RPD Limit**      **Qual**

nC6 to nC12	22.43	0.48	24.15	0	92.9	75 - 125	24.63	9.36	20
>nC12 to nC28	26.75	0.48	24.15	0	111	75 - 125	27.56	2.99	20
<i>Surr: 2-Fluorobiphenyl</i>	2.43	0	2.415	0	101	70 - 130	2.389	1.67	20
<i>Surr: Trifluoromethyl benzene</i>	2.55	0	2.415	0	106	70 - 130	2.553	0.118	20

**The following samples were analyzed in this batch:**

HS20030619-02	HS20030619-03	HS20030619-04	HS20030619-05
HS20030619-06	HS20030619-07	HS20030619-08	HS20030619-09
HS20030619-10	HS20030619-12		



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151787 ( 0 )		Instrument: FID-12		Method: LOW-LEVEL TEXAS TPH BY TX1005						
<b>MBLK</b>	Sample ID: <b>MBLK-151787</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 17:10</b>					
Client ID:		Run ID: <b>FID-12_358504</b>		SeqNo: <b>5520592</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	< 0.20	0.50								
>nC12 to nC28	< 0.20	0.50								
>nC28 to nC35	< 0.20	0.50								
Total Petroleum Hydrocarbon	< 0.20	0.50								
Surr: 2-Fluorobiphenyl	1.936	0	2.5	0	77.5	70 - 130				
Surr: Trifluoromethyl benzene	2.31	0	2.5	0	92.4	70 - 130				
<b>LCS</b>	Sample ID: <b>LCS-151787</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 17:40</b>					
Client ID:		Run ID: <b>FID-12_358504</b>		SeqNo: <b>5520593</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	20.43	0.50	25	0	81.7	75 - 125				
>nC12 to nC28	24.35	0.50	25	0	97.4	75 - 125				
Surr: 2-Fluorobiphenyl	2.281	0	2.5	0	91.2	70 - 130				
Surr: Trifluoromethyl benzene	2.412	0	2.5	0	96.5	70 - 130				
<b>LCSD</b>	Sample ID: <b>LCSD-151787</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 18:10</b>					
Client ID:		Run ID: <b>FID-12_358504</b>		SeqNo: <b>5520594</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	19.96	0.50	25	0	79.9	75 - 125	20.43	2.3	20	
>nC12 to nC28	23.88	0.50	25	0	95.5	75 - 125	24.35	1.96	20	
Surr: 2-Fluorobiphenyl	2.28	0	2.5	0	91.2	70 - 130	2.281	0.0255	20	
Surr: Trifluoromethyl benzene	2.401	0	2.5	0	96.0	70 - 130	2.412	0.47	20	
<b>MS</b>	Sample ID: <b>HS20030619-13MS</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 19:10</b>					
Client ID: <b>WG-1620-MW97A-20200312</b>		Run ID: <b>FID-12_358504</b>		SeqNo: <b>5520615</b>	PrepDate: <b>17-Mar-2020</b>	DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
nC6 to nC12	19.86	0.49	24.4	0	81.4	75 - 125				
>nC12 to nC28	22.87	0.49	24.4	0	93.7	75 - 125				
Surr: 2-Fluorobiphenyl	2.002	0	2.44	0	82.1	70 - 130				
Surr: Trifluoromethyl benzene	2.177	0	2.44	0	89.2	70 - 130				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

**Batch ID:** 151787 ( 0 )      **Instrument:** FID-12      **Method:** LOW-LEVEL TEXAS TPH BY TX1005

**MSD**      Sample ID: **HS20030619-13MSD**      Units: **mg/L**      Analysis Date: **18-Mar-2020 19:40**  
 Client ID: **WG-1620-MW97A-20200312**      Run ID: **FID-12\_358504**      SeqNo: **5520616**      PrepDate: **17-Mar-2020**      DF: **1**  
 Analyte      Result      MQL      SPK Val      SPK Ref Value      %REC      Control Limit      RPD Ref Value      %RPD      RPD Limit Qual

nC6 to nC12	21.86	0.49	24.54	0	89.1	75 - 125	19.86	9.62	20
>nC12 to nC28	26.03	0.49	24.54	0	106	75 - 125	22.87	12.9	20
Surr: 2-Fluorobiphenyl	2.344	0	2.454	0	95.5	70 - 130	2.002	15.7	20
Surr: Trifluoromethyl benzene	2.479	0	2.454	0	101	70 - 130	2.177	13	20

The following samples were analyzed in this batch: 

HS20030619-13	HS20030619-14	HS20030619-15	HS20030619-16
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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151763 ( 0 )		Instrument: ICPMS05			Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-151763</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 17:32</b>					
Client ID:		Run ID: <b>ICPMS05_358398</b>			SeqNo: <b>5519290</b>		PrepDate: <b>17-Mar-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	< 0.000400	0.00200								
<b>LCS</b>	Sample ID: <b>LCS-151763</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 17:34</b>					
Client ID:		Run ID: <b>ICPMS05_358398</b>			SeqNo: <b>5519291</b>		PrepDate: <b>17-Mar-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.04791	0.00200	0.05	0	95.8	80 - 120				
<b>MS</b>	Sample ID: <b>HS20030614-05MS</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 18:13</b>					
Client ID:		Run ID: <b>ICPMS05_358398</b>			SeqNo: <b>5519307</b>		PrepDate: <b>17-Mar-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.04448	0.00200	0.05	0.000142	88.7	80 - 120				
<b>MSD</b>	Sample ID: <b>HS20030614-05MSD</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 18:16</b>					
Client ID:		Run ID: <b>ICPMS05_358398</b>			SeqNo: <b>5519308</b>		PrepDate: <b>17-Mar-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.04389	0.00200	0.05	0.000142	87.5	80 - 120	0.04448	1.33	20	
<b>PDS</b>	Sample ID: <b>HS20030614-05PDS</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 18:18</b>					
Client ID:		Run ID: <b>ICPMS05_358398</b>			SeqNo: <b>5519309</b>		PrepDate: <b>17-Mar-2020</b>		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	0.1094	0.00200	0.1	0.000142	109	75 - 125				
<b>SD</b>	Sample ID: <b>HS20030614-05SD</b>	Units: <b>mg/L</b>			Analysis Date: <b>18-Mar-2020 18:11</b>					
Client ID:		Run ID: <b>ICPMS05_358398</b>			SeqNo: <b>5519306</b>		PrepDate: <b>17-Mar-2020</b>		DF: <b>5</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D	RPD Limit	Qual
Arsenic	< 0.00200	0.0100					0.000142	0	10	

The following samples were analyzed in this batch: HS20030619-02 HS20030619-03 HS20030619-04 HS20030619-05  
 HS20030619-06

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151813 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-151813</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Mar-2020 16:14</b>					
Client ID:	Run ID: <b>ICPMS05_358593</b>	SeqNo: <b>5523720</b>		PrepDate: <b>18-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	< 0.000400	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-151813</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Mar-2020 16:16</b>					
Client ID:	Run ID: <b>ICPMS05_358593</b>	SeqNo: <b>5523721</b>		PrepDate: <b>18-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.04715	0.00200	0.05	0	94.3	80 - 120			
<b>MS</b>	Sample ID: <b>HS20030654-04MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Mar-2020 17:49</b>					
Client ID:	Run ID: <b>ICPMS05_358593</b>	SeqNo: <b>5523737</b>		PrepDate: <b>18-Mar-2020</b>		DF: <b>5</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.04826	0.0100	0.05	0.000622	95.3	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20030654-04MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Mar-2020 17:51</b>					
Client ID:	Run ID: <b>ICPMS05_358593</b>	SeqNo: <b>5523738</b>		PrepDate: <b>18-Mar-2020</b>		DF: <b>5</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.04948	0.0100	0.05	0.000622	97.7	80 - 120	0.04826	2.49 20	
<b>PDS</b>	Sample ID: <b>HS20030654-04PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Mar-2020 17:53</b>					
Client ID:	Run ID: <b>ICPMS05_358593</b>	SeqNo: <b>5523739</b>		PrepDate: <b>18-Mar-2020</b>		DF: <b>5</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.6009	0.0100	0.5	0	120	75 - 125			
<b>SD</b>	Sample ID: <b>HS20030654-04SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>20-Mar-2020 17:46</b>					
Client ID:	Run ID: <b>ICPMS05_358593</b>	SeqNo: <b>5523736</b>		PrepDate: <b>18-Mar-2020</b>		DF: <b>25</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D %D Limit Qual	
Arsenic	< 0.0100	0.0500					0.000622	0 10	

The following samples were analyzed in this batch:

HS20030619-07	HS20030619-08	HS20030619-09	HS20030619-10
HS20030619-12	HS20030619-13	HS20030619-14	HS20030619-15
HS20030619-16			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151716 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-151716	Units: ug/L			Analysis Date: 16-Mar-2020 13:32					
Client ID:	Run ID: SV-7_358216	SeqNo: 5515267	PrepDate: 16-Mar-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	< 0.021	0.20								
2,4-Dimethylphenol	< 0.040	0.20								
2,4-Dinitrotoluene	< 0.058	0.20								
2,6-Dinitrotoluene	< 0.042	0.20								
2-Chloronaphthalene	< 0.021	0.20								
2-Methylnaphthalene	< 0.019	0.10								
4,6-Dinitro-2-methylphenol	< 0.020	0.20								
4-Nitrophenol	< 0.047	1.0								
Acenaphthene	< 0.027	0.10								
Acenaphthylene	< 0.015	0.10								
Anthracene	< 0.014	0.10								
Benz(a)anthracene	< 0.050	0.10								
Benzo(a)pyrene	< 0.020	0.10								
Bis(2-chloroethoxy)methane	< 0.030	0.20								
Bis(2-ethylhexyl)phthalate	< 0.037	0.20								
Chrysene	< 0.021	0.10								
Dibenzofuran	< 0.020	0.10								
Di-n-butyl phthalate	< 0.020	0.20								
Fluoranthene	< 0.010	0.10								
Fluorene	< 0.030	0.10								
Naphthalene	< 0.020	0.10								
Nitrobenzene	< 0.024	0.20								
N-Nitrosodiphenylamine	< 0.025	0.20								
Pentachlorophenol	< 0.079	0.20								
Phenanthrene	< 0.021	0.10								
Phenol	< 0.035	0.20								
Pyrene	< 0.019	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	4.07	0.20	5	0	81.4	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	3.367	0.20	5	0	67.3	40 - 125				
<i>Surr: 2-Fluorophenol</i>	3.058	0.20	5	0	61.2	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	4.326	0.20	5	0	86.5	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	3.169	0.20	5	0	63.4	41 - 120				
<i>Surr: Phenol-d6</i>	3.349	0.20	5	0	67.0	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151716 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-151716	Units: ug/L			Analysis Date: 16-Mar-2020 12:16					
Client ID:	Run ID: SV-7_358216	SeqNo: 5515265	PrepDate: 16-Mar-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	4.21	0.20	5	0	84.2	39 - 127				
2,4-Dimethylphenol	4.506	0.20	5	0	90.1	35 - 120				
2,4-Dinitrotoluene	5.519	0.20	5	0	110	50 - 122				
2,6-Dinitrotoluene	5.528	0.20	5	0	111	50 - 120				
2-Chloronaphthalene	5.308	0.20	5	0	106	50 - 120				
2-Methylnaphthalene	5.216	0.10	5	0	104	50 - 120				
4,6-Dinitro-2-methylphenol	4.232	0.20	5	0	84.6	25 - 121				
4-Nitrophenol	4.723	1.0	5	0	94.5	30 - 130				
Acenaphthene	4.719	0.10	5	0	94.4	45 - 120				
Acenaphthylene	4.688	0.10	5	0	93.8	47 - 120				
Anthracene	4.816	0.10	5	0	96.3	45 - 120				
Benz(a)anthracene	5.312	0.10	5	0	106	40 - 120				
Benzo(a)pyrene	5.285	0.10	5	0	106	45 - 120				
Bis(2-chloroethoxy)methane	5.043	0.20	5	0	101	45 - 120				
Bis(2-ethylhexyl)phthalate	6.344	0.20	5	0	127	40 - 139				
Chrysene	4.782	0.10	5	0	95.6	43 - 120				
Dibenzofuran	4.968	0.10	5	0	99.4	50 - 120				
Di-n-butyl phthalate	5.755	0.20	5	0	115	45 - 123				
Fluoranthene	5.205	0.10	5	0	104	45 - 125				
Fluorene	5.174	0.10	5	0	103	49 - 120				
Naphthalene	4.795	0.10	5	0	95.9	45 - 120				
Nitrobenzene	4.375	0.20	5	0	87.5	44 - 120				
N-Nitrosodiphenylamine	4.465	0.20	5	0	89.3	40 - 125				
Pentachlorophenol	4.311	0.20	5	0	86.2	19 - 121				
Phenanthrene	4.815	0.10	5	0	96.3	45 - 121				
Phenol	5.116	0.20	5	0	102	20 - 124				
Pyrene	5.307	0.10	5	0	106	40 - 130				
Surr: 2,4,6-Tribromophenol	5.449	0.20	5	0	109	34 - 129				
Surr: 2-Fluorobiphenyl	3.867	0.20	5	0	77.3	40 - 125				
Surr: 2-Fluorophenol	3.617	0.20	5	0	72.3	20 - 120				
Surr: 4-Terphenyl-d14	5.306	0.20	5	0	106	40 - 135				
Surr: Nitrobenzene-d5	3.465	0.20	5	0	69.3	41 - 120				
Surr: Phenol-d6	4.106	0.20	5	0	82.1	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

Batch ID: 151716 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-151716		Units: ug/L		Analysis Date: 16-Mar-2020 12:35				
Client ID:		Run ID: SV-7_358216		SeqNo: 5515266		PrepDate: 16-Mar-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Diphenylhydrazine	4.205	0.20	5	0	84.1	39 - 127	4.21	0.114	20	
2,4-Dimethylphenol	4.426	0.20	5	0	88.5	35 - 120	4.506	1.8	20	
2,4-Dinitrotoluene	5.542	0.20	5	0	111	50 - 122	5.519	0.403	20	
2,6-Dinitrotoluene	5.689	0.20	5	0	114	50 - 120	5.528	2.87	20	
2-Chloronaphthalene	5.269	0.20	5	0	105	50 - 120	5.308	0.732	20	
2-Methylnaphthalene	5.229	0.10	5	0	105	50 - 120	5.216	0.249	20	
4,6-Dinitro-2-methylphenol	4.262	0.20	5	0	85.2	25 - 121	4.232	0.702	30	
4-Nitrophenol	4.721	1.0	5	0	94.4	30 - 130	4.723	0.0308	20	
Acenaphthene	4.724	0.10	5	0	94.5	45 - 120	4.719	0.102	20	
Acenaphthylene	4.75	0.10	5	0	95.0	47 - 120	4.688	1.31	20	
Anthracene	4.858	0.10	5	0	97.2	45 - 120	4.816	0.869	20	
Benz(a)anthracene	5.279	0.10	5	0	106	40 - 120	5.312	0.633	20	
Benzo(a)pyrene	5.402	0.10	5	0	108	45 - 120	5.285	2.19	20	
Bis(2-chloroethoxy)methane	4.997	0.20	5	0	99.9	45 - 120	5.043	0.922	20	
Bis(2-ethylhexyl)phthalate	6.297	0.20	5	0	126	40 - 139	6.344	0.747	20	
Chrysene	4.735	0.10	5	0	94.7	43 - 120	4.782	0.99	20	
Dibenzofuran	5.019	0.10	5	0	100	50 - 120	4.968	1.01	20	
Di-n-butyl phthalate	5.883	0.20	5	0	118	45 - 123	5.755	2.19	20	
Fluoranthene	5.237	0.10	5	0	105	45 - 125	5.205	0.616	20	
Fluorene	5.177	0.10	5	0	104	49 - 120	5.174	0.0563	20	
Naphthalene	4.761	0.10	5	0	95.2	45 - 120	4.795	0.718	20	
Nitrobenzene	4.464	0.20	5	0	89.3	44 - 120	4.375	2.02	20	
N-Nitrosodiphenylamine	4.485	0.20	5	0	89.7	40 - 125	4.465	0.452	20	
Pentachlorophenol	4.428	0.20	5	0	88.6	19 - 121	4.311	2.66	20	
Phenanthrene	4.88	0.10	5	0	97.6	45 - 121	4.815	1.35	20	
Phenol	5.144	0.20	5	0	103	20 - 124	5.116	0.544	20	
Pyrene	5.254	0.10	5	0	105	40 - 130	5.307	1.01	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>5.379</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>108</i>	<i>34 - 129</i>	<i>5.449</i>	<i>1.3</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.919</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>78.4</i>	<i>40 - 125</i>	<i>3.867</i>	<i>1.34</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>3.698</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>74.0</i>	<i>20 - 120</i>	<i>3.617</i>	<i>2.24</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>5.131</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>103</i>	<i>40 - 135</i>	<i>5.306</i>	<i>3.35</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>3.449</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.0</i>	<i>41 - 120</i>	<i>3.465</i>	<i>0.452</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>4.168</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>83.4</i>	<i>20 - 120</i>	<i>4.106</i>	<i>1.5</i>	<i>20</i>	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

<b>Batch ID:</b> 151716 ( 0 )	<b>Instrument:</b> SV-7	<b>Method:</b> LOW-LEVEL SEMIVOLATILES BY 8270D		
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The following samples were analyzed in this batch:

HS20030619-02	HS20030619-03	HS20030619-04	HS20030619-05
HS20030619-06	HS20030619-07	HS20030619-08	HS20030619-09
HS20030619-10	HS20030619-12	HS20030619-13	HS20030619-14
HS20030619-15	HS20030619-16		



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

**Batch ID:** R358243 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200314</b>			Units: <b>ug/L</b>		Analysis Date: <b>14-Mar-2020 11:00</b>			
Client ID:		Run ID: <b>VOA4_358243</b>			SeqNo: <b>5514256</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	< 0.20	1.0								
Benzene	< 0.20	1.0								
Chlorobenzene	< 0.30	1.0								
Ethylbenzene	< 0.30	1.0								
Methylene chloride	< 1.0	2.0								
Toluene	< 0.20	1.0								
Xylenes, Total	< 0.30	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	52.52	1.0	50	0	105	70 - 123				
<i>Surr: 4-Bromofluorobenzene</i>	47.01	1.0	50	0	94.0	82 - 115				
<i>Surr: Dibromofluoromethane</i>	50.2	1.0	50	0	100	73 - 126				
<i>Surr: Toluene-d8</i>	48.89	1.0	50	0	97.8	81 - 120				

<b>LCS</b>		Sample ID: <b>VLCSW-200314</b>			Units: <b>ug/L</b>		Analysis Date: <b>14-Mar-2020 10:11</b>			
Client ID:		Run ID: <b>VOA4_358243</b>			SeqNo: <b>5514255</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	19.35	1.0	20	0	96.7	70 - 124				
Benzene	19.67	1.0	20	0	98.4	74 - 120				
Chlorobenzene	19.67	1.0	20	0	98.4	76 - 113				
Ethylbenzene	20.16	1.0	20	0	101	77 - 117				
Methylene chloride	19.96	2.0	20	0	99.8	70 - 127				
Toluene	19.96	1.0	20	0	99.8	77 - 118				
Xylenes, Total	61.63	1.0	60	0	103	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	49.34	1.0	50	0	98.7	70 - 130				
<i>Surr: 4-Bromofluorobenzene</i>	49.66	1.0	50	0	99.3	82 - 115				
<i>Surr: Dibromofluoromethane</i>	50.66	1.0	50	0	101	73 - 126				
<i>Surr: Toluene-d8</i>	50.8	1.0	50	0	102	81 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

**Batch ID:** R358243 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20030599-05MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>14-Mar-2020 14:20</b>			
Client ID:		Run ID: <b>VOA4_358243</b>			SeqNo: <b>5514264</b>		PrepDate:		DF: <b>100</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	2420	100	2000	632.1	89.4	70 - 127				
Benzene	2933	100	2000	969.3	98.2	70 - 127				
Chlorobenzene	1887	100	2000	0	94.4	70 - 114				
Ethylbenzene	1921	100	2000	0	96.0	70 - 124				
Methylene chloride	1828	200	2000	0	91.4	70 - 128				
Toluene	1960	100	2000	34.94	96.2	70 - 123				
Xylenes, Total	5878	100	6000	0	98.0	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>4941</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>98.8</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>4950</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>99.0</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>4933</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>98.7</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>4956</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>99.1</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20030599-05MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>14-Mar-2020 14:45</b>			
Client ID:		Run ID: <b>VOA4_358243</b>			SeqNo: <b>5514265</b>		PrepDate:		DF: <b>100</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	2451	100	2000	632.1	91.0	70 - 127	2420	1.29	20	
Benzene	2824	100	2000	969.3	92.7	70 - 127	2933	3.78	20	
Chlorobenzene	1923	100	2000	0	96.1	70 - 114	1887	1.85	20	
Ethylbenzene	1956	100	2000	0	97.8	70 - 124	1921	1.85	20	
Methylene chloride	1790	200	2000	0	89.5	70 - 128	1828	2.13	20	
Toluene	1948	100	2000	34.94	95.6	70 - 123	1960	0.622	20	
Xylenes, Total	6166	100	6000	0	103	70 - 130	5878	4.78	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>4858</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>97.2</i>	<i>70 - 126</i>	<i>4941</i>	<i>1.68</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>5133</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>103</i>	<i>81 - 113</i>	<i>4950</i>	<i>3.62</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>4820</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>96.4</i>	<i>77 - 123</i>	<i>4933</i>	<i>2.31</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>4990</i>	<i>100</i>	<i>5000</i>	<i>0</i>	<i>99.8</i>	<i>82 - 127</i>	<i>4956</i>	<i>0.691</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20030619-02    HS20030619-03    HS20030619-04    HS20030619-05  
 HS20030619-06

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

**Batch ID:** R358244 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200314</b>			Units: <b>ug/L</b>		Analysis Date: <b>14-Mar-2020 23:24</b>			
Client ID:		Run ID: <b>VOA4_358244</b>			SeqNo: <b>5514283</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	< 0.20	1.0								
Benzene	< 0.20	1.0								
Chlorobenzene	< 0.30	1.0								
Ethylbenzene	< 0.30	1.0								
Methylene chloride	< 1.0	2.0								
Toluene	< 0.20	1.0								
Xylenes, Total	< 0.30	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.27</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.31</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>94.6</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>51.08</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200314</b>			Units: <b>ug/L</b>		Analysis Date: <b>14-Mar-2020 22:35</b>			
Client ID:		Run ID: <b>VOA4_358244</b>			SeqNo: <b>5514282</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.47	1.0	20	0	92.4	70 - 124				
Benzene	19.05	1.0	20	0	95.3	74 - 120				
Chlorobenzene	19.21	1.0	20	0	96.1	76 - 113				
Ethylbenzene	19.21	1.0	20	0	96.1	77 - 117				
Methylene chloride	19.03	2.0	20	0	95.1	70 - 127				
Toluene	19.36	1.0	20	0	96.8	77 - 118				
Xylenes, Total	58.72	1.0	60	0	97.9	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.71</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.66</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>50.17</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>50.4</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QC BATCH REPORT**

**Batch ID:** R358244 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20030619-07MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Mar-2020 00:37</b>			
Client ID: <b>WG-1620-MW84A-20200312</b>		Run ID: <b>VOA4_358244</b>			SeqNo: <b>5514286</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.46	1.0	20	0	92.3	70 - 127				
Benzene	19.3	1.0	20	0	96.5	70 - 127				
Chlorobenzene	19.36	1.0	20	0	96.8	70 - 114				
Ethylbenzene	19.95	1.0	20	0	99.7	70 - 124				
Methylene chloride	18.37	2.0	20	0	91.8	70 - 128				
Toluene	19.89	1.0	20	0	99.4	70 - 123				
Xylenes, Total	60.5	1.0	60	0	101	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.72</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.9</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.8</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.95</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.54</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20030619-07MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>15-Mar-2020 01:01</b>			
Client ID: <b>WG-1620-MW84A-20200312</b>		Run ID: <b>VOA4_358244</b>			SeqNo: <b>5514287</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.37	1.0	20	0	91.9	70 - 127	18.46	0.492	20	
Benzene	19.26	1.0	20	0	96.3	70 - 127	19.3	0.213	20	
Chlorobenzene	18.78	1.0	20	0	93.9	70 - 114	19.36	3.05	20	
Ethylbenzene	19.35	1.0	20	0	96.8	70 - 124	19.95	3.02	20	
Methylene chloride	17.61	2.0	20	0	88.1	70 - 128	18.37	4.2	20	
Toluene	18.89	1.0	20	0	94.4	70 - 123	19.89	5.16	20	
Xylenes, Total	60.07	1.0	60	0	100	70 - 130	60.5	0.704	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.82</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.6</i>	<i>70 - 126</i>	<i>48.72</i>	<i>0.19</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>50.43</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>81 - 113</i>	<i>48.9</i>	<i>3.09</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>48.21</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.4</i>	<i>77 - 123</i>	<i>48.95</i>	<i>1.51</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>49.43</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.9</i>	<i>82 - 127</i>	<i>50.54</i>	<i>2.23</i>	<i>20</i>	

The following samples were analyzed in this batch:

HS20030619-01	HS20030619-07	HS20030619-08	HS20030619-09
HS20030619-10	HS20030619-11	HS20030619-12	HS20030619-13
HS20030619-14	HS20030619-15	HS20030619-16	

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030619

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
Arkansas	19-028-0	27-Mar-2020
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231 V009	22-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20030619

Date/Time Received: 13-Mar-2020 14:00
Received by: PS

Checklist completed by: Asad Chaudhry
eSignature
Date: 13-Mar-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 17-Mar-2020

Matrices: GW, Water

Carrier name: ALS Courier

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

2 Page(s)
COC IDs:221279, 221277

Temperature(s)/Thermometer(s): 1.2°C, 2.9°C, 1.7°C, 1.4°C UC/C IR 25
Cooler(s)/Kit(s): 45544, 45564, 45666, 433355
Date/Time sample(s) sent to storage: 03/13/2020 17:20

- Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



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Holland, MI  
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# Chain of Custody Form

Page      of     

COC ID: 221279

## HS20030619

Golder Associates Inc.  
Houston TX-Wood Preserving Works

in, WV  
8

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ALS Project Manager:

Customer Information		Project Information		
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A
Work Order		Project Number	1620-14-Rev0 SR 92688	B
Company Name	Golder Associates Inc.	Bill To Company	Union Pacific Railroad- A/P	C
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E
	Suite 4004		Stop 0780	F
City/State/Zip	Round Rock, TX 78684	City/State/Zip	Omaha NE 681790750	G
Phone	(512) 671-3434	Phone		H
Fax	(512) 671-3446	Fax		I
e-Mail Address	Eric_Matzner@golder.com	e-Mail Address		J

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TB (G-02) 2019 2020 03 12	3-12-20	16:30	Water	1.8	2	X										
2	WG-1620-MW 996 2020 03 11	3-11-20	15:35	Groundwater	12.0	29	X	X	X	X							
3	WG-1620-MW 545 2020 03 11	3-11-20	14:20	G-W	12.8	9	X	X	X	X							
4	WG-1620-MW 94A 2020 03 12	3-12-20	15:45	G-W	12.8	9	X	X	X	X							
5	WG-1620-MW 93B 2020 03 12	3-12-20	14:35	G-W	12.8	9	X	X	X	X							
6	WG-1620-MW 92B 2020 03 12	3-12-20	13:20	G-W	12.8	9	X	X	X	X							
7	WG-1620-MW 84A 2020 03 12	3-12-20	12:15	G-W	12.8	9	X	X	X	X							
8	WG-1620-MW 70C 2020 03 12	3-12-20	10:35	G-W	12.8	9	X	X	X	X							
9	WG-1620-MW 91A 2020 03 12	3-12-20	9:20	G-W	12.8	9	X	X	X	X							
10	WG-1620-FD (1) 2020 03 12	3-12-20	-	G-W	12.8	9	X	X	X	X							

Sampler(s) Please Print & Sign: *Tim McSpadden*      Shipment Method: \_\_\_\_\_      Required Turnaround Time: (Check Box)  5-10 WK Log  3-5 WK Log  2-3 Days  24 Hour      Results Due Date: \_\_\_\_\_

Relinquished by: *[Signature]*      Date: *3-13-20*      Time: *13:15*      Received by: *[Signature]*      Notes: UPRR HWPW 1620-14

Relinquished by: *[Signature]*      Date: *3-13-20*      Time: *14:00*      Received by (Laboratory): *[Signature]*      Cooler ID: *45544*      Cooler Temp.: *1.2*      QC Package: (Check One Box Below)  IIR- Certified  IIR- Not Certified

Logged by (Laboratory): *[Signature]*      Date: \_\_\_\_\_      Time: \_\_\_\_\_      Checked by (Laboratory): *[Signature]*      Cooler ID: *45564*      Cooler Temp.: *2.9*      QC Package:  IIR- Certified  IIR- Not Certified

Cooler ID: *45660*      Cooler Temp.: *1.7*      QC Package:  IIR- Certified  IIR- Not Certified

Cooler ID: *43355*      Cooler Temp.: *1.4*      QC Package:  IIR- Certified  IIR- Not Certified

Preservative Key: 1-HCl 2-HNO<sub>3</sub> 3-H<sub>2</sub>SO<sub>4</sub> 4-NaOH 5-Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> 6-NaHSO<sub>4</sub> 7-Other 8-4°C 9-5035

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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# Chain of Custody Form

Page      of     

COC ID: 221277

## HS20030619

Golder Associates Inc.  
Houston TX-Wood Preserving Works



n, WV

Customer Information		Project Information		ALS Project Manager:	
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260_LL_W (5632526 VOC Site Specific)
Work Order		Project Number	1620-14-Rev0 SR 92088	B	TX1005_W Low (5643233 TPH TX1005)
Company Name	Golder Associates Inc.	Bill To Company	Union Pacific Railroad- A/P	C	8270_LOW_W (5632532 SemiVolatiles Site specific)
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 Metals - As)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E	
	Suite 4004		Stop 0750	F	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	G	
Phone	(512) 671-3434	Phone		H	
Fax	(512) 671-3446	Fax		I	
e-Mail Address	Eric_Matzner@golder.com	e-Mail Address		J	

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WG-1620-TB (61-621720-18) <sup>20</sup> 857	3-12-20	12:36	Water	1.8	2	X										
2	WG-1620-MW76B-20200312	3-12-20	10:35	Groundwater	1.2.8	29	X	X	X	X							
3	" - MW97A - "	3-12-20	11:47	↓	↓	↓	↓	↓	↓	↓							
4	" - MW98A - "	3-12-20	12:54	↓	↓	↓	↓	↓	↓	↓							
5	" - MW98B - "	3-12-20	13:53	↓	↓	↓	↓	↓	↓	↓							
6	" - MW50B - "	3-12-20	15:20	↓	↓	↓	↓	↓	↓	↓							
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>Sarah Balke</i>		Shipment Method	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> 1-2 Business Days <input type="checkbox"/> 3-5 Business Days <input type="checkbox"/> 7-10 Business Days <input type="checkbox"/> Other		Results Due Date:
Relinquished by: <i>T. Haddock</i>	Date: 3-15-20	Time: 13:15	Received by: <i>[Signature]</i>	Notes: UPRR HWPV 1620-14	
Relinquished to: <i>[Signature]</i>	Date: 3-13-20	Time: 1400	Received by (Laboratory): <i>[Signature]</i>	Cooler ID	Cooler Temp.
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	QC Package: (Check One Box Below)	
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035				<input type="checkbox"/> Ice/Insulated	<input checked="" type="checkbox"/> 100% Clean
				<input type="checkbox"/> Ice/Insulated/CO <sub>2</sub> Purge	<input type="checkbox"/> 100% Clean
				<input type="checkbox"/> Ice/Insulated/CO <sub>2</sub>	<input type="checkbox"/> 100% Clean

- note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
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F: +1 281 530 5887

March 27, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20030961**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 3 sample(s) on Mar 20, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read 'Dane J. Wacasey'.

Generated By: JUMOKE.LAWAL  
Dane J. Wacasey

---

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

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**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey

Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 03/27/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20030961			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 152024,152161,R358905			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?	X				
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			1
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 03/27/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20030961			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 152024,152161,R358905			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
S1	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
S2	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
S3	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
S4	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
S5	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section)					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
S6	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
S7	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
S8	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
S9	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
S10	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
S11	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
S12	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
S13	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
S14	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
S15	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
S16	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 03/27/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20030961
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 152024,152161,R358905

ER# <sup>5</sup>	Description
------------------	-------------

1	Batch 152024, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.
---	---

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20030961

**SAMPLE SUMMARY**

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Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20030961-01	WQ-1620-TB01-20200320	Water	CG-021720 -183	20-Mar-2020 11:00	20-Mar-2020 11:50	<input type="checkbox"/>
HS20030961-02	WG-1620-MW47A-20200320	Groundwater		20-Mar-2020 10:30	20-Mar-2020 11:50	<input type="checkbox"/>
HS20030961-03	WG-1620-MW60AR-20200320	Groundwater		20-Mar-2020 08:55	20-Mar-2020 11:50	<input type="checkbox"/>



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB01-20200320  
 Collection Date: 20-Mar-2020 11:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20030961  
 Lab ID:HS20030961-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>					Analyst: AKP
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	24-Mar-2020 22:58
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	24-Mar-2020 22:58
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	24-Mar-2020 22:58
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	24-Mar-2020 22:58
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	24-Mar-2020 22:58
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	24-Mar-2020 22:58
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	24-Mar-2020 22:58
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>104</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>24-Mar-2020 22:58</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.8</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>24-Mar-2020 22:58</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>24-Mar-2020 22:58</i>
<i>Surr: Toluene-d8</i>	<i>104</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>24-Mar-2020 22:58</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW47A-20200320  
 Collection Date: 20-Mar-2020 10:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20030961  
 Lab ID:HS20030961-02  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	25-Mar-2020 05:30
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	25-Mar-2020 05:30
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	25-Mar-2020 05:30
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	25-Mar-2020 05:30
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	25-Mar-2020 05:30
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	25-Mar-2020 05:30
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	25-Mar-2020 05:30
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>103</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:30</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>93.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:30</i>
<i>Surr: Dibromofluoromethane</i>	<i>101</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:30</i>
<i>Surr: Toluene-d8</i>	<i>105</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:30</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW47A-20200320  
 Collection Date: 20-Mar-2020 10:30

**ANALYTICAL REPORT**  
 WorkOrder:HS20030961  
 Lab ID:HS20030961-02  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	24-Mar-2020 10:50
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	24-Mar-2020 10:50
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	24-Mar-2020 10:50
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	24-Mar-2020 10:50
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	24-Mar-2020 10:50
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	24-Mar-2020 10:50
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	24-Mar-2020 10:50
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	24-Mar-2020 10:50
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	24-Mar-2020 10:50
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	24-Mar-2020 10:50
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	24-Mar-2020 10:50
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	24-Mar-2020 10:50
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	24-Mar-2020 10:50
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	24-Mar-2020 10:50
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00011</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	24-Mar-2020 10:50
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	24-Mar-2020 10:50
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	24-Mar-2020 10:50
<b>Di-n-butyl phthalate</b>	<b>0.000076</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	24-Mar-2020 10:50
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	24-Mar-2020 10:50
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	24-Mar-2020 10:50
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	24-Mar-2020 10:50
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	24-Mar-2020 10:50
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	24-Mar-2020 10:50
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	24-Mar-2020 10:50
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	24-Mar-2020 10:50
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	24-Mar-2020 10:50
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	24-Mar-2020 10:50
<i>Surr: 2,4,6-Tribromophenol</i>	96.9			34-129	%REC	1	24-Mar-2020 10:50
<i>Surr: 2-Fluorobiphenyl</i>	71.7			40-125	%REC	1	24-Mar-2020 10:50
<i>Surr: 2-Fluorophenol</i>	58.4			20-120	%REC	1	24-Mar-2020 10:50
<i>Surr: 4-Terphenyl-d14</i>	89.6			40-135	%REC	1	24-Mar-2020 10:50
<i>Surr: Nitrobenzene-d5</i>	56.7			41-120	%REC	1	24-Mar-2020 10:50
<i>Surr: Phenol-d6</i>	61.7			20-120	%REC	1	24-Mar-2020 10:50
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.000566</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	26-Mar-2020 19:34

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW60AR-20200320  
 Collection Date: 20-Mar-2020 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20030961  
 Lab ID:HS20030961-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	25-Mar-2020 05:54
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	25-Mar-2020 05:54
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	25-Mar-2020 05:54
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	25-Mar-2020 05:54
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	25-Mar-2020 05:54
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	25-Mar-2020 05:54
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	25-Mar-2020 05:54
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>104</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:54</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:54</i>
<i>Surr: Dibromofluoromethane</i>	<i>96.9</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:54</i>
<i>Surr: Toluene-d8</i>	<i>103</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>25-Mar-2020 05:54</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW60AR-20200320  
 Collection Date: 20-Mar-2020 08:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20030961  
 Lab ID:HS20030961-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 23-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	24-Mar-2020 11:28
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	24-Mar-2020 11:28
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	24-Mar-2020 11:28
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	24-Mar-2020 11:28
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	24-Mar-2020 11:28
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	24-Mar-2020 11:28
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	24-Mar-2020 11:28
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	24-Mar-2020 11:28
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	24-Mar-2020 11:28
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	24-Mar-2020 11:28
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	24-Mar-2020 11:28
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	24-Mar-2020 11:28
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	24-Mar-2020 11:28
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	24-Mar-2020 11:28
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.000063</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	24-Mar-2020 11:28
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	24-Mar-2020 11:28
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	24-Mar-2020 11:28
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	24-Mar-2020 11:28
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	24-Mar-2020 11:28
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	24-Mar-2020 11:28
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	24-Mar-2020 11:28
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	24-Mar-2020 11:28
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	24-Mar-2020 11:28
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	24-Mar-2020 11:28
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	24-Mar-2020 11:28
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	24-Mar-2020 11:28
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	24-Mar-2020 11:28
<i>Surr: 2,4,6-Tribromophenol</i>	89.1			34-129	%REC	1	24-Mar-2020 11:28
<i>Surr: 2-Fluorobiphenyl</i>	60.2			40-125	%REC	1	24-Mar-2020 11:28
<i>Surr: 2-Fluorophenol</i>	48.9			20-120	%REC	1	24-Mar-2020 11:28
<i>Surr: 4-Terphenyl-d14</i>	77.0			40-135	%REC	1	24-Mar-2020 11:28
<i>Surr: Nitrobenzene-d5</i>	48.7			41-120	%REC	1	24-Mar-2020 11:28
<i>Surr: Phenol-d6</i>	47.3			20-120	%REC	1	24-Mar-2020 11:28
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00189</b>	J	<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	26-Mar-2020 19:36

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**Batch ID:** 152024      **Start Date:** 23 Mar 2020 14:30      **End Date:** 23 Mar 2020 17:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030961-02		1000 (mL)	1 (mL)	0.001
HS20030961-03		1000 (mL)	1 (mL)	0.001

**Batch ID:** 152161      **Start Date:** 26 Mar 2020 08:47      **End Date:** 26 Mar 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030961-02		10 (mL)	10 (mL)	1
HS20030961-03		10 (mL)	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID:</b> 152024 ( 0 )		<b>Test Name :</b> LOW-LEVEL SEMIVOLATILES BY 8270D			<b>Matrix:</b> Groundwater	
HS20030961-02	WG-1620-MW47A-20200320	20 Mar 2020 10:30		23 Mar 2020 15:37	24 Mar 2020 10:50	1
HS20030961-03	WG-1620-MW60AR-20200320	20 Mar 2020 08:55		23 Mar 2020 15:37	24 Mar 2020 11:28	1
<b>Batch ID:</b> 152161 ( 0 )		<b>Test Name :</b> ICP-MS METALS BY SW6020A			<b>Matrix:</b> Groundwater	
HS20030961-02	WG-1620-MW47A-20200320	20 Mar 2020 10:30		26 Mar 2020 13:00	26 Mar 2020 19:34	1
HS20030961-03	WG-1620-MW60AR-20200320	20 Mar 2020 08:55		26 Mar 2020 13:00	26 Mar 2020 19:36	1
<b>Batch ID:</b> R358905 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Groundwater	
HS20030961-02	WG-1620-MW47A-20200320	20 Mar 2020 10:30			25 Mar 2020 05:30	1
HS20030961-03	WG-1620-MW60AR-20200320	20 Mar 2020 08:55			25 Mar 2020 05:54	1
<b>Batch ID:</b> R358905 ( 0 )		<b>Test Name :</b> LOW LEVEL VOLATILES BY SW8260C			<b>Matrix:</b> Water	
HS20030961-01	WQ-1620-TB01-20200320	20 Mar 2020 11:00			24 Mar 2020 22:58	1

WorkOrder: HS20030961  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200



WorkOrder: HS20030961  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20030961  
 InstrumentID: VOA4  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00061	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00057	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00058	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00048	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00070	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00050	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00070	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QC BATCH REPORT**

Batch ID: 152161 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-152161</b>	Units: <b>mg/L</b>		Analysis Date: <b>26-Mar-2020 19:13</b>					
Client ID:	Run ID: <b>ICPMS05_358931</b>	SeqNo: <b>5532745</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	< 0.000400	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-152161</b>	Units: <b>mg/L</b>		Analysis Date: <b>26-Mar-2020 19:15</b>					
Client ID:	Run ID: <b>ICPMS05_358931</b>	SeqNo: <b>5532746</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.04894	0.00200	0.05	0	97.9	80 - 120			
<b>MS</b>	Sample ID: <b>HS20030965-04MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>26-Mar-2020 19:22</b>					
Client ID:	Run ID: <b>ICPMS05_358931</b>	SeqNo: <b>5532749</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.05528	0.00200	0.05	0.004711	101	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20030965-04MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>26-Mar-2020 19:25</b>					
Client ID:	Run ID: <b>ICPMS05_358931</b>	SeqNo: <b>5532750</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.05392	0.00200	0.05	0.004711	98.4	80 - 120	0.05528	2.48 20	
<b>PDS</b>	Sample ID: <b>HS20030965-04PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>26-Mar-2020 19:27</b>					
Client ID:	Run ID: <b>ICPMS05_358931</b>	SeqNo: <b>5532751</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.106	0.00200	0.1	0.004711	101	75 - 125			
<b>SD</b>	Sample ID: <b>HS20030965-04SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>26-Mar-2020 19:20</b>					
Client ID:	Run ID: <b>ICPMS05_358931</b>	SeqNo: <b>5532748</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>5</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D %D Limit Qual	
Arsenic	0.004019	0.0100					0.004711	0 10 J	

The following samples were analyzed in this batch: HS20030961-02 HS20030961-03

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QC BATCH REPORT**

Batch ID: 152024 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-152024	Units: ug/L			Analysis Date: 24-Mar-2020 09:53					
Client ID:	Run ID: SV-7_358753	SeqNo: 5526601	PrepDate: 23-Mar-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	< 0.021	0.20								
2,4-Dimethylphenol	< 0.040	0.20								
2,4-Dinitrotoluene	< 0.058	0.20								
2,6-Dinitrotoluene	< 0.042	0.20								
2-Chloronaphthalene	< 0.021	0.20								
2-Methylnaphthalene	< 0.019	0.10								
4,6-Dinitro-2-methylphenol	< 0.020	0.20								
4-Nitrophenol	< 0.047	1.0								
Acenaphthene	< 0.027	0.10								
Acenaphthylene	< 0.015	0.10								
Anthracene	< 0.014	0.10								
Benz(a)anthracene	< 0.050	0.10								
Benzo(a)pyrene	< 0.020	0.10								
Bis(2-chloroethoxy)methane	< 0.030	0.20								
Bis(2-ethylhexyl)phthalate	< 0.037	0.20								
Chrysene	< 0.021	0.10								
Dibenzofuran	< 0.020	0.10								
Di-n-butyl phthalate	< 0.020	0.20								
Fluoranthene	< 0.010	0.10								
Fluorene	< 0.030	0.10								
Naphthalene	< 0.020	0.10								
Nitrobenzene	< 0.024	0.20								
N-Nitrosodiphenylamine	< 0.025	0.20								
Pentachlorophenol	< 0.079	0.20								
Phenanthrene	< 0.021	0.10								
Phenol	< 0.035	0.20								
Pyrene	< 0.019	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.818</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>76.4</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.671</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>73.4</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>3.241</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>64.8</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.592</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>91.8</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>3.079</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>61.6</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.448</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>69.0</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QC BATCH REPORT**

Batch ID: 152024 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-152024	Units: ug/L			Analysis Date: 24-Mar-2020 10:12					
Client ID:	Run ID: SV-7_358753	SeqNo: 5526602		PrepDate: 23-Mar-2020		DF: 1				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.044	0.20	5	0	60.9	39 - 127				
2,4-Dimethylphenol	3.518	0.20	5	0	70.4	35 - 120				
2,4-Dinitrotoluene	4.266	0.20	5	0	85.3	50 - 122				
2,6-Dinitrotoluene	4.362	0.20	5	0	87.2	50 - 120				
2-Chloronaphthalene	4.343	0.20	5	0	86.9	50 - 120				
2-Methylnaphthalene	3.829	0.10	5	0	76.6	50 - 120				
4,6-Dinitro-2-methylphenol	3.498	0.20	5	0	70.0	25 - 121				
4-Nitrophenol	2.946	1.0	5	0	58.9	30 - 130				
Acenaphthene	3.623	0.10	5	0	72.5	45 - 120				
Acenaphthylene	3.667	0.10	5	0	73.3	47 - 120				
Anthracene	3.781	0.10	5	0	75.6	45 - 120				
Benz(a)anthracene	3.921	0.10	5	0	78.4	40 - 120				
Benzo(a)pyrene	3.898	0.10	5	0	78.0	45 - 120				
Bis(2-chloroethoxy)methane	3.614	0.20	5	0	72.3	45 - 120				
Bis(2-ethylhexyl)phthalate	3.616	0.20	5	0	72.3	40 - 139				
Chrysene	3.587	0.10	5	0	71.7	43 - 120				
Dibenzofuran	3.894	0.10	5	0	77.9	50 - 120				
Di-n-butyl phthalate	3.806	0.20	5	0	76.1	45 - 123				
Fluoranthene	3.938	0.10	5	0	78.8	45 - 125				
Fluorene	3.911	0.10	5	0	78.2	49 - 120				
Naphthalene	3.624	0.10	5	0	72.5	45 - 120				
Nitrobenzene	3.199	0.20	5	0	64.0	44 - 120				
N-Nitrosodiphenylamine	3.714	0.20	5	0	74.3	40 - 125				
Pentachlorophenol	1.418	0.20	5	0	28.4	19 - 121				
Phenanthrene	3.672	0.10	5	0	73.4	45 - 121				
Phenol	3.565	0.20	5	0	71.3	20 - 124				
Pyrene	4.122	0.10	5	0	82.4	40 - 130				
<i>Surr: 2,4,6-Tribromophenol</i>	<i>4.236</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>84.7</i>	<i>34 - 129</i>				
<i>Surr: 2-Fluorobiphenyl</i>	<i>3.409</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>68.2</i>	<i>40 - 125</i>				
<i>Surr: 2-Fluorophenol</i>	<i>2.949</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>59.0</i>	<i>20 - 120</i>				
<i>Surr: 4-Terphenyl-d14</i>	<i>4.003</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>80.1</i>	<i>40 - 135</i>				
<i>Surr: Nitrobenzene-d5</i>	<i>2.711</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>54.2</i>	<i>41 - 120</i>				
<i>Surr: Phenol-d6</i>	<i>3.075</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>61.5</i>	<i>20 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QC BATCH REPORT**

Batch ID: 152024 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCSD		Sample ID: LCSD-152024		Units: ug/L		Analysis Date: 24-Mar-2020 11:09				
Client ID:		Run ID: SV-7_358753		SeqNo: 5526759		PrepDate: 23-Mar-2020		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.078	0.20	5	0	61.6	39 - 127	3.044	1.11	20	
2,4-Dimethylphenol	3.536	0.20	5	0	70.7	35 - 120	3.518	0.508	20	
2,4-Dinitrotoluene	4.103	0.20	5	0	82.1	50 - 122	4.266	3.91	20	
2,6-Dinitrotoluene	4.141	0.20	5	0	82.8	50 - 120	4.362	5.2	20	
2-Chloronaphthalene	4.317	0.20	5	0	86.3	50 - 120	4.343	0.588	20	
2-Methylnaphthalene	3.833	0.10	5	0	76.7	50 - 120	3.829	0.114	20	
4,6-Dinitro-2-methylphenol	3.767	0.20	5	0	75.3	25 - 121	3.498	7.42	30	
4-Nitrophenol	3.196	1.0	5	0	63.9	30 - 130	2.946	8.14	20	
Acenaphthene	3.518	0.10	5	0	70.4	45 - 120	3.623	2.94	20	
Acenaphthylene	3.588	0.10	5	0	71.8	47 - 120	3.667	2.18	20	
Anthracene	3.775	0.10	5	0	75.5	45 - 120	3.781	0.165	20	
Benz(a)anthracene	3.835	0.10	5	0	76.7	40 - 120	3.921	2.22	20	
Benzo(a)pyrene	3.734	0.10	5	0	74.7	45 - 120	3.898	4.3	20	
Bis(2-chloroethoxy)methane	3.687	0.20	5	0	73.7	45 - 120	3.614	2.01	20	
Bis(2-ethylhexyl)phthalate	3.596	0.20	5	0	71.9	40 - 139	3.616	0.554	20	
Chrysene	3.449	0.10	5	0	69.0	43 - 120	3.587	3.93	20	
Dibenzofuran	3.833	0.10	5	0	76.7	50 - 120	3.894	1.59	20	
Di-n-butyl phthalate	3.871	0.20	5	0	77.4	45 - 123	3.806	1.69	20	
Fluoranthene	4.045	0.10	5	0	80.9	45 - 125	3.938	2.67	20	
Fluorene	3.875	0.10	5	0	77.5	49 - 120	3.911	0.945	20	
Naphthalene	3.65	0.10	5	0	73.0	45 - 120	3.624	0.72	20	
Nitrobenzene	3.414	0.20	5	0	68.3	44 - 120	3.199	6.51	20	
N-Nitrosodiphenylamine	3.737	0.20	5	0	74.7	40 - 125	3.714	0.598	20	
Pentachlorophenol	1.54	0.20	5	0	30.8	19 - 121	1.418	8.25	20	
Phenanthrene	3.71	0.10	5	0	74.2	45 - 121	3.672	1.03	20	
Phenol	3.505	0.20	5	0	70.1	20 - 124	3.565	1.71	20	
Pyrene	3.879	0.10	5	0	77.6	40 - 130	4.122	6.05	20	
Surr: 2,4,6-Tribromophenol	4.294	0.20	5	0	85.9	34 - 129	4.236	1.37	20	
Surr: 2-Fluorobiphenyl	3.184	0.20	5	0	63.7	40 - 125	3.409	6.83	20	
Surr: 2-Fluorophenol	2.98	0.20	5	0	59.6	20 - 120	2.949	1.05	20	
Surr: 4-Terphenyl-d14	3.844	0.20	5	0	76.9	40 - 135	4.003	4.06	20	
Surr: Nitrobenzene-d5	2.828	0.20	5	0	56.6	41 - 120	2.711	4.25	20	
Surr: Phenol-d6	3.023	0.20	5	0	60.5	20 - 120	3.075	1.71	20	

The following samples were analyzed in this batch: HS20030961-02      HS20030961-03

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QC BATCH REPORT**

**Batch ID:** R358905 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200324</b>			Units: <b>ug/L</b>		Analysis Date: <b>24-Mar-2020 20:29</b>			
Client ID:		Run ID: <b>VOA4_358905</b>			SeqNo: <b>5530354</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	< 0.20	1.0								
Benzene	< 0.20	1.0								
Chlorobenzene	< 0.30	1.0								
Ethylbenzene	< 0.30	1.0								
Methylene chloride	< 1.0	2.0								
Toluene	< 0.20	1.0								
Xylenes, Total	< 0.30	1.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.63</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 123</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>46.77</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.5</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>49.16</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>51.52</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>81 - 120</i>				

<b>LCS</b>		Sample ID: <b>VLCSW-200324</b>			Units: <b>ug/L</b>		Analysis Date: <b>24-Mar-2020 19:39</b>			
Client ID:		Run ID: <b>VOA4_358905</b>			SeqNo: <b>5530353</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	18.16	1.0	20	0	90.8	70 - 124				
Benzene	18.54	1.0	20	0	92.7	74 - 120				
Chlorobenzene	19.29	1.0	20	0	96.5	76 - 113				
Ethylbenzene	19.8	1.0	20	0	99.0	77 - 117				
Methylene chloride	18.31	2.0	20	0	91.6	70 - 127				
Toluene	19.66	1.0	20	0	98.3	77 - 118				
Xylenes, Total	61.08	1.0	60	0	102	75 - 122				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.14</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>98.3</i>	<i>70 - 130</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.3</i>	<i>82 - 115</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.16</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>96.3</i>	<i>73 - 126</i>				
<i>Surr: Toluene-d8</i>	<i>51.1</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QC BATCH REPORT**

**Batch ID:** R358905 ( 0 )      **Instrument:** VOA4      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MS</b>		Sample ID: <b>HS20030969-04MS</b>			Units: <b>ug/L</b>		Analysis Date: <b>24-Mar-2020 23:47</b>			
Client ID:		Run ID: <b>VOA4_358905</b>			SeqNo: <b>5530362</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.67	1.0	20	0	88.3	70 - 127				
Benzene	18.29	1.0	20	0	91.5	70 - 127				
Chlorobenzene	18.92	1.0	20	0	94.6	70 - 114				
Ethylbenzene	19.84	1.0	20	0	99.2	70 - 124				
Methylene chloride	16.71	2.0	20	0	83.6	70 - 128				
Toluene	19.19	1.0	20	0	95.9	70 - 123				
Xylenes, Total	60.72	1.0	60	0	101	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>48.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.8</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.61</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.2</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>47.63</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.3</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>50.49</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: <b>HS20030969-04MSD</b>			Units: <b>ug/L</b>		Analysis Date: <b>25-Mar-2020 00:11</b>			
Client ID:		Run ID: <b>VOA4_358905</b>			SeqNo: <b>5530363</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Dichloroethane	17.47	1.0	20	0	87.3	70 - 127	17.67	1.16	20	
Benzene	17.79	1.0	20	0	88.9	70 - 127	18.29	2.8	20	
Chlorobenzene	18.79	1.0	20	0	93.9	70 - 114	18.92	0.69	20	
Ethylbenzene	19.53	1.0	20	0	97.7	70 - 124	19.84	1.55	20	
Methylene chloride	16.28	2.0	20	0	81.4	70 - 128	16.71	2.63	20	
Toluene	19.09	1.0	20	0	95.4	70 - 123	19.19	0.533	20	
Xylenes, Total	59.99	1.0	60	0	100.0	70 - 130	60.72	1.22	20	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>49.88</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>70 - 126</i>	<i>48.88</i>	<i>2.01</i>	<i>20</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>49.71</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.4</i>	<i>81 - 113</i>	<i>49.61</i>	<i>0.185</i>	<i>20</i>	
<i>Surr: Dibromofluoromethane</i>	<i>46.83</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>93.7</i>	<i>77 - 123</i>	<i>47.63</i>	<i>1.69</i>	<i>20</i>	
<i>Surr: Toluene-d8</i>	<i>50.32</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>101</i>	<i>82 - 127</i>	<i>50.49</i>	<i>0.333</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20030961-01      HS20030961-02      HS20030961-03



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030961

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231 V009	22-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020

Sample Receipt Checklist

Client Name: PBW
Work Order: HS20030961

Date/Time Received: 20-Mar-2020 11:50
Received by: PS

Checklist completed by: Jared R. Makan
eSignature
Date: 20-Mar-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 23-Mar-2020

Matrices: Water

Carrier name: ALS Courier

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:221276

Temperature(s)/Thermometer(s): 1.0°C/1.0°C UC/C IR11
Cooler(s)/Kit(s): 43906
Date/Time sample(s) sent to storage: 03/20/2020 19:00
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by:

Login Notes:

Client Contacted: Date Contacted: Person Contacted:

Contacted By: Regarding:

Comments:

Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Everett, WA  
+1 425 356 2600

Fort Collins, CO  
+1 970 490 1511

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

Page      of     

COC ID: 221276

Houston, TX  
+1 281 530 5656

Middletown, PA  
+1 717 944 5541

Spring City, PA  
+1 610 948 4903

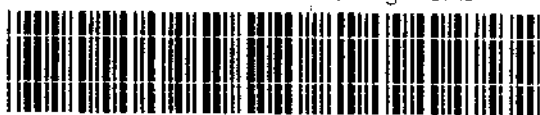
Salt Lake City, UT  
+1 801 266 7700

South Charleston, WV  
+1 304 356 3168

York, PA  
+1 717 505 5280

Customer Information		ALS Project Manager:		ALS Work Order #:	
Purchase Order	UPRR/Kevin Petaruburs	Project Name	Houston TX-Wood Preserving Works	Parameter/Method Request for Analysis	
Work Order		Project Number	1620-14-Rev0 SR 92288	A	5250 LL_W (5632528 VOC Site Specific)
Company Name	Golder Associates Inc.	Bill To Company	Union Pacific Railroad- A/P	B	<del>TX1605_W_Low (5643203 TPH TX1605)</del> <b>NO TPH</b>
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	C	2270 LOW_W (5632530 SemiVolatiles Site specific)
Address	2201 Double Creek Drive	Address	1400 Douglas Street	D	IOP_TW (5636002 Metals - As)
	Suite 4004		Stop 0750	E	
City/State/Zip	Round Rock, TX 78664	City/State/Zip	Omaha NE 681790750	F	
Phone	(512) 671-3434	Phone		G	
Fax	(512) 671-3446	Fax		H	
e-Mail Address	Eric_Matzner@golder.com	e-Mail Address		I	
				J	

**HS20030961**  
Golder Associates Inc.  
Houston TX-Wood Preserving Works




No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WC-1620-TB <b>CC-021720183</b>	3-20-20	11:00	Water	1.0	2	X										
2	WC-1620-MW <b>47A20200320</b>	3-20-20	10:30	Groundwater	1.2.0	19	X	X	X	X							
3	WC-1620-MW <b>60AR20200320</b>	3-20-20	8:55	GW	1,2,8	9	X		X	X							
4																	
5																	
6																	
7																	
8																	
9																	
10																	

Sampler(s) Please Print & Sign <i>Kevin Petaruburs</i>		Shipment Method	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> 2 Business Days <input type="checkbox"/> 5 Business Days <input type="checkbox"/> 7 Business Days <input type="checkbox"/> 10 Business Days		Results Due Date:
Relinquished by: <i>Kevin Petaruburs</i>	Date: 3-20-20	Time: 11:25	Received by: <i>Eric Matzner</i>	Notes: UPRR HWPW 1620-14	
Relinquished by: <i>Eric Matzner</i>	Date: 3-20-20	Time: 11:50	Received by (Laboratory): <i>Eric Matzner</i>	Cooler ID: 43906	Cooler Temp: UC 10 12.11 66.0
Logged by (Laboratory):	Date:	Time:	Checked by (Laboratory):	QC Package: (Check One Box Below) <input type="checkbox"/> 1-11-11-11-11 <input checked="" type="checkbox"/> 1-11-11-11-11-11 <input type="checkbox"/> 1-11-11-11-11-11-11 <input type="checkbox"/> Other	

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
3. The Chain of Custody is a legal document. All information must be completed accurately.

Copyright 2011 by ALS Environmental.

 <p><b>ALS</b>          10450 Standlin Rd., Suite 210          Houston, Texas 77099          Tel. +1 281 530 5658          Fax. +1 281 530 5887</p>	<b>CUSTODY SEAL</b>		Seal Broken By:
	Date: 5-26-20	Time: 11:36	Date:
	Name: T. M. [Signature]	[Signature]	Date: 3-20-20
Company: [Signature]	[Signature]		



---

10450 Stancliff Rd. Suite 210  
Houston, TX 77099  
T: +1 281 530 5656  
F: +1 281 530 5887

March 31, 2020

Eric Matzner  
Golder Associates Inc.  
2201 Double Creek Drive  
Suite 4004  
Round Rock, TX 78664

Work Order: **HS20030819**

Laboratory Results for: **Houston TX-Wood Preserving Works**

Dear Eric,

ALS Environmental received 8 sample(s) on Mar 18, 2020 for the analysis presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested. Results are expressed as "as received" unless otherwise noted.

QC sample results for this data met EPA or laboratory specifications except as noted in the Case Narrative or as noted with qualifiers in the QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained by ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

If you have any questions regarding this report, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read 'Dane J. Wacasey'.

Generated By: DANE.WACASEY  
Dane J. Wacasey

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**TRRP Laboratory Data  
Package Cover Page**

This data package consists of all or some of the following as applicable:

This signature page, the laboratory review checklist, and the following reportable data:

- R1 Field chain-of-custody documentation;
- R2 Sample identification cross-reference;
- R3 Test reports (analytical data sheets) for each environmental sample that includes:
  - a) Items consistent with NELAC Chapter 5,
  - b) dilution factors,
  - c) preparation methods,
  - d) cleanup methods, and
  - e) if required for the project, tentatively identified compounds (TICs).
- R4 Surrogate recovery data including:
  - a) Calculated recovery (%R), and
  - b) The laboratory's surrogate QC limits.
- R5 Test reports/summary forms for blank samples;
- R6 Test reports/summary forms for laboratory control samples (LCSs) including:
  - a) LCS spiking amounts,
  - b) Calculated %R for each analyte, and
  - c) The laboratory's LCS QC limits.
- R7 Test reports for project matrix spike/matrix spike duplicates (MS/MSDs) including:
  - a) Samples associated with the MS/MSD clearly identified,
  - b) MS/MSD spiking amounts,
  - c) Concentration of each MS/MSD analyte measured in the parent and spiked samples,
  - d) Calculated %Rs and relative percent differences (RPDs), and
  - e) The laboratory's MS/MSD QC limits.
- R8 Laboratory analytical duplicate (if applicable) recovery and precision:
  - a) the amount of analyte measured in the duplicate,
  - b) the calculated RPD, and
  - c) the laboratory's QC limits for analytical duplicates.
- R9 List of method quantitation limits (MQLs) and detectability check sample results for each analyte for each method and matrix.
- R10 Other problems or anomalies.  
The Exception Report for each "No" or "Not Reviewed (NR)" item in Laboratory Review Checklist and for each analyte, matrix, and method for which the laboratory does not hold NELAC accreditation under the Texas Laboratory Accreditation Program.

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**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

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**TRRP Laboratory Data  
Package Cover Page**

Release Statement: I am responsible for the release of this laboratory data package. This laboratory is NELAC accredited under the Texas Laboratory Accreditation Program for all the methods, analytes and matrices reported in this data package except as noted in the Exception Reports. The data have been reviewed and are technically compliant with the requirements of the methods used, except where noted by the laboratory in the attached exception reports. By my signature below, I affirm to the best of my knowledge, all problems/anomalies, observed by the laboratory have been identified by the laboratory in the Laboratory Review Checklist, and no information affecting the quality of the data has been knowingly withheld.

Check, if applicable:  [NA] This laboratory meets an exception under 30 TAC §25.6 and was last inspected by  TCEQ or  \_\_\_\_\_ on (enter date of last inspection). Any findings affecting the data in this laboratory data package are noted in the Exception Reports herein. The official signing the cover page of the report in which these data are used is responsible for releasing this data package and is by signature affirming the above release statement is true.



Dane J. Wacasey



Laboratory Review Checklist: Reportable Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 03/30/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20030819			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 151876,152162,R358799			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
<b>R1</b>	OI	<b>Chain-of-custody (C-O-C)</b>					
		Did samples meet the laboratory's standard conditions of sample acceptability upon receipt?	X				
		Were all departures from standard conditions described in an exception report?	X				
<b>R2</b>	OI	<b>Sample and quality control (QC) identification</b>					
		Are all field sample ID numbers cross-referenced to the laboratory ID numbers?	X				
		Are all laboratory ID numbers cross-referenced to the corresponding QC data?	X				
<b>R3</b>	OI	<b>Test reports</b>					
		Were all samples prepared and analyzed within holding times?	X				
		Other than those results < MQL, were all other raw values bracketed by calibration standards?	X				
		Were calculations checked by a peer or supervisor?	X				
		Were all analyte identifications checked by a peer or supervisor?	X				
		Were sample detection limits reported for all analytes not detected?	X				
		Were all results for soil and sediment samples reported on a dry weight basis?			X		
		Were % moisture (or solids) reported for all soil and sediment samples?			X		
		Were bulk soils/solids samples for volatile analysis extracted with methanol per SW-846 Method 5035?			X		
		If required for the project, TICs reported?			X		
<b>R4</b>	O	<b>Surrogate recovery data</b>					
		Were surrogates added prior to extraction?	X				
		Were surrogate percent recoveries in all samples within the laboratory QC limits?	X				
<b>R5</b>	OI	<b>Test reports/summary forms for blank samples</b>					
		Were appropriate type(s) of blanks analyzed?	X				
		Were blanks analyzed at the appropriate frequency?	X				
		Were method blanks taken through the entire analytical process, including preparation and, if applicable, cleanup procedures?	X				
		Were blank concentrations < MQL?	X				
<b>R6</b>	OI	<b>Laboratory control samples (LCS):</b>					
		Were all COCs included in the LCS?	X				
		Was each LCS taken through the entire analytical procedure, including prep and cleanup steps?	X				
		Were LCSs analyzed at the required frequency?	X				
		Were LCS (and LCSD, if applicable) %Rs within the laboratory QC limits?	X				
		Does the detectability data document the laboratory's capability to detect the COCs at the MDL used to calculate the SDLs?	X				
		Was the LCSD RPD within QC limits?	X				
<b>R7</b>	OI	<b>Matrix spike (MS) and matrix spike duplicate (MSD) data</b>					
		Were the project/method specified analytes included in the MS and MSD?	X				
		Were MS/MSD analyzed at the appropriate frequency?		X			1
		Were MS (and MSD, if applicable) %Rs within the laboratory QC limits?	X				
		Were MS/MSD RPDs within laboratory QC limits?	X				
<b>R8</b>	OI	<b>Analytical duplicate data</b>					
		Were appropriate analytical duplicates analyzed for each matrix?			X		
		Were analytical duplicates analyzed at the appropriate frequency?			X		
		Were RPDs or relative standard deviations within the laboratory QC limits?			X		
<b>R9</b>	OI	<b>Method quantitation limits (MQLs):</b>					
		Are the MQLs for each method analyte included in the laboratory data package?	X				
		Do the MQLs correspond to the concentration of the lowest non-zero calibration standard?	X				
		Are unadjusted MQLs and DCSs included in the laboratory data package?	X				
<b>R10</b>	OI	<b>Other problems/anomalies</b>					
		Are all known problems/anomalies/special conditions noted in this LRC and ER?	X				
		Were all necessary corrective actions performed for the reported data?	X				
		Was applicable and available technology used to lower the SDL and minimize the matrix interference effects on the sample results?	X				
		Is the laboratory NELAC-accredited under the Texas Laboratory Program for the analytes, matrices and methods associated with this laboratory data package?	X				

Laboratory Review Checklist: Supporting Data							
Laboratory Name: ALS Laboratory Group				LRC Date: 03/30/2020			
Project Name: Houston TX-Wood Preserving Works				Laboratory Job Number: HS20030819			
Reviewer Name: Dane Wacasey				Prep Batch Number(s): 151876,152162,R358799			
# <sup>1</sup>	A <sup>2</sup>	Description	Yes	No	NA <sup>3</sup>	NR <sup>4</sup>	ER# <sup>5</sup>
S1	OI	<b>Initial calibration (ICAL)</b>					
		Were response factors and/or relative response factors for each analyte within QC limits?	X				
		Were percent RSDs or correlation coefficient criteria met?	X				
		Was the number of standards recommended in the method used for all analytes?	X				
		Were all points generated between the lowest and highest standard used to calculate the curve?	X				
		Are ICAL data available for all instruments used?	X				
		Has the initial calibration curve been verified using an appropriate second source standard?	X				
S2	OI	<b>Initial and continuing calibration verification (ICCV and CCV) and continuing calibration blank (CCB)</b>					
		Was the CCV analyzed at the method-required frequency?	X				
		Were percent differences for each analyte within the method-required QC limits?	X				
		Was the ICAL curve verified for each analyte?	X				
		Was the absolute value of the analyte concentration in the inorganic CCB < MDL?	X				
S3	O	<b>Mass spectral tuning:</b>					
		Was the appropriate compound for the method used for tuning?	X				
		Were ion abundance data within the method-required QC limits?	X				
S4	O	<b>Internal standards (IS):</b>					
		Were IS area counts and retention times within the method-required QC limits?	X				
S5	OI	<b>Raw data</b> (NELAC section 1 appendix A glossary, and section 5.12 or ISO/IEC 17025 section					
		Were the raw data (for example, chromatograms, spectral data) reviewed by an analyst?	X				
		Were data associated with manual integrations flagged on the raw data?	X				
S6	O	<b>Dual column confirmation</b>					
		Did dual column confirmation results meet the method-required QC?			X		
S7	O	<b>Tentatively identified compounds (TICs):</b>					
		If TICs were requested, were the mass spectra and TIC data subject to appropriate checks?			X		
S8	I	<b>Interference Check Sample (ICS) results:</b>					
		Were percent recoveries within method QC limits?	X				
S9	I	<b>Serial dilutions, post digestion spikes, and method of standard additions</b>					
		Were percent differences, recoveries, and the linearity within the QC limits specified in the method?	X				
S10	OI	<b>Method detection limit (MDL) studies</b>					
		Was a MDL study performed for each reported analyte?	X				
		Is the MDL either adjusted or supported by the analysis of DCSs?	X				
S11	OI	<b>Proficiency test reports:</b>					
		Was the laboratory's performance acceptable on the applicable proficiency tests or evaluation studies?	X				
S12	OI	<b>Standards documentation</b>					
		Are all standards used in the analyses NIST-traceable or obtained from other appropriate sources?	X				
S13	OI	<b>Compound/analyte identification procedures</b>					
		Are the procedures for compound/analyte identification documented?	X				
S14	OI	<b>Demonstration of analyst competency (DOC)</b>					
		Was DOC conducted consistent with NELAC Chapter 5C or ISO/IEC 4?	X				
		Is documentation of the analyst's competency up-to-date and on file?	X				
S15	OI	<b>Verification/validation documentation for methods</b> (NELAC Chap 5 or ISO/IEC 17025 Section 5)					
		Are all the methods used to generate the data documented, verified, and validated, where applicable?	X				
S16	OI	<b>Laboratory standard operating procedures (SOPs):</b>					
		Are laboratory SOPs current and on file for each method performed?	X				

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.

O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);

NA = Not Applicable;

NR = Not Reviewed;

R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Laboratory Review Checklist: Exception Reports**

Laboratory Name: ALS Laboratory Group	LRC Date: 03/30/2020
Project Name: Houston TX-Wood Preserving Works	Laboratory Job Number: HS20030819
Reviewer Name: Dane Wacasey	Prep Batch Number(s): 151876,152162,R358799

ER# <sup>5</sup>	Description
	Batch 151876, Semivolatile Organics Method SW8270, LCS/LCSD were analyzed and reported in lieu of an MS/MSD for this batch.

Items identified by the letter "R" must be included in the laboratory data package submitted in the TRRP-required report(s). Items identified by the letter "S" should be retained and made available upon request for the appropriate retention period.  
O = Organic Analyses; I = Inorganic Analyses (and general chemistry, when applicable);  
NA = Not Applicable;  
NR = Not Reviewed;  
R# = Exception Report identification number (an Exception Report should be completed for an item if "NR" or "No" is checked).

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**Work Order:** HS20030819

**SAMPLE SUMMARY**

Lab Samp ID	Client Sample ID	Matrix	TagNo	Collection Date	Date Received	Hold
HS20030819-01	WQ-1620-TB01-20200317	Water	CG 021720 -179	17-Mar-2020 15:59	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-02	WG-1620-MW88A-20200317	Groundwater		17-Mar-2020 12:55	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-03	WG-1620-MW61B-20200317	Groundwater		17-Mar-2020 11:45	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-04	WG-1620-MW96B-20200317	Groundwater		17-Mar-2020 10:25	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-05	WG-1620-MW95A-20200317	Groundwater		17-Mar-2020 09:10	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-06	WG-1620-MW60B-20200317	Groundwater		17-Mar-2020 15:15	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-07	WG-1620-MW88B-20200317	Groundwater		17-Mar-2020 13:50	18-Mar-2020 16:35	<input type="checkbox"/>
HS20030819-08	WG-1620-FD02-20200317	Groundwater		17-Mar-2020 00:00	18-Mar-2020 16:35	<input type="checkbox"/>

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WQ-1620-TB01-20200317  
 Collection Date: 17-Mar-2020 15:59

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-01  
 Matrix:Water

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 19:53
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 19:53
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 19:53
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 19:53
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 19:53
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 19:53
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 19:53
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>100</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 19:53</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.3</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 19:53</i>
<i>Surr: Dibromofluoromethane</i>	<i>97.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 19:53</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 19:53</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88A-20200317  
 Collection Date: 17-Mar-2020 12:55

**ANALYTICAL REPORT**

WorkOrder:HS20030819  
 Lab ID:HS20030819-02  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 20:39
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 20:39
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 20:39
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 20:39
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 20:39
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 20:39
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 20:39
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>100</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:39</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.6</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:39</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:39</i>
<i>Surr: Toluene-d8</i>	<i>102</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:39</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88A-20200317  
 Collection Date: 17-Mar-2020 12:55

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-02  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 18:27
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 18:27
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 18:27
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 18:27
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 18:27
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 18:27
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 18:27
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 18:27
<b>Acenaphthene</b>	<b>0.0015</b>		<b>0.000027</b>	<b>0.00010</b>	<b>mg/L</b>	1	19-Mar-2020 18:27
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 18:27
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 18:27
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 18:27
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 18:27
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 18:27
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00010</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	19-Mar-2020 18:27
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 18:27
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 18:27
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 18:27
<b>Fluoranthene</b>	<b>0.00025</b>		<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	19-Mar-2020 18:27
<b>Fluorene</b>	<b>0.00026</b>		<b>0.000030</b>	<b>0.00010</b>	<b>mg/L</b>	1	19-Mar-2020 18:27
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 18:27
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 18:27
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 18:27
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 18:27
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 18:27
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 18:27
<b>Pyrene</b>	<b>0.00053</b>		<b>0.000019</b>	<b>0.00010</b>	<b>mg/L</b>	1	19-Mar-2020 18:27
<i>Surr: 2,4,6-Tribromophenol</i>	97.6			34-129	%REC	1	19-Mar-2020 18:27
<i>Surr: 2-Fluorobiphenyl</i>	68.6			40-125	%REC	1	19-Mar-2020 18:27
<i>Surr: 2-Fluorophenol</i>	64.7			20-120	%REC	1	19-Mar-2020 18:27
<i>Surr: 4-Terphenyl-d14</i>	104			40-135	%REC	1	19-Mar-2020 18:27
<i>Surr: Nitrobenzene-d5</i>	51.5			41-120	%REC	1	19-Mar-2020 18:27
<i>Surr: Phenol-d6</i>	58.4			20-120	%REC	1	19-Mar-2020 18:27
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00355</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Mar-2020 13:41

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW61B-20200317  
 Collection Date: 17-Mar-2020 11:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:11
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:11
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:11
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:11
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 22:11
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:11
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:11
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>99.1</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:11</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.5</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:11</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.8</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:11</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:11</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW61B-20200317  
 Collection Date: 17-Mar-2020 11:45

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-03  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 18:46
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 18:46
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 18:46
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 18:46
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 18:46
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 18:46
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 18:46
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 18:46
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	19-Mar-2020 18:46
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 18:46
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 18:46
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 18:46
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 18:46
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 18:46
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00015</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	19-Mar-2020 18:46
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 18:46
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 18:46
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 18:46
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	19-Mar-2020 18:46
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	19-Mar-2020 18:46
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 18:46
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 18:46
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 18:46
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 18:46
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 18:46
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 18:46
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 18:46
<i>Surr: 2,4,6-Tribromophenol</i>	<i>64.0</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 18:46</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>62.5</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 18:46</i>
<i>Surr: 2-Fluorophenol</i>	<i>50.3</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 18:46</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>79.6</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 18:46</i>
<i>Surr: Nitrobenzene-d5</i>	<i>46.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 18:46</i>
<i>Surr: Phenol-d6</i>	<i>44.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 18:46</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00461</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Mar-2020 14:44

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW96B-20200317  
 Collection Date: 17-Mar-2020 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-04  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:34
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:34
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:34
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:34
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 22:34
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:34
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:34
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:34</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.2</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:34</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:34</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:34</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW96B-20200317  
 Collection Date: 17-Mar-2020 10:25

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-04  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 19:05
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 19:05
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 19:05
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 19:05
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 19:05
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 19:05
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 19:05
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 19:05
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	19-Mar-2020 19:05
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 19:05
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 19:05
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 19:05
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:05
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 19:05
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00035</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	19-Mar-2020 19:05
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 19:05
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:05
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 19:05
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	19-Mar-2020 19:05
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	19-Mar-2020 19:05
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:05
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 19:05
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 19:05
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 19:05
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 19:05
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 19:05
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 19:05
<i>Surr: 2,4,6-Tribromophenol</i>	<i>47.2</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 19:05</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>62.3</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 19:05</i>
<i>Surr: 2-Fluorophenol</i>	<i>48.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 19:05</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>86.2</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 19:05</i>
<i>Surr: Nitrobenzene-d5</i>	<i>45.0</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 19:05</i>
<i>Surr: Phenol-d6</i>	<i>41.8</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 19:05</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00312</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Mar-2020 14:46

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW95A-20200317  
 Collection Date: 17-Mar-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-05  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:57
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:57
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:57
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:57
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 22:57
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 22:57
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 22:57
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>100</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:57</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>95.9</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:57</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.6</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:57</i>
<i>Surr: Toluene-d8</i>	<i>103</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 22:57</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW95A-20200317  
 Collection Date: 17-Mar-2020 09:10

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-05  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 19:24
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 19:24
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 19:24
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 19:24
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 19:24
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 19:24
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 19:24
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 19:24
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	19-Mar-2020 19:24
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 19:24
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 19:24
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 19:24
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:24
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 19:24
Bis(2-ethylhexyl)phthalate	< 0.000037		0.000037	0.00020	mg/L	1	19-Mar-2020 19:24
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 19:24
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:24
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 19:24
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	19-Mar-2020 19:24
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	19-Mar-2020 19:24
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:24
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 19:24
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 19:24
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 19:24
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 19:24
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 19:24
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 19:24
Surr: 2,4,6-Tribromophenol	60.4			34-129	%REC	1	19-Mar-2020 19:24
Surr: 2-Fluorobiphenyl	61.4			40-125	%REC	1	19-Mar-2020 19:24
Surr: 2-Fluorophenol	57.8			20-120	%REC	1	19-Mar-2020 19:24
Surr: 4-Terphenyl-d14	82.2			40-135	%REC	1	19-Mar-2020 19:24
Surr: Nitrobenzene-d5	48.9			41-120	%REC	1	19-Mar-2020 19:24
Surr: Phenol-d6	45.1			20-120	%REC	1	19-Mar-2020 19:24
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
Arsenic	0.000977	J	0.000400	0.00200	mg/L	1	27-Mar-2020 14:48

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW60B-20200317  
 Collection Date: 17-Mar-2020 15:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-06  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 23:20
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 23:20
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 23:20
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 23:20
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 23:20
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 23:20
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 23:20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>96.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:20</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:20</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:20</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW60B-20200317  
 Collection Date: 17-Mar-2020 15:15

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-06  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 19:43
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 19:43
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 19:43
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 19:43
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 19:43
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 19:43
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 19:43
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 19:43
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	19-Mar-2020 19:43
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 19:43
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 19:43
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 19:43
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:43
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 19:43
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00092</b>		<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	19-Mar-2020 19:43
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 19:43
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:43
<b>Di-n-butyl phthalate</b>	<b>0.000047</b>	J	<b>0.000020</b>	<b>0.00020</b>	<b>mg/L</b>	1	19-Mar-2020 19:43
<b>Fluoranthene</b>	<b>0.000015</b>	J	<b>0.000010</b>	<b>0.00010</b>	<b>mg/L</b>	1	19-Mar-2020 19:43
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	19-Mar-2020 19:43
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 19:43
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 19:43
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 19:43
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 19:43
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 19:43
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 19:43
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 19:43
<i>Surr: 2,4,6-Tribromophenol</i>	74.8			34-129	%REC	1	19-Mar-2020 19:43
<i>Surr: 2-Fluorobiphenyl</i>	50.1			40-125	%REC	1	19-Mar-2020 19:43
<i>Surr: 2-Fluorophenol</i>	47.8			20-120	%REC	1	19-Mar-2020 19:43
<i>Surr: 4-Terphenyl-d14</i>	82.6			40-135	%REC	1	19-Mar-2020 19:43
<i>Surr: Nitrobenzene-d5</i>	41.6			41-120	%REC	1	19-Mar-2020 19:43
<i>Surr: Phenol-d6</i>	43.0			20-120	%REC	1	19-Mar-2020 19:43
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00234</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Mar-2020 14:51

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88B-20200317  
 Collection Date: 17-Mar-2020 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-07  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 23:43
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 23:43
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 23:43
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 23:43
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 23:43
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 23:43
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 23:43
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>102</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:43</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.0</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:43</i>
<i>Surr: Dibromofluoromethane</i>	<i>100</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:43</i>
<i>Surr: Toluene-d8</i>	<i>101</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 23:43</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.



Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-MW88B-20200317  
 Collection Date: 17-Mar-2020 13:50

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-07  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 20:02
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 20:02
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 20:02
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 20:02
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 20:02
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 20:02
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 20:02
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 20:02
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	19-Mar-2020 20:02
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 20:02
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 20:02
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 20:02
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 20:02
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 20:02
Bis(2-ethylhexyl)phthalate	< 0.000037		0.000037	0.00020	mg/L	1	19-Mar-2020 20:02
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 20:02
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 20:02
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 20:02
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	19-Mar-2020 20:02
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	19-Mar-2020 20:02
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 20:02
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 20:02
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 20:02
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 20:02
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 20:02
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 20:02
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 20:02
Surr: 2,4,6-Tribromophenol	59.9			34-129	%REC	1	19-Mar-2020 20:02
Surr: 2-Fluorobiphenyl	63.9			40-125	%REC	1	19-Mar-2020 20:02
Surr: 2-Fluorophenol	55.5			20-120	%REC	1	19-Mar-2020 20:02
Surr: 4-Terphenyl-d14	93.3			40-135	%REC	1	19-Mar-2020 20:02
Surr: Nitrobenzene-d5	52.0			41-120	%REC	1	19-Mar-2020 20:02
Surr: Phenol-d6	46.3			20-120	%REC	1	19-Mar-2020 20:02
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
Arsenic	0.00166	J	0.000400	0.00200	mg/L	1	27-Mar-2020 14:53

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD02-20200317  
 Collection Date: 17-Mar-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-08  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW LEVEL VOLATILES BY SW8260C</b>		<b>Method:SW8260</b>		Analyst: AKP			
1,2-Dichloroethane	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 20:16
Benzene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 20:16
Chlorobenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 20:16
Ethylbenzene	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 20:16
Methylene chloride	< 0.0010		0.0010	0.0020	mg/L	1	23-Mar-2020 20:16
Toluene	< 0.00020		0.00020	0.0010	mg/L	1	23-Mar-2020 20:16
Xylenes, Total	< 0.00030		0.00030	0.0010	mg/L	1	23-Mar-2020 20:16
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>101</i>			<i>70-126</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:16</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>97.1</i>			<i>81-113</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:16</i>
<i>Surr: Dibromofluoromethane</i>	<i>98.3</i>			<i>77-123</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:16</i>
<i>Surr: Toluene-d8</i>	<i>100</i>			<i>82-127</i>	<i>%REC</i>	<i>1</i>	<i>23-Mar-2020 20:16</i>

Note: See Qualifiers Page for a list of qualifiers and their explanation.

Client: Golder Associates Inc.  
 Project: Houston TX-Wood Preserving Works  
 Sample ID: WG-1620-FD02-20200317  
 Collection Date: 17-Mar-2020 00:00

**ANALYTICAL REPORT**  
 WorkOrder:HS20030819  
 Lab ID:HS20030819-08  
 Matrix:Groundwater

ANALYSES	RESULT	QUAL	SDL	MQL	UNITS	DILUTION FACTOR	DATE ANALYZED
<b>LOW-LEVEL SEMIVOLATILES BY 8270D</b>		<b>Method:SW8270</b>		Prep:SW3510 / 19-Mar-2020		Analyst: GEY	
1,2-Diphenylhydrazine	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 20:22
2,4-Dimethylphenol	< 0.000040		0.000040	0.00020	mg/L	1	19-Mar-2020 20:22
2,4-Dinitrotoluene	< 0.000058		0.000058	0.00020	mg/L	1	19-Mar-2020 20:22
2,6-Dinitrotoluene	< 0.000042		0.000042	0.00020	mg/L	1	19-Mar-2020 20:22
2-Chloronaphthalene	< 0.000021		0.000021	0.00020	mg/L	1	19-Mar-2020 20:22
2-Methylnaphthalene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 20:22
4,6-Dinitro-2-methylphenol	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 20:22
4-Nitrophenol	< 0.000047		0.000047	0.0010	mg/L	1	19-Mar-2020 20:22
Acenaphthene	< 0.000027		0.000027	0.00010	mg/L	1	19-Mar-2020 20:22
Acenaphthylene	< 0.000015		0.000015	0.00010	mg/L	1	19-Mar-2020 20:22
Anthracene	< 0.000014		0.000014	0.00010	mg/L	1	19-Mar-2020 20:22
Benz(a)anthracene	< 0.000050		0.000050	0.00010	mg/L	1	19-Mar-2020 20:22
Benzo(a)pyrene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 20:22
Bis(2-chloroethoxy)methane	< 0.000030		0.000030	0.00020	mg/L	1	19-Mar-2020 20:22
<b>Bis(2-ethylhexyl)phthalate</b>	<b>0.00013</b>	J	<b>0.000037</b>	<b>0.00020</b>	<b>mg/L</b>	1	19-Mar-2020 20:22
Chrysene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 20:22
Dibenzofuran	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 20:22
Di-n-butyl phthalate	< 0.000020		0.000020	0.00020	mg/L	1	19-Mar-2020 20:22
Fluoranthene	< 0.000010		0.000010	0.00010	mg/L	1	19-Mar-2020 20:22
Fluorene	< 0.000030		0.000030	0.00010	mg/L	1	19-Mar-2020 20:22
Naphthalene	< 0.000020		0.000020	0.00010	mg/L	1	19-Mar-2020 20:22
Nitrobenzene	< 0.000024		0.000024	0.00020	mg/L	1	19-Mar-2020 20:22
N-Nitrosodiphenylamine	< 0.000025		0.000025	0.00020	mg/L	1	19-Mar-2020 20:22
Pentachlorophenol	< 0.000079		0.000079	0.00020	mg/L	1	19-Mar-2020 20:22
Phenanthrene	< 0.000021		0.000021	0.00010	mg/L	1	19-Mar-2020 20:22
Phenol	< 0.000035		0.000035	0.00020	mg/L	1	19-Mar-2020 20:22
Pyrene	< 0.000019		0.000019	0.00010	mg/L	1	19-Mar-2020 20:22
<i>Surr: 2,4,6-Tribromophenol</i>	<i>57.8</i>			<i>34-129</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 20:22</i>
<i>Surr: 2-Fluorobiphenyl</i>	<i>59.0</i>			<i>40-125</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 20:22</i>
<i>Surr: 2-Fluorophenol</i>	<i>48.5</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 20:22</i>
<i>Surr: 4-Terphenyl-d14</i>	<i>89.9</i>			<i>40-135</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 20:22</i>
<i>Surr: Nitrobenzene-d5</i>	<i>47.3</i>			<i>41-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 20:22</i>
<i>Surr: Phenol-d6</i>	<i>44.4</i>			<i>20-120</i>	<i>%REC</i>	<i>1</i>	<i>19-Mar-2020 20:22</i>
<b>ICP-MS METALS BY SW6020A</b>		<b>Method:SW6020</b>		Prep:SW3010A / 26-Mar-2020		Analyst: JHD	
<b>Arsenic</b>	<b>0.00311</b>		<b>0.000400</b>	<b>0.00200</b>	<b>mg/L</b>	1	27-Mar-2020 14:55

Note: See Qualifiers Page for a list of qualifiers and their explanation.

## Weight / Prep Log

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**Batch ID:** 151876      **Start Date:** 19 Mar 2020 07:00      **End Date:** 19 Mar 2020 12:30  
**Method:** SV AQ SEP FUN EXTRACT-LOWLEV - 3510C      **Prep Code:** 3510\_B\_LOW

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030819-02	1	1000 (mL)	1 (mL)	0.001
HS20030819-03	1	1000 (mL)	1 (mL)	0.001
HS20030819-04	1	1000 (mL)	1 (mL)	0.001
HS20030819-05	1	1000 (mL)	1 (mL)	0.001
HS20030819-06	1	1000 (mL)	1 (mL)	0.001
HS20030819-07	1	1000 (mL)	1 (mL)	0.001
HS20030819-08	1	1000 (mL)	1 (mL)	0.001

**Batch ID:** 152162      **Start Date:** 26 Mar 2020 08:57      **End Date:** 26 Mar 2020 13:00  
**Method:** WATER - SW3010A      **Prep Code:** 3010A

Sample ID	Container	Sample Wt/Vol	Final Volume	Prep Factor
HS20030819-02		10 (mL)	10 (mL)	1
HS20030819-03		10 (mL)	10 (mL)	1
HS20030819-04		10 (mL)	10 (mL)	1
HS20030819-05		10 (mL)	10 (mL)	1
HS20030819-06		10 (mL)	10 (mL)	1
HS20030819-07		10 (mL)	10 (mL)	1
HS20030819-08		10 (mL)	10 (mL)	1

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**DATES REPORT**

Sample ID	Client Samp ID	Collection Date	Leachate Date	Prep Date	Analysis Date	DF
<b>Batch ID: 151876 ( 0 )</b>		<b>Test Name : LOW-LEVEL SEMIVOLATILES BY 8270D</b>			<b>Matrix: Groundwater</b>	
HS20030819-02	WG-1620-MW88A-20200317	17 Mar 2020 12:55		19 Mar 2020 10:38	19 Mar 2020 18:27	1
HS20030819-03	WG-1620-MW61B-20200317	17 Mar 2020 11:45		19 Mar 2020 10:38	19 Mar 2020 18:46	1
HS20030819-04	WG-1620-MW96B-20200317	17 Mar 2020 10:25		19 Mar 2020 10:38	19 Mar 2020 19:05	1
HS20030819-05	WG-1620-MW95A-20200317	17 Mar 2020 09:10		19 Mar 2020 10:38	19 Mar 2020 19:24	1
HS20030819-06	WG-1620-MW60B-20200317	17 Mar 2020 15:15		19 Mar 2020 10:38	19 Mar 2020 19:43	1
HS20030819-07	WG-1620-MW88B-20200317	17 Mar 2020 13:50		19 Mar 2020 10:38	19 Mar 2020 20:02	1
HS20030819-08	WG-1620-FD02-20200317	17 Mar 2020 00:00		19 Mar 2020 10:38	19 Mar 2020 20:22	1
<b>Batch ID: 152162 ( 0 )</b>		<b>Test Name : ICP-MS METALS BY SW6020A</b>			<b>Matrix: Groundwater</b>	
HS20030819-02	WG-1620-MW88A-20200317	17 Mar 2020 12:55		26 Mar 2020 13:00	27 Mar 2020 13:41	1
HS20030819-03	WG-1620-MW61B-20200317	17 Mar 2020 11:45		26 Mar 2020 13:00	27 Mar 2020 14:44	1
HS20030819-04	WG-1620-MW96B-20200317	17 Mar 2020 10:25		26 Mar 2020 13:00	27 Mar 2020 14:46	1
HS20030819-05	WG-1620-MW95A-20200317	17 Mar 2020 09:10		26 Mar 2020 13:00	27 Mar 2020 14:48	1
HS20030819-06	WG-1620-MW60B-20200317	17 Mar 2020 15:15		26 Mar 2020 13:00	27 Mar 2020 14:51	1
HS20030819-07	WG-1620-MW88B-20200317	17 Mar 2020 13:50		26 Mar 2020 13:00	27 Mar 2020 14:53	1
HS20030819-08	WG-1620-FD02-20200317	17 Mar 2020 00:00		26 Mar 2020 13:00	27 Mar 2020 14:55	1
<b>Batch ID: R358799 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Groundwater</b>	
HS20030819-02	WG-1620-MW88A-20200317	17 Mar 2020 12:55			23 Mar 2020 20:39	1
HS20030819-03	WG-1620-MW61B-20200317	17 Mar 2020 11:45			23 Mar 2020 22:11	1
HS20030819-04	WG-1620-MW96B-20200317	17 Mar 2020 10:25			23 Mar 2020 22:34	1
HS20030819-05	WG-1620-MW95A-20200317	17 Mar 2020 09:10			23 Mar 2020 22:57	1
HS20030819-06	WG-1620-MW60B-20200317	17 Mar 2020 15:15			23 Mar 2020 23:20	1
HS20030819-07	WG-1620-MW88B-20200317	17 Mar 2020 13:50			23 Mar 2020 23:43	1
HS20030819-08	WG-1620-FD02-20200317	17 Mar 2020 00:00			23 Mar 2020 20:16	1
<b>Batch ID: R358799 ( 0 )</b>		<b>Test Name : LOW LEVEL VOLATILES BY SW8260C</b>			<b>Matrix: Water</b>	
HS20030819-01	WQ-1620-TB01-20200317	17 Mar 2020 15:59			23 Mar 2020 19:53	1

WorkOrder: HS20030819  
 InstrumentID: ICPMS05  
 Test Code: ICP\_TW  
 Test Number: SW6020  
 Test Name: ICP-MS Metals by SW6020A

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous      **Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	Arsenic	7440-38-2	0.000500	0.000499	0.000400	0.00200

WorkOrder: HS20030819  
 InstrumentID: SV-7  
 Test Code: 8270\_LOW\_W  
 Test Number: SW8270  
 Test Name: Low-Level Semivolatiles by 8270D

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Diphenylhydrazine	122-66-7	0.00010	0.00013	0.000021	0.00020
A	2,4-Dimethylphenol	105-67-9	0.00010	0.000081	0.000040	0.00020
A	2,4-Dinitrotoluene	121-14-2	0.00010	0.00010	0.000058	0.00020
A	2,6-Dinitrotoluene	606-20-2	0.00010	0.00010	0.000042	0.00020
A	2-Chloronaphthalene	91-58-7	0.00010	0.00012	0.000021	0.00020
A	2-Methylnaphthalene	91-57-6	0.000050	0.000052	0.000019	0.00010
A	4,6-Dinitro-2-methylphenol	534-52-1	0.00010	0.000060	0.000020	0.00020
A	4-Nitrophenol	100-02-7	0.00010	0.000069	0.000047	0.0010
A	Acenaphthene	83-32-9	0.000050	0.000050	0.000027	0.00010
A	Acenaphthylene	208-96-8	0.000050	0.000046	0.000015	0.00010
A	Anthracene	120-12-7	0.000050	0.000046	0.000014	0.00010
A	Benz(a)anthracene	56-55-3	0.000050	0.000057	0.000050	0.00010
A	Benzo(a)pyrene	50-32-8	0.000050	0.000047	0.000020	0.00010
A	Bis(2-chloroethoxy)methane	111-91-1	0.00010	0.00012	0.000030	0.00020
A	Bis(2-ethylhexyl)phthalate	117-81-7	0.00010	0.00014	0.000037	0.00020
A	Chrysene	218-01-9	0.000050	0.000049	0.000021	0.00010
A	Dibenzofuran	132-64-9	0.000050	0.000050	0.000020	0.00010
A	Di-n-butyl phthalate	84-74-2	0.00010	0.00012	0.000020	0.00020
A	Fluoranthene	206-44-0	0.000050	0.000053	0.000010	0.00010
A	Fluorene	86-73-7	0.000050	0.000050	0.000030	0.00010
A	Naphthalene	91-20-3	0.000050	0.000048	0.000020	0.00010
A	Nitrobenzene	98-95-3	0.00010	0.00014	0.000024	0.00020
A	N-Nitrosodiphenylamine	86-30-6	0.00010	0.00012	0.000025	0.00020
A	Pentachlorophenol	87-86-5	0.00010	0.000057	0.000079	0.00020
A	Phenanthrene	85-01-8	0.000050	0.000047	0.000021	0.00010
A	Phenol	108-95-2	0.00010	0.00012	0.000035	0.00020
A	Pyrene	129-00-0	0.000050	0.000051	0.000019	0.00010
S	2,4,6-Tribromophenol	118-79-6	0	0	0	0.00020
S	2-Fluorobiphenyl	321-60-8	0	0	0	0.00020
S	2-Fluorophenol	367-12-4	0	0	0	0.00020
S	4-Terphenyl-d14	1718-51-0	0	0	0	0.00020
S	Nitrobenzene-d5	4165-60-0	0	0	0	0.00020
S	Phenol-d6	13127-88-3	0	0	0	0.00020

WorkOrder: HS20030819  
 InstrumentID: VOA2  
 Test Code: 8260\_LL\_W  
 Test Number: SW8260  
 Test Name: Low Level Volatiles by SW8260C

**METHOD DETECTION /  
 REPORTING LIMITS**

**Matrix:** Aqueous

**Units:** mg/L

Type	Analyte	CAS	DCS Spike	DCS	MDL	PQL
A	1,2-Dichloroethane	107-06-2	0.00050	0.00062	0.00020	0.0010
A	Benzene	71-43-2	0.00050	0.00059	0.00020	0.0010
A	Chlorobenzene	108-90-7	0.00050	0.00071	0.00030	0.0010
A	Ethylbenzene	100-41-4	0.00050	0.00063	0.00030	0.0010
A	Methylene chloride	75-09-2	0.0010	0.00075	0.0010	0.0020
A	Toluene	108-88-3	0.00050	0.00069	0.00020	0.0010
A	Xylenes, Total	1330-20-7	0.00050	0.00059	0.00030	0.0010
S	1,2-Dichloroethane-d4	17060-07-0	0	0	0	0.0010
S	4-Bromofluorobenzene	460-00-4	0	0	0	0.0010
S	Dibromofluoromethane	1868-53-7	0	0	0	0.0010
S	Toluene-d8	2037-26-5	0	0	0	0.0010



**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QC BATCH REPORT**

Batch ID: 152162 ( 0 )		Instrument: ICPMS05		Method: ICP-MS METALS BY SW6020A					
<b>MBLK</b>	Sample ID: <b>MBLK-152162</b>	Units: <b>mg/L</b>		Analysis Date: <b>27-Mar-2020 13:36</b>					
Client ID:	Run ID: <b>ICPMS05_359030</b>	SeqNo: <b>5534467</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	< 0.000400	0.00200							
<b>LCS</b>	Sample ID: <b>LCS-152162</b>	Units: <b>mg/L</b>		Analysis Date: <b>27-Mar-2020 13:39</b>					
Client ID:	Run ID: <b>ICPMS05_359030</b>	SeqNo: <b>5534468</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.05069	0.00200	0.05	0	101	80 - 120			
<b>MS</b>	Sample ID: <b>HS20031033-01MS</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Mar-2020 11:14</b>					
Client ID:	Run ID: <b>ICPMS05_359110</b>	SeqNo: <b>5536564</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.1729	0.00200	0.05	0.1254	95.0	80 - 120			
<b>MSD</b>	Sample ID: <b>HS20031033-01MSD</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Mar-2020 11:17</b>					
Client ID:	Run ID: <b>ICPMS05_359110</b>	SeqNo: <b>5536565</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.1842	0.00200	0.05	0.1254	118	80 - 120	0.1729	6.31 20	
<b>PDS</b>	Sample ID: <b>HS20031033-01PDS</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Mar-2020 11:19</b>					
Client ID:	Run ID: <b>ICPMS05_359110</b>	SeqNo: <b>5536566</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>1</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	RPD %RPD Limit Qual	
Arsenic	0.2352	0.00200	0.1	0.1254	110	75 - 125			
<b>SD</b>	Sample ID: <b>HS20031033-01SD</b>	Units: <b>mg/L</b>		Analysis Date: <b>30-Mar-2020 11:12</b>					
Client ID:	Run ID: <b>ICPMS05_359110</b>	SeqNo: <b>5536563</b>		PrepDate: <b>26-Mar-2020</b>		DF: <b>5</b>			
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%D %D Limit Qual	
Arsenic	0.1248	0.0100					0.1254	0.469 10	
The following samples were analyzed in this batch:									
	HS20030819-02	HS20030819-03	HS20030819-04	HS20030819-05	HS20030819-06	HS20030819-07	HS20030819-08		

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QC BATCH REPORT**

Batch ID: 151876 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
MBLK	Sample ID: MBLK-151876	Units: ug/L			Analysis Date: 19-Mar-2020 13:39					
Client ID:	Run ID: SV-7_358509	SeqNo: 5520890	PrepDate: 19-Mar-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	< 0.021	0.20								
2,4-Dimethylphenol	< 0.040	0.20								
2,4-Dinitrotoluene	< 0.058	0.20								
2,6-Dinitrotoluene	< 0.042	0.20								
2-Chloronaphthalene	< 0.021	0.20								
2-Methylnaphthalene	< 0.019	0.10								
4,6-Dinitro-2-methylphenol	< 0.020	0.20								
4-Nitrophenol	< 0.047	1.0								
Acenaphthene	< 0.027	0.10								
Acenaphthylene	< 0.015	0.10								
Anthracene	< 0.014	0.10								
Benz(a)anthracene	< 0.050	0.10								
Benzo(a)pyrene	< 0.020	0.10								
Bis(2-chloroethoxy)methane	< 0.030	0.20								
Bis(2-ethylhexyl)phthalate	< 0.037	0.20								
Chrysene	< 0.021	0.10								
Dibenzofuran	< 0.020	0.10								
Di-n-butyl phthalate	< 0.020	0.20								
Fluoranthene	< 0.010	0.10								
Fluorene	< 0.030	0.10								
Naphthalene	< 0.020	0.10								
Nitrobenzene	< 0.024	0.20								
N-Nitrosodiphenylamine	< 0.025	0.20								
Pentachlorophenol	< 0.079	0.20								
Phenanthrene	< 0.021	0.10								
Phenol	< 0.035	0.20								
Pyrene	< 0.019	0.10								
<i>Surr: 2,4,6-Tribromophenol</i>	2.949	0.20	5	0	59.0	34 - 129				
<i>Surr: 2-Fluorobiphenyl</i>	2.962	0.20	5	0	59.2	40 - 125				
<i>Surr: 2-Fluorophenol</i>	2.88	0.20	5	0	57.6	20 - 120				
<i>Surr: 4-Terphenyl-d14</i>	3.57	0.20	5	0	71.4	40 - 135				
<i>Surr: Nitrobenzene-d5</i>	2.498	0.20	5	0	50.0	41 - 120				
<i>Surr: Phenol-d6</i>	2.798	0.20	5	0	56.0	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QC BATCH REPORT**

Batch ID: 151876 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
LCS	Sample ID: LCS-151876	Units: ug/L			Analysis Date: 19-Mar-2020 15:15					
Client ID:	Run ID: SV-7_358509	SeqNo: 5520894	PrepDate: 19-Mar-2020	DF: 1						
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	3.006	0.20	5	0	60.1	39 - 127				
2,4-Dimethylphenol	3.069	0.20	5	0	61.4	35 - 120				
2,4-Dinitrotoluene	3.767	0.20	5	0	75.3	50 - 122				
2,6-Dinitrotoluene	3.698	0.20	5	0	74.0	50 - 120				
2-Chloronaphthalene	4.153	0.20	5	0	83.1	50 - 120				
2-Methylnaphthalene	3.43	0.10	5	0	68.6	50 - 120				
4,6-Dinitro-2-methylphenol	3.658	0.20	5	0	73.2	25 - 121				
4-Nitrophenol	2.531	1.0	5	0	50.6	30 - 130				
Acenaphthene	3.225	0.10	5	0	64.5	45 - 120				
Acenaphthylene	3.269	0.10	5	0	65.4	47 - 120				
Anthracene	3.472	0.10	5	0	69.4	45 - 120				
Benz(a)anthracene	3.507	0.10	5	0	70.1	40 - 120				
Benzo(a)pyrene	3.777	0.10	5	0	75.5	45 - 120				
Bis(2-chloroethoxy)methane	3.186	0.20	5	0	63.7	45 - 120				
Bis(2-ethylhexyl)phthalate	3.478	0.20	5	0	69.6	40 - 139				
Chrysene	3.344	0.10	5	0	66.9	43 - 120				
Dibenzofuran	3.454	0.10	5	0	69.1	50 - 120				
Di-n-butyl phthalate	3.575	0.20	5	0	71.5	45 - 123				
Fluoranthene	3.531	0.10	5	0	70.6	45 - 125				
Fluorene	3.427	0.10	5	0	68.5	49 - 120				
Naphthalene	3.329	0.10	5	0	66.6	45 - 120				
Nitrobenzene	2.977	0.20	5	0	59.5	44 - 120				
N-Nitrosodiphenylamine	3.366	0.20	5	0	67.3	40 - 125				
Pentachlorophenol	1.871	0.20	5	0	37.4	19 - 121				
Phenanthrene	3.392	0.10	5	0	67.8	45 - 121				
Phenol	3.406	0.20	5	0	68.1	20 - 124				
Pyrene	3.501	0.10	5	0	70.0	40 - 130				
Surr: 2,4,6-Tribromophenol	3.613	0.20	5	0	72.3	34 - 129				
Surr: 2-Fluorobiphenyl	2.764	0.20	5	0	55.3	40 - 125				
Surr: 2-Fluorophenol	2.851	0.20	5	0	57.0	20 - 120				
Surr: 4-Terphenyl-d14	3.276	0.20	5	0	65.5	40 - 135				
Surr: Nitrobenzene-d5	2.453	0.20	5	0	49.1	41 - 120				
Surr: Phenol-d6	2.853	0.20	5	0	57.1	20 - 120				

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QC BATCH REPORT**

Batch ID: 151876 ( 0 )		Instrument: SV-7		Method: LOW-LEVEL SEMIVOLATILES BY 8270D						
<b>LCSD</b>	Sample ID: <b>LCSD-151876</b>	Units: <b>ug/L</b>			Analysis Date: <b>19-Mar-2020 14:18</b>					
Client ID:	Run ID: <b>SV-7_358509</b>	SeqNo: <b>5520891</b>		PrepDate: <b>19-Mar-2020</b>		DF: <b>1</b>				
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
1,2-Diphenylhydrazine	2.865	0.20	5	0	57.3	39 - 127	3.006	4.83	20	
2,4-Dimethylphenol	3.104	0.20	5	0	62.1	35 - 120	3.069	1.1	20	
2,4-Dinitrotoluene	3.649	0.20	5	0	73.0	50 - 122	3.767	3.18	20	
2,6-Dinitrotoluene	3.677	0.20	5	0	73.5	50 - 120	3.698	0.567	20	
2-Chloronaphthalene	3.78	0.20	5	0	75.6	50 - 120	4.153	9.39	20	
2-Methylnaphthalene	3.4	0.10	5	0	68.0	50 - 120	3.43	0.881	20	
4,6-Dinitro-2-methylphenol	3.494	0.20	5	0	69.9	25 - 121	3.658	4.58	30	
4-Nitrophenol	2.988	1.0	5	0	59.8	30 - 130	2.531	16.5	20	
Acenaphthene	3.232	0.10	5	0	64.6	45 - 120	3.225	0.23	20	
Acenaphthylene	3.295	0.10	5	0	65.9	47 - 120	3.269	0.783	20	
Anthracene	3.376	0.10	5	0	67.5	45 - 120	3.472	2.8	20	
Benz(a)anthracene	3.682	0.10	5	0	73.6	40 - 120	3.507	4.86	20	
Benzo(a)pyrene	3.802	0.10	5	0	76.0	45 - 120	3.777	0.658	20	
Bis(2-chloroethoxy)methane	3.188	0.20	5	0	63.8	45 - 120	3.186	0.0664	20	
Bis(2-ethylhexyl)phthalate	3.54	0.20	5	0	70.8	40 - 139	3.478	1.76	20	
Chrysene	3.571	0.10	5	0	71.4	43 - 120	3.344	6.57	20	
Dibenzofuran	3.429	0.10	5	0	68.6	50 - 120	3.454	0.747	20	
Di-n-butyl phthalate	3.482	0.20	5	0	69.6	45 - 123	3.575	2.64	20	
Fluoranthene	3.413	0.10	5	0	68.3	45 - 125	3.531	3.4	20	
Fluorene	3.408	0.10	5	0	68.2	49 - 120	3.427	0.573	20	
Naphthalene	3.336	0.10	5	0	66.7	45 - 120	3.329	0.225	20	
Nitrobenzene	2.957	0.20	5	0	59.1	44 - 120	2.977	0.66	20	
N-Nitrosodiphenylamine	3.264	0.20	5	0	65.3	40 - 125	3.366	3.1	20	
Pentachlorophenol	1.812	0.20	5	0	36.2	19 - 121	1.871	3.23	20	
Phenanthrene	3.322	0.10	5	0	66.4	45 - 121	3.392	2.07	20	
Phenol	3.274	0.20	5	0	65.5	20 - 124	3.406	3.96	20	
Pyrene	3.554	0.10	5	0	71.1	40 - 130	3.501	1.51	20	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>3.602</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>72.0</i>	<i>34 - 129</i>	<i>3.613</i>	<i>0.299</i>	<i>20</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>2.776</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>55.5</i>	<i>40 - 125</i>	<i>2.764</i>	<i>0.445</i>	<i>20</i>	
<i>Surr: 2-Fluorophenol</i>	<i>2.733</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>54.7</i>	<i>20 - 120</i>	<i>2.851</i>	<i>4.25</i>	<i>20</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>3.381</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>67.6</i>	<i>40 - 135</i>	<i>3.276</i>	<i>3.15</i>	<i>20</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>2.448</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>49.0</i>	<i>41 - 120</i>	<i>2.453</i>	<i>0.189</i>	<i>20</i>	
<i>Surr: Phenol-d6</i>	<i>2.753</i>	<i>0.20</i>	<i>5</i>	<i>0</i>	<i>55.1</i>	<i>20 - 120</i>	<i>2.853</i>	<i>3.58</i>	<i>20</i>	

The following samples were analyzed in this batch: HS20030819-02 HS20030819-03 HS20030819-04 HS20030819-05  
 HS20030819-06 HS20030819-07 HS20030819-08

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QC BATCH REPORT**

**Batch ID:** R358799 ( 0 )      **Instrument:** VOA2      **Method:** LOW LEVEL VOLATILES BY SW8260C

<b>MBLK</b>		Sample ID: <b>VBLKW-200323</b>		Units: <b>ug/L</b>		Analysis Date: <b>23-Mar-2020 19:22</b>			
Client ID:		Run ID: <b>VOA2_358799</b>		SeqNo: <b>5527631</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	< 0.20	1.0							
Benzene	< 0.20	1.0							
Chlorobenzene	< 0.30	1.0							
Ethylbenzene	< 0.30	1.0							
Methylene chloride	< 1.0	2.0							
Toluene	< 0.20	1.0							
Xylenes, Total	< 0.30	1.0							
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>51.61</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>103</i>	<i>70 - 123</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>47.84</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>95.7</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>49.76</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.5</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>51.01</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>102</i>	<i>81 - 120</i>			

<b>LCS</b>		Sample ID: <b>VLCSW-203023</b>		Units: <b>ug/L</b>		Analysis Date: <b>23-Mar-2020 18:36</b>			
Client ID:		Run ID: <b>VOA2_358799</b>		SeqNo: <b>5527630</b>		PrepDate:		DF: <b>1</b>	
Analyte	Result	MLQ	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	19.11	1.0	20	0	95.6	70 - 124			
Benzene	18.46	1.0	20	0	92.3	74 - 120			
Chlorobenzene	17.94	1.0	20	0	89.7	76 - 113			
Ethylbenzene	18.37	1.0	20	0	91.9	77 - 117			
Methylene chloride	17.55	2.0	20	0	87.7	70 - 127			
Toluene	18.17	1.0	20	0	90.9	77 - 118			
Xylenes, Total	55.92	1.0	60	0	93.2	75 - 122			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>53.36</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>107</i>	<i>70 - 130</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.6</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.2</i>	<i>82 - 115</i>			
<i>Surr: Dibromofluoromethane</i>	<i>48.48</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.0</i>	<i>73 - 126</i>			
<i>Surr: Toluene-d8</i>	<i>49.98</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100.0</i>	<i>81 - 120</i>			

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QC BATCH REPORT**

Batch ID: R358799 ( 0 )		Instrument: VOA2		Method: LOW LEVEL VOLATILES BY SW8260C						
<b>MS</b>		Sample ID: HS20030819-02MS		Units: ug/L		Analysis Date: 23-Mar-2020 21:02				
Client ID: WG-1620-MW88A-20200317		Run ID: VOA2_358799		SeqNo: 5527635		PrepDate:		DF: 1		
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual	
1,2-Dichloroethane	17.84	1.0	20	0	89.2	70 - 127				
Benzene	18.14	1.0	20	0	90.7	70 - 127				
Chlorobenzene	17.35	1.0	20	0	86.8	70 - 114				
Ethylbenzene	18.32	1.0	20	0	91.6	70 - 124				
Methylene chloride	16.47	2.0	20	0	82.3	70 - 128				
Toluene	17.75	1.0	20	0	88.8	70 - 123				
Xylenes, Total	54.81	1.0	60	0	91.4	70 - 130				
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>52.67</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>70 - 126</i>				
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.7</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.4</i>	<i>81 - 113</i>				
<i>Surr: Dibromofluoromethane</i>	<i>48.94</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.9</i>	<i>77 - 123</i>				
<i>Surr: Toluene-d8</i>	<i>49.9</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>99.8</i>	<i>82 - 127</i>				

<b>MSD</b>		Sample ID: HS20030819-02MSD		Units: ug/L		Analysis Date: 23-Mar-2020 21:25			
Client ID: WG-1620-MW88A-20200317		Run ID: VOA2_358799		SeqNo: 5527636		PrepDate:		DF: 1	
Analyte	Result	MQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit Qual
1,2-Dichloroethane	17.75	1.0	20	0	88.8	70 - 127	17.84	0.511	20
Benzene	17.84	1.0	20	0	89.2	70 - 127	18.14	1.62	20
Chlorobenzene	17.38	1.0	20	0	86.9	70 - 114	17.35	0.187	20
Ethylbenzene	18.24	1.0	20	0	91.2	70 - 124	18.32	0.447	20
Methylene chloride	16.37	2.0	20	0	81.9	70 - 128	16.47	0.587	20
Toluene	17.8	1.0	20	0	89.0	70 - 123	17.75	0.297	20
Xylenes, Total	54.38	1.0	60	0	90.6	70 - 130	54.81	0.78	20
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>52.39</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>105</i>	<i>70 - 126</i>	<i>52.67</i>	<i>0.533</i>	<i>20</i>
<i>Surr: 4-Bromofluorobenzene</i>	<i>48.55</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.1</i>	<i>81 - 113</i>	<i>48.7</i>	<i>0.314</i>	<i>20</i>
<i>Surr: Dibromofluoromethane</i>	<i>48.89</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>97.8</i>	<i>77 - 123</i>	<i>48.94</i>	<i>0.106</i>	<i>20</i>
<i>Surr: Toluene-d8</i>	<i>50.09</i>	<i>1.0</i>	<i>50</i>	<i>0</i>	<i>100</i>	<i>82 - 127</i>	<i>49.9</i>	<i>0.39</i>	<i>20</i>

The following samples were analyzed in this batch: HS20030819-01 HS20030819-02 HS20030819-03 HS20030819-04  
 HS20030819-05 HS20030819-06 HS20030819-07 HS20030819-08

**Client:** Golder Associates Inc.  
**Project:** Houston TX-Wood Preserving Works  
**WorkOrder:** HS20030819

**QUALIFIERS,  
ACRONYMS, UNITS**

<b>Qualifier</b>	<b>Description</b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
M	Manually integrated, see raw data for justification
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL/SDL

<b>Acronym</b>	<b>Description</b>
DCS	Detectability Check Study
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
SDL	Sample Detection Limit
TRRP	Texas Risk Reduction Program

<b>Unit Reported</b>	<b>Description</b>
mg/L	Milligrams per Liter

**CERTIFICATIONS,ACCREDITATIONS & LICENSES**

<b>Agency</b>	<b>Number</b>	<b>Expire Date</b>
California	2919, 2019-2020	30-Apr-2020
Dept of Defense	ANAB L2231 V009	22-Dec-2021
Florida	E87611-28	30-Jun-2020
Illinois	2000322019-2	09-May-2020
Kansas	E-10352 2019-2020	31-Jul-2020
Kentucky	123043, 2019-2020	30-Apr-2020
Louisiana	03087, 2019-2020	30-Jun-2020
Maryland	343, 2019-2020	30-Jun-2020
North Carolina	624-2020	31-Dec-2020
North Dakota	R-193 2019-2020	30-Apr-2020
Oklahoma	2019-067	31-Aug-2020
Texas	T104704231-19-25	30-Apr-2020



Sample Receipt Checklist

Client Name: PBW
Work Order: HS20030819

Date/Time Received: 18-Mar-2020 16:35
Received by: DDG

Checklist completed by: Jared R. Makan
eSignature
Date: 18-Mar-2020

Reviewed by: Dane J. Wacasey
eSignature
Date: 19-Mar-2020

Matrices: Water

Carrier name: ALS Courier

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on shipping container/cooler? Yes [ ] No [ ] Not Present [checked]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
VOA/TX1005/TX1006 Solids in hermetically sealed vials? Yes [ ] No [ ] Not Present [checked]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Samplers name present on COC? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [ ] No [checked]
All samples received within holding time? Yes [checked] No [ ]
Container/Temp Blank temperature in compliance? Yes [checked] No [ ]

1 Page(s)
COC IDs:221274

Temperature(s)/Thermometer(s): 1.3°C, 1.0°C UC/C IR11
Cooler(s)/Kit(s): 45498, 43911
Date/Time sample(s) sent to storage: 03/18/2020 20:45
Water - VOA vials have zero headspace? Yes [checked] No [ ] No VOA vials submitted [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] N/A [ ]
pH adjusted? Yes [ ] No [checked] N/A [ ]
pH adjusted by:

Login Notes: Limited sample volume for MW 60B - approximately 750ml of sample in each SVOC amber bottle.

Client Contacted: Date Contacted: Person Contacted:
Contacted By: Regarding:
Comments:
Corrective Action:



Cincinnati, OH  
+1 513 733 5336

Everett, WA  
+1 425 356 2600

Fort Collins, CO  
+1 970 490 1511

Holland, MI  
+1 616 399 6070

# Chain of Custody Form

## HS20030819

Golder Associates Inc.  
Houston TX-Wood Preserving Works

Page      of     

COC ID: 221274



Customer Information		Project Information		ALS Project Manager:											
Purchase Order	UPRR/Kevin Peterburs	Project Name	Houston TX-Wood Preserving Works	A	8260 LI_W (5632528 VOC Site Specific)										
Work Order		Project Number	1620-14-Rev0 SR 92688	B	TX1005_W Low (5643233 TPH TX1005)										
Company Name	Golder Associates Inc.	Bill To Company	Union Pacific Railroad- A/P	C	3270 LOW_W (5632533 SemiVolatiles Site specific)										
Send Report To	Eric Matzner	Invoice Attn	Accounts Payable	D	ICP_TW (5636002 Metals - As)										
Address	2201 Double Creek Drive	Address	1400 Douglas Street	E											
	Suite 4004		Stop 0750	F											
City/State/Zip	Round Rock, TX 78864	City/State/Zip	Omaha NE 681790750	G											
Phone	(512) 671-3434	Phone		H											
Fax	(512) 671-3446	Fax		I											
e-Mail Address	Eric_Matzner@golder.com	e-Mail Address		J											

No.	Sample Description	Date	Time	Matrix	Pres.	# Bottles	A	B	C	D	E	F	G	H	I	J	Hold
1	WQ-1620-TB CG-021720 17A	3-17-2020	15:59	Water	1.8	2	X										
2	WQ-1620-MW 88A 20200317	3-17-2020	12:55	Groundwater	1.28	9	X	X	X	X							
3	WG 1620 MW 61B 2020 0317	3-17-2020	11:45	GW	1.28	9	X	X	X	X							
4	WG 1620 MW 96B 2020 0317	3-17-2020	10:25	GW	1.28	9	X	X	X	X							
5	WG 1620 MW 95A 2020 0317	3-17-2020	9:10	GW	1.28	9	X	X	X	X							
6	WG 1620 MW 60B 2020 0317	3-17-2020	15:15	GW	1.28	9	X	X	X	X							
7	WG 1620 MW 88B 2020 0317	3-17-2020	13:50	GW	1.28	9	X	X	X	X							
8	WG 1620 F02 2020 0317	3-17-2020	-	GW	1.28	9	X	X	X	X							
9																	
10																	

Sampler(s) Please Print & Sign <i>TIM McSpadden</i> <i>T. McSpadden</i>		Shipment Method	Required Turnaround Time: (Check Box) <input checked="" type="checkbox"/> STD 18 HR DEL <input type="checkbox"/> 5 HR DEL <input type="checkbox"/> 2 HR DEL <input type="checkbox"/> 24 HR				Results Due Date:			
Relinquished by: <i>T. McSpadden</i>	Date: 3-18-20 Time: 16:00	Received by: <i>D.S.</i>	Notes: UPRR HWPW 1620-14							
Relinquished by: <i>D.S.</i>	Date: 3-18-20 Time: 16:35	Received by (Laboratory): <i>D.S.</i>	Cooler ID	Cooler Temp.	QC Package: (Check One Box Below)					
Logged by (Laboratory):	Date:	Checked by (Laboratory):	45498	1.3	<input type="checkbox"/> Level II (CFC)	<input checked="" type="checkbox"/> ITRP Checker				
			43911	1.0	<input type="checkbox"/> Level III (CFC/Procedure)	<input type="checkbox"/> ITRP Level II				
					<input type="checkbox"/> Level IV (CFC/PCDA)	<input type="checkbox"/> ITRP Level I				
Preservative Key: 1-HCl 2-HNO <sub>3</sub> 3-H <sub>2</sub> SO <sub>4</sub> 4-NaOH 5-Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> 6-NaHSO <sub>4</sub> 7-Other 8-4°C 9-5035										

Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.  
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.  
 3. The Chain of Custody is a legal document. All information must be completed accurately.

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**ATTACHMENT C**

# Groundwater Sampling Records



**GROUNDWATER SAMPLING RECORD**

Project Number: **19119232** Project Name: **UPRR - HWPW - SMU** Date: **7-2-19**  
 Sampling Location (well ID, etc.): **MW 11B** Starting Water Level (ft. BMP): **5.21**  
 Sample Number: **WG-1620-MW11B-20190702** Casing Stickup (ft.):  
 Sampled by: **JTB** WL (ft. BMP): **5.21** (ft. BGL):  
 Measuring Point (MP) of Well: **TOC** - Steel or **AVC** TD (ft. BMP): (ft. BGL):  
 Screened Interval (ft. BGL): - Ft. water: Casing Dia. (In ID):  
 Filter Pack Interval (ft. BGL): - 1X Casing Vol (gal.): 3X (gal):

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: **Peristaltic Pump** / SS Pump / Bailer Sampling: **Peristaltic Pump** / SS Pump / Bailer  
 Disposal of Discharged Water: **55-gallon drum**


INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: **SOUNST** Other: -  
 Multi Meter: **HORIBA**  
 Field Calibration: **AUTO CAL**  
 Filter / Filter Size: -

<b>SAMPLING MEASUREMENTS</b>										
Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L/m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
0801										
0812	5.39		2	22.4	0.79	1120	6.92	-42	16	milky
0817	5.43		↓	22.1	0.62	1150	6.97	-37	12	↓
0822	5.41		↓	22.1	0.61	1160	6.96	-39	12	↓

WL (ft. BMP) at End of Purge: Sample Intake Depth (ft. BMP):

<b>SAMPLE INVENTORY</b>					Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
Time	Bottles Collected		No.				
	Volume	Composition (G, P)					
0835	16	G	2	N	-	SVOCs	

Comments:


**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446





**GROUNDWATER SAMPLING RECORD**

Project Number: **19119232** Project Name: **VPRR-HWPW-SMDU** Date: **7-1-19**  
 Sampling Location (well ID, etc.): **MW07** Starting Water Level (ft. BMP): **3.93**  
 Sample Number: **WG-1620-MW07-20190701** Casing Stickup (ft.):  
 Sampled by: **JTB** WL (ft. BMP): **3.93** (ft. BGL):  
 Measuring Point (MP) of Well: **TOC** Steel or **PVC** TD (ft. BMP): (ft. BGL):  
 Screened Interval (ft. BGL): Ft. water: Casing Dia. (In ID):  
 Filter Pack Interval (ft. BGL): 1X Casing Vol (gal.): 3X (gal):

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: **Peristaltic Pump** / SS Pump / Bailer Sampling: **Peristaltic Pump** / SS Pump / Bailer  
 Disposal of Discharged Water: **55-gallon drum**

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: **SOLINST** Other:   
 Multi Meter: **HORIBA**  
 Field Calibration: **AUTO CAL**  
 Filter / Filter Size:

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol (gal. or L)	Purge Rate (gal. or L/m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
1623										
1632	4.26		2	22.7	0.34	1340	6.86	-41	16	neutral
1638	4.29		↓	22.3	0.26	1310	6.92	-35	11	↓
1643	4.22		↓	22.4	0.26	1310	6.91	-36	12	↓

WL (ft. BMP) at End of Purge: Sample Intake Depth (ft. BMP):

**SAMPLE INVENTORY**

Bottles Collected					Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
Time	Volume	Composition (G, P)	No.				
1655	1L	G	2	N	-	SVOCs	

Comments:  
**WG-1620-FD02-20190701**  
**1655-Sample Time**

**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446









**GROUNDWATER SAMPLING RECORD**

PAGE 1 of 1

Project Number: 1911 9232 Project Name: UPRR - HWPW - SMU Date: 7-2-19  
 Sampling Location (well ID, etc.): MW 01A Starting Water Level (ft. BMP): 2.85  
 Sample Number: WG-1620-MW01A-20190702 Casing Stickup (ft.): \_\_\_\_\_  
 Sampled by: YTB WL (ft. BMP): 2.85 (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: (TOC) - Steel or (PVC) TD (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): - Ft. water: \_\_\_\_\_ Casing Dia. (In ID): \_\_\_\_\_  
 Filter Pack Interval (ft. BGL): - 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer Sampling: Peristaltic Pump / SS Pump / Bailer  
 Disposal of Discharged Water: 55-gallon drum

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: SAMST Other: -  
 Multi Meter: HORIBA  
 Field Calibration: AUTO CAL  
 Filter / Filter Size: -

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L/m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
1132										
1143	3.07	2	2	21.7	0.59	1150	6.89	-26	7.1	rel. turbid
1148	3.08		↓	21.7	0.39	1140	6.81	-25	5.2	↓
1153	3.08		↓	21.7	0.42	1140	6.82	-27	6.1	↓

WL (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): \_\_\_\_\_

**SAMPLE INVENTORY**

Time	Bottles Collected				Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.				
1210	1L	G	2		N	-	SVOCs

Comments:  
WG-1620-FDOT-20190702  
1210 - Sample Time

**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446







**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.21</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-9-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.21</u>	ft. BGL
MW ID	<u>MW56A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW56A-20190709</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1009				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1019	<u>0.2</u>	<u>/</u>	<u>4.42</u>	<u>21.9</u>	<u>7.21</u>	<u>1140</u>	<u>0.82</u>	<u>-62</u>	<u>5.1</u>
1024	<u>↓</u>	<u>/</u>	<u>4.43</u>	<u>21.8</u>	<u>7.17</u>	<u>1170</u>	<u>0.64</u>	<u>-54</u>	<u>6.7</u>
1029	<u>↓</u>	<u>/</u>	<u>4.42</u>	<u>21.8</u>	<u>7.16</u>	<u>1160</u>	<u>0.67</u>	<u>-56</u>	<u>6.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1040	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1040	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.91</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-19-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>7.91</u>	ft. BGL
MW ID	<u>MW81B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW81B - 20190708</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0916</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0926</u>	<u>.2</u>	<u>/</u>	<u>8.12</u>	<u>22.6</u>	<u>7.31</u>	<u>1040</u>	<u>0.91</u>	<u>-79</u>	<u>5.2</u>
<u>0931</u>	<u>↓</u>	<u>/</u>	<u>8.13</u>	<u>22.3</u>	<u>7.25</u>	<u>1060</u>	<u>0.72</u>	<u>-76</u>	<u>4.1</u>
<u>0936</u>	<u>↓</u>	<u>/</u>	<u>8.12</u>	<u>22.4</u>	<u>7.26</u>	<u>1060</u>	<u>0.71</u>	<u>-77</u>	<u>4.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0945</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0945</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>20.39</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>07-19-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>20.39</u> ft. BGL
MW ID	<u>MWSIC</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MWSIC-20190709</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0822</u>	(L/min)	(L)	(ft)	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0832</u>	<u>.2</u>	<u>/</u>	<u>20.56</u>	<u>21.9</u>	<u>7.13</u>	<u>1430</u>	<u>0.67</u>	<u>-76</u>	<u>4.7</u>
<u>0837</u>	<u>↓</u>	<u>/</u>	<u>20.58</u>	<u>21.6</u>	<u>7.10</u>	<u>1420</u>	<u>0.51</u>	<u>-74</u>	<u>7.9</u>
<u>0842</u>	<u>↓</u>	<u>/</u>	<u>20.58</u>	<u>21.7</u>	<u>7.11</u>	<u>1410</u>	<u>0.52</u>	<u>-77</u>	<u>7.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0855</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0855</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.29</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-9-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>6.29</u> ft. BGL
MW ID	<u>MW51A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW51A-20190709</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0652				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0702	<u>.2</u>	<u>/</u>	<u>6.41</u>	<u>21.2</u>	<u>7.21</u>	<u>1760</u>	<u>1.31</u>	<u>-91</u>	<u>9.2</u>
0707	<u>↓</u>	<u>/</u>	<u>6.42</u>	<u>21.2</u>	<u>7.25</u>	<u>1710</u>	<u>1.11</u>	<u>-85</u>	<u>8.6</u>
0712	<u>↓</u>	<u>/</u>	<u>6.41</u>	<u>21.1</u>	<u>7.26</u>	<u>1720</u>	<u>1.07</u>	<u>-86</u>	<u>8.9</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0725</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0725</u>	<u>40ML / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.23</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.23</u> ft. BGL
MW ID	<u>MWISB</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MWISB-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1729</u>	(L/min)	(L)	(ft)	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1739</u>	<u>.2</u>	<u>/</u>	<u>9.42</u>	<u>22.4</u>	<u>6.74</u>	<u>960</u>	<u>0.76</u>	<u>-46</u>	<u>8.2</u>
<u>1744</u>	<u>↓</u>	<u>/</u>	<u>9.44</u>	<u>22.2</u>	<u>6.80</u>	<u>910</u>	<u>0.71</u>	<u>-40</u>	<u>6.5</u>
<u>1749</u>	<u>↓</u>	<u>/</u>	<u>9.43</u>	<u>22.3</u>	<u>6.81</u>	<u>920</u>	<u>0.72</u>	<u>-41</u>	<u>6.8</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1800</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1800</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FB01-20190710  
1815 - SAMPLE TIME

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>22.38</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>22.38</u> ft. BGL
MW ID	<u>MWISC</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1632				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1642	<u>.2</u>	<u>/</u>	<u>22.61</u>	<u>22.6</u>	<u>6.91</u>	<u>820</u>	<u>1.06</u>	<u>-71</u>	<u>6.2</u>
1647	<u>↓</u>	<u>/</u>	<u>22.63</u>	<u>22.3</u>	<u>6.93</u>	<u>870</u>	<u>0.80</u>	<u>-61</u>	<u>4.7</u>
1652	<u>↓</u>	<u>/</u>	<u>22.64</u>	<u>22.4</u>	<u>6.93</u>	<u>860</u>	<u>0.81</u>	<u>-62</u>	<u>5.2</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1705</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1705</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.03</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>9.03</u> ft. BGL
MW ID	<u>MW15A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW15A-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1536</u>	<u>1.2</u>	<u>/</u>	<u>9.42</u>	<u>22.7</u>	<u>6.91</u>	<u>1190</u>	<u>0.52</u>	<u>-71</u>	<u>10</u>
<u>1551</u>	<u>↓</u>	<u>/</u>	<u>9.44</u>	<u>22.3</u>	<u>6.81</u>	<u>1150</u>	<u>0.49</u>	<u>-62</u>	<u>8.2</u>
<u>1556</u>	<u>↓</u>	<u>/</u>	<u>9.45</u>	<u>22.4</u>	<u>6.82</u>	<u>1160</u>	<u>0.51</u>	<u>-67</u>	<u>9.3</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1610</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1610</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.04</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>9.04</u> ft. BGL
MW ID	<u>MW20A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW20A-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1439				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1449	<u>2</u>	<u>/</u>	<u>9.41</u>	<u>22.4</u>	<u>6.61</u>	<u>1690</u>	<u>0.67</u>	<u>-101</u>	<u>3.7</u>
1454	<u>↓</u>	<u>/</u>	<u>9.42</u>	<u>22.5</u>	<u>6.62</u>	<u>1660</u>	<u>0.54</u>	<u>-101</u>	<u>3.2</u>
1459	<u>↓</u>	<u>/</u>	<u>9.42</u>	<u>22.1</u>	<u>6.62</u>	<u>1670</u>	<u>0.53</u>	<u>-92</u>	<u>3.4</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1510</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1510</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>22.14</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>22.14</u>	ft. BGL
MW ID	<u>MW17C</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW17C-20190710</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1251</u>	<u>1.2</u>	<u>/</u>	<u>22.39</u>	<u>21.3</u>	<u>6.71</u>	<u>1170</u>	<u>0.79</u>	<u>-29</u>	<u>4.1</u>
<u>1304</u>	<u>↓</u>	<u>/</u>	<u>22.46</u>	<u>22.9</u>	<u>6.74</u>	<u>1160</u>	<u>0.74</u>	<u>-32</u>	<u>3.1</u>
<u>1311</u>	<u>↓</u>	<u>/</u>	<u>22.47</u>	<u>22.8</u>	<u>6.77</u>	<u>1160</u>	<u>0.71</u>	<u>-33</u>	<u>2.9</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>22.14</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>22.14</u>	ft. BGL
MW ID	<u>MW17</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW17-20190710</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1341				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1351	<u>0.2</u>	<u>/</u>	<u>22.41</u>	<u>22.6</u>	<u>6.90</u>	<u>1320</u>	<u>0.91</u>	<u>-61</u>	<u>9.1</u>
1356	<u>↓</u>	<u>/</u>	<u>22.43</u>	<u>22.5</u>	<u>6.93</u>	<u>1340</u>	<u>0.71</u>	<u>-52</u>	<u>8.4</u>
1401	<u>↓</u>	<u>/</u>	<u>22.43</u>	<u>22.5</u>	<u>6.93</u>	<u>1340</u>	<u>0.72</u>	<u>-51</u>	<u>8.6</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1415</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1415</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>24.58</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>24.58</u> ft. BGL
MW ID	<u>MW19C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW19C-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1154				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1204	<u>0.2</u>	<u>/</u>	<u>24.81</u>	<u>21.6</u>	<u>6.70</u>	<u>1260</u>	<u>0.79</u>	<u>-102</u>	<u>7.2</u>
1209	<u>↓</u>	<u>/</u>	<u>24.82</u>	<u>21.8</u>	<u>6.71</u>	<u>1240</u>	<u>0.72</u>	<u>-101</u>	<u>6.2</u>
1214	<u>↓</u>	<u>/</u>	<u>24.81</u>	<u>21.9</u>	<u>6.71</u>	<u>1230</u>	<u>0.74</u>	<u>-102</u>	<u>6.7</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1225</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1225</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.28</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>15.28</u> ft. BGL
MW ID	<u>MW72B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW72B-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1103				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1113	<u>.2</u>	<u>/</u>	<u>15.51</u>	<u>25.6</u>	<u>6.79</u>	<u>1730</u>	<u>0.86</u>	<u>-103</u>	<u>12</u>
1118	<u>↓</u>	<u>/</u>	<u>15.56</u>	<u>25.2</u>	<u>6.80</u>	<u>1710</u>	<u>0.86</u>	<u>-103</u>	<u>11</u>
1123	<u>↓</u>	<u>/</u>	<u>15.57</u>	<u>25.1</u>	<u>6.81</u>	<u>1710</u>	<u>0.79</u>	<u>-101</u>	<u>11</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1135</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1135</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>13.71</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>13.71</u> ft. BGL
MW ID	<u>MW57B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW57B-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1009</u>	<u>1.2</u>	<u>/</u>	<u>14.03</u>	<u>21.6</u>	<u>6.69</u>	<u>1270</u>	<u>1.42</u>	<u>-31</u>	<u>14</u>
<u>1024</u>	<u>↓</u>	<u>/</u>	<u>14.05</u>	<u>21.7</u>	<u>6.65</u>	<u>1270</u>	<u>1.02</u>	<u>-25</u>	<u>11</u>
<u>1029</u>	<u>↓</u>	<u>/</u>	<u>14.05</u>	<u>21.7</u>	<u>6.67</u>	<u>1260</u>	<u>1.01</u>	<u>-26</u>	<u>11</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1040</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1040</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>13.11</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>13.11</u>	ft. BGL
MW ID	<u>MW57A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW57A-20190710</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0924				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0934	<u>0.2</u>	<u>/</u>	<u>13.41</u>	<u>23.1</u>	<u>6.71</u>	<u>1470</u>	<u>0.79</u>	<u>-91</u>	<u>9.1</u>
0939	<u>↓</u>	<u>/</u>	<u>13.42</u>	<u>23.2</u>	<u>6.56</u>	<u>1430</u>	<u>0.59</u>	<u>-91</u>	<u>7.1</u>
0944	<u>↓</u>	<u>/</u>	<u>13.42</u>	<u>23.2</u>	<u>6.51</u>	<u>1430</u>	<u>0.57</u>	<u>-90</u>	<u>8.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0955	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
0955	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>12.29</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>12.29</u> ft. BGL
MW ID	<u>MWS2A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0821</u>	<u>1.2</u>	<u>/</u>	<u>12.61</u>	<u>22.9</u>	<u>6.64</u>	<u>1460</u>	<u>1.41</u>	<u>-77</u>	<u>3.7</u>
<u>0836</u>	<u>↓</u>	<u>/</u>	<u>12.63</u>	<u>22.9</u>	<u>6.71</u>	<u>1420</u>	<u>1.11</u>	<u>-72</u>	<u>3.2</u>
<u>0841</u>	<u>↓</u>	<u>/</u>	<u>12.63</u>	<u>22.8</u>	<u>6.81</u>	<u>1440</u>	<u>1.09</u>	<u>-71</u>	<u>4.6</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0855</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0855</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>23.08</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>23.08</u> ft. BGL
MW ID	<u>MW18C-</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW18C-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0734				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0744	<u>2</u>	<u>/</u>	<u>23.39</u>	<u>22.7</u>	<u>6.71</u>	<u>1340</u>	<u>0.79</u>	<u>-101</u>	<u>7.3</u>
0749	<u>↓</u>	<u>/</u>	<u>23.41</u>	<u>22.4</u>	<u>6.64</u>	<u>1320</u>	<u>0.65</u>	<u>-93</u>	<u>6.7</u>
0754	<u>↓</u>	<u>/</u>	<u>23.41</u>	<u>22.4</u>	<u>6.67</u>	<u>1310</u>	<u>0.67</u>	<u>-93</u>	<u>7.4</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0805</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0805</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>17.51</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-10-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>17.51</u> ft. BGL
MW ID	<u>MW18A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW18A-20190719</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0642</u>	<u>1.2</u>	<u>/</u>	<u>17.89</u>	<u>22.4</u>	<u>6.79</u>	<u>1570</u>	<u>0.90</u>	<u>-41</u>	<u>5.2</u>
<u>0657</u>	<u>↓</u>	<u>/</u>	<u>17.92</u>	<u>22.1</u>	<u>6.73</u>	<u>1550</u>	<u>0.71</u>	<u>-35</u>	<u>6.1</u>
<u>0702</u>	<u>↓</u>	<u>/</u>	<u>17.92</u>	<u>22.1</u>	<u>6.74</u>	<u>1540</u>	<u>0.71</u>	<u>-36</u>	<u>6.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G/P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0720</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.09</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>5.09</u> ft. BGL
MW ID	<u>MW64A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW64A-20190711</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1521</u>	<u>1.2</u>	<u>/</u>	<u>5.26</u>	<u>21.6</u>	<u>6.91</u>	<u>1540</u>	<u>0.61</u>	<u>-62</u>	<u>5.6</u>
<u>1536</u>	<u>↓</u>	<u>/</u>	<u>5.29</u>	<u>21.7</u>	<u>6.93</u>	<u>1570</u>	<u>0.51</u>	<u>-50</u>	<u>4.1</u>
<u>1541</u>	<u>↓</u>	<u>/</u>	<u>5.29</u>	<u>21.7</u>	<u>6.93</u>	<u>1560</u>	<u>0.52</u>	<u>-51</u>	<u>4.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1550</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1550</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.56</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.56</u>	ft. BGL
MW ID	<u>MW05</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW05-20190711</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1417</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1427</u>	<u>.2</u>	<u>/</u>	<u>6.77</u>	<u>22.6</u>	<u>6.91</u>	<u>580</u>	<u>0.69</u>	<u>-39</u>	<u>5.6</u>
<u>1432</u>	<u>↓</u>	<u>/</u>	<u>6.78</u>	<u>22.4</u>	<u>6.95</u>	<u>590</u>	<u>0.42</u>	<u>-33</u>	<u>2.7</u>
<u>1437</u>	<u>↓</u>	<u>/</u>	<u>6.78</u>	<u>22.3</u>	<u>6.96</u>	<u>590</u>	<u>0.41</u>	<u>-34</u>	<u>3.4</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1450</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1450</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.31</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.31</u>	ft. BGL
MW ID	<u>P11</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-P11 -20190711</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1323</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1333</u>	<u>1.2</u>	<u>/</u>	<u>5.49</u>	<u>22.7</u>	<u>6.97</u>	<u>1160</u>	<u>0.71</u>	<u>-134</u>	<u>7.9</u>
<u>1338</u>	<u>↓</u>	<u>/</u>	<u>5.51</u>	<u>22.5</u>	<u>6.96</u>	<u>1140</u>	<u>0.76</u>	<u>-131</u>	<u>6.7</u>
<u>1343</u>	<u>↓</u>	<u>/</u>	<u>5.51</u>	<u>22.6</u>	<u>6.96</u>	<u>1140</u>	<u>0.77</u>	<u>-132</u>	<u>6.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1355</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1355</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.42</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.42</u> ft. BGL
MW ID	<u>MW42B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW42B-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1228</u>	<u>1.2</u>	<u>/</u>	<u>7.61</u>	<u>23.2</u>	<u>6.79</u>	<u>1610</u>	<u>0.89</u>	<u>-21</u>	<u>8.2</u>
<u>1243</u>	<u>↓</u>	<u>/</u>	<u>7.62</u>	<u>23.1</u>	<u>6.76</u>	<u>1550</u>	<u>0.78</u>	<u>-16</u>	<u>9.1</u>
<u>1248</u>	<u>↓</u>	<u>/</u>	<u>7.60</u>	<u>23.1</u>	<u>6.77</u>	<u>1540</u>	<u>0.79</u>	<u>-17</u>	<u>9.2</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1300</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1300</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.28</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>7.28</u>	ft. BGL
MW ID	<u>MW40B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620 MW40B-20190711</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1126</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1136</u>	<u>.2</u>	<u>/</u>	<u>7.42</u>	<u>22.5</u>	<u>6.96</u>	<u>1650</u>	<u>0.81</u>	<u>-79</u>	<u>8.6</u>
<u>1141</u>	<u>↓</u>	<u>/</u>	<u>7.44</u>	<u>22.3</u>	<u>6.98</u>	<u>1630</u>	<u>0.70</u>	<u>-77</u>	<u>7.1</u>
<u>1146</u>	<u>↓</u>	<u>/</u>	<u>7.44</u>	<u>22.4</u>	<u>6.98</u>	<u>1640</u>	<u>0.71</u>	<u>-77</u>	<u>7.1</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1200</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1200</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.02</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.02</u>	ft. BGL
MW ID	<u>MW12A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW12A-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1021				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1031	<u>1.2</u>	<u>/</u>	<u>8.22</u>	<u>22.6</u>	<u>6.79</u>	<u>490</u>	<u>1.06</u>	<u>-31</u>	<u>8.1</u>
1036	<u>↓</u>	<u>/</u>	<u>8.21</u>	<u>22.6</u>	<u>6.77</u>	<u>460</u>	<u>0.82</u>	<u>-26</u>	<u>6.1</u>
1041	<u>↓</u>	<u>/</u>	<u>8.21</u>	<u>22.7</u>	<u>6.76</u>	<u>470</u>	<u>0.81</u>	<u>-27</u>	<u>6.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1055</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1055</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>23.73</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>23.73</u>	ft. BGL
MW ID	<u>MW12C</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW12C-20190711</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0929	(L/min)	(L)	(ft)	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0939	<u>.2</u>	<u>/</u>	<u>23.91</u>	<u>21.3</u>	<u>6.91</u>	<u>810</u>	<u>0.91</u>	<u>-39</u>	<u>7.4</u>
0944	<u>↓</u>	<u>/</u>	<u>23.92</u>	<u>21.1</u>	<u>6.86</u>	<u>820</u>	<u>0.70</u>	<u>-60</u>	<u>8.2</u>
0949	<u>↓</u>	<u>/</u>	<u>23.91</u>	<u>21.2</u>	<u>6.87</u>	<u>830</u>	<u>0.71</u>	<u>-61</u>	<u>8.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1000</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1000</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.47</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.47</u>	ft. BGL
MW ID	<u>MW39B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620 MW39B - 20190711</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0843				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0852	<u>1.2</u>	<u>/</u>	<u>7.68</u>	<u>22.4</u>	<u>6.91</u>	<u>1270</u>	<u>0.51</u>	<u>-91</u>	<u>5.2</u>
0858	<u>↓</u>	<u>/</u>	<u>7.69</u>	<u>22.4</u>	<u>6.85</u>	<u>1280</u>	<u>0.54</u>	<u>-86</u>	<u>4.1</u>
0903	<u>↓</u>	<u>/</u>	<u>7.68</u>	<u>22.3</u>	<u>6.86</u>	<u>1280</u>	<u>0.56</u>	<u>-87</u>	<u>4.7</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0915</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0915</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>10.63</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>10.63</u> ft. BGL
MW ID	<u>MW13</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW13-20190711</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0742</u>	<u>2</u>	<u>/</u>	<u>10.81</u>	<u>21.6</u>	<u>6.93</u>	<u>750</u>	<u>0.59</u>	<u>-52</u>	<u>13</u>
<u>0757</u>	<u>↓</u>	<u>/</u>	<u>10.84</u>	<u>21.9</u>	<u>6.94</u>	<u>730</u>	<u>0.54</u>	<u>-50</u>	<u>11</u>
<u>0802</u>	<u>↓</u>	<u>/</u>	<u>10.83</u>	<u>21.9</u>	<u>6.94</u>	<u>740</u>	<u>0.56</u>	<u>-51</u>	<u>11</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0815</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0815</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.57</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-11-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.57</u> ft. BGL
MW ID	<u>MW14</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW14-20190710</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0647</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0657</u>	<u>1.2</u>	<u>/</u>	<u>9.73</u>	<u>22.4</u>	<u>6.71</u>	<u>1090</u>	<u>1.31</u>	<u>-76</u>	<u>16</u>
<u>0702</u>	<u>↓</u>	<u>/</u>	<u>9.74</u>	<u>22.4</u>	<u>6.64</u>	<u>1070</u>	<u>1.01</u>	<u>-70</u>	<u>11</u>
<u>0707</u>	<u>↓</u>	<u>/</u>	<u>9.73</u>	<u>22.5</u>	<u>6.65</u>	<u>1060</u>	<u>1.02</u>	<u>-71</u>	<u>12</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0720</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0720</u>	<u>40mL/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVPCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.22</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-12-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.22</u>	ft. BGL
MW ID	<u>TW41B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-TW41B-20190712</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1009				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1019	<u>2</u>	<u>/</u>	<u>8.43</u>	<u>22.6</u>	<u>6.67</u>	<u>1410</u>	<u>1.13</u>	<u>-106</u>	<u>8.6</u>
1024	<u>↓</u>	<u>/</u>	<u>8.44</u>	<u>22.3</u>	<u>6.56</u>	<u>1430</u>	<u>0.91</u>	<u>-109</u>	<u>7.7</u>
1029	<u>↓</u>	<u>/</u>	<u>8.44</u>	<u>22.3</u>	<u>6.57</u>	<u>1440</u>	<u>0.92</u>	<u>-109</u>	<u>7.9</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1040	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1040	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.02</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-12-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.02</u> ft. BGL
MW ID	<u>MW04</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW04-20190712</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0856				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0906	<u>2</u>	<u>/</u>	<u>6.31</u>	<u>22.3</u>	<u>6.87</u>	<u>720</u>	<u>0.71</u>	<u>-77</u>	<u>6.2</u>
0911	<u>↓</u>	<u>/</u>	<u>6.32</u>	<u>22.4</u>	<u>6.84</u>	<u>730</u>	<u>0.52</u>	<u>-73</u>	<u>5.1</u>
0916	<u>↓</u>	<u>/</u>	<u>6.31</u>	<u>22.2</u>	<u>6.81</u>	<u>740</u>	<u>0.51</u>	<u>-74</u>	<u>5.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0925	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0925	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.57</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-12-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>5.57</u> ft. BGL
MW ID	<u>MW03</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW03-20190712</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0802				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0812	<u>0.2</u>	<u>/</u>	<u>5.81</u>	<u>21.7</u>	<u>7.12</u>	<u>560</u>	<u>0.74</u>	<u>-34</u>	<u>7.7</u>
0817	<u>↓</u>	<u>/</u>	<u>5.86</u>	<u>21.8</u>	<u>7.06</u>	<u>570</u>	<u>0.62</u>	<u>-31</u>	<u>7.2</u>
0822	<u>↓</u>	<u>/</u>	<u>5.87</u>	<u>21.8</u>	<u>7.07</u>	<u>570</u>	<u>0.62</u>	<u>-31</u>	<u>6.6</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0835	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
0835	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.21</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-12-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.21</u>	ft. BGL
MW ID	<u>MW09</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW09 -20190712</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0704</u>	<u>(L/min)</u>	<u>(L)</u>	<u>(ft)</u>	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0714</u>	<u>.2</u>	<u>/</u>	<u>6.46</u>	<u>22.9</u>	<u>6.77</u>	<u>660</u>	<u>1.39</u>	<u>-21</u>	<u>5.2</u>
<u>0719</u>	<u>↓</u>	<u>/</u>	<u>6.48</u>	<u>23.0</u>	<u>6.75</u>	<u>630</u>	<u>1.02</u>	<u>-19</u>	<u>4.2</u>
<u>0724</u>	<u>↓</u>	<u>/</u>	<u>6.49</u>	<u>23.1</u>	<u>6.74</u>	<u>640</u>	<u>1.03</u>	<u>-19</u>	<u>4.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0735</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0735</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>3.96</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>3.96</u> ft. BGL
MW ID	<u>MW63B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW63B-20190716</u>		Water Quality	MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1723</u>	<u>2</u>	<u>/</u>	<u>4.29</u>	<u>22.4</u>	<u>6.81</u>	<u>1640</u>	<u>1.31</u>	<u>-46</u>	<u>9.3</u>
<u>1738</u>	<u>↓</u>	<u>/</u>	<u>4.31</u>	<u>22.6</u>	<u>6.79</u>	<u>1630</u>	<u>1.01</u>	<u>-43</u>	<u>7.4</u>
<u>1743</u>	<u>↓</u>	<u>/</u>	<u>4.31</u>	<u>22.6</u>	<u>6.79</u>	<u>1630</u>	<u>1.02</u>	<u>-49</u>	<u>7.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1755</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1755</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.28</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.52</u> ft. BGL
MW ID	<u>MW28A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW28A-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1629				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1639	<u>2</u>	<u>/</u>	<u>6.81</u>	<u>21.9</u>	<u>6.67</u>	<u>1340</u>	<u>0.91</u>	<u>-106</u>	<u>7.4</u>
1644	<u>↓</u>	<u>/</u>	<u>6.82</u>	<u>21.8</u>	<u>6.61</u>	<u>1320</u>	<u>0.76</u>	<u>-101</u>	<u>5.1</u>
1649	<u>↓</u>	<u>/</u>	<u>6.82</u>	<u>21.8</u>	<u>6.62</u>	<u>1310</u>	<u>0.77</u>	<u>-102</u>	<u>5.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1700</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1700</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.03</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>15.03</u> ft. BGL
MW ID	<u>MW28C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW28C-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1543</u>	(L/min)	(L)	(ft)	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1553</u>	<u>.2</u>	<u>/</u>	<u>15.29</u>	<u>22.3</u>	<u>6.92</u>	<u>1460</u>	<u>0.79</u>	<u>-101</u>	<u>3.2</u>
<u>1558</u>	<u>↓</u>	<u>/</u>	<u>15.32</u>	<u>22.1</u>	<u>6.88</u>	<u>1420</u>	<u>0.60</u>	<u>-96</u>	<u>4.9</u>
<u>1603</u>	<u>↓</u>	<u>/</u>	<u>15.32</u>	<u>22.1</u>	<u>6.89</u>	<u>1430</u>	<u>0.62</u>	<u>-93</u>	<u>5.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1615</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1615</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

MS/MSD

Field Team Leader JOHN BRAYTON name

John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.32</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>8.32</u> ft. BGL
MW ID	<u>MW36A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW36A-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1448</u>	<u>1.2</u>	<u>/</u>	<u>8.72</u>	<u>22.6</u>	<u>6.71</u>	<u>1410</u>	<u>0.79</u>	<u>-17</u>	<u>5.9</u>
<u>1458</u>	<u>↓</u>	<u>/</u>	<u>8.74</u>	<u>22.4</u>	<u>6.72</u>	<u>1340</u>	<u>0.61</u>	<u>-13</u>	<u>4.1</u>
<u>1508</u>	<u>↓</u>	<u>/</u>	<u>8.74</u>	<u>22.4</u>	<u>6.74</u>	<u>1360</u>	<u>0.62</u>	<u>-14</u>	<u>4.6</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1520</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1520</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>1.86</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>1.86</u> ft. BGL
MW ID	<u>MW36B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW36B-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1338				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1348	<u>.2</u>	<u>/</u>	<u>2.37</u>	<u>21.7</u>	<u>6.71</u>	<u>1160</u>	<u>0.67</u>	<u>-46</u>	<u>27</u>
1353	<u>↓</u>	<u>/</u>	<u>2.39</u>	<u>21.8</u>	<u>6.61</u>	<u>1190</u>	<u>0.53</u>	<u>-42</u>	<u>20</u>
1358	<u>↓</u>	<u>/</u>	<u>2.40</u>	<u>21.8</u>	<u>6.63</u>	<u>1190</u>	<u>0.56</u>	<u>-41</u>	<u>21</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1420</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1420</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.77</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>6.77</u> ft. BGL
MW ID	<u>MW25A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW25A-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HDR15A</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1237				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1247	<u>.2</u>	<u>/</u>	<u>7.03</u>	<u>22.3</u>	<u>6.94</u>	<u>1520</u>	<u>0.57</u>	<u>-36</u>	<u>17</u>
1252	<u>↓</u>	<u>/</u>	<u>7.06</u>	<u>22.1</u>	<u>6.92</u>	<u>1470</u>	<u>0.48</u>	<u>-29</u>	<u>12</u>
1257	<u>↓</u>	<u>/</u>	<u>7.07</u>	<u>22.1</u>	<u>6.91</u>	<u>1460</u>	<u>0.49</u>	<u>-31</u>	<u>12</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1310</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1310</u>	<u>40mL/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.84</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>15.84</u> ft. BGL
MW ID	<u>MW25C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW25C-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1139				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1149	<u>.2</u>	<u>/</u>	<u>16.09</u>	<u>23.6</u>	<u>6.81</u>	<u>1160</u>	<u>0.71</u>	<u>-47</u>	<u>6.2</u>
1154	<u>↓</u>	<u>/</u>	<u>16.11</u>	<u>23.7</u>	<u>6.76</u>	<u>1190</u>	<u>0.56</u>	<u>-42</u>	<u>5.1</u>
1159	<u>↓</u>	<u>/</u>	<u>16.12</u>	<u>23.7</u>	<u>6.77</u>	<u>1190</u>	<u>0.58</u>	<u>-43</u>	<u>5.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1215</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1215</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.27</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>15.27</u> ft. BGL
MW ID	<u>MW54C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW54C-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1043				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1053	<u>1.2</u>	<u>/</u>	<u>15.61</u>	<u>21.6</u>	<u>6.79</u>	<u>1610</u>	<u>1.21</u>	<u>-81</u>	<u>4.6</u>
1058	<u>↓</u>	<u>/</u>	<u>15.62</u>	<u>21.6</u>	<u>6.72</u>	<u>1540</u>	<u>0.86</u>	<u>-77</u>	<u>2.7</u>
1103	<u>↓</u>	<u>/</u>	<u>15.62</u>	<u>21.9</u>	<u>6.71</u>	<u>1510</u>	<u>0.89</u>	<u>-71</u>	<u>2.6</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1115</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1115</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.86</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>15.86</u> ft. BGL
MW ID	<u>MW53C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW53C-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
0946									
0956	<u>1.2</u>	<u>/</u>	<u>16.11</u>	<u>22.6</u>	<u>6.91</u>	<u>1560</u>	<u>0.89</u>	<u>-22</u>	<u>9.2</u>
1001	<u>↓</u>	<u>/</u>	<u>16.16</u>	<u>22.3</u>	<u>6.84</u>	<u>1530</u>	<u>0.76</u>	<u>-21</u>	<u>7.2</u>
1006	<u>↓</u>	<u>/</u>	<u>16.18</u>	<u>22.4</u>	<u>6.86</u>	<u>1540</u>	<u>0.71</u>	<u>-21</u>	<u>7.6</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1015</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1015</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.34</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.34</u> ft. BGL
MW ID	<u>MW62B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW62B-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0757				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0807	<u>1.2</u>	<u>/</u>	<u>4.71</u>	<u>21.6</u>	<u>6.71</u>	<u>1120</u>	<u>1.13</u>	<u>-26</u>	<u>7.4</u>
0812	<u>↓</u>	<u>/</u>	<u>4.73</u>	<u>21.3</u>	<u>6.73</u>	<u>1170</u>	<u>0.92</u>	<u>-30</u>	<u>6.0</u>
0817	<u>↓</u>	<u>/</u>	<u>4.74</u>	<u>21.4</u>	<u>6.74</u>	<u>1160</u>	<u>0.91</u>	<u>-31</u>	<u>6.1</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0825</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0825</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader

JOHN BRAYTON  
name

John Brayton  
signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>23.11</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-16-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>23.11</u> ft. BGL
MW ID	<u>MW21C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW21C-20190716</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0656				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0706	<u>2</u>	<u>/</u>	<u>23.34</u>	<u>21.6</u>	<u>7.07</u>	<u>810</u>	<u>0.92</u>	<u>-51</u>	<u>8.6</u>
0711	<u>↓</u>	<u>/</u>	<u>23.39</u>	<u>21.7</u>	<u>7.01</u>	<u>820</u>	<u>0.71</u>	<u>-44</u>	<u>7.0</u>
0716	<u>↓</u>	<u>/</u>	<u>23.39</u>	<u>21.7</u>	<u>7.02</u>	<u>820</u>	<u>0.72</u>	<u>-46</u>	<u>7.1</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0735</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0735</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FDD1-20190716

0735-SAMPLE TIME

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.09</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.09</u>	ft. BGL
MW ID	<u>MW33A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW33A-20190717</u>		Water Quality	MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
1712	0.2	/	5.33	21.3	7.12	780	1.26	-31	4.2
1722	↓	/	5.32	21.0	7.08	710	1.01	-25	3.3
1727	↓	/	5.32	21.1	7.09	710	1.02	-26	3.6
1732									

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1750	60mL	P	1	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HNO <sub>3</sub>	METALS
1750	40mL / 1L	G/G	5	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HCL / NONE	VOCS / SVOCs

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FD02-20190717

1750-SAMPLE TIME



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>0.62</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>0.62</u>	ft. BGL
MW ID	<u>MW71B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW71B-20190717</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1606</u>	<u>1.2</u>	<u>/</u>	<u>1.1</u>	<u>22.7</u>	<u>6.91</u>	<u>510</u>	<u>0.46</u>	<u>-76</u>	<u>11</u>
<u>1621</u>	<u>↓</u>	<u>/</u>	<u>1.10</u>	<u>22.5</u>	<u>6.94</u>	<u>520</u>	<u>0.33</u>	<u>-71</u>	<u>7.9</u>
<u>1624</u>	<u>↓</u>	<u>/</u>	<u>1.10</u>	<u>22.5</u>	<u>6.93</u>	<u>510</u>	<u>0.34</u>	<u>-71</u>	<u>8.2</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1640</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1640</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.83</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>15.83</u> ft. BGL
MW ID	<u>MW87C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW87C-20190717</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1509</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1519</u>	<u>.2</u>	<u>/</u>	<u>16.12</u>	<u>22.4</u>	<u>7.46</u>	<u>1670</u>	<u>0.92</u>	<u>-92</u>	<u>10</u>
<u>1524</u>	<u>↓</u>	<u>/</u>	<u>16.13</u>	<u>22.6</u>	<u>7.61</u>	<u>1630</u>	<u>0.74</u>	<u>-76</u>	<u>8.2</u>
<u>1529</u>	<u>↓</u>	<u>/</u>	<u>16.13</u>	<u>22.3</u>	<u>7.61</u>	<u>1640</u>	<u>0.71</u>	<u>-71</u>	<u>9.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1540</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1540</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.67</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>8.67</u> ft. BGL
MW ID	<u>MW44A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW44A-20190717</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1329				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1339	<u>.2</u>	<u>/</u>	<u>8.96</u>	<u>22.7</u>	<u>6.62</u>	<u>910</u>	<u>0.99</u>	<u>-27</u>	<u>8.1</u>
1344	<u>↓</u>	<u>/</u>	<u>8.94</u>	<u>22.9</u>	<u>6.79</u>	<u>930</u>	<u>0.84</u>	<u>-21</u>	<u>6.2</u>
1349	<u>↓</u>	<u>/</u>	<u>8.94</u>	<u>22.4</u>	<u>6.56</u>	<u>920</u>	<u>0.86</u>	<u>-22</u>	<u>7.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1400</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1400</u>	<u>40mL/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVPCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.06</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.06</u> ft. BGL
MW ID	<u>MW59B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW59B-20190717</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1236</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1246</u>	<u>0.2</u>	<u>/</u>	<u>9.39</u>	<u>22.3</u>	<u>6.81</u>	<u>1260</u>	<u>0.71</u>	<u>-34</u>	<u>7.4</u>
<u>1251</u>	<u>↓</u>	<u>/</u>	<u>9.41</u>	<u>22.4</u>	<u>6.77</u>	<u>1250</u>	<u>0.22</u>	<u>-29</u>	<u>6.1</u>
<u>1256</u>	<u>↓</u>	<u>/</u>	<u>9.42</u>	<u>22.1</u>	<u>6.79</u>	<u>1240</u>	<u>0.16</u>	<u>-31</u>	<u>6.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1310</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1310</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BEAYTON name John Beayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.92</u> ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.92</u> ft. BGL
MW ID	<u>MW59A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620 - 2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1147				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1157	<u>2</u>	<u>/</u>	<u>9.23</u>	<u>23.2</u>	<u>6.96</u>	<u>1460</u>	<u>-72</u>	<u>0.51</u>	<u>8.1</u>
1202	<u>↓</u>	<u>/</u>	<u>9.24</u>	<u>23.1</u>	<u>6.92</u>	<u>1470</u>	<u>-68</u>	<u>0.42</u>	<u>7.1</u>
1207	<u>↓</u>	<u>/</u>	<u>9.23</u>	<u>23.1</u>	<u>6.93</u>	<u>1470</u>	<u>-69</u>	<u>0.42</u>	<u>7.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1220</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1220</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.23</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>16.23</u> ft. BGL
MW ID	<u>MW48C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW48C-20190717</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1048				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1058	<u>.2</u>	<u>/</u>	<u>16.47</u>	<u>21.4</u>	<u>6.79</u>	<u>1090</u>	<u>0.77</u>	<u>-31</u>	<u>16</u>
1103	<u>↓</u>	<u>/</u>	<u>16.51</u>	<u>21.7</u>	<u>6.77</u>	<u>1130</u>	<u>0.72</u>	<u>-25</u>	<u>10</u>
1108	<u>↓</u>	<u>/</u>	<u>16.52</u>	<u>21.9</u>	<u>6.72</u>	<u>1160</u>	<u>0.74</u>	<u>-26</u>	<u>10</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1120</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1120</u>	<u>40mL/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.43</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>16.43</u> ft. BGL
MW ID	<u>MW47C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW47C-20190717</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0956				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1002	<u>.2</u>	<u>/</u>	<u>16.78</u>	<u>22.5</u>	<u>6.74</u>	<u>960</u>	<u>1.31</u>	<u>-7</u>	<u>9.2</u>
1011	<u>↓</u>	<u>/</u>	<u>16.80</u>	<u>22.4</u>	<u>6.72</u>	<u>910</u>	<u>1.16</u>	<u>-6</u>	<u>7.4</u>
1016	<u>↓</u>	<u>/</u>	<u>16.81</u>	<u>22.4</u>	<u>6.71</u>	<u>920</u>	<u>1.13</u>	<u>-6</u>	<u>8.1</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1025</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1025</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.71</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.71</u>	ft. BGL
MW ID	<u>MW69A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW69A-20190717</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0901				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0911	<u>0.2</u>	<u>/</u>	<u>9.97</u>	<u>22.6</u>	<u>6.94</u>	<u>1730</u>	<u>1.13</u>	<u>-37</u>	<u>8.6</u>
0916	<u>↓</u>	<u>/</u>	<u>10.01</u>	<u>22.5</u>	<u>6.88</u>	<u>1640</u>	<u>0.95</u>	<u>-44</u>	<u>7.7</u>
0921	<u>↓</u>	<u>/</u>	<u>10.02</u>	<u>22.4</u>	<u>6.89</u>	<u>1620</u>	<u>0.96</u>	<u>-47</u>	<u>7.1</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0935	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0935	<u>40mL / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.52</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>5.52</u> ft. BGL
MW ID	<u>MW60A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620 MW60A - 20190717</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0802				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0812	<u>.2</u>	<u>/</u>	<u>5.81</u>	<u>22.3</u>	<u>6.98</u>	<u>1260</u>	<u>0.92</u>	<u>-77</u>	<u>9.1</u>
0817	<u>↓</u>	<u>/</u>	<u>5.83</u>	<u>22.4</u>	<u>6.97</u>	<u>1230</u>	<u>0.70</u>	<u>-76</u>	<u>8.4</u>
0822	<u>↓</u>	<u>/</u>	<u>5.84</u>	<u>22.4</u>	<u>6.96</u>	<u>1240</u>	<u>0.71</u>	<u>-76</u>	<u>8.2</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0835</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0835</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.96</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>6.96</u>	ft. BGL
MW ID	<u>MW61A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW61A-20190717</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0711				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0721	<u>0.2</u>	<u>/</u>	<u>7.16</u>	<u>21.7</u>	<u>6.71</u>	<u>1320</u>	<u>0.74</u>	<u>-106</u>	<u>8.1</u>
0726	<u>↓</u>	<u>/</u>	<u>7.17</u>	<u>21.6</u>	<u>6.73</u>	<u>1350</u>	<u>0.56</u>	<u>-101</u>	<u>7.7</u>
0731	<u>↓</u>	<u>/</u>	<u>7.17</u>	<u>21.7</u>	<u>6.74</u>	<u>1360</u>	<u>0.57</u>	<u>-102</u>	<u>7.6</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0750	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
0750	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

ms/msd Field Team Leader JOHN BRAYTON John Brayton

name signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.31</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-17-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.31</u>	ft. BGL
MW ID	<u>MW26A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW26A-20190717</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1816</u>	(L/min)	(L)	(ft)	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1826</u>	<u>.2</u>	<u>/</u>	<u>6.49</u>	<u>21.4</u>	<u>6.71</u>	<u>1740</u>	<u>1.39</u>	<u>-107</u>	<u>17</u>
<u>1831</u>	<u>↓</u>	<u>/</u>	<u>6.51</u>	<u>21.8</u>	<u>6.57</u>	<u>1710</u>	<u>1.03</u>	<u>-106</u>	<u>13</u>
<u>1836</u>	<u>↓</u>	<u>/</u>	<u>6.51</u>	<u>21.7</u>	<u>6.56</u>	<u>1720</u>	<u>1.04</u>	<u>-106</u>	<u>14</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1845</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1845</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**

 Page 1 of 1

Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>15.61</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>15.61</u>	ft. BGL
MW ID	<u>MW27C</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW27C-20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
1731									
1741	<u>1.2</u>	<u>/</u>	<u>15.92</u>	<u>22.2</u>	<u>6.51</u>	<u>1230</u>	<u>0.74</u>	<u>-44</u>	<u>6.8</u>
1744	<u>↓</u>	<u>/</u>	<u>15.94</u>	<u>22.3</u>	<u>6.41</u>	<u>1270</u>	<u>0.70</u>	<u>-42</u>	<u>5.1</u>
1751	<u>↓</u>	<u>/</u>	<u>15.95</u>	<u>22.3</u>	<u>6.42</u>	<u>1260</u>	<u>0.71</u>	<u>-43</u>	<u>5.2</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1805</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1805</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with &lt;1 foot of stable draw down.

 Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader

JOHN BRAYTON  
name

signature





# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.21</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.21</u>	ft. BGL
MW ID	<u>MW27A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW27A-20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1641				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1651	<u>2</u>	<u>/</u>	<u>7.47</u>	<u>21.7</u>	<u>7.06</u>	<u>1420</u>	<u>0.74</u>	<u>-21</u>	<u>4.1</u>
1656	<u>↓</u>	<u>/</u>	<u>7.49</u>	<u>21.6</u>	<u>7.09</u>	<u>1430</u>	<u>0.67</u>	<u>-22</u>	<u>9.6</u>
1701	<u>↓</u>	<u>/</u>	<u>7.49</u>	<u>21.6</u>	<u>7.08</u>	<u>1430</u>	<u>0.66</u>	<u>-22</u>	<u>9.7</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1715</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1715</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCs/SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>3.26</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>3.24</u>	ft. BGL
MW ID	<u>MW388</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW388-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1539				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1549	<u>.2</u>	<u>/</u>	<u>3.47</u>	<u>22.7</u>	<u>6.74</u>	<u>1340</u>	<u>0.79</u>	<u>-101</u>	<u>9.2</u>
1554	<u>↓</u>	<u>/</u>	<u>3.51</u>	<u>22.5</u>	<u>6.75</u>	<u>1320</u>	<u>0.51</u>	<u>-92</u>	<u>8.7</u>
1559	<u>↓</u>	<u>/</u>	<u>3.52</u>	<u>22.6</u>	<u>6.76</u>	<u>1330</u>	<u>0.52</u>	<u>-93</u>	<u>8.6</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1610</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1610</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader

JOHN BRAYTON  
name

signature

John Brayton

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.74</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.74</u>	ft. BGL
MW ID	<u>MW89B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW89B-20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1444</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1454</u>	<u>.2</u>	<u>/</u>	<u>7.93</u>	<u>22.7</u>	<u>6.81</u>	<u>1420</u>	<u>1.09</u>	<u>-84</u>	<u>8.6</u>
<u>1459</u>	<u>↓</u>	<u>/</u>	<u>7.96</u>	<u>22.6</u>	<u>6.77</u>	<u>1430</u>	<u>1.01</u>	<u>-80</u>	<u>7.1</u>
<u>1504</u>	<u>↓</u>	<u>/</u>	<u>7.97</u>	<u>22.6</u>	<u>6.76</u>	<u>1430</u>	<u>1.02</u>	<u>-81</u>	<u>7.3</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1515</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1515</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**

Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.13</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.13</u> ft. BGL
MW ID	<u>MW90B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW90B-20190719</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
1357									
1407	<u>2</u>	<u>/</u>	<u>4.31</u>	<u>23.2</u>	<u>6.93</u>	<u>1470</u>	<u>1.41</u>	<u>-41</u>	<u>7.4</u>
1412	<u>↓</u>	<u>/</u>	<u>4.32</u>	<u>23.1</u>	<u>6.92</u>	<u>1420</u>	<u>0.99</u>	<u>-42</u>	<u>6.5</u>
1417	<u>↓</u>	<u>/</u>	<u>4.31</u>	<u>23.2</u>	<u>6.94</u>	<u>1410</u>	<u>0.96</u>	<u>-42</u>	<u>6.7</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1425	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1425	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with &lt;1 foot of stable draw down.

 Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader

JOHN BRAYTON  
name

signature





**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.62</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.62</u>	ft. BGL
MW ID	<u>MW35B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620 MW35B -20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1256				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1306	<u>2</u>	<u>/</u>	<u>5.86</u>	<u>21.4</u>	<u>6.74</u>	<u>1940</u>	<u>0.81</u>	<u>-39</u>	<u>7.9</u>
1311	<u>↓</u>	<u>/</u>	<u>5.89</u>	<u>21.3</u>	<u>6.72</u>	<u>1920</u>	<u>0.71</u>	<u>-41</u>	<u>6.1</u>
1316	<u>↓</u>	<u>/</u>	<u>5.88</u>	<u>21.1</u>	<u>6.71</u>	<u>1910</u>	<u>0.72</u>	<u>-42</u>	<u>6.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1330	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1330	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.92</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>7.92</u>	ft. BGL
MW ID	<u>MW35A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW35A-20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1208				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1218	<u>2</u>	<u>/</u>	<u>8.16</u>	<u>21.4</u>	<u>6.64</u>	<u>1710</u>	<u>0.69</u>	<u>-41</u>	<u>6.7</u>
1223	<u>↓</u>	<u>/</u>	<u>8.17</u>	<u>21.6</u>	<u>6.60</u>	<u>1640</u>	<u>0.52</u>	<u>-33</u>	<u>5.4</u>
1228	<u>↓</u>	<u>/</u>	<u>8.16</u>	<u>21.7</u>	<u>6.61</u>	<u>1630</u>	<u>0.59</u>	<u>-34</u>	<u>5.9</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1240</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1240</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>17.17</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>17.17</u>	ft. BGL
MW ID	<u>MW83C</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW83C-20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1059				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1109	<u>0.2</u>	<u>/</u>	<u>17.39</u>	<u>23.4</u>	<u>7.16</u>	<u>1260</u>	<u>0.91</u>	<u>-28</u>	<u>4.9</u>
1114	<u>↓</u>	<u>/</u>	<u>17.41</u>	<u>23.2</u>	<u>7.12</u>	<u>1230</u>	<u>0.73</u>	<u>-25</u>	<u>7.1</u>
1119	<u>↓</u>	<u>/</u>	<u>17.42</u>	<u>23.1</u>	<u>7.13</u>	<u>1240</u>	<u>0.74</u>	<u>-26</u>	<u>7.2</u>
Purging was completed based on:			<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)						

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1130	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1130	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BEAYTON name John Beayton signature

# Groundwater Sample Collection



Project/Phase	19119232	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	6.11	ft. BMP
Site Location	UPRR - HWPW		<input type="checkbox"/> Decon between locations	Casing Stickup	-	ft.
Date	7-18-19	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	6.11	ft. BGL
MW ID	MW 83B	Location	Other _____	Total MW Depth	-	ft. BGL
Sample ID	WG-1620 MW 83B - 20190718	Water Quality		MW Diameter	2.0	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	HORIBA	MW Volume	-	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	U-50	Pump Intake Depth	-	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1008				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1018	2	/	6.34	22.1	6.74	1740	0.71	-61	16
1023	↓	/	6.36	22.1	6.72	1770	0.61	-61	12
1028	↓	/	6.36	22.2	6.71	1760	0.62	-62	13

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1040	60ML	P	1	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	HNO <sub>3</sub>	METALS
1040	40ML / 1L	G/G	5	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HCL / NONE	VOCs / SVOCs

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.39</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>16.39</u> ft. BGL
MW ID	<u>MW68C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW68C-20190718</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Watterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0913				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0923	<u>1.2</u>	<u>/</u>	<u>16.67</u>	<u>21.7</u>	<u>7.19</u>	<u>1560</u>	<u>1.16</u>	<u>-23</u>	<u>4.1</u>
0928	<u>↓</u>	<u>/</u>	<u>16.66</u>	<u>21.6</u>	<u>7.19</u>	<u>1520</u>	<u>0.92</u>	<u>-20</u>	<u>3.2</u>
0933	<u>↓</u>	<u>/</u>	<u>16.67</u>	<u>21.6</u>	<u>7.21</u>	<u>1530</u>	<u>0.91</u>	<u>-21</u>	<u>3.4</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0945</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> Unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0945</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> Unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.78</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.78</u>	ft. BGL
MW ID	<u>MW68B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW68B-20190718</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Watterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0816</u>	<u>(L/min)</u>	<u>(L)</u>	<u>(ft)</u>	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0826</u>	<u>1.2</u>	<u>/</u>	<u>5.96</u>	<u>22.6</u>	<u>6.71</u>	<u>940</u>	<u>0.86</u>	<u>-23</u>	<u>4.7</u>
<u>0831</u>	<u>↓</u>	<u>/</u>	<u>5.98</u>	<u>22.4</u>	<u>6.63</u>	<u>940</u>	<u>0.70</u>	<u>-22</u>	<u>4.9</u>
<u>0836</u>	<u>↓</u>	<u>/</u>	<u>5.98</u>	<u>22.4</u>	<u>6.64</u>	<u>930</u>	<u>0.71</u>	<u>-22</u>	<u>6.1</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0855</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>0855</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FDD3-20190718  
0855-SAMPLE TIME

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.01</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-18-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.01</u> ft. BGL
MW ID	<u>MW08A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW08A-20190718</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0721				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0731	<u>2</u>	<u>/</u>	<u>5.32</u>	<u>21.4</u>	<u>6.94</u>	<u>1040</u>	<u>0.79</u>	<u>-22</u>	<u>6.4</u>
0736	<u>↓</u>	<u>/</u>	<u>5.36</u>	<u>21.6</u>	<u>6.91</u>	<u>1010</u>	<u>0.61</u>	<u>-19</u>	<u>5.6</u>
0741	<u>↓</u>	<u>/</u>	<u>5.37</u>	<u>21.7</u>	<u>6.92</u>	<u>1020</u>	<u>0.62</u>	<u>-21</u>	<u>5.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0755	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
0755	<u>40ml / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.68</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.68</u> ft. BGL
MW ID	<u>MW79A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW79A-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1838</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1848</u>	<u>.2</u>	<u>/</u>	<u>9.89</u>	<u>23.4</u>	<u>6.94</u>	<u>860</u>	<u>1.06</u>	<u>-32</u>	<u>16</u>
<u>1853</u>	<u>↓</u>	<u>/</u>	<u>9.91</u>	<u>23.1</u>	<u>6.91</u>	<u>850</u>	<u>0.84</u>	<u>-32</u>	<u>13</u>
<u>1858</u>	<u>↓</u>	<u>/</u>	<u>9.91</u>	<u>23.1</u>	<u>6.91</u>	<u>810</u>	<u>0.87</u>	<u>-31</u>	<u>14</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1910</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1910</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.59</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>8.59</u> ft. BGL
MW ID	<u>MW74B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW74B-2019030</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1747				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1757	<u>.2</u>	<u>/</u>	<u>8.91</u>	<u>24.2</u>	<u>6.91</u>	<u>1370</u>	<u>0.67</u>	<u>-72</u>	<u>16</u>
1802	<u>↓</u>	<u>/</u>	<u>8.89</u>	<u>24.1</u>	<u>6.93</u>	<u>1330</u>	<u>0.52</u>	<u>-62</u>	<u>11</u>
1807	<u>↓</u>	<u>/</u>	<u>8.89</u>	<u>24.1</u>	<u>6.97</u>	<u>1340</u>	<u>0.53</u>	<u>-61</u>	<u>12</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1820</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1820</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>20.18</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>20.18</u>	ft. BGL
MW ID	<u>MW76C</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW76C-20190130</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1658				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1708	<u>.2</u>	<u>/</u>	<u>20.39</u>	<u>22.7</u>	<u>7.16</u>	<u>1920</u>	<u>0.59</u>	<u>-81</u>	<u>4.2</u>
1713	<u>↓</u>	<u>/</u>	<u>20.41</u>	<u>22.6</u>	<u>7.21</u>	<u>1920</u>	<u>0.42</u>	<u>-76</u>	<u>5.1</u>
1718	<u>↓</u>	<u>/</u>	<u>20.40</u>	<u>22.6</u>	<u>7.22</u>	<u>1930</u>	<u>0.41</u>	<u>-77</u>	<u>5.6</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1730	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1730	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>19.73</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>19.73</u> ft. BGL
MW ID	<u>MWB6C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MWB6C-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1601</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1611</u>	<u>2</u>	<u>/</u>	<u>19.96</u>	<u>22.7</u>	<u>7.29</u>	<u>1060</u>	<u>0.81</u>	<u>-47</u>	<u>6.9</u>
<u>1614</u>	<u>↓</u>	<u>/</u>	<u>19.98</u>	<u>22.6</u>	<u>7.22</u>	<u>1020</u>	<u>0.85</u>	<u>-42</u>	<u>7.1</u>
<u>1621</u>	<u>↓</u>	<u>/</u>	<u>19.98</u>	<u>22.6</u>	<u>7.21</u>	<u>1030</u>	<u>0.86</u>	<u>-42</u>	<u>5.6</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1640</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1640</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FD04-20190730  
1640-SAMPLE TIME

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>21.78</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>21.78</u> ft. BGL
MW ID	<u>MW85C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW85C-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1507</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1517</u>	<u>1.2</u>	<u>/</u>	<u>21.93</u>	<u>22.5</u>	<u>7.49</u>	<u>1340</u>	<u>1.09</u>	<u>-29</u>	<u>3.1</u>
<u>1522</u>	<u>↓</u>	<u>/</u>	<u>21.94</u>	<u>22.4</u>	<u>7.62</u>	<u>1360</u>	<u>1.04</u>	<u>-25</u>	<u>3.7</u>
<u>1527</u>	<u>↓</u>	<u>/</u>	<u>21.94</u>	<u>22.4</u>	<u>7.61</u>	<u>1370</u>	<u>1.06</u>	<u>-26</u>	<u>5.6</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

1535  
1535

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1540</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1540</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.11</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.11</u>	ft. BGL
MW ID	<u>MW77A</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW77A-20190730</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate	Cumulative Purge Volume	Depth to Water	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1122</u>	<u>(L/min)</u>	<u>(L)</u>	<u>(ft)</u>	<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1132</u>	<u>.2</u>	<u>/</u>	<u>6.33</u>	<u>22.3</u>	<u>6.91</u>	<u>1230</u>	<u>0.74</u>	<u>-51</u>	<u>6.7</u>
<u>1137</u>	<u>↓</u>	<u>/</u>	<u>6.32</u>	<u>22.2</u>	<u>6.94</u>	<u>1270</u>	<u>0.61</u>	<u>-44</u>	<u>5.2</u>
<u>1142</u>	<u>↓</u>	<u>/</u>	<u>6.32</u>	<u>22.1</u>	<u>6.96</u>	<u>1260</u>	<u>0.62</u>	<u>-46</u>	<u>5.1</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1155</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1155</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>11.02</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>11.02</u> ft. BGL
MW ID	<u>MW80B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW80B-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1037				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1047	<u>2</u>	<u>/</u>	<u>11.34</u>	<u>22.6</u>	<u>7.11</u>	<u>1060</u>	<u>0.42</u>	<u>-47</u>	<u>4.1</u>
1052	<u>↓</u>	<u>/</u>	<u>11.37</u>	<u>22.8</u>	<u>7.08</u>	<u>1020</u>	<u>0.41</u>	<u>-41</u>	<u>3.7</u>
1057	<u>↓</u>	<u>/</u>	<u>11.36</u>	<u>22.9</u>	<u>7.06</u>	<u>1030</u>	<u>0.41</u>	<u>-42</u>	<u>3.4</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1110</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1110</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.03</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.03</u> ft. BGL
MW ID	<u>MW82B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW82B-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0940				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0950	<u>2</u>	<u>/</u>	<u>4.27</u>	<u>22.5</u>	<u>6.94</u>	<u>1940</u>	<u>0.91</u>	<u>-21</u>	<u>19</u>
0955	<u>↓</u>	<u>/</u>	<u>4.29</u>	<u>22.5</u>	<u>6.92</u>	<u>1960</u>	<u>0.75</u>	<u>-15</u>	<u>15</u>
1000	<u>↓</u>	<u>/</u>	<u>4.30</u>	<u>22.4</u>	<u>6.91</u>	<u>1970</u>	<u>0.77</u>	<u>-16</u>	<u>16</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1010</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1010</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

ms/msd

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.09</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.09</u>	ft. BGL
MW ID	<u>MW 84B</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW84B-20190830</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
0843	2	/	7.31	22.7	7.11	560	0.56	-107	8.6
0858	↓	/	7.32	22.6	7.07	520	0.40	-103	5.1
0903	↓	/	7.31	22.6	7.06	530	0.41	-104	6.2

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0915	60ML	P	1	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HNO <sub>3</sub>	METALS
0915	40ML / 1L	G / G	5	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HCL / NONE	VOCs / SVOCs

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.92</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>5.92</u> ft. BGL
MW ID	<u>MW33BR</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW33BR-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0746				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0756	<u>.2</u>	<u>/</u>	<u>6.21</u>	<u>22.7</u>	<u>6.91</u>	<u>1570</u>	<u>0.61</u>	<u>-27</u>	<u>8.2</u>
0801	<u>↓</u>	<u>/</u>	<u>6.22</u>	<u>22.8</u>	<u>6.85</u>	<u>1590</u>	<u>0.63</u>	<u>-21</u>	<u>7.4</u>
0806	<u>↓</u>	<u>/</u>	<u>6.21</u>	<u>22.8</u>	<u>6.86</u>	<u>1590</u>	<u>0.62</u>	<u>-22</u>	<u>7.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0820</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0820</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.41</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-30-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.41</u> ft. BGL
MW ID	<u>MW32AR</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW32AR-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0651				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0701	<u>1.2</u>	<u>/</u>	<u>6.74</u>	<u>21.5</u>	<u>6.77</u>	<u>1910</u>	<u>0.63</u>	<u>-71</u>	<u>4.2</u>
0706	<u>↓</u>	<u>/</u>	<u>6.75</u>	<u>21.6</u>	<u>6.75</u>	<u>1830</u>	<u>0.55</u>	<u>-66</u>	<u>5.1</u>
0711	<u>↓</u>	<u>/</u>	<u>6.74</u>	<u>21.6</u>	<u>6.74</u>	<u>1820</u>	<u>0.56</u>	<u>-67</u>	<u>5.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0725</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0725</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	19119232	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	11.21	ft. BMP
Site Location	UPRR - HWPW		<input type="checkbox"/> Decon between locations	Casing Stickup	-	ft.
Date	7-31-19	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	11.21	ft. BGL
MW ID	MW49A	Location	Other _____	Total MW Depth	-	ft. BGL
Sample ID	WG-162D-MW49A-20190731	Water Quality		MW Diameter	2.0	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	HORIBA	MW Volume	-	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	U-50	Pump Intake Depth	-	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1714				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1726	2	/	11.47	22.7	6.91	1260	1.13	-36	21
1731	↓	/	11.51	22.5	6.83	1220	0.90	-41	16
1736	↓	/	11.51	22.4	6.84	1230	0.91	-41	17

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1745	60ML	P	1	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	HNO <sub>3</sub>	METALS
1745	40ML / 1L	G/G	5	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HCL / NONE	VOCs / SVOCs

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.09</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.09</u> ft. BGL
MW ID	<u>MW67B</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-1620-MW67B-20190731</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1601				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1617	<u>2</u>	<u>/</u>	<u>9.28</u>	<u>22.5</u>	<u>7.21</u>	<u>430</u>	<u>0.91</u>	<u>-79</u>	<u>8.2</u>
1622	<u>↓</u>	<u>/</u>	<u>9.27</u>	<u>22.4</u>	<u>7.18</u>	<u>410</u>	<u>0.73</u>	<u>-76</u>	<u>6.1</u>
1629	<u>↓</u>	<u>/</u>	<u>9.28</u>	<u>22.4</u>	<u>7.17</u>	<u>390</u>	<u>0.74</u>	<u>-77</u>	<u>6.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1645</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1645</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

MS/MSD Field Team Leader JOHN BRAYTON John Brayton  
name signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>24.13</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>24.13</u> ft. BGL
MW ID	<u>MW88C</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW88C-20190731</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1516</u>	<u>1.2</u>	<u>/</u>	<u>24.39</u>	<u>22.6</u>	<u>7.39</u>	<u>1270</u>	<u>0.59</u>	<u>-91</u>	<u>5.7</u>
<u>1531</u>	<u>↓</u>	<u>/</u>	<u>24.41</u>	<u>22.6</u>	<u>7.41</u>	<u>1270</u>	<u>0.46</u>	<u>-86</u>	<u>4.7</u>
<u>1536</u>	<u>↓</u>	<u>/</u>	<u>24.42</u>	<u>22.9</u>	<u>7.34</u>	<u>1260</u>	<u>0.47</u>	<u>-86</u>	<u>5.1</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G/P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1545</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1545</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.41</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.41</u>	ft. BGL
MW ID	<u>MW22 BR</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620-MW22 BR-20190731</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1321				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1331	<u>1.2</u>	<u>/</u>	<u>4.71</u>	<u>21.6</u>	<u>6.91</u>	<u>1260</u>	<u>0.91</u>	<u>-61</u>	<u>10</u>
1336	<u>↓</u>	<u>/</u>	<u>4.72</u>	<u>21.6</u>	<u>6.93</u>	<u>1230</u>	<u>0.74</u>	<u>-55</u>	<u>8.6</u>
1341	<u>↓</u>	<u>/</u>	<u>4.71</u>	<u>21.7</u>	<u>6.82</u>	<u>1240</u>	<u>0.71</u>	<u>-56</u>	<u>9.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
1350	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1350	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>3.96</u>	ft. BMP
Site Location	<u>UPRR - HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>3.96</u>	ft. BGL
MW ID	<u>MW22PR</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620 MW22PR 2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1232				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1242	<u>.2</u>	<u>/</u>	<u>4.23</u>	<u>22.4</u>	<u>6.59</u>	<u>1260</u>	<u>0.57</u>	<u>-43</u>	<u>16</u>
1247	<u>↓</u>	<u>/</u>	<u>4.24</u>	<u>22.2</u>	<u>6.53</u>	<u>1270</u>	<u>0.44</u>	<u>-41</u>	<u>12</u>
1252	<u>↓</u>	<u>/</u>	<u>4.22</u>	<u>22.1</u>	<u>6.52</u>	<u>1280</u>	<u>0.47</u>	<u>-41</u>	<u>13</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1305</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1305</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>81.31</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>81.31</u>	ft. BGL
MW ID	<u>MW59D</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-1620 MW59D-20190731</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1112</u>	<u>1.2</u>	<u>/</u>	<u>81.37</u>	<u>21.7</u>	<u>6.96</u>	<u>1120</u>	<u>0.93</u>	<u>-77</u>	<u>5.9</u>
<u>1127</u>	<u>↓</u>	<u>/</u>	<u>81.56</u>	<u>21.6</u>	<u>6.94</u>	<u>1140</u>	<u>0.70</u>	<u>-67</u>	<u>4.3</u>
<u>1132</u>	<u>↓</u>	<u>/</u>	<u>81.56</u>	<u>21.7</u>	<u>6.94</u>	<u>1150</u>	<u>0.71</u>	<u>-69</u>	<u>5.7</u>
Purging was completed based on:		<input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)							

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1150</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1150</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FDOS-20190731

1150 SAMPLE TIME



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>82.99</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>82.99</u> ft. BGL
MW ID	<u>MW66D</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW66D-20190731</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1013				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1023	<u>.2</u>	<u>/</u>	<u>83.18</u>	<u>22.6</u>	<u>7.11</u>	<u>11130</u>	<u>0.94</u>	<u>-86</u>	<u>6.7</u>
1028	<u>↓</u>	<u>/</u>	<u>83.19</u>	<u>22.7</u>	<u>7.11</u>	<u>1170</u>	<u>0.94</u>	<u>-80</u>	<u>6.1</u>
1033	<u>↓</u>	<u>/</u>	<u>83.18</u>	<u>22.7</u>	<u>7.09</u>	<u>1110</u>	<u>0.86</u>	<u>-81</u>	<u>5.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>82.71</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>82.71</u> ft. BGL
MW ID	<u>MW65D</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW65D-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0909				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0919	<u>.2</u>	<u>/</u>	<u>82.96</u>	<u>21.6</u>	<u>6.92</u>	<u>1290</u>	<u>0.89</u>	<u>-86</u>	<u>9.2</u>
0924	<u>↓</u>	<u>/</u>	<u>82.97</u>	<u>21.7</u>	<u>6.94</u>	<u>1250</u>	<u>0.74</u>	<u>-91</u>	<u>8.6</u>
0929	<u>↓</u>	<u>/</u>	<u>82.96</u>	<u>21.6</u>	<u>6.98</u>	<u>1260</u>	<u>0.77</u>	<u>-90</u>	<u>8.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0945</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

MS/MSD Field Team Leader JOHN BRAYTON John Brayton  
name signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.13</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>4.13</u> ft. BGL
MW ID	<u>MW38A</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW38A-20190731</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0802				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0812	<u>.2</u>	<u>/</u>	<u>4.32</u>	<u>22.6</u>	<u>6.47</u>	<u>1060</u>	<u>0.92</u>	<u>-74</u>	<u>27</u>
0817	<u>↓</u>	<u>/</u>	<u>4.33</u>	<u>22.5</u>	<u>6.48</u>	<u>1020</u>	<u>0.72</u>	<u>-66</u>	<u>21</u>
0822	<u>↓</u>	<u>/</u>	<u>4.33</u>	<u>22.4</u>	<u>6.42</u>	<u>1030</u>	<u>0.73</u>	<u>-62</u>	<u>22</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0835</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0835</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>82.91</u> ft. BMP
Site Location	<u>UPRR-HWPPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>7-31-19</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	Depth to Water	<u>82.91</u> ft. BGL
MW ID	<u>MW 36D-</u>	Location	Other _____	Total MW Depth	<u>-</u> ft. BGL
Sample ID	<u>WG-162D-MW36D-20190730</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0709				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0719	<u>.2</u>	<u>/</u>	<u>83.16</u>	<u>21.7</u>	<u>6.79</u>	<u>1340</u>	<u>1.06</u>	<u>-51</u>	<u>6.2</u>
0724	<u>↓</u>	<u>/</u>	<u>83.19</u>	<u>21.5</u>	<u>6.71</u>	<u>1310</u>	<u>0.92</u>	<u>-47</u>	<u>6.7</u>
0729	<u>↓</u>	<u>/</u>	<u>83.19</u>	<u>21.6</u>	<u>6.72</u>	<u>1310</u>	<u>0.93</u>	<u>-49</u>	<u>6.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
0740	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0740	<u>40ML / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
signature



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<b>GROUNDWATER SAMPLING RECORD</b>		PAGE ___ of ___
Project Number: 19119232	Project Name: UPRR HWPW	Date: 1-14-2020
Sampling Location (well ID, etc.): MW-01A	Starting Water Level (ft. BMP): 2.71	
Sample Number: WG-1620 MW01A 20200114	Casing Stickup (ft.):	
Sampled by: Tim McSpadden	WL (ft. BMP):	(ft. BGL):
Measuring Point (MP) of Well: TOC - Steel or PVC	TD (ft. BMP): 19.90	(ft. BGL):
Screened Interval (ft. BGL):	Ft. water:	Casing Dia. (In ID): 4"
Filter Pack Interval (ft. BGL):	1X Casing Vol (gal.):	3X (gal.):

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox-Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Horan Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: pH-4.01 NTU-0 COND 4.45  
 Filter / Filter Size:

SAMPLING MEASUREMENTS										
Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:55	2.71	.2	.2	24.17	4.54	.844	6.17	+78	26.7	Clear/m
14:05	2.80	.3	.2	24.01	2.16	.969	6.06	-67	28.4	
14:10	2.89	.4	.2	23.71	0.50	1.03	6.09	-52	11.2	
14:15	2.91	.5	.2	24.50	2.29	1.03	6.14	-26	9.9	
14:20	3.02	.6	.2	24.58	1.01	1.04	6.14	-25	9.6	
14:25	3.09	.7	.2	24.74	1.53	1.05	6.13	-21	9.6	
14:30	3.17	.8	.2	24.90	1.94	1.06	6.15	-16	8.6	
14:35	3.20	.9	.2	25.6	1.91	1.06	6.15	-16	8.4	

WL (ft. BMP) at End of Purge: 3.20 Sample Intake Depth (ft. BMP): 16.

SAMPLE INVENTORY						
Time	Bottles Collected			Filtration (Y / G)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (C P)	No.			
14:35			2		None	

Comments:  
 Duplicate, Field blank  
 WG-1620 FD01 20200114  
 WG-1620 FB02 20200114

**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: **19119232** Project Name: **UPRR MW/PW** Date: **1-14-2020**  
 Sampling Location (well ID, etc.): **MW-02** Starting Water Level (ft. BMP): **2.64**  
 Sample Number: **WG-1620 MW-02 20200114** Casing Stickup (ft.): **2'**  
 Sampled by: **T. M. McSpadden** WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or **(PVC)** TD (ft. BMP): **20.20** (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): **2'**  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): **Low Flow - Dedicated Tubing/Equipment**

Cleaning Equipment: **DI/Alconox Rinse**

Purge: **Peristaltic Pump / SS Pump / Bailer / Bladder**

Sampling: **Peristaltic Pump / SS Pump / Bailer / Bladder**

Disposal of Discharged Water: **Purge H<sub>2</sub>O drum 1-7-2020**

## INSTRUMENTS (Indicate make, model, I.D.)

Water Level: **P-211 Horon**

Other: **Geo Pump P 1102**

Multi Meter: **Humbel C1103**

Field Calibration: **PH-4.01 NTU-0 Comp 4.49**

Filter / Filter Size:

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft. BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µS/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:55	2.64	.2	.2	23.20	3.94	.605	6.57	-81	33.5	Clear/no
13:05	3.39	.3	.2	23.29	1.82	.605	6.44	-83	30.2	
13:10	3.47	.4	.2	22.91	0.93	.596	6.10	-85	28.7	
13:15	3.51	.5	.2	22.70	0.91	.592	6.16	-84	27.8	
13:20	3.55	.6	.2	22.32	0.59	.578	5.98	-79	40.4	
13:25	3.59	.7	.2	22.28	0.61	.578	5.99	-79	42.6	
13:30	3.62	.8	.2	22.24	0.60	.578	5.96	-77	40.0	

WL (ft. BMP) at End of Purge: **3.62**

Sample Intake Depth (ft. BMP): **17.0**

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
13:30			2		Heat	

Comments:



**GOLDER**

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 Round Rock, Texas 78664

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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_\_ of \_\_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-21-2020

Sample Number: WG-1620MW0320200121 Starting Water Level (ft. BMP): 3.05

Sampling Location (well ID, etc.): mw03 Casing Stickup (ft.): 2'

Sampled by: Tim MCSadden Starting Water Level (ft. BGL): \_\_\_\_\_

Measuring Point (MP) of Well: TBC Total Depth (ft. BMP): 18.90 soft bottom

Screened Interval (ft. BGL): \_\_\_\_\_ Casing Diameter (In ID): 2"

Filter Pack Interval (ft. BGL): \_\_\_\_\_ Casing Volume (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

METHODS (describe): Low flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.D.)

Water Level: P2111 Horos Thermometer: u<sup>6</sup>

pH Meter: Hanna CI103 Field Calibration: PA 4.0

Conductivity Meter: " " Field Calibration: \_\_\_\_\_

Filter / Filter Size: 10 microns Other: \_\_\_\_\_

## SAMPLING MEASUREMENTS

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
				± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	
14:20	3.05	.2	.2	18.08	6.82	.779	3.20	34	76.8	clear
14:30	3.41	.3	.2	18.10	6.72	.779	1.08	41	70.4	
14:35	3.34	.4	.2	18.09	6.63	.783	.25	53	44.9	
14:40	3.28	.5	.2	17.91	6.60	.776	.15	57	53.3	
14:45	3.30	.6	.2	17.93	6.60	.773	0	60	33.0	
14:50	3.28	.7	.2	17.71	6.56	.774	0	63	39.9	
14:55	3.29	.8	.2	17.67	6.55	.774	0	64	41.2	
15:00	3.31	.9	.2	17.58	6.55	.776	0	65	47.8	

Water Level (ft. BMP) at End of Purge: 3.30 Sample Intake Depth (ft. BMP): 15.0

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
1500			6	Y		

Comments:



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(832) 916-3690



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-21-2020  
 Sample Number: W61620 MW03 20200121 Starting Water Level (ft. BMP): 4.61  
 Sampling Location (well ID, etc.): MW 04 Casing Stickup (ft.): 1'  
 Sampled by: Tim McSpadden Starting Water Level (ft. BGL):  
 Measuring Point (MP) of Well: Total Depth (ft. BMP): 21.38 soft bottom  
 Screened Interval (ft. BGL): Casing Diameter (In ID):  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment:  
 Purging: Sampling:  
 Disposal of Discharged Water:

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: P2111 Horva Thermometer: " "  
 pH Meter: Horiba C1103 Field Calibration: PH 4.0  
 Conductivity Meter: " " Field Calibration:  
 Filter / Filter Size: Other:

**SAMPLING MEASUREMENTS**


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
15:20	4.61	.2	.2	17.98	6.74	.626	3.19	67	51.8	
15:30	4.96	.3	.2	18.09	6.66	-.628	2.22	71	52.5	
15:35	5.00	.4	.2	18.17	6.62	-.628	1.01	76	55.0	
15:40	5.08	.5	.2	18.16	6.59	-.626	1.66	80	45.3	
15:45	5.10	.6	.2	18.13	6.52	-.626	1.40	84	42.1	
15:50	5.10	.7	.2	18.60	6.50	-.621	1.23	96	56.5	
15:55	5.12	.8	.2	18.72	6.45	-.621	1.23	97	37.0	
16:00	5.12	.9	.2							

Water Level (ft. BMP) at End of Purge: 5.12 Sample Intake Depth (ft. BMP): 18.0

**SAMPLE INVENTORY**

Bottles Collected				Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
Time	Volume	Composition (G/P)	No.			
16:00			6			

Comments:



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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.69</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.69</u>	ft. BGL
MW ID	<u>MW05</u>	Location	Other _____	Total MW Depth	<u>27.30</u>	ft. BGL
Sample ID	<u>WG-1620-MW05-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1422				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1432	<u>.2</u>	<u>/</u>	<u>4.92</u>	<u>21.7</u>	<u>6.94</u>	<u>690</u>	<u>0.86</u>	<u>-52</u>	<u>6.7</u>
1437	<u>↓</u>	<u>/</u>	<u>4.91</u>	<u>21.6</u>	<u>6.93</u>	<u>660</u>	<u>0.59</u>	<u>-44</u>	<u>4.1</u>
1442	<u>↓</u>	<u>/</u>	<u>4.92</u>	<u>21.6</u>	<u>6.92</u>	<u>670</u>	<u>0.57</u>	<u>-46</u>	<u>4.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1455	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1455	<u>40mL / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: VPRR HWPW Date: 1-13-2020  
 Sampling Location (well ID, etc.): MW 07 Starting Water Level (ft. BMP): 4.29  
 Sample Number: AWG-1620 MW07 2020 0113 Casing Stickup (ft.): 3'  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC Steel or (FVC) TD (ft. BMP): 24.90 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 4"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE**

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe) Low Flow - Dedicated Tubing/Equipment

Cleaning Equipment: DI/Alconox Rinse

Purge: Peristaltic Pump / SS Pump / Bailer / Bladder

Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder

Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: P2111 Heron

Other: Geo Pump P1103

Multi Meter: Homba C1103

Field Calibration: pH-3.94 NTU-0 conc 449

Filter / Filter Size: \_\_\_\_\_

**SAMPLING MEASUREMENTS**

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:30	4.29	.2	.2	17.71	9.11	252	6.56	149	2.50	Clear/no
12:40	4.41	.3	.2	17.79	5.91	.821	7.30	110	512	
12:45	4.43	.4	.2	19.02	4.97	.814	7.16	116	59.5	
12:50	4.44	.5	.2	19.30	4.86	.813	7.15	117	45.4	
12:55	4.46	.6	.2	19.34	4.84	.813	7.13	118	33.3	
13:00	4.46	.7	.2	19.31	4.85	.813	7.11	119	29.4	
13:05	4.57	.8	.2	19.65	4.57	.811	7.06	121	15.9	
13:10	4.52	.9	.2	19.61	4.60	.812	7.05	120	15.4	

WL (ft. BMP) at End of Purge: 4.52

Sample Intake Depth (ft. BMP): 21.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
13:15			2		Heat	

Comments:



**GOLDER**

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 Round Rock, Texas 78664

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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-13-2020  
 Sampling Location (well ID, etc.): MW 08 Starting Water Level (ft. BMP): 4.45  
 Sample Number: WG-1620 MW 08 20200113 Casing Stickup (ft.): 2.5'  
 Sampled by: T.M. McSpadden WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TOC - Steel or FVC TD (ft. BMP): 25.10 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID): 4"  
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O down 1-7-2020

**INSTRUMENTS** (Indicate make, model, I.d.)

Water Level: P2111 Horon Other: Geo Pump P1102  
 Multi Meter: Honke C1103  
 Field Calibration: PH-392 NTU-0 cond 449  
 Filter / Filter Size:     

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
11:00	4.45	.2	.2	20.32	4.74	.612	7.22	100	33.7	Clear/No
11:10	4.61	.3	.2	20.16	4.63	.610	7.20	102	26.4	
11:05	4.63	.4	.2	19.98	4.54	.609	7.16	104	19.9	
11:10	4.65	.5	.2	19.83	4.40	.607	7.18	105	12.1	
11:15	4.65	.6	.2	19.90	3.90	.603	7.14	109	6.1	
11:20	4.64	.7	.2	19.69	3.86	.604	7.13	110	4.9	
11:25	4.65	.8	.2	19.66	3.78	.604	7.12	111	2.1	
11:30	4.65	.9	.2	19.59	3.62	.605	7.10	112	0.1	

WL (ft. BMP) at End of Purge: 4.65 Sample Intake Depth (ft. BMP): 22

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (C/P)	No.			
11:30			2		Neat	

Comments:  
Field blank collected @ MW-08

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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.11</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.11</u>	ft. BGL
MW ID	<u>MW09</u>	Location	Other _____	Total MW Depth	<u>25.40</u>	ft. BGL
Sample ID	<u>WG-1620-MW09-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
0826									
0836	<u>2</u>	<u>/</u>	<u>4.31</u>	<u>23.6</u>	<u>6.71</u>	<u>660</u>	<u>1.29</u>	<u>-26</u>	<u>4.7</u>
0841	<u>↓</u>	<u>/</u>	<u>4.33</u>	<u>23.2</u>	<u>6.64</u>	<u>620</u>	<u>1.06</u>	<u>-22</u>	<u>3.1</u>
0846	<u>↓</u>	<u>/</u>	<u>4.33</u>	<u>23.4</u>	<u>6.67</u>	<u>630</u>	<u>1.16</u>	<u>-23</u>	<u>3.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0900	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0900	<u>40ml / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR Houston Date: 1-14-2020  
 Sampling Location (well ID, etc.): MW 10A Starting Water Level (ft. BMP): 4.67  
 Sample Number: WG 1620 MW 10A 20200119 Casing Stickup (ft.): 3'  
 Sampled by: TIM MCS podden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC Steel or PVC TD (ft. BMP): 25.55 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 4"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow, Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge 170 drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P 2111 Heron Other: Geo Pump P1102  
 Multi Meter: Hanna C1103  
 Field Calibration: Ph-4.01 NTU-0 Comp 4.49  
 Filter / Filter Size: \_\_\_\_\_

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
11:45	4.67	.2	.2	22.74	2.35	.947	6.24	-21	0.0	
11:55	4.96	.3	.2	22.76	2.18	.944	6.26	71	0.0	
12:00	5.19	.4	.2	22.75	2.21	.942	6.27	18	0.0	
12:05	5.24	.5	.2	22.67	2.04	.940	6.28	40	0.0	
12:10	5.29	.6	.2	22.83	1.98	.945	6.29	52	0.0	
12:15	5.31	.7	.2	22.89	1.95	.945	6.29	54	0.0	
12:20	5.33	.8	.2	22.92	1.96	.944	6.30	55	0.0	
12:25	5.34	.9	.2	22.94	1.92	.944	6.30	57	0.0	

WL (ft. BMP) at End of Purge: 5.34 Sample Intake Depth (ft. BMP): 2.2

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
12:25			2		heat	

Comments:



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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: **19119232** Project Name: **UPRR HWPU** Date: **1-14-2020**  
 Sampling Location (well ID, etc.): **MW-103** Starting Water Level (ft. BMP): **4.90**  
 Sample Number: **WG 1620 MW 103 20200114** Casing Stickup (ft.): **3.**  
 Sampled by: **TIM McSpadden** WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or **PVC** TD (ft. BMP): **46.52** (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): **4"**  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): **Low Flow** - Dedicated Tubing/Equipment

Cleaning Equipment: **DI/Alconox Rinse**

Purge: **Peristaltic Pump** / SS Pump / Bailer / Bladder

Sampling: **Peristaltic Pump** / SS Pump / Bailer / Bladder

Disposal of Discharged Water: **Purge water drum 1-7-2020**

INSTRUMENTS (Indicate make, model, I.d.)

Water Level: **P2111 HANON**

Other: **Geo Pump P1102**

Multi Meter: **HANNA C1103**

Field Calibration: **PH-4.01 NTC-0 CAL0449**

Filter / Filter Size:

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
11:00	4.90	.2	.2	22.07	1.94	.990	6.28	-90	17.5	clear/no
11:10	5.15	.3	.2	22.26	0.19	1.06	6.05	-92	0.0	
11:15	5.17	.4	.2	23.30	0.21	1.08	5.90	-94	0.0	
11:20	5.19	.5	.2	23.26	0.05	1.07	5.90	-94	0.0	
11:25	5.19	.6	.2	23.05	0.00	1.06	5.91	-95	0.0	
11:30	5.21	.7	.2	22.78	0.00	1.06	5.92	-97	0.0	
11:35	5.23	.8	.2	22.74	0.00	1.07	5.90	-96	0.0	

WL (ft. BMP) at End of Purge: **5.23**

Sample Intake Depth (ft. BMP): **43.**

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (C, P)	No.			
11:35			2		Meat	

Comments:



**GOLDER**

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Q

16

**GROUNDWATER SAMPLING RECORD** PAGE \_\_\_\_ of \_\_\_\_

Project Number: 19119232 Project Name: UPRR HWPu Date: 1-14-2020  
 Sampling Location (well ID, etc.): MW 11A Starting Water Level (ft. BMP): 5.11  
 Sample Number: WG 1620 MW 11A 20200114 Casing Stickup (ft.): 3'  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 24.05 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 4"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P-2111 Heron Other: Geo pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: PH-401 NTU-0 COND 4.49  
 Filter / Filter Size: \_\_\_\_\_

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
9:55	5.11	.2	.2	21.27	0.40	1.03	6.09	-81	0.0	
10:05	5.22	.3	.2	21.22	0.91	0.921	6.14	-19	0.0	
10:10	5.51	.4	.2	21.30	2.56	.896	6.19	56	11.7	
10:15	5.55	.5	.2	21.27	2.62	.893	6.20	63	10.9	
10:20	5.58	.6	.2	21.37	2.47	.884	6.22	89	7.3	
10:25	5.58	.7	.2	21.41	2.38	.881	6.22	93	3.9	
10:30	5.59	.8	.2	21.44	2.42	.881	6.22	93	4.5	
10:35	5.59	.9	.2	21.49	2.29	.880	6.22	95	7.4	

WL (ft. BMP) at End of Purge: 5.59 Sample Intake Depth (ft. BMP): 21

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (GP)	No.			
10:35			2		Heat	

Comments:  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-14-2020  
 Sampling Location (well ID, etc.): MW-113 Starting Water Level (ft. BMP): 5.30  
 Sample Number: WG 1620 MW 113 2020 014 Casing Stickup (ft.): 3'  
 Sampled by: Tim McSpedden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 46.80 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 44  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Hanna C1103  
 Field Calibration: Ph 4.01 NTU=0 Comp. 449  
 Filter / Filter Size: -

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:45	5.30	.2	.2	20.92	2.08	1.04	5.20	-71	0.0	Clear/No
8:55	5.62	.5	.2	21.09	0.57	1.03	5.37	-94	0.0	
9:00	5.73	.9	.2	21.12	0.41	1.03	5.47	-102	0.0	
9:05	5.81	.5	.2	21.01	0.00	1.02	5.78	-101	0.0	
9:10	5.86	.6	.2	21.06	0.00	1.02	5.88	-101	0.0	
9:15	5.91	.7	.2	21.09	0.00	1.02	5.90	-101	0.0	
9:20	5.97	.8	.2	21.12	0.00	1.02	5.92	-101	0.0	

WL (ft. BMP) at End of Purge: 5.97 Sample Intake Depth (ft. BMP): 41.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
9:25			2		Neat	

Comments: \_\_\_\_\_



**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.04</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.04</u>	ft. BGL
MW ID	<u>MW12A</u>	Location	Other _____	Total MW Depth	<u>30.35</u>	ft. BGL
Sample ID	<u>WG-1620-MW12A-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1612</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1622</u>	<u>.3</u>	<u>/</u>	<u>6.29</u>	<u>23.6</u>	<u>6.79</u>	<u>670</u>	<u>1.21</u>	<u>-37</u>	<u>2.4</u>
<u>1627</u>	<u>↓</u>	<u>/</u>	<u>6.31</u>	<u>23.5</u>	<u>6.72</u>	<u>610</u>	<u>1.03</u>	<u>-33</u>	<u>5.9</u>
<u>1632</u>	<u>↓</u>	<u>/</u>	<u>6.32</u>	<u>23.4</u>	<u>6.71</u>	<u>590</u>	<u>1.04</u>	<u>-34</u>	<u>6.1</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1645</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1645</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HarpW Date: 1-23-2020  
 Sample Number: WG 1620 MW 12B 2020023 Starting Water Level (ft. BMP): 5.66  
 Sampling Location (well ID, etc.): MW 12B Casing Stickup (ft.): 4'  
 Sampled by: TIM McSpodden Starting Water Level (ft. BGL):  
 Measuring Point (MP) of Well: Total Depth (ft. BMP): 46.01  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment:  
 Purging: Sampling:  
 Disposal of Discharged Water:

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: P2111 Herau Hi Water Thermometer: " "  
 pH Meter: Hanna CI103 Field Calibration: PH 4.0 PH 4.0  
 Conductivity Meter: " " Field Calibration:  
 Filter / Filter Size: 10 micron Other: Oreo pump D1102

**SAMPLING MEASUREMENTS**


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
13:20	5.66	.2	.2	19.65	7.04	1.09	0.91	-89	67.7	Clear
13:30	6.71	.3	.2	19.32	7.03	1.10	0.25	-89	0.20	
13:35	6.75	.4	.2	19.24	7.00	1.10	0.13	-89	51.7	
13:40	6.84	.5	.2	19.20	7.02	1.09	0	-89	49.7	
13:45	6.88	.6	.2	19.13	7.01	1.08	0	-89	52.6	
13:50	6.93	.7	.2	19.10	7.01	1.08	0	-89	52.9	
13:55	6.99	.8	.2	19.02	7.00	1.08	0	-89	52.1	
14:00	6.99	.9	.2	18.99	7.01	1.08	0	-88	51.4	

Water Level (ft. BMP) at End of Purge: 6.99 Sample Intake Depth (ft. BMP): 20.

**SAMPLE INVENTORY**

Bottles Collected				Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
Time	Volume	Composition (G.P)	No.			
14:00			6	10 micron		

Comments:



**Golder Associates Inc.**  
 11231 Richmond Avenue, Suite D104  
 Houston, TX 77082  
 (832) 916-3690



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>24.61</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>24.61</u>	ft. BGL
MW ID	<u>MW-12C</u>	Location	Other _____	Total MW Depth	<u>75.60</u>	ft. BGL
Sample ID	<u>WG-162D-MW12C-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1519</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1529</u>	<u>.2</u>	<u>/</u>	<u>24.87</u>	<u>21.3</u>	<u>6.92</u>	<u>840</u>	<u>1.01</u>	<u>-86</u>	<u>6.7</u>
<u>1534</u>	<u>↓</u>	<u>/</u>	<u>24.91</u>	<u>21.1</u>	<u>6.83</u>	<u>860</u>	<u>0.83</u>	<u>-77</u>	<u>8.1</u>
<u>1539</u>	<u>↓</u>	<u>/</u>	<u>24.90</u>	<u>21.1</u>	<u>6.84</u>	<u>860</u>	<u>0.81</u>	<u>-79</u>	<u>9.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1555</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1555</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.42</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.42</u>	ft. BGL
MW ID	<u>MW13</u>	Location	Other _____	Total MW Depth	<u>26.10</u>	ft. BGL
Sample ID	<u>WG-1620-MW13-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0739				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0749	<u>.2</u>	<u>/</u>	<u>9.67</u>	<u>21.2</u>	<u>6.92</u>	<u>710</u>	<u>0.96</u>	<u>-62</u>	<u>12</u>
0754	<u>↓</u>	<u>/</u>	<u>9.71</u>	<u>21.2</u>	<u>6.94</u>	<u>720</u>	<u>0.71</u>	<u>-56</u>	<u>18</u>
0759	<u>↓</u>	<u>/</u>	<u>9.72</u>	<u>21.3</u>	<u>6.97</u>	<u>720</u>	<u>0.72</u>	<u>-57</u>	<u>17</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0810	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0810	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.39</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.39</u>	ft. BGL
MW ID	<u>MW14</u>	Location	Other _____	Total MW Depth	<u>44.80</u>	ft. BGL
Sample ID	<u>WG-1620-MW14 - 20190114</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1546				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1556	<u>2</u>	<u>/</u>	<u>8.67</u>	<u>22.4</u>	<u>6.67</u>	<u>1040</u>	<u>1.17</u>	<u>-67</u>	<u>21</u>
1601	<u>↓</u>	<u>/</u>	<u>8.71</u>	<u>22.3</u>	<u>6.61</u>	<u>1050</u>	<u>1.11</u>	<u>-61</u>	<u>15</u>
1606	<u>↓</u>	<u>/</u>	<u>8.72</u>	<u>22.3</u>	<u>6.59</u>	<u>1030</u>	<u>1.09</u>	<u>-62</u>	<u>16</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1615	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1615	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton

name signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.57</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.57</u>	ft. BGL
MW ID	<u>MWISA</u>	Location	Other _____	Total MW Depth	<u>29.90</u>	ft. BGL
Sample ID	<u>WG-162D-MWISA-20190114</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1313				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1322	<u>.2</u>	<u>/</u>	<u>8.79</u>	<u>21.6</u>	<u>6.86</u>	<u>1160</u>	<u>0.91</u>	<u>-76</u>	<u>13</u>
1326	<u>↓</u>	<u>/</u>	<u>8.81</u>	<u>21.4</u>	<u>6.77</u>	<u>1130</u>	<u>0.77</u>	<u>-71</u>	<u>10</u>
1333	<u>↓</u>	<u>/</u>	<u>8.81</u>	<u>21.7</u>	<u>6.79</u>	<u>1120</u>	<u>0.74</u>	<u>-73</u>	<u>11</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1345</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1345</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
 name John Brayton  
 signature [Signature]



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.39</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.39</u>	ft. BGL
MW ID	<u>MWISB</u>	Location	Other _____	Total MW Depth	<u>40.70</u>	ft. BGL
Sample ID	<u>WG-1620-MWISB-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
1358									
1408	<u>.2</u>	<u>/</u>	<u>8.61</u>	<u>22.6</u>	<u>6.91</u>	<u>930</u>	<u>0.71</u>	<u>-51</u>	<u>6.7</u>
1413	<u>↓</u>	<u>/</u>	<u>8.62</u>	<u>22.4</u>	<u>6.86</u>	<u>940</u>	<u>0.53</u>	<u>-55</u>	<u>7.4</u>
1418	<u>↓</u>	<u>/</u>	<u>8.61</u>	<u>22.1</u>	<u>6.88</u>	<u>910</u>	<u>0.56</u>	<u>-56</u>	<u>7.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1430</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1430</u>	<u>40ML / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>23.12</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>23.12</u>	ft. BGL
MW ID	<u>MWISC</u>	Location	Other _____	Total MW Depth	<u>75.25</u>	ft. BGL
Sample ID	<u>WG-162D-MWISG20190114</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>7</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>7</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1443				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1453	<u>.2</u>	/	<u>23.31</u>	<u>21.6</u>	<u>6.79</u>	<u>860</u>	<u>1.12</u>	<u>-74</u>	<u>8.6</u>
1458	↓	/	<u>23.32</u>	<u>21.1</u>	<u>6.81</u>	<u>810</u>	<u>0.93</u>	<u>-62</u>	<u>6.8</u>
1503	↓	/	<u>23.32</u>	<u>21.3</u>	<u>6.82</u>	<u>820</u>	<u>0.92</u>	<u>-61</u>	<u>7.1</u>

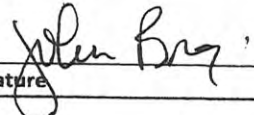
Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1520	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1520	<u>40ML / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON 

name signature



# Groundwater Sample Collection



Project/Phase	19119232	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	8.23	ft. BMP
Site Location	UPRR-HWPW		<input type="checkbox"/> Decon between locations	Casing Stickup	-	ft.
Date	1-14-20	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	8.23	ft. BGL
MW ID	MW7	Location	Other _____	Total MW Depth	36.95	ft. BGL
Sample ID	WG-1620-MW7-2019	Water Quality		MW Diameter	2.0	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	HORIBA	MW Volume	-	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	U-50	Pump Intake Depth	-	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1637				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1647	2	/	8.42	21.6	7.11	1440	0.52	-61	16
1652	↓	/	8.46	21.7	7.06	1420	0.45	-52	11
1657	↓	/	8.47	21.9	7.04	1410	0.46	-51	12

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1710	60ml	P	1	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	HNO <sub>3</sub>	METALS
1710	40ml / 1L	G/G		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	HCL / NONE	VOCS / SVOCS

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>23.21</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>23.21</u>	ft. BGL
MW ID	<u>MW7C</u>	Location	Other _____	Total MW Depth	<u>72.15</u>	ft. BGL
Sample ID	<u>WG-162D-MW7C-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Watterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0735				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0745	<u>2</u>	<u>/</u>	<u>23.39</u>	<u>22.6</u>	<u>6.92</u>	<u>1270</u>	<u>1.12</u>	<u>-46</u>	<u>11</u>
0754	<u>↓</u>	<u>/</u>	<u>23.41</u>	<u>22.2</u>	<u>6.96</u>	<u>1260</u>	<u>0.84</u>	<u>-36</u>	<u>8.2</u>
0759	<u>↓</u>	<u>/</u>	<u>23.41</u>	<u>22.1</u>	<u>6.97</u>	<u>1260</u>	<u>0.83</u>	<u>-39</u>	<u>9.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0815	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
0815	<u>40ML / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



# GROUNDWATER SAMPLING RECORD

PAGE 1 of \_\_\_\_

Project Number: 19119232 Project Name: HWPW Date: 1-8-2020  
 Sampling Location (well ID, etc.): MW-18A Starting Water Level (ft. BMP): 17.71  
 Sample Number: WC-1620-MW18A 20200108 Casing Stickup (ft.): 18"  
 Sampled by: TIM McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 35.65 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow / Dedicated Tubing / Equipment  
 Cleaning Equipment: D1/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: D2111 Heron Other: Geo pump P1102  
 Multi Meter: Hanba C1103  
 Field Calibration: PH-4.04 NTU-0 COND.450  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft. B <sup>1</sup> OC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
10:30	17.71	.2	.2	24.03	11.41	1.88	6.78	-1.55	0.0	
10:40	19.23	.3	.2	24.12	0.52	1.80	6.80	-1.54	0.0	
10:45	19.29	.4	.2	24.00	0.04	1.65	6.81	-1.46	973	
10:50	19.32	.5	.2	23.96	0.00	1.56	6.77	-1.45	1000	
10:55	19.36	.6	.2	23.94	0.00	1.38	6.78	-1.37	1000	
11:00	19.40	.7	.2	23.96	0.00	1.02	6.83	-1.26	990	
11:05	19.42	.8	.2	24.12	0.00	1.952	6.82	-1.25	978	
11:10	19.43	.9	.2	24.15	0.00	1.933	6.83	-1.25	986	

WL (ft. BMP) at End of Purge: 19.43 Sample Intake Depth (ft. BMP): 32.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
11:10	2.5		9			

Comments:


**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

**GROUNDWATER SAMPLING RECORD**

Project Number: HWPW19119232 Project Name: HWPW Date: 1-8-2020  
 Sampling Location (well ID, etc.): MW-18C Starting Water Level (ft. BMP): 23.54  
 Sample Number: WG-1620-MW18C 20200108 Casing Stickup (ft.): 14'  
 Sampled by: T.M. McSpadden WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 61.80 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge Hydro drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: D 2111 Horow Other: Geo Pum P1102  
 Multi Meter: Humba C1103  
 Field Calibration: Ph: 4.04 NTU: 0 COND: 450  
 Filter / Filter Size: W m 002

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
9:30	23.54	.2	.2	19.99	1.98	3.26	11.83	-196	253	clear/no
9:40	25.01	.3	.2	19.78	0.25	3.26	11.88	-203	226	
9:45	25.04	.4	.2	20.33	0.39	3.26	11.90	-214	186	
9:50	25.09	.5	.2	20.38	0.40	3.27	11.91	-219	181	
9:55	25.13	.6	.2	20.75	0.16	3.28	11.90	-220	167	
10:00	25.14	.7	.2	21.21	0.12	3.30	11.90	-224	193	
10:05	25.16	.8	.2	21.53	0.01	3.30	11.89	-230	116	
10:10	25.17	.9	.2	21.58	0.00	3.30	11.89	-230	108	

WL (ft. BMP) at End of Purge: 25.17 Sample Intake Depth (ft. BMP): 60.

**SAMPLE INVENTORY**

Bottles Collected				Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
Time	Volume	Composition (G/P)	No.			
10:10	2.0		9			

Comments:  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_


**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>25.71</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>1.0</u>	ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>24.71</u>	ft. BGL
MW ID	<u>MW-19C</u>	Location	Other _____	Total MW Depth	<u>75.15</u>	ft. BGL
Sample ID	<u>WG-162D-MW19C-2020009</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1016				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1021	<u>1.2</u>	<u>/</u>	<u>26.06</u>	<u>21.9</u>	<u>6.67</u>	<u>1500</u>	<u>2.42</u>	<u>-111</u>	<u>17</u>
1026	<u>↓</u>	<u>/</u>	<u>26.12</u>	<u>21.8</u>	<u>6.71</u>	<u>1530</u>	<u>2.06</u>	<u>-107</u>	<u>13</u>
1031	<u>↓</u>	<u>/</u>	<u>26.13</u>	<u>21.8</u>	<u>6.72</u>	<u>1540</u>	<u>2.07</u>	<u>-108</u>	<u>14</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm) / 45µm	(type)	(quality control sample, other)
1045	<u>60ml</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1045	<u>40ml / 1L</u>	<u>G/G</u>	<u>5</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.06</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.06</u> ft. BGL
MW ID	<u>MW20A</u>	Location	Other _____	Total MW Depth	<u>22.80</u> ft. BGL
Sample ID	<u>WG-162D-MW20A-20190114</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
<u>1219</u>									
<u>1228</u>	<u>.2</u>		<u>9.26</u>	<u>21.6</u>	<u>6.74</u>	<u>1440</u>	<u>0.96</u>	<u>-91</u>	<u>5.1</u>
<u>1234</u>	<u>↓</u>		<u>9.31</u>	<u>21.2</u>	<u>6.76</u>	<u>1410</u>	<u>0.73</u>	<u>-86</u>	<u>3.8</u>
<u>1239</u>	<u>↓</u>		<u>9.32</u>	<u>21.3</u>	<u>6.74</u>	<u>1420</u>	<u>0.74</u>	<u>-86</u>	<u>4.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1255</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1255</u>	<u>40ML / IL</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>23.46</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>—</u>	ft.
Date	_____	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>23.46</u>	ft. BGL
MW ID	<u>MW21C</u>	Location	Other _____	Total MW Depth	<u>74.60</u>	ft. BGL
Sample ID	<u>WG-1620-MW21C-20190114</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0922				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0932	<u>.2</u>	<u>/</u>	<u>23.61</u>	<u>21.3</u>	<u>7.11</u>	<u>910</u>	<u>0.96</u>	<u>-36</u>	<u>10</u>
0937	<u>↓</u>	<u>/</u>	<u>23.64</u>	<u>21.2</u>	<u>7.10</u>	<u>850</u>	<u>0.76</u>	<u>-31</u>	<u>7.4</u>
0942	<u>↓</u>	<u>/</u>	<u>23.64</u>	<u>21.3</u>	<u>7.06</u>	<u>860</u>	<u>0.79</u>	<u>-31</u>	<u>8.6</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1000</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1000</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-1620-FD02-20190114

**GROUNDWATER SAMPLING RECORD**

Project Number: 9119232 Project Name: UPRR HWPW Date: 2-10-2020  
 Sample Number: WG-1620 MW22AR 20200210 Starting Water Level (ft. BMP): 1.55  
 Sampling Location (well ID, etc.): MW-22AR Casing Stickup (ft.): 0  
 Sampled by: TIM McSpadden Starting Water Level (ft. BGL):  
 Measuring Point (MP) of Well: Total Depth (ft. BMP): 19.80  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment:  
 Purging: Sampling:  
 Disposal of Discharged Water:

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: H. Oil Heron 017720 Thermometer: C1113  
 pH Meter: Hanna C1113 Field Calibration:  
 Conductivity Meter: C1113 Field Calibration:  
 Filter / Filter Size: 10 micron Other: Geo Pump P1119

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L / m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
9:00	1.55	.2	.2	22.33	6.50	1.20	16.40	-33	0.0	LT Brown
9:10	2.36	.3	.2	22.49	6.56	1.24	1.38	-72	910	
9:15	3.01	.4	.2	22.53	6.57	1.23	0.11	-71	852	
9:20	3.34	.5	.2	22.56	6.60	1.09	0.00	-60	818	
9:25	3.47	.6	.2	22.67	6.61	1.04	0.00	-52	648	
9:30	3.53	.7	.2	22.72	6.62	1.03	0.00	-50	611	
9:35	3.67	.8	.2	22.80	6.63	1.01	0.00	-44	502	
9:40	3.72	.9	.2	22.84	6.63	0.993	0.00	-40	219	

Water Level (ft. BMP) at End of Purge: 3.74 Sample Intake Depth (ft. BMP): 17.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
10:00			6			

Comments:  
WG-1620 MW22AR 20200210 10:00  
WG-1620 FB17 20200210  
Trip Blank 10



**Golder Associates Inc.**  
 11231 Richmond Avenue, Suite D104  
 Houston, TX 77082  
 (832) 916-3690



# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: \_\_\_\_\_ Date: 2-10-2020

Sample Number: WG-1620 MW 22 BR 20200210 Starting Water Level (ft. BMP): 1.44

Sampling Location (well ID, etc.): MW-22BR Casing Stickup (ft.): 0

Sampled by: TIM McSpadden Starting Water Level (ft. BGL): 1.44

Measuring Point (MP) of Well: \_\_\_\_\_ Total Depth (ft. BMP): 37.80

Screened Interval (ft. BGL): \_\_\_\_\_ Casing Diameter (In ID): 2"

Filter Pack Interval (ft. BGL): \_\_\_\_\_ Casing Volume (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

METHODS (describe): Low Flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: H. O17 Heron 017720 Thermometer: C-1113

pH Meter: HoriBa C1113 Field Calibration: \_\_\_\_\_

Conductivity Meter: C-1113 Field Calibration: \_\_\_\_\_

Filter / Filter Size: 10 Micron Other: Geo pump P1119

## SAMPLING MEASUREMENTS

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
10:20	1.44	.2	.2	23.08	6.48	1.24	3.07	-50	135	
10:30	1.72	.3	.2	23.12	6.44	1.25	1.23	-61	124	
10:35	1.68	.4	.2	23.01	6.33	1.34	0.00	-53	21.0	
10:40	1.66	.5	.2	23.00	6.32	1.25	0.00	-52	20.2	
10:45	1.62	.6	.2	23.02	6.32	1.35	0.00	-50	13.3	
10:50	1.66	.7	.2	23.04	6.33	1.36	0.00	-53	11.0	
10:55	1.67	.8	.2	23.01	6.33	1.35	0.00	-57	10.9	
11:00	1.65	.9	.2	23.03	6.33	1.35	0.00	-58	13.2	

Water Level (ft. BMP) at End of Purge: 1.65 Sample Intake Depth (ft. BMP): 34.0

## SAMPLE INVENTORY

Bottles Collected				Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
Time	Volume	Composition (G, P)	No.			
11:00		G	6			

Comments:



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**GROUNDWATER SAMPLING RECORD**

PAGE 1 of     

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW-23C Starting Water Level (ft. BMP): 26.59  
 Sample Number: WG-1620 MW23C 20200109 Casing Stickup (ft.):       
 Sampled by: JIM MASPADDA WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TWC - Steel or FVC TD (ft. BMP): 76.80 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID):       
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge into drum 1-7-2020

**INSTRUMENTS** (Indicate make, model, I.d.)  
 Water Level: D2111 Merom Other: Geo pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: 0-4.08 NTU = 0 COND 445  
 Filter / Filter Size: 10-micron Alfa

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L/m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:59	26.59	.2	.2	23.11	0.82	.475	9.92	-41	64.7	Clear/No
9:10	26.71	.3	.2	23.08	0.72	.473	9.95	-39	69.8	
9:15	26.72	.4	.2	23.08	.45	.472	9.95	-39	58.8	
9:20	26.74	.5	.2	23.16	0	.469	10.02	-42	50.3	
9:25	26.75	.6	.2	23.22	0	.469	10.02	-42	47.2	
9:30	26.75	.7	.2	23.24	0	.469	10.03	-44	47.1	
9:35	26.77	.8	.2	23.24	0	.469	10.04	-47	47.4	

WL (ft. BMP) at End of Purge: 26.77 Sample Intake Depth (ft. BMP): 70.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (P)	No.			
9:40	2.0		6			

Comments: No tubing in well

 **GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.01</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.01</u> ft. BGL
MW ID	<u>MW25A</u>	Location	Other _____	Total MW Depth	<u>28.80</u> ft. BGL
Sample ID	<u>WG-162D-MW25A-20190115</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1049				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1059	<u>.2</u>	<u>/</u>	<u>8.26</u>	<u>22.2</u>	<u>6.91</u>	<u>1440</u>	<u>0.79</u>	<u>-42</u>	<u>26</u>
1104	<u>↓</u>	<u>/</u>	<u>8.31</u>	<u>22.3</u>	<u>6.96</u>	<u>1420</u>	<u>0.64</u>	<u>-36</u>	<u>20</u>
1109	<u>↓</u>	<u>/</u>	<u>8.32</u>	<u>22.4</u>	<u>6.84</u>	<u>1410</u>	<u>0.67</u>	<u>-37</u>	<u>21</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1120</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1120</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>1/2</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.72</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>16.72</u>	ft. BGL
MW ID	<u>MW25C</u>	Location	Other _____	Total MW Depth	<u>74.05</u>	ft. BGL
Sample ID	<u>WG-162D-MW25C-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1519				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1529	<u>.2</u>	<u>/</u>	<u>16.93</u>	<u>23.4</u>	<u>6.77</u>	<u>1110</u>	<u>0.91</u>	<u>-42</u>	<u>6.7</u>
1534	<u>↓</u>	<u>/</u>	<u>16.96</u>	<u>23.1</u>	<u>6.73</u>	<u>1120</u>	<u>0.71</u>	<u>-47</u>	<u>4.8</u>
1539	<u>↓</u>	<u>/</u>	<u>16.97</u>	<u>23.1</u>	<u>6.72</u>	<u>1130</u>	<u>0.71</u>	<u>-46</u>	<u>5.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1550	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1550	<u>40ML / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-16-2020  
 Sampling Location (well ID, etc.): MW 26A Starting Water Level (ft. BMP): 3.91  
 Sample Number: WG 1620 MW 26A 20200116 Casing Stickup (ft.): 0  
 Sampled by: T M McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 19.90 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P 2111 Heron Other: Geo pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: Ph. 4.0  
 Filter / Filter Size: 10 Micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
15:15	3.91	.2	.2	22.38	4.76	1.47	6.42	-87	0	Brown/nd
15:25	4.05	.3	.2	22.43	1.20	1.63	6.40	-87	0	prilled
15:30	4.04	.4	.2	22.13	0	2.38	6.08	-71	706	feeding
15:35	4.00	.5	.2	22.04	0	2.40	6.06	-71	1000	up out
15:40	4.02	.6	.2	22.32	0	1.25	6.25	-72	226	df well
15:45	4.04	.7	.2	22.29	0	1.23	6.25	-72	345	to check
15:50	4.09	.8	.2	22.27	0	1.23	6.25	-72	308	the depth
15:55	4.12	.9	.2	22.22	0	1.23	6.24	-71	263	

WL (ft. BMP) at End of Purge: 4.12 Sample Intake Depth (ft. BMP): 16.0

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
16:00			6			

Comments: \_\_\_\_\_



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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.22</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.22</u>	ft. BGL
MW ID	<u>MW27A</u>	Location	Other _____	Total MW Depth	<u>29.85</u>	ft. BGL
Sample ID	<u>WG-162D-MW27A-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1607				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1617	<u>.2</u>	<u>/</u>	<u>7.43</u>	<u>21.6</u>	<u>6.77</u>	<u>1060</u>	<u>1.14</u>	<u>-36</u>	<u>5.2</u>
1622	<u>↓</u>	<u>/</u>	<u>7.44</u>	<u>21.7</u>	<u>6.75</u>	<u>1080</u>	<u>0.92</u>	<u>-31</u>	<u>4.4</u>
1627	<u>↓</u>	<u>/</u>	<u>7.43</u>	<u>21.9</u>	<u>6.74</u>	<u>1090</u>	<u>0.93</u>	<u>-32</u>	<u>4.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1640	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1640	<u>40ML / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.93</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>16.93</u>	ft. BGL
MW ID	<u>MW27C</u>	Location	Other _____	Total MW Depth	<u>73.60</u>	ft. BGL
Sample ID	<u>WG-162D-MW27C-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Watterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1649				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1659	<u>.2</u>	<u>/</u>	<u>17.26</u>	<u>22.7</u>	<u>6.96</u>	<u>1140</u>	<u>1.09</u>	<u>-51</u>	<u>6.7</u>
1704	<u>↓</u>	<u>/</u>	<u>17.29</u>	<u>22.4</u>	<u>6.92</u>	<u>1120</u>	<u>0.79</u>	<u>-46</u>	<u>5.3</u>
1709	<u>↓</u>	<u>/</u>	<u>17.29</u>	<u>22.4</u>	<u>6.91</u>	<u>1120</u>	<u>0.74</u>	<u>-47</u>	<u>5.1</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1720</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1720</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-16-2020  
 Sampling Location (well ID, etc.): MW-28A Starting Water Level (ft. BMP): 5.01  
 Sample Number: WG 1620 MW 28A 20200116 Casing Stickup (ft.): 0  
 Sampled by: TJM McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 25.60 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge thru drum 1-7-2020

**INSTRUMENTS** (Indicate make, model, I.d.)

Water Level: P 2111 Heron Other: Geo pump P1102  
 Multi Meter: Homba C 1103  
 Field Calibration: PH-4.0  
 Filter / Filter Size: 0

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:30	5.01	.2	.2	22.97	3.01	1.47	5.29	-23	160	clear/no
8:40	5.09	.3	.2	22.37	3.89	1.47	5.61	-46	99.1	
8:45	5.12	.4	.2	22.41	5.22	1.45	5.77	-46	50.1	
8:50	5.17	.5	.2	22.58	5.47	1.45	5.83	-44	34.8	
8:55	5.22	.6	.2	23.25	5.4	1.43	5.97	-40	13.2	
9:00	5.26	.7	.2	23.43	5.63	1.41	6.14	-38	1.5	
9:05	5.30	.8	.2	23.38	5.57	1.41	6.17	-38	2.9	
9:10	5.32	.9	.2	23.43	5.77	1.41	6.18	-38	3.5	

WL (ft. BMP) at End of Purge: 5.32 Sample Intake Depth (ft. BMP): 23.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y / <input checked="" type="radio"/> N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (C/P)	No.			
9:10			6			

Comments:



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MS MSD

**GROUNDWATER SAMPLING RECORD** PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-16-2020  
 Sampling Location (well ID, etc.): TIM McSpadden Starting Water Level (ft. BMP): 15.45  
 Sample Number: mw. 28C Casing Stickup (ft.):  
 Sampled by: WG-1620 MW28C 20200116 WL (ft. BMP): (ft. BGL):  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 87.30 (ft. BGL):  
 Screened Interval (ft. BGL): Ft. water: Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): 1X Casing Vol (gal.): 3X (gal):

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge 1/2 drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: PH-4.0  
 Filter / Filter Size:

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
9:35	15.45	.2	.2	23.92	1.32	0.186	7.58	-159	15.3	Clear/no
9:45	16.61	.3	.2	23.02	0.92	0.180	7.73	-176	14.9	
9:50	17.12	.4	.2	23.87	0.21	0.168	8.04	-191	10.7	
9:55	17.22	.5	.2	23.96	0.0	0.163	8.23	-200	6.2	
10:00	17.57	.6	.2	23.76	0.0	0.164	8.38	-206	0.5	
10:05	17.73	.7	.2	23.72	0.0	0.164	8.46	-206	0	
10:10	17.91	.8	.2	23.71	0.0	0.163	8.56	-208	0	

WL (ft. BMP) at End of Purge: Sample Intake Depth (ft. BMP):

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
10:15			6			

Comments:  
 WG-1620 MW28C MS 20200116  
 WG-1620 MW28C MSD 20200116

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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19110232 Project Name: VPRR HW Pan Date: 1-20-2020  
 Sampling Location (well ID, etc.): MW-32AR Starting Water Level (ft. BMP): 3.86  
 Sample Number: WG 1620 MW 32AR 20200120 Casing Stickup (ft.): 0  
 Sampled by: Jim McSpalden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 18.11 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): low flow Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge Hydro Drum 1-20-20

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Horva Other: Geo pump P1102  
 Multi Meter: Hanba C1103  
 Field Calibration: PK-40  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:30	3.86	.2	.2	17.52	11.89	.946	6.88	3	240	Clear/no
8:40	4.10	.3	.2	18.14	3.14	.944	6.85	-4	2.55	
8:45	4.06	.4	.2	18.65	.85	.962	6.74	-47	496	
8:50	4.04	.5	.2	18.35	.19	.977	6.63	-60	606	
8:55	4.04	.6	.2	18.24	.10	.982	6.62	-61	600	
9:00	4.04	.7	.2	18.29	0	.984	6.52	-73	453	
9:05	4.05	.8	.2	18.49	0	1.02	6.51	-77	346	
9:10	4:03	.9	.2							

WL (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): 15

**SAMPLE INVENTORY** 4.03

Bottles Collected					Filtration	Preservation	Remarks
Time	Volume	Composition	G P	No.	(Y) N	(type)	(quality control sample, other)
9:10				6			

Comments: \_\_\_\_\_


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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-20-2020  
 Sampling Location (well ID, etc.): MW = 32B Starting Water Level (ft. BMP): 4.50  
 Sample Number: WA 1620 MW 32B 20200120 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpedde WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - (Steel) or PVC TD (ft. BMP): 36.45 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe) Low Flow - Dedicated Tubing/Equipment

Cleaning Equipment: DI/Alconox Rinse

Purge: Peristaltic Pump / SS Pump / Bailer / Bladder

Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder

Disposal of Discharged Water: Purge Hydro drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P2111 Horon

Other: Geo Pump P1102

Multi Meter: Humbler C1103

Field Calibration: Ph. 4.0

Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
9:35	4.50	.2	.2	16.84	4.09	.697	6.79	-23	121	clear/no
9:45	6.55	.3	.2	16.69	1.49	.701	6.78	-30	115	
9:50	6.79	.4	.2	16.94	.08	.706	6.72	-34	105	
9:55	7.01	.5	.2	16.89	0	.708	6.77	-34	101	
10:00	7.11	.6	.2	17.15	0	.689	6.77	-39	101	
10:05	7.14	.7	.2	16.99	0	.690	6.77	-39	102	
10:10	7.17	.8	.2	16.92	0	.691	6.76	-38	102	
10:15	7.20	.9	.2	15.88	0	.684	6.76	-38	936	

WL (ft. BMP) at End of Purge: 7.20

Sample Intake Depth (ft. BMP): 31

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
10:15			6			

Comments:



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Duplicate

<b>GROUNDWATER SAMPLING RECORD</b>		PAGE ___ of ___
Project Number: 19119232	Project Name: UPRR MWPN	Date: 1-20-2020
Sampling Location (well ID, etc.): MW-33A	Starting Water Level (ft. BMP): 4.41	
Sample Number: WG 1620 MW 33A 20200120	Casing Stickup (ft.): 0	
Sampled by: Tim McSpadden	WL (ft. BMP): (ft. BGL):	
Measuring Point (MP) of Well: TOC - Steel or PVC	TD (ft. BMP): 25.20 (ft. BGL):	
Screened Interval (ft. BGL):	Ft. water: Casing Dia. (In ID): 2 1/2	
Filter Pack Interval (ft. BGL):	1X Casing Vol (gal.): 3X (gal.):	

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: Q1/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge to drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: Ph. 4.0-  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:00	4.41	.2	.2	20.87	1.90	1.36	6.65	-21	140	clear/m
12:10	4.58	.3	.2	21.27	.35	1.36	6.68	-31	140	
12:15	4.62	.4	.2	21.65	0	1.37	6.64	-44	79.2	
12:20	4.61	.5	.2	22.11	0	1.39	6.50	-54	80.4	
12:25	4.64	.6	.2	22.21	0	1.37	6.49	-56	56.2	
12:30	4.62	.7	.2	22.56	0	1.38	6.47	-59	46.9	
12:35	4.58	.8	.2	22.57	0	1.38	6.47	-61	48.7	
12:40	4.56	.9	.2	22.92	0	1.39	6.47	-61	14.8	

WL (ft. BMP) at End of Purge: 4.56 Sample Intake Depth (ft. BMP): 22

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
12:45			12			

Comments:  
 Duplicate  
 WG-1620 FDO5 20200120

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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.06</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.06</u>	ft. BGL
MW ID	<u>MW33BR</u>	Location	Other _____	Total MW Depth	<u>38.05</u>	ft. BGL
Sample ID	<u>WG-162D-MW33BR-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1424</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1434</u>	<u>.2</u>		<u>5.22</u>	<u>21.4</u>	<u>6.91</u>	<u>1560</u>	<u>0.74</u>	<u>-31</u>	<u>6.7</u>
<u>1439</u>	<u>↓</u>		<u>5.26</u>	<u>21.4</u>	<u>6.92</u>	<u>1520</u>	<u>0.56</u>	<u>-26</u>	<u>4.7</u>
<u>1444</u>	<u>↓</u>		<u>5.27</u>	<u>21.6</u>	<u>6.92</u>	<u>1530</u>	<u>0.59</u>	<u>-27</u>	<u>5.1</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1500</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1500</u>	<u>40ML / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_\_ of \_\_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-10-2020  
 Sampling Location (well ID, etc.): 3 MW 35A Starting Water Level (ft. BMP): 4.42  
 Sample Number: WG 1620 MW 35A 20200110 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 28.25 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment

Cleaning Equipment: DI/Alconox Rinse

Purge: Peristaltic Pump / SS Pump / Bailer / Bladder

Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder

Disposal of Discharged Water: Purge to drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.D.)

Water Level: P2111 Heron

Other: Geo Pump P1102

Multi Meter: Hanna C1103

Field Calibration: pH-7.99. NTU-0 cond 4.49

Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
10:40	4.42	.2	.2	23.54	1.99	2.17	6.41	-33	47.1	
10:50	4.91	.3	.2	23.63	1.84	2.32	6.25	-39	34.9	
10:55	4.89	.4	.2	23.65	1.22	2.33	6.23	-42	35.6	
11:00	4.98	.5	.2	23.58	0	2.34	6.20	-48	27.2	
11:05	4.90	.6	.2	23.52	0	2.34	6.19	-51	24.4	
11:10	4.92	.7	.2	23.52	0	2.33	6.19	-54	22.9	
11:15	4.92	.8	.2	23.52	0	2.29	6.20	-55	22.5	

WL (ft. BMP) at End of Purge: 4.0 4.92

Sample Intake Depth (ft. BMP): 2.5

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y) N	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G) P	No.			
11:20			6			

Comments:



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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPLW Date: 1-10-2020  
 Sampling Location (well ID, etc.): MW-35B Starting Water Level (ft. BMP): 4.65  
 Sample Number: WG-1620 MW35B 2020 0110 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 42.95 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P-2111 Heron Other: Geo Pump P1102  
 Multi Meter: Hanna C1103  
 Field Calibration: PH-3.99 NTU-0 COND 4.10  
 Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:00	4.65	.2	.2	23.17	.92	4.64	6.57	-88	0.0	Brown / yes
12:10	5.31	.3	.2	23.17	0	1.71	6.50	-64	0.0	clear/no
12:15	5.49	.4	.2	23.05	0	1.16	6.55	-79	247	clear/no
12:20	5.58	.5	.2	23.25	0	1.33	6.38	-84	16.2	
12:25	5.59	.6	.2	23.29	0	1.36	6.38	-85	15.3	
12:30	5.60	.7	.2	23.37	0	1.40	6.38	-85	11.9	
12:35	5.61	.8	.2	23.41	0	1.40	6.38	-85	11.2	

WL (ft. BMP) at End of Purge: 5.61 Sample Intake Depth (ft. BMP): 33

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
12:45			6			

Comments:  
Duplicate  
WG-1620 F001 2020 0110

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 Round Rock, Texas 78664  
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Stop work 12:30

# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW 36A Starting Water Level (ft. BMP): 8.83  
 Sample Number: WG 1620 MW 36A 20200109 Casing Stickup (ft.): 0  
 Sampled by: TIM McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 27.35 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2 1/2  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): low flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H2O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Heron Other: Geo pump P1102  
 Multi Meter: Homba C1103  
 Field Calibration: PK-4-05 NTU-0 comp 445  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
15:15	8.83	.2	.2	23.60	11.0	.963	7.03	12	0	Clear/nt
15:25	9.71	.3	.2	23.61	14.85	.965	7.01	11	0	
15:30	10.82	.4	.2	23.82	1.41	.821	6.97	42	183	
15:35	1.08	.5	.2	23.84	2.20	.799	6.97	45	139	
15:40	1.22	.6	.2	23.89	2.36	.791	6.97	49	103	
15:45	1.37	.7	.2	23.88	2.28	.790	6.97	50	90.3	
15:50	1.39	.8	.2	23.84	2.23	.791	6.97	50	88.2	
15:55	1.42	.9	.2	23.86	2.18	.790	6.97	51	89.1	

WL (ft. BMP) at End of Purge: 11.42 Sample Intake Depth (ft. BMP): 24.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (GP)	No.			
16:00	2.5		6			

Comments: \_\_\_\_\_



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# GROUNDWATER SAMPLING RECORD

PAGE 1 of 1

Project Number: 19119232 Project Name: URR HWPH Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW-36B Starting Water Level (ft. BMP): .39  
 Sample Number: WG-1620 MW36B-20200109 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 42.85 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: 101/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.D.)

Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Humba C1103  
 Field Calibration: Ph 4.06 ATU-0 COMB 445  
 Filter / Filter Size: 10 micron Filter

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L/m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
14:00	0.39	.2	.2	22.61	2.82	.153	7.91	41	13.4	
14:10	1.51	.3	.2	22.50	2.14	.152	7.84	47	8.2	
14:15	2.21	.4	.2	22.40	2.02	.161	7.76	53	8.1	
14:20	2.68	.5	.2	22.55	1.64	.149	7.51	64	8.1	
14:25	2.71	.6	.2	22.56	1.56	.149	7.45	67	7.6	
14:30	2.86	.7	.2	22.79	1.49	.149	7.34	83	8.3	
14:35	3.27	.8	.2	22.93	1.46	.149	7.32	73	8.9	

WL (ft. BMP) at End of Purge: 3.27

Sample Intake Depth (ft. BMP): 38

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y) N	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (GP)	No.			
14:40	2.0		6			

Comments:



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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>82.06</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>82.06</u>	ft. BGL
MW ID	<u>MW36D</u>	Location	Other _____	Total MW Depth	<u>110.10</u>	ft. BGL
Sample ID	<u>WG-162D-MW36D-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C) <input type="checkbox"/> NM	pH <input type="checkbox"/> NM	Conductivity (µS/m or mS/cm) <input type="checkbox"/> not measured	Dissolved Oxygen (mg/L) <input type="checkbox"/> not measured	Redox Potential (mV) <input type="checkbox"/> not measured	Turbidity (NTU) <input type="checkbox"/> NM
1652									
1702	<u>.2</u>	<u>/</u>	<u>82.29</u>	<u>21.9</u>	<u>6.71</u>	<u>1360</u>	<u>1.39</u>	<u>-73</u>	<u>6.7</u>
1707	<u>↓</u>	<u>/</u>	<u>82.31</u>	<u>21.4</u>	<u>6.72</u>	<u>1320</u>	<u>1.09</u>	<u>-70</u>	<u>7.2</u>
1712	<u>↓</u>	<u>/</u>	<u>82.32</u>	<u>21.3</u>	<u>6.74</u>	<u>1330</u>	<u>1.12</u>	<u>-71</u>	<u>7.1</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1725</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1725</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HWPLW Date: 1-21-2020

Sample Number: MW 38A Starting Water Level (ft. BMP): 1.51

Sampling Location (well ID, etc.): WG-1620 MW 38A 20200121 Casing Stickup (ft.): 0

Sampled by: TIM McSpadden Starting Water Level (ft. BGL):

Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 37.60

Screened Interval (ft. BGL): Casing Diameter (In ID): 2"

Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe): Low Flow

Cleaning Equipment:

Purging: Sampling:

Disposal of Discharged Water:

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: P2111 Heron Thermometer: " "

pH Meter: Hanna C1103 Field Calibration: Ph. 4.0

Conductivity Meter: " " Field Calibration:

Filter / Filter Size: Other:

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L / m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
10:30	1.51	.2	.2	19.08	7.34	.507	5.05	8	42.2	clay/no
10:40	1.65	.3	.2	19.42	7.34	.499	2.87	18	42.0	
10:45	1.69	.4	.2	18.27	7.18	.606	.88	32	8.9	
10:50	1.68	.5	.2	18.17	7.02	.769	.31	26	8.9	
10:55	1.66	.6	.2	18.52	6.89	.805	.11	14	8.2	
11:00	1.67	.7	.2	18.27	6.88	.803	.02	14	8.1	
11:05	1.69	.8	.2	18.21	6.90	.805	0	12	7.9	
11:10										

Water Level (ft. BMP) at End of Purge: Sample Intake Depth (ft. BMP):

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
11:10			6			

Comments:



**Golder Associates Inc.**  
 11231 Richmond Avenue, Suite D104  
 Houston, TX 77082  
 (832) 916-3690



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119252 Project Name: UPRR HWPW Date: 1-21-2026  
 Sample Number: MW 38B Starting Water Level (ft. BMP): 3.79  
 Sampling Location (well ID, etc.): WG-1620MW38B20240121 Casing Stickup (ft.): 3'  
 Sampled by: TIM MCSadden Starting Water Level (ft. BGL):       
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 35.70 soft bottom  
 Screened Interval (ft. BGL):      Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL):      Casing Volume (gal.):     

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment:       
 Purging:      Sampling:       
 Disposal of Discharged Water:     

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: P2111 Horaw Thermometer:       
 pH Meter: Hanna CI103 Field Calibration: Ph. 4.0  
 Conductivity Meter:      Field Calibration:       
 Filter / Filter Size: 10 micron Other:     

**SAMPLING MEASUREMENTS**


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
12:00	3.79	.2	.2	18.75	7.04	1.33	3.20	-60	254	
12:10	4.08	.3	.2	19.28	6.59	1.32	.69	-53	264	
12:15	3.98	.4	.2	19.18	6.33	1.29	0	-44	193	
12:20	3.99	.5	.2	18.69	6.28	1.29	0	-45	149	
12:25	3.99	.6	.2	18.86	6.26	1.31	0.39	-45	145	
12:30	3.97	.7	.2	18.79	6.24	1.30	1.20	-45	116	
12:35	3.98	.8	.2	18.68	6.23	1.30	.34	-46	88.9	
13:40	3.97	.9	.2	18.59	6.22	1.30	.45	-46	73.8	

Water Level (ft. BMP) at End of Purge: 3.97 Sample Intake Depth (ft. BMP): 31

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
13:45			6			

Comments:     



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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.02</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.02</u>	ft. BGL
MW ID	<u>MW39B</u>	Location	Other _____	Total MW Depth	<u>41.25</u>	ft. BGL
Sample ID	<u>WG-1620-MW39B-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1703				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1713	<u>.2</u>	<u>/</u>	<u>6.31</u>	<u>21.8</u>	<u>6.81</u>	<u>1260</u>	<u>0.79</u>	<u>-91</u>	<u>8.2</u>
1718	<u>↓</u>	<u>/</u>	<u>6.34</u>	<u>21.7</u>	<u>6.82</u>	<u>1230</u>	<u>0.64</u>	<u>-80</u>	<u>7.7</u>
1723	<u>↓</u>	<u>/</u>	<u>6.35</u>	<u>21.7</u>	<u>6.82</u>	<u>1240</u>	<u>0.67</u>	<u>-81</u>	<u>7.8</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1735	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1735	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.17</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.17</u> ft. BGL
MW ID	<u>MW40B</u>	Location	Other _____	Total MW Depth	<u>42.55</u> ft. BGL
Sample ID	<u>WG-1620-MW40B-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1131				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1141	<u>.3</u>	<u>/</u>	<u>6.29</u>	<u>22.9</u>	<u>6.91</u>	<u>1620</u>	<u>0.74</u>	<u>-72</u>	<u>4.7</u>
1146	<u>↓</u>	<u>/</u>	<u>6.36</u>	<u>22.4</u>	<u>6.85</u>	<u>1620</u>	<u>0.61</u>	<u>-70</u>	<u>5.8</u>
1151	<u>↓</u>	<u>/</u>	<u>6.35</u>	<u>22.7</u>	<u>6.86</u>	<u>1610</u>	<u>0.63</u>	<u>-71</u>	<u>5.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1210</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1210</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-23-2020

Sample Number: W6 1620 MW 41B 20200123 Starting Water Level (ft. BMP): 4.20

Sampling Location (well ID, etc.): MW 41B Casing Stickup (ft.): 41

Sampled by: TIM MCSpeckle Starting Water Level (ft. BGL): \_\_\_\_\_

Measuring Point (MP) of Well: \_\_\_\_\_ Total Depth (ft. BMP): 42.63

Screened Interval (ft. BGL): \_\_\_\_\_ Casing Diameter (In ID): 6.2" PVC

Filter Pack Interval (ft. BGL): \_\_\_\_\_ Casing Volume (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

METHODS (describe): Low Flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P2111 Horon Hi Water Thermometer: v 11

pH Meter: Hanna CI103 Field Calibration: PH 4.0 Ph 7.0

Conductivity Meter: v 4 Field Calibration: \_\_\_\_\_

Filter / Filter Size: \_\_\_\_\_ Other: Geo Amy 1102

## SAMPLING MEASUREMENTS


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
12:00	4.20	.2	.2	18.29	6.96	1.28	2.16	-130	13.4	Clear
12:10	4.40	.3	.2	18.23	6.90	1.28	1.16	-136	10.2	
12:15	4.44	.4	.2	18.19	6.87	1.28	.75	-135	6.4	
12:20	4.46	.5	.2	18.23	6.85	1.28	1.41	+132	4.6	
12:25	4.48	.6	.2	18.25	6.80	1.28	0.11	-124	8.2	
12:30	4.52	.7	.2	18.27	6.80	1.28	0	-124	9.4	
12:35	4.55	.8	.2	18.29	6.80	1.28	0	-124	8.2	

Water Level (ft. BMP) at End of Purge: 4.55 Sample Intake Depth (ft. BMP): 35.6

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (GP)	No.			
12:40			6			

Comments: \_\_\_\_\_



**Golder Associates Inc.**  
 11231 Richmond Avenue, Suite D104  
 Houston, TX 77082  
 (832) 916-3690



# Groundwater Sample Collection

Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.97</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.97</u> ft. BGL
MW ID	<u>MW42B</u>	Location	Other _____	Total MW Depth	<u>43.80</u> ft. BGL
Sample ID	<u>WG-162D-MW42B-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1232</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1242</u>	<u>.2</u>	<u>/</u>	<u>7.19</u>	<u>21.9</u>	<u>6.77</u>	<u>1450</u>	<u>0.96</u>	<u>-41</u>	<u>7.2</u>
<u>1247</u>	<u>↓</u>	<u>/</u>	<u>7.22</u>	<u>21.6</u>	<u>6.75</u>	<u>1470</u>	<u>0.81</u>	<u>-36</u>	<u>6.2</u>
<u>1252</u>	<u>↓</u>	<u>/</u>	<u>7.21</u>	<u>21.7</u>	<u>6.74</u>	<u>1470</u>	<u>0.84</u>	<u>-39</u>	<u>6.9</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1310</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1310</u>	<u>40mL / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: HWPW UPRR Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW 44A Starting Water Level (ft. BMP): 10.18  
 Sample Number: WG 1620 MW44A 20200109 Casing Stickup (ft.): 0  
 Sampled by: TIM MCSPODDEN WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or (PVC) TD (ft. BMP): 28.05 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: D2111 Herou Other: Geo Pump P-1102  
 Multi Meter: Humbro C1103  
 Field Calibration: PH-4.08 NTU-0 Cond-445  
 Filter / Filter Size: 10 micron Polte

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:00	10.18	.2	.2	23.38	0.88	.814	7.21	-95	22.2	
13:10	10.34	.3	.2	23.35	0.63	.813	7.19	-95	18.7	
13:15	10.38	.4	.2	23.28	0.16	.806	7.09	-93	15.5	
13:20	10.39	.5	.2	23.25	0.41	.795	7.02	-91	14.4	
13:25	10.41	.6	.2	23.26	0.41	.795	7.03	-91	14.0	

WL (ft. BMP) at End of Purge: 10.41 Sample Intake Depth (ft. BMP): 25

**SAMPLE INVENTORY**

Time	Bottles Collected			No.	Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (Y/P)					
13:30	2.0			6	Y		

Comments:



**GOLDER**

2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.81</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>16.81</u>	ft. BGL
MW ID	<u>MW44C</u>	Location	Other _____	Total MW Depth	<u>47.80</u>	ft. BGL
Sample ID	<u>WG-162D-MW44C-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1241</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1251</u>	<u>2</u>	<u>/</u>	<u>17.06</u>	<u>21.6</u>	<u>6.91</u>	<u>1140</u>	<u>0.91</u>	<u>-41</u>	<u>6.2</u>
<u>1256</u>	<u>↓</u>	<u>/</u>	<u>17.11</u>	<u>21.4</u>	<u>6.82</u>	<u>1120</u>	<u>0.71</u>	<u>-33</u>	<u>5.7</u>
<u>1301</u>	<u>↓</u>	<u>/</u>	<u>17.12</u>	<u>21.7</u>	<u>6.84</u>	<u>1110</u>	<u>0.72</u>	<u>-34</u>	<u>5.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1315</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1315</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HW PW Date: 1-20-2020  
 Sampling Location (well ID, etc.): MW 45C Starting Water Level (ft. BMP): 16.01  
 Sample Number: WG 1620M MW 45C 20200120 Casing Stickup (ft.): \_\_\_\_\_  
 Sampled by: J.M. McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC Steel or PVC TD (ft. BMP): 70.60 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: NI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.D.)  
 Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Hanna C1103  
 Field Calibration: PH-4.0  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
15:50	16.01	.2	.2	21.85	15.19	.245	8.09	-9	52.2	clear/w
16:00	16.71	.3	.2	21.70	15.52	.245	8.02	-3	49.7	
16:05	16.79	.4	.2	21.66	15.35	.253	7.80	1	47.2	
16:10	16.66	.5	.2	21.60	15.22	.257	7.73	2	49.1	
16:15	16.61	.6	.2	21.36	15.15	.246	7.64	6	33.4	
16:20	16.63	.7	.2	20.38	15.00	.234	7.54	11	16.7	
16:25	16.66	.8	.2	20.23	15.05	.233	7.50	2	15.1	
16:30	16.74	.9	.2	20.95	15.00	.227	7.57	-10	39.4	

WL (ft. BMP) at End of Purge: 16.74 Sample Intake Depth (ft. BMP): 60.0

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y) (N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
16:30			6			

Comments: \_\_\_\_\_



**GOLDER**  
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 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>16.34</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>16.34</u>	ft. BGL
MW ID	<u>MW46C</u>	Location	Other _____	Total MW Depth	<u>72.80</u>	ft. BGL
Sample ID	<u>WG-162D-MW46C-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0952				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1002	<u>2</u>	<u>/</u>	<u>16.56</u>	<u>23.4</u>	<u>6.74</u>	<u>1160</u>	<u>1.09</u>	<u>-31</u>	<u>26</u>
1007	<u>↓</u>	<u>/</u>	<u>16.57</u>	<u>23.5</u>	<u>6.72</u>	<u>1170</u>	<u>1.06</u>	<u>-25</u>	<u>18</u>
1012	<u>↓</u>	<u>/</u>	<u>16.57</u>	<u>23.4</u>	<u>6.71</u>	<u>1170</u>	<u>1.07</u>	<u>-26</u>	<u>19</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1025	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1025	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



**Groundwater Sample Collection**



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>17.28</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	_____ ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	_____ ft. BGL
MW ID	<u>MW 47C</u>	Location	Other _____	Total MW Depth	<u>68.85</u> ft. BGL
Sample ID	<u>WG-1620-MW47C-20200116</u>	Water Quality		MW Diameter	_____ inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	_____ gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	_____ ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>10303</u>	<u>1.2</u>	<u>/</u>	<u>17.46</u>	<u>22.7</u>	<u>6.91</u>	<u>1260</u>	<u>1.39</u>	<u>-27</u>	<u>23</u>
<u>1048</u>	<u>↓</u>	<u>/</u>	<u>17.47</u>	<u>22.4</u>	<u>6.96</u>	<u>1240</u>	<u>1.07</u>	<u>-36</u>	<u>12</u>
<u>1053</u>	<u>↓</u>	<u>/</u>	<u>17.47</u>	<u>22.5</u>	<u>6.94</u>	<u>1240</u>	<u>1.08</u>	<u>-36</u>	<u>13</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1105</u>	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1105</u>	<u>40ml / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>17.09</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	_____ ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	_____ ft. BGL
MW ID	<u>MW 48C</u>	Location	Other _____	Total MW Depth	<u>72.50</u> ft. BGL
Sample ID	<u>WG-1620-MW48C-20200110</u>	Water Quality		MW Diameter	_____ inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	_____ gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	_____ ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0947</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0957</u>	<u>1.2</u>	<u>/</u>	<u>17.29</u>	<u>22.8</u>	<u>7.11</u>	<u>980</u>	<u>2.34</u>	<u>-96</u>	<u>16</u>
<u>1002</u>	<u>↓</u>	<u>/</u>	<u>17.31</u>	<u>22.9</u>	<u>7.13</u>	<u>960</u>	<u>1.96</u>	<u>-91</u>	<u>17</u>
<u>1007</u>	<u>↓</u>	<u>/</u>	<u>17.31</u>	<u>22.8</u>	<u>7.11</u>	<u>960</u>	<u>1.95</u>	<u>-92</u>	<u>17</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1020</u>	<u>60ml</u>	<u>P</u>	<u>1</u>	<input checked="" type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1020</u>	<u>40ml / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

**GROUNDWATER SAMPLING RECORD** PAGE 1 of     

Project Number: 19119232 Project Name: VPRR HWPW Date: 1-7-2020  
 Sampling Location (well ID, etc.): MW-49A Starting Water Level (ft. BMP): 11.12  
 Sample Number: WG-1620-MW49A-20200107 Casing Stickup (ft.): 0  
 Sampled by: TIM MCSADDEN WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 30.18 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge to drum.

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: D2111 Horou Other: Portable sampler P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: PH 3.99 +VU-0 cond 4.52  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:30	11.12	.2	.2	21.40	9.64	1.00	6.33	-37	286	Clear/no
12:40	11.41	.3	.2	22.79	6.20	1.00	6.57	-86	224	
12:45	11:44	.4	.2	22.90	6.60	1.01	6.60	-94	188	
12:50	11:48	.5	.2	22.99	5.37	1.01	6.63	-98	165	
12:55	11:52	.6	.2	23.18	4.89	1.00	6.63	-102	142	
13:00	11:54	.7	.2	23.73	0.29	.979	6.65	-100	124	
13:05	11:56	.8	.2	23.91	0.21	.963	6.65	-101	91.4	
13:10	11:59	.9	.2	24.69	0.13	.845	6.67	-99	31.2	

WL (ft. BMP) at End of Purge: 11.59 Sample Intake Depth (ft. BMP): 27.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (C/P)	No.			
13:10	1.5		9			

Comments:



**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-7-2020  
 Sampling Location (well ID, etc.): MW-49B Starting Water Level (ft. BMP): 11.57  
 Sample Number: WG-1620-MW49B-20200107 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 33.05 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum

**INSTRUMENTS** (Indicate make, model, I.D.)

Water Level: D214 Horu Other: Portable Sampler P-1102  
 Multi Meter: Horiba C1103  
 Field Calibration: Ph-3.99 NTU -0 comp 4.52  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:45	11.51	.2	.2	25.64	4.20	.708	6.91	-76	81.2	
13:55	11.76	.3	.2	25.59	4.10	.703	6.91	-76	85.4	
14:00	11.83	.4	.2	24.89	3.62	.714	6.88	-70	69.2	
14:05	11.93	.5	.2	24.90	3.41	.715	6.87	-69	63.2	
14:10	11.85	.6	.2	24.86	3.51	.714	6.86	-68	60.2	
14:15	11.86	.7	.2	24.88	3.06	.716	6.81	-65	74.2	
14:20	11.86	.8	.2	24.90	3.00	.716	6.80	-64	72.7	
14:25	11.86	.9	.2	24.87	2.73	.716	6.76	-61	58.9	

WL (ft. BMP) at End of Purge: 11.86 Sample Intake Depth (ft. BMP): 15.0

**SAMPLE INVENTORY**

Bottles Collected				Filtration	Preservation	Remarks
Time	Volume	Composition (S/P)	No.	(Y/N)	(type)	(quality control sample, other)
14:30	1.5		9			

Comments:



**GOLDER**

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20.09



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>2.09</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.09</u>	ft. BGL
MW ID	<u>MWS0A</u>	Location	Other _____	Total MW Depth	<u>24.80</u>	ft. BGL
Sample ID	<u>WG-1620-MWS0A-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1643</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1648</u>	<u>1.2</u>	<u>/</u>	<u>7.31</u>	<u>21.6</u>	<u>7.11</u>	<u>1160</u>	<u>0.92</u>	<u>-61</u>	<u>2.2</u>
<u>1703</u>	<u>↓</u>	<u>/</u>	<u>7.32</u>	<u>21.5</u>	<u>7.12</u>	<u>1130</u>	<u>0.81</u>	<u>-52</u>	<u>8.6</u>
<u>1708</u>	<u>↓</u>	<u>/</u>	<u>7.32</u>	<u>21.4</u>	<u>7.06</u>	<u>1120</u>	<u>0.74</u>	<u>-51</u>	<u>8.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1715</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1715</u>	<u>40ML/1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.47</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.47</u> ft. BGL
MW ID	<u>MWSIA</u>	Location	Other _____	Total MW Depth	<u>24.90</u> ft. BGL
Sample ID	<u>WG-1620-MWSIA-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>-</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>-</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1116</u>	<u>1.2</u>	<u>/</u>	<u>8.72</u>	<u>20.4</u>	<u>7.12</u>	<u>1630</u>	<u>1.14</u>	<u>-71</u>	<u>9.2</u>
<u>1131</u>	<u>↓</u>	<u>/</u>	<u>8.75</u>	<u>20.6</u>	<u>7.13</u>	<u>1640</u>	<u>0.96</u>	<u>-72</u>	<u>7.1</u>
<u>1136</u>	<u>↓</u>	<u>/</u>	<u>8.76</u>	<u>20.7</u>	<u>7.17</u>	<u>1670</u>	<u>0.93</u>	<u>-72</u>	<u>6.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1150</u>	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1150</u>	<u>40ml / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>20.34</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>20.34</u> ft. BGL
MW ID	<u>MWSIC</u>	Location	Other _____	Total MW Depth	<u>72.80</u> ft. BGL
Sample ID	<u>WG-1620-MWSIC-2019</u>	Water Quality		MW Diameter	<u>20</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1207				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1217	<u>.2</u>	<u>/</u>	<u>20.62</u>	<u>21.3</u>	<u>7.11</u>	<u>1460</u>	<u>0.81</u>	<u>-67</u>	<u>6.1</u>
1223	<u>↓</u>	<u>/</u>	<u>20.61</u>	<u>21.1</u>	<u>7.05</u>	<u>1420</u>	<u>0.51</u>	<u>-60</u>	<u>7.9</u>
1227	<u>↓</u>	<u>/</u>	<u>20.61</u>	<u>21.2</u>	<u>7.06</u>	<u>1430</u>	<u>0.47</u>	<u>-61</u>	<u>7.7</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1240</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1240</u>	<u>40ML/1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL/NONE</u>	<u>VOCS/SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW-53C Starting Water Level (ft. BMP): 16.72  
 Sample Number: WG-1620 MW53C 20200109 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC Steel or PVC TD (ft. BMP): 71.05 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): \_\_\_\_\_  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment

Cleaning Equipment: Q1/Alconox Rinse

Purge: Peristaltic Pump / SS Pump / Bailer / Bladder

Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder

Disposal of Discharged Water: Purge 170 drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: D2111 Heron

Other: Geo Pump P1102

Multi Meter: Hanba C1103

Field Calibration: D-4.08 NTU-0 AND 445

Filter / Filter Size: 10 micron All

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
10:35	16.72	.2	.2	23.64	1.72	.224	9.01	-123	11.1	
10:45	16.99	.3	.2	23.57	0.72	.221	8.84	-129	40.1	
10:50	17.01	.4	.2	23.64	0.17	.208	8.64	-123	24.6	
10:55	17.01	.5	.2	23.75	0	.210	8.24	-104	14.9	
11:00	17.04	.6	.2	23.81	0	.213	8.16	-84	15.5	
11:05	17.06	.7	.2	23.80	0	.220	8.04	-70	17.0	
11:10	17.08	.8	.2	23.82	0	.237	7.92	-57	17.3	
11:15	17.08	.9	.2	23.87	0	.276	7.79	-48	19.9	

WL (ft. BMP) at End of Purge: 17:08

Sample Intake Depth (ft. BMP): 65.

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (C, P)	No.			
11:15	1.5		6			

Comments:



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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPLW Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW 54C Starting Water Level (ft. BMP): 15.59  
 Sample Number: WG 1620 MW 54C 20200109 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - (Steel) or PVC TD (ft. BMP): 72.05 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59  
 METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

## INSTRUMENTS (Indicate make, model, l.d.)

Water Level: 02111 Heron Other: Geo Pump P1102  
 Multi Meter: Hanna C1103  
 Field Calibration: 4.08 NTU-0 COND 445  
 Filter / Filter Size: 10 micro Brite

## SAMPLING MEASUREMENTS


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
11:55	5.59	1.2	1.2	24.62	1.28	1.21	7.23	.95	42.3	Clear/nd
12:05	6.72	1.3	1.2	24.57	1.91	1.21	7.17	.94	41.4	
12:10	6.76	1.4	1.2	23.43	1.13	1.51	6.85	.88	300	
12:15	6.78	1.5	1.2	23.27	0.00	1.52	6.88	-84	155	
12:20	6.81	1.6	1.2	23.28	0.00	1.52	6.86	-82	135	
12:25	6.83	1.7	1.2	23.53	0.00	1.52	6.86	-83	126	
12:30	6.83	1.7	1.2	23.34	0.00	1.52	6.88	-83	125	

WL (ft. BMP) at End of Purge: 16.83 Sample Intake Depth (ft. BMP): 69.

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y) N	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
12:35	1.5		6			

Comments: \_\_\_\_\_


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# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: HWPU Date: 1-8-2020  
 Sampling Location (well ID, etc.): M.W. 57A Starting Water Level (ft. BMP): 14.46  
 Sample Number: FAW-57WG-1620 MW57A20200108 Casing Stickup (ft.): 3'  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 25.90 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2'  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O down 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: D 2111 Herox Other: Geo Pump P1101  
 Multi Meter: Horiba E1103  
 Field Calibration: D-4.04 NTU 0.0 COND 450  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
14:15	14.46	.2	.2	22.49	0.12	.700	6.77	-87	123	
14:25	14.53	.3	.2	22.41	0.01	.777	6.81	-69	68.3	
14:30	14.57	.4	.2	22.48	0.0	.770	6.82	-90	70.4	
14:35	14.66	.5	.2	22.84	0.0	.782	6.82	-96	71.4	
14:40	14.68	.6	.2	22.91	0.0	.783	6.82	-96	72.9	
14:45	14.71	.7	.2	23.11	0.0	.788	6.82	-103	53.1	
14:50	14.73	.8	.2	23.16	0.0	.788	6.81	-103	49.3	

WL (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): 2.2

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
14:55	2.0		6			

Comments:  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_


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# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: HWPW Date: 1-8-2020  
 Sampling Location (well ID, etc.): MW-57B Starting Water Level (ft. BMP): 13.74  
 Sample Number: WG 1620 MW 57B 2020 0108 Casing Stickup (ft.): 3'  
 Sampled by: Tim McSpadden WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 42.85 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID):       
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge to drum 1-7-2020

**INSTRUMENTS** (Indicate make, model, l.d.)

Water Level: D 2111 Horan Other: Geo pump P 1103  
 Multi Meter: Horiba C1103  
 Field Calibration: D-4.04 NTU-0.0 COMP 450  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:45	13.74	.2	.2	23.31	1.86	.996	7.32	-60	0.0	
12:55	14.70	.3	.2	23.15	1.35	.999	7.32	-65	0.0	
13:00	14.74	.4	.2	23.12	1.80	1.00	7.28	-69	628	
13:05	14.74	.5	.2	23.09	.72	1.00	7.28	-68	541	
13:10	14.76	.6	.2	23.00	.61	1.01	7.27	-66	467	
13:20	14.88	.7	.2	23.08	.51	1.01	7.27	-60	212	
13:25	14.95	.8	.2	23.01	.45	1.01	7.26	-58	240	
13:30	14.96	.9	.2	23.03	0.0	1.00	7.25	-57	92.4	

WL (ft. BMP) at End of Purge: 14.96 Sample Intake Depth (ft. BMP): 35

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, Y)	No.			
13:30	2.5		9			

Comments:     



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# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: HWPW Date: 1-8-2020  
 Sampling Location (well ID, etc.): MW-58A Starting Water Level (ft. BMP): 17.71  
 Sample Number: WG 1620 MW58A 20200108 Casing Stickup (ft.): 0  
 Sampled by: TIM McSpodde WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or (PVC) TD (ft. BMP): 28.60 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge 150 drum 1-7-2020

**INSTRUMENTS** (Indicate make, model, I.d.)  
 Water Level: D2111 Heron Other: Geo pump P1102  
 Multi Meter: Hanna C1103  
 Field Calibration: Ph-4.04 NTU-0 COND 450  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
11:30	17.71	.2	.2	25.16	5.06	648	6.86	-98	47.5	
11:40	15.12	.3	.2	24.96	5.51	650	6.84	-99	43.8	
11:45	15.15	.3	.2	24.98	5.13	637	6.82	-97	32.6	
11:50	15.17	.4	.2	24.87	5.74	637	6.80	-95	30.1	
11:55	15.19	.5	.2	24.67	5.75	635	6.81	-97	30.3	
12:00	15.22	.6	.2	24.69	5.71	633	6.80	-96	35.9	
12:05	15.21	.7	.2	24.67	5.61	622	6.80	-94	36.9	
12:10	15.23	.8	.2	24.69	5.57	632	6.80	-95	42.9	

WL (ft. BMP) at End of Purge: 15.23 Sample Intake Depth (ft. BMP): 25.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G.P)	No.			
12:15	2.5		9			

Comments: \_\_\_\_\_



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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.21</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date		Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.21</u>	ft. BGL
MW ID	<u>MW59A</u>	Location	Other _____	Total MW Depth	<u>21.10</u>	ft. BGL
Sample ID	<u>WG-1620-MW59A-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0729				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0739	<u>.2</u>	<u>/</u>	<u>9.47</u>	<u>22.7</u>	<u>6.96</u>	<u>1460</u>	<u>0.67</u>	<u>-71</u>	<u>7.4</u>
0744	<u>↓</u>	<u>/</u>	<u>9.51</u>	<u>22.3</u>	<u>6.92</u>	<u>1420</u>	<u>0.50</u>	<u>-61</u>	<u>7.8</u>
0749	<u>↓</u>	<u>/</u>	<u>9.52</u>	<u>22.4</u>	<u>6.91</u>	<u>1410</u>	<u>0.51</u>	<u>-62</u>	<u>7.4</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0805	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0805	<u>40ml / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.17</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.17</u>	ft. BGL
MW ID	<u>MW59B</u>	Location	Other _____	Total MW Depth	<u>32.90</u>	ft. BGL
Sample ID	<u>WG-1620-MW59B-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0827				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0837	<u>.2</u>	<u>/</u>	<u>9.36</u>	<u>22.3</u>	<u>6.79</u>	<u>1280</u>	<u>0.36</u>	<u>-29</u>	<u>6.2</u>
0842	<u>↓</u>	<u>/</u>	<u>9.37</u>	<u>22.1</u>	<u>6.71</u>	<u>1250</u>	<u>0.28</u>	<u>-25</u>	<u>5.1</u>
0847	<u>↓</u>	<u>/</u>	<u>9.36</u>	<u>22.1</u>	<u>6.72</u>	<u>1260</u>	<u>0.29</u>	<u>-26</u>	<u>5.3</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>0900</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0900</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>83.06</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>83.06</u>	ft. BGL
MW ID	<u>MWS9D</u>	Location	Other _____	Total MW Depth	<u>117.80</u>	ft. BGL
Sample ID	<u>WG-162D-MWS9D-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1316				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1326	<u>.2</u>	<u>/</u>	<u>83.19</u>	<u>21.6</u>	<u>7.06</u>	<u>1140</u>	<u>0.78</u>	<u>-91</u>	<u>4.7</u>
1331	<u>↓</u>	<u>/</u>	<u>83.21</u>	<u>21.3</u>	<u>7.03</u>	<u>1120</u>	<u>0.61</u>	<u>-86</u>	<u>2.6</u>
1336	<u>↓</u>	<u>/</u>	<u>83.22</u>	<u>21.2</u>	<u>7.03</u>	<u>1110</u>	<u>0.62</u>	<u>-86</u>	<u>3.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1350	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1350	<u>40ML / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

WG-162D-PDO3-20190116  
1350-Damplitude

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.87</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	_____ ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	_____ ft. BGL
MW ID	<u>MW 61A</u>	Location	Other _____	Total MW Depth	<u>21.95</u> ft. BGL
Sample ID	<u>WG-1620-MW61A-20200116</u>	Water Quality		MW Diameter	_____ inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	_____ gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	_____ ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1117</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1127</u>	<u>1.2</u>	<u>/</u>	<u>5.39</u>	<u>22.9</u>	<u>6.72</u>	<u>1170</u>	<u>1.42</u>	<u>-37</u>	<u>6.7</u>
<u>1132</u>	<u>↓</u>	<u>/</u>	<u>5.41</u>	<u>22.6</u>	<u>6.69</u>	<u>1130</u>	<u>0.74</u>	<u>-36</u>	<u>7.2</u>
<u>1137</u>	<u>↓</u>	<u>/</u>	<u>5.42</u>	<u>22.7</u>	<u>6.68</u>	<u>1120</u>	<u>0.73</u>	<u>-36</u>	<u>7.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>1200</u>	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1200</u>	<u>40ml / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

ms/msd

Field Team Leader JOHN BRAYTON name

John Brayton signature



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HWPu Date: 1-27-2024  
 Sample Number: WG-1620 m w 623 2020 0127 Starting Water Level (ft. BMP): 3.05  
 Sampling Location (well ID, etc.): MW 623 Casing Stickup (ft.): 3'  
 Sampled by: TIM McSpalden Starting Water Level (ft. BGL):       
 Measuring Point (MP) of Well:      Total Depth (ft. BMP): 35.55  
 Screened Interval (ft. BGL):      Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL):      Casing Volume (gal.):     

**QUALITY ASSURANCE**

METHODS (describe):       
 Cleaning Equipment:       
 Purging:      Sampling:       
 Disposal of Discharged Water:     

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: 01-7720 Golder Thermometer: " "  
 pH Meter: Hanna CI106 Field Calibration: PH-4.0 PH 4.0  
 Conductivity Meter: " " Field Calibration:       
 Filter / Filter Size: 10 micron Filter Other: Geo Pump P 1108

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
				± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
<u>8:45</u>	<u>3.05</u>	<u>.2</u>	<u>.2</u>	<u>19.52</u>	<u>6.45</u>	<u>1.33</u>	<u>7.51</u>	<u>-2</u>	<u>89.1</u>	
<u>8:55</u>	<u>3.41</u>	<u>.3</u>	<u>.2</u>	<u>19.71</u>	<u>6.56</u>	<u>1.34</u>	<u>2.06</u>	<u>-72</u>	<u>79.8</u>	
<u>9:00</u>	<u>3.39</u>	<u>.4</u>	<u>.2</u>	<u>19.60</u>	<u>6.57</u>	<u>1.34</u>	<u>1.21</u>	<u>-100</u>	<u>72.0</u>	
<u>9:05</u>	<u>3.40</u>	<u>.5</u>	<u>.2</u>	<u>19.79</u>	<u>6.52</u>	<u>1.34</u>	<u>0.17</u>	<u>-130</u>	<u>39.7</u>	
<u>9:10</u>	<u>3.37</u>	<u>.6</u>	<u>.2</u>	<u>17.72</u>	<u>6.48</u>	<u>1.34</u>	<u>0.10</u>	<u>-129</u>	<u>34.4</u>	
<u>9:15</u>	<u>3.35</u>	<u>.7</u>	<u>.2</u>	<u>17.65</u>	<u>6.50</u>	<u>1.34</u>	<u>0.14</u>	<u>-131</u>	<u>32.1</u>	
<u>9:20</u>	<u>3.37</u>	<u>.8</u>	<u>.2</u>	<u>17.67</u>	<u>6.49</u>	<u>1.34</u>	<u>0.08</u>	<u>-130</u>	<u>32.2</u>	

Water Level (ft. BMP) at End of Purge: 3.37 Sample Intake Depth (ft. BMP): 3-32

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
<u>9:25</u>			<u>6</u>	<u>Y</u>		

Comments:



Golder Associates Inc.  
 11231 Richmond Avenue, Suite D104  
 Houston, TX 77082  
 (832) 916-3690

**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-16-2020  
 Sampling Location (well ID, etc.): MW 63B Starting Water Level (ft. BMP): 4.90  
 Sample Number: WG-1620 MW 63B 20200116 Casing Stickup (ft.): 0  
 Sampled by: Tim Mespedden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 36.27 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID) 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge Hydro drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Horos Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: Ph-4.0  
 Filter / Filter Size: \_\_\_\_\_

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
14:15	4.90	.2	.2	24.48	3.12	1.79	6.69	-110	2.0	
14:25	5.81	.3	.2	24.60	0.90	1.80	6.61	-112	4.9	
14:30	6.57	.4	.2	24.42	1.63	1.82	6.48	-110	0.4	
14:35	6.91	.5	.2	24.26	3.11	1.81	6.45	-101	0	
14:40	7.15	.6	.2	24.45	2.88	1.80	6.46	-103	0	
14:45	7.29	.7	.2	24.46	2.84	1.78	6.46	-104	0	

WL (ft. BMP) at End of Purge: 7.29 Sample Intake Depth (ft. BMP): 32

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G P)	No.			
14:45			6			

Comments:  
 \_\_\_\_\_  
 \_\_\_\_\_



**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119 232 Project Name: UPPR HWPW Date: 1-27-20  
 Sample Number: W0-1620 MW 64A 20200127 Starting Water Level (ft. BMP): 4.91  
 Sampling Location (well ID, etc.): MW 64A Casing Stickup (ft.): 3'  
 Sampled by: T.M. McSpadden Starting Water Level (ft. BGL):  
 Measuring Point (MP) of Well: Total Depth (ft. BMP): 22.20  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe):  
 Cleaning Equipment:  
 Purging: Sampling:  
 Disposal of Discharged Water:

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: 01-7720 Thermometer: " "  
 pH Meter: Hanna CI106 Field Calibration:  
 Conductivity Meter: " " Field Calibration:  
 Filter / Filter Size: 10 micron Other: Geo Pump P1108

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L /m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
10:35	4.91	.2	.2	17.04	6.49	482	13.16	118	0	
10:45	8.67	.3	.2	8.91	6.71	1.65	3.09	120	0	
10:50	8.72	.4	.2	8.84	6.70	1.66	2.81	122	964	
10:55	8.75	.5	.2	8.76	6.70	1.66	2.76	124	711	
11:00	8.77	.6	.2	8.72	6.69	1.66	2.93	124	734	
11:05	8.79	.7	.2	8.58	6.70	1.65	2.51	125	575	
11:10	8.82	.8	.2	8.46	6.70	1.65	2.01	125	531	

Water Level (ft. BMP) at End of Purge: 8.82 Sample Intake Depth (ft. BMP): 17

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y) (N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G.P)	No.			
11:15			6			

Comments:



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 Houston, TX 77082  
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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>83.29</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>83.29</u>	ft. BGL
MW ID	<u>MW65D</u>	Location	Other _____	Total MW Depth	<u>109.75</u>	ft. BGL
Sample ID	<u>WG-1620-MW65D-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1538				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1548	<u>.2</u>	<u>/</u>	<u>83.46</u>	<u>21.6</u>	<u>6.91</u>	<u>1240</u>	<u>1.13</u>	<u>-91</u>	<u>9.1</u>
1553	<u>↓</u>	<u>/</u>	<u>83.47</u>	<u>21.4</u>	<u>6.92</u>	<u>1260</u>	<u>0.94</u>	<u>-83</u>	<u>5.6</u>
1558	<u>↓</u>	<u>/</u>	<u>83.47</u>	<u>21.4</u>	<u>6.94</u>	<u>1270</u>	<u>0.91</u>	<u>-82</u>	<u>6.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1620</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1620</u>	<u>40ML / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

MS / MSD

Field Team Leader JOHN BRAYTON name

John Brayton signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>81.26</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-16-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>81.26</u>	ft. BGL
MW ID	<u>MW-66D</u>	Location	Other _____	Total MW Depth	<u>103.05</u>	ft. BGL
Sample ID	<u>WG-162D-mw66D-20190116</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1426				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1432	<u>.2</u>	<u>/</u>	<u>81.43</u>	<u>21.6</u>	<u>7.07</u>	<u>1160</u>	<u>1.21</u>	<u>-71</u>	<u>6.7</u>
1441	<u>↓</u>	<u>/</u>	<u>81.44</u>	<u>21.8</u>	<u>7.01</u>	<u>1160</u>	<u>0.92</u>	<u>-65</u>	<u>9.1</u>
1446	<u>↓</u>	<u>/</u>	<u>81.44</u>	<u>21.9</u>	<u>7.02</u>	<u>1170</u>	<u>0.93</u>	<u>-67</u>	<u>8.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1500	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1500	<u>40ML / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.83</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.82</u> ft. BGL
MW ID	<u>MW67B</u>	Location	Other _____	Total MW Depth	<u>39.50</u> ft. BGL
Sample ID	<u>WG-162D-MW67B-20190115</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0843				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0853	<u>.2</u>	<u>/</u>	<u>9.06</u>	<u>21.6</u>	<u>7.16</u>	<u>510</u>	<u>0.71</u>	<u>-96</u>	<u>7.6</u>
0858	<u>↓</u>	<u>/</u>	<u>9.08</u>	<u>21.9</u>	<u>7.12</u>	<u>460</u>	<u>0.54</u>	<u>-91</u>	<u>6.8</u>
0903	<u>↓</u>	<u>/</u>	<u>9.08</u>	<u>22.2</u>	<u>7.11</u>	<u>470</u>	<u>0.56</u>	<u>-92</u>	<u>7.4</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0930	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
0930	<u>40ML / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

MS/MSD



# GROUNDWATER SAMPLING RECORD

Project Number: \_\_\_\_\_ Project Name: \_\_\_\_\_ Date: **1-23-2020**

Sample Number: **W61620 MW 68B 2020 0123** Starting Water Level (ft. BMP): **1.20**

Sampling Location (well ID, etc.): **MW 68B** Casing Stickup (ft.): **0**

Sampled by: **Tim McSpadden** Starting Water Level (ft. BGL): **1.20**

Measuring Point (MP) of Well: \_\_\_\_\_ Total Depth (ft. BMP): **37.40**

Screened Interval (ft. BGL): \_\_\_\_\_ Casing Diameter (In ID): **2"**

Filter Pack Interval (ft. BGL): \_\_\_\_\_ Casing Volume (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

METHODS (describe): **Low Flow**

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: **P2161 Horaw Hi water** Thermometer: **h 1'**

pH Meter: **Homba C1103** Field Calibration: **PH 4.0 PH 4.0**

Conductivity Meter: **h** Field Calibration: \_\_\_\_\_

Filter / Filter Size: \_\_\_\_\_ Other: **Geo Pump P 1102**

## SAMPLING MEASUREMENTS


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
14:45	1.20	.2	.2	21.41	6.85	1.89	9.02	+55	59.1	
14:55	2.18	.3	.2	21.47	6.84	1.89	8.20	-52	57.6	
15:00	2.20	.4	.2	21.44	6.85	1.89	5.01	-55	55.0	
15:05	2.23	.5	.2	21.43	6.85	1.89	2.42	-56	52.1	
15:10	2.28	.6	.2	21.45	6.85	1.89	0	-56	53.9	
15:15	2.33	.7	.2	21.44	6.85	1.89	0	-56	54.7	
15:20	2.41	.8	.2	21.37	6.84	1.89	0	-55	53.7	
15:25	2.44	.9	.2	21.41	6.85	1.89	0	-56	53.1	

Water Level (ft. BMP) at End of Purge: **2.44** Sample Intake Depth (ft. BMP): **24.0**

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
<del>15:15</del>						
15:25		G	6	Y		

Comments: **Power Supply error Homba  
Printed circuit board failure  
Change batteries to fix the error**



**Golder Associates Inc.**  
11231 Richmond Avenue, Suite D104  
Houston, TX 77082  
(832) 916-3690

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 1919232 Project Name: UPRR HWPH Date: 1-20-2020  
 Sampling Location (well ID, etc.): MW-7013 Starting Water Level (ft. BMP): 6.55  
 Sample Number: WGI620 MW 7013 20200120 Casing Stickup (ft.): 0  
 Sampled by: TIM MCSpeider WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 35.30 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59  
 METHODS (describe): Low Flow Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge into drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P2111 Hera Other: Cred pump P1102  
 Multi Meter: Homba C1103  
 Field Calibration: Ph. 4.0  
 Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
10:50	6.55	.2	.2	20.16	1.34	2.69	6.52	9	81.4	clear/nd
11:00	7.31	.3	.2	20.64	0.78	2.61	6.53	1	92.9	
11:05	7.49	.4	.2	20.44	1.03	2.61	6.56	3	84.9	
11:10	7.64	.5	.2	20.49	0	2.61	6.57	-5	79.9	
11:15	7.79	.6	.2	20.86	0	2.61	6.58	-12	43.1	
11:20	7.86	.7	.2	20.93	0	2.62	6.58	-14	35.9	
11:25	7.92	.8	.2	20.90	0	2.61	6.58	-14	34.7	
11:30	8.00	.9	.2	20.67	0	2.61	6.59	-13	22.6	

WL (ft. BMP) at End of Purge: 8.0 Sample Intake Depth (ft. BMP): 31.0

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y) (N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G) (P)	No.			
11:30			6			

Comments: \_\_\_\_\_


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 Phone: (512) 671-3434 Fax: (512) 671-3446

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>1.12</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-15-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>1.12</u>	ft. BGL
MW ID	<u>MW71B</u>	Location	Other _____	Total MW Depth	<u>36.80</u>	ft. BGL
Sample ID	<u>WG-162D-MW71B-20190115</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1328				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1338	<u>.2</u>	<u>/</u>	<u>1.29</u>	<u>22.7</u>	<u>6.96</u>	<u>610</u>	<u>0.59</u>	<u>-72</u>	<u>7.2</u>
1343	<u>↓</u>	<u>/</u>	<u>1.34</u>	<u>22.3</u>	<u>6.92</u>	<u>570</u>	<u>0.41</u>	<u>-66</u>	<u>7.4</u>
1348	<u>↓</u>	<u>/</u>	<u>1.35</u>	<u>22.4</u>	<u>6.91</u>	<u>580</u>	<u>0.42</u>	<u>-61</u>	<u>7.4</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1400</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1400</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-9-2020  
 Sampling Location (well ID, etc.): MW-72B Starting Water Level (ft. BMP): 14.78  
 Sample Number: WG 1620 MW 72B 20200109 Casing Stickup (ft.):       
 Sampled by: Tim McSpadden WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TOC - (Steel) or PVC TD (ft. BMP): 40.31 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H2O drum 1-7-2020

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: P2111 Hema Other: Geo pump 1102  
 Multi Meter: Hema C1103  
 Field Calibration: PH. 4.8 NTC-0 comp. 4/18  
 Filter / Filter Size: 10 microp

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:00	14.78	.2	.2	22.43	.46	1.75	6.36	-39	96.6	clear/no
8:10	15.58	.3	.2	22.41	.31	1.74	6.46	-40	99.9	
8:15	15.77	.4	.2	22.46	.11	1.71	6.72	-33	70.8	
8:20	15.84	.5	.2	22.52	.06	1.70	6.76	-32	65.8	
8:25	15.99	.6	.2	22.65	0.0	1.66	6.88	-30	33.4	
8:30	16.21	.7	.2	22.70	0	1.64	6.90	-30	32.8	
8:35	16.29	.8	.2	22.79	0	1.59	6.94	-31	32.5	

WL (ft. BMP) at End of Purge: 16.29 Sample Intake Depth (ft. BMP): 35

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
8:35	2.0		6			

Comments:

NO tubing in well



**GOLDER**

2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.29</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-17-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.29</u>	ft. BGL
MW ID	<u>MW 74B</u>	Location	Other _____	Total MW Depth	<u>35.10</u>	ft. BGL
Sample ID	<u>WG-162D-MW74B-20190117</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>0946</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>0956</u>	<u>.2</u>	<u>/</u>	<u>8.47</u>	<u>23.1</u>	<u>6.89</u>	<u>1380</u>	<u>0.79</u>	<u>-71</u>	<u>16</u>
<u>1061</u>	<u>↓</u>	<u>/</u>	<u>8.49</u>	<u>23.1</u>	<u>6.85</u>	<u>1360</u>	<u>0.65</u>	<u>-72</u>	<u>12</u>
<u>1006</u>	<u>↓</u>	<u>/</u>	<u>8.49</u>	<u>23.2</u>	<u>6.84</u>	<u>1350</u>	<u>0.67</u>	<u>-72</u>	<u>13</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1020</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1020</u>	<u>40mL / 1L</u>	<u>G / G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>8.53</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-17-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>8.52</u>	ft. BGL
MW ID	<u>MW75B</u>	Location	Other _____	Total MW Depth	<u>37.25</u>	ft. BGL
Sample ID	<u>WG-162D-MW75B-20190117</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0742				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0753	<u>2</u>	<u>/</u>	<u>8.79</u>	<u>21.7</u>	<u>6.79</u>	<u>1170</u>	<u>0.74</u>	<u>-42</u>	<u>5.2</u>
0757	<u>↓</u>	<u>/</u>	<u>8.81</u>	<u>21.3</u>	<u>6.76</u>	<u>1140</u>	<u>0.63</u>	<u>-41</u>	<u>4.4</u>
0802	<u>↓</u>	<u>/</u>	<u>8.80</u>	<u>21.2</u>	<u>6.76</u>	<u>1130</u>	<u>0.67</u>	<u>-41</u>	<u>4.7</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0815	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0815	<u>40ml / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton

name signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>20.68</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>20.68</u> ft. BGL
MW ID	<u>MW76C</u>	Location	Other _____	Total MW Depth	<u>70.40</u> ft. BGL
Sample ID	<u>WG-162D-MW76C-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1444				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1454	<u>.2</u>	<u>/</u>	<u>20.91</u>	<u>22.6</u>	<u>7.06</u>	<u>1940</u>	<u>0.74</u>	<u>-77</u>	<u>5.6</u>
1459	<u>↓</u>	<u>/</u>	<u>20.94</u>	<u>22.3</u>	<u>7.12</u>	<u>1920</u>	<u>0.51</u>	<u>-70</u>	<u>7.1</u>
1504	<u>↓</u>	<u>/</u>	<u>20.93</u>	<u>22.4</u>	<u>7.13</u>	<u>1910</u>	<u>0.52</u>	<u>-71</u>	<u>7.2</u>
Purging was completed based on: <input checked="" type="checkbox"/> stabilization of water quality parameters <input type="checkbox"/> removal of three well volumes <input type="checkbox"/> removal of at least one half well volume (low yield well)									

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1520	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1520	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>6.81</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>6.81</u> ft. BGL
MW ID	<u>MW 77A</u>	Location	Other _____	Total MW Depth	<u>22.80</u> ft. BGL
Sample ID	<u>WG-162D-MW77A-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
<u>1352</u>				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
<u>1402</u>	<u>.3</u>	<u>/</u>	<u>6.97</u>	<u>21.4</u>	<u>6.92</u>	<u>1270</u>	<u>0.79</u>	<u>-42</u>	<u>9.2</u>
<u>1407</u>	<u>↓</u>	<u>/</u>	<u>7.01</u>	<u>21.3</u>	<u>6.91</u>	<u>1240</u>	<u>0.66</u>	<u>-39</u>	<u>7.4</u>
<u>1412</u>	<u>↓</u>	<u>/</u>	<u>7.02</u>	<u>21.6</u>	<u>6.91</u>	<u>1250</u>	<u>0.69</u>	<u>-39</u>	<u>8.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1425</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1425</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.96</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>H7-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.96</u>	ft. BGL
MW ID	<u>MW78A</u>	Location	Other _____	Total MW Depth	<u>24.85</u>	ft. BGL
Sample ID	<u>WG-162D-MW78A-20190117</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1053				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1103	<u>.2</u>	<u>/</u>	<u>8.21</u>	<u>21.3</u>	<u>6.79</u>	<u>1240</u>	<u>1.46</u>	<u>-71</u>	<u>6.7</u>
1108	<u>↓</u>	<u>/</u>	<u>8.24</u>	<u>21.2</u>	<u>6.72</u>	<u>1260</u>	<u>1.31</u>	<u>-65</u>	<u>5.1</u>
1113	<u>↓</u>	<u>/</u>	<u>8.25</u>	<u>21.6</u>	<u>6.71</u>	<u>1230</u>	<u>1.26</u>	<u>-67</u>	<u>5.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1125</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1125</u>	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>9.81</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-17-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>9.81</u>	ft. BGL
MW ID	<u>MW79A</u>	Location	Other _____	Total MW Depth	<u>26.45</u>	ft. BGL
Sample ID	<u>WG-162D-MW79A-20190117</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0849				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0859	<u>.2</u>	<u>/</u>	<u>10.06</u>	<u>21.6</u>	<u>6.91</u>	<u>720</u>	<u>0.61</u>	<u>-31</u>	<u>16</u>
0904	<u>↓</u>	<u>/</u>	<u>10.08</u>	<u>21.5</u>	<u>6.84</u>	<u>760</u>	<u>0.46</u>	<u>-25</u>	<u>17</u>
0909	<u>↓</u>	<u>/</u>	<u>10.09</u>	<u>21.5</u>	<u>6.86</u>	<u>770</u>	<u>0.47</u>	<u>-26</u>	<u>17</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration	Preserved	Notes and Observations
Time	Volume	Composition (G / P)	No.	(10µm / 45µm)	(type)	(quality control sample, other)
<u>0920</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>0920</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UPRR HW/PW Date: 1-7-2020  
 Sampling Location (well ID, etc.): MW-803 Starting Water Level (ft. BMP): 10.87  
 Sample Number: WG-1620-MW 803 2020.01.07 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP):      (ft. BGL):       
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 37.74 (ft. BGL):       
 Screened Interval (ft. BGL):      Ft. water:      Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL):      1X Casing Vol (gal.):      3X (gal.):     

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum

INSTRUMENTS (Indicate make, model, l.d.)  
 Water Level: D2111 Horva Other:       
 Multi Meter: Hanna C1103  
 Field Calibration: pH-3.99 NTC-0 Cond. 4.52  
 Filter / Filter Size:     

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
15:15	10.87	.2	.2	24.86	11.31	.949	7.18	-7	20.8	
15:25	11.11	.3	.2	25.00	3.75	.946	7.17	-5	19.6	
15:30	11.14	.4	.2	25.17	3.37	.945	7.17	-6	18.1	
15:35	11.21	.5	.2	25.23	3.21	.946	7.18	-5	12.5	
15:40	11.24	.6	.2	25.25	3.11	.945	7.18	-5	10.1	
15:45	11.31	.7	.2	25.38	2.97	.943	7.19	-3	9.3	
15:50	11.33	.8	.2	25.42	2.94	.942	7.19	-4	4.8	
15:55	11:33	.9	.2	25.42	2.01	.942	7.19	-3	9.8	
16:00	11:35	.10	.2	25.40	2.91	.912	7.19	-5	4.5	

WL (ft. BMP) at End of Purge: 11.35 Sample Intake Depth (ft. BMP): 34.0

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
16:00	2.0		9			

Comments:     



**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446

# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>7.12</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>7.12</u> ft. BGL
MW ID	<u>MW81B</u>	Location	Other _____	Total MW Depth	<u>33.80</u> ft. BGL
Sample ID	<u>WG-1620-MW81B-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1548				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1558	<u>.2</u>	<u>/</u>	<u>7.22</u>	<u>21.9</u>	<u>7.21</u>	<u>1060</u>	<u>0.91</u>	<u>-62</u>	<u>8.1</u>
1603	<u>↓</u>	<u>/</u>	<u>7.27</u>	<u>21.6</u>	<u>7.15</u>	<u>1080</u>	<u>0.77</u>	<u>-60</u>	<u>6.7</u>
1608	<u>↓</u>	<u>/</u>	<u>7.28</u>	<u>21.6</u>	<u>7.16</u>	<u>1090</u>	<u>0.79</u>	<u>-61</u>	<u>7.1</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1620</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
<u>1620</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

MS / MSD

**GROUNDWATER SAMPLING RECORD** PAGE \_\_\_\_ of \_\_\_\_

Project Number: 19119232 Project Name: UPRR HW Pu Date: 1-21-2020  
 Sampling Location (well ID, etc.): MW 82B Starting Water Level (ft. BMP): 2.65  
 Sample Number: WG 1620 MW 82B 20200121 Casing Stickup (ft.): \_\_\_\_\_  
 Sampled by: T.M. MCS padden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 34.90 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Horco Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: PK. 4.0  
 Filter / Filter Size: \_\_\_\_\_

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:30	2.65	.2	.2	16.87	0.73	2.38	5.60	-61	65.3	
8:40	2.94	.3	.2	18.24	0.60	2.35	5.88	-67	84.0	
8:45	2.88	.4	.2	18.94	0.44	2.33	6.10	-76	93.8	
8:50	2.81	.5	.2	18.40	0	2.34	6.23	-64	73.0	
8:55	2.80	.6	.2	18.79	0	2.27	6.34	-58	6.3	
9:00	2.82	.7	.2	18.87	0	2.26	6.35	-58	5.2	
9:05	2.83	.8	.2	18.93	0	2.25	6.36	-60	5.1	
9:10	2.81	.9	.2	19.03	0	2.22	6.37	-61	3.9	

WL (ft. BMP) at End of Purge: 2.81 Sample Intake Depth (ft. BMP): 30.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
9:10			18			

Comments:  
WG 1620 MW 82B MS 2020 0121  
WG 1620 MW 82B MSD 2020 0121

**GOLDER**  
 2201 Double Creek Dr., Suite 4004  
 Round Rock, Texas 78664  
 Phone: (512) 671-3434 Fax: (512) 671-3446







Jan

<b>GROUNDWATER SAMPLING RECORD</b>		PAGE <u>1</u> of <u>    </u>
Project Number: <u>19119232</u>	Project Name: <u>UPRR HWPW</u>	Date: <u>1-10-2020</u>
Sampling Location (well ID, etc.): <u>MW 83C</u>	Starting Water Level (ft. BMP): <u>17.21</u>	
Sample Number: <u>WG 1620 MW 83C 2020 0110</u>	Casing Stickup (ft.): <u>0</u>	
Sampled by: <u>T.M. McSpadden</u>	WL (ft. BMP): <u>    </u>	(ft. BGL): <u>    </u>
Measuring Point (MP) of Well: <u>TOC</u> Steel or <u>PVC</u>	TD (ft. BMP): <u>67.90</u>	(ft. BGL): <u>    </u>
Screened Interval (ft. BGL): <u>    </u>	Ft. water: <u>    </u>	Casing Dia. (In ID): <u>2"</u>
Filter Pack Interval (ft. BGL): <u>    </u>	1X Casing Vol (gal.): <u>    </u>	3X (gal.): <u>    </u>

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16    4": 0.65    5.25": 1.12    6": 1.47    6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Altronex Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder    Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge 140 drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.D.)  
 Water Level: R2111 Heron    Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: PH-3.99 NTV-0 cond - 4.49  
 Filter / Filter Size: 10 microns

**SAMPLING MEASUREMENTS**    Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
8:15	17.71	.2	.2	23.56	2.10	1.23	6.26	-100	67.9	clear/no
8:25	17.99	.3	.2	23.75	1.94	1.22	6.88	-133	66.9	
8:30	18.0	.4	.2	23.80	1.52	1.21	7.13	-147	59.2	
8:35	17.99	.5	.2	23.98	1.06	1.17	7.95	-178	30.2	
8:40	17.97	.6	.2	24.07	0	1.17	8.28	-188	25.8	
8:45	17.99	.7	.2	24.10	0	1.16	8.47	-201	23.2	
8:50	17.99	.8	.2	24.12	0	1.16	8.52	-209	23.7	
8:55	17.97	.9	.2	24.09	0	1.16	8.60	-209	20.9	

WL (ft. BMP) at End of Purge: 17.97    Sample Intake Depth (ft. BMP): 63

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
9:00			6			

Comments: Tubing 3' in well, used tools to fish out tubing.



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**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-16-2020  
 Sampling Location (well ID, etc.): MW 84B Starting Water Level (ft. BMP): 4.33  
 Sample Number: W61620 MW84B 20200116 Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 46.40 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

**INSTRUMENTS** (Indicate make, model, I.D.)

Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Hanba C1103  
 Field Calibration: Ph. 4.0  
 Filter / Filter Size: \_\_\_\_\_

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:15	4.33	.2	.2	23.77	1.05	1.64	6.92	-15	9.0	clear/lo
13:25	4.67	.3	.2	23.37	.28	1.69	6.52	37	0	
13:30	4.68	.4	.2	23.27	.02	1.71	6.46	66	0	
13:35	4.70	.5	.2	23.26	.17	1.71	6.44	75	0	
13:40	4.73	.6	.2	23.24	.26	1.71	6.41	53	0	
13:45	4.73	.7	.2	23.22	.21	1.71	6.40	-9	0	

WL (ft. BMP) at End of Purge: 4.73 Sample Intake Depth (ft. BMP): 36

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G,P)	No.			
13:45			6			

Comments:  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_


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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>22.27</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-9-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>22.27</u>	ft. BGL
MW ID	<u>MW-85C</u>	Location	Other _____	Total MW Depth	<u>70.10</u>	ft. BGL
Sample ID	<u>WG-1620-MW85C-2019</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1303				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1313	<u>2</u>	<u>/</u>	<u>22.47</u>	<u>22.6</u>	<u>7.52</u>	<u>1360</u>	<u>0.96</u>	<u>-21</u>	<u>6.7</u>
1318	<u>↓</u>	<u>/</u>	<u>22.51</u>	<u>22.4</u>	<u>7.46</u>	<u>1360</u>	<u>0.73</u>	<u>-16</u>	<u>5.2</u>
1323	<u>↓</u>	<u>/</u>	<u>22.52</u>	<u>22.1</u>	<u>7.47</u>	<u>1390</u>	<u>0.76</u>	<u>-17</u>	<u>4.3</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1335	<u>60ml</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
1335	<u>40ml / 1L</u>	<u>G/G</u>		<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature



# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>21.09</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-17-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>21.09</u>	ft. BGL
MW ID	<u>MW86C</u>	Location	Other _____	Total MW Depth	<u>70.25</u>	ft. BGL
Sample ID	<u>WG-1620-MW86C-20190117</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
0642				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
0652	<u>2</u>	<u>/</u>	<u>21.32</u>	<u>21.6</u>	<u>7.12</u>	<u>1060</u>	<u>0.92</u>	<u>-36</u>	<u>4.7</u>
0657	<u>↓</u>	<u>/</u>	<u>21.37</u>	<u>21.2</u>	<u>7.11</u>	<u>1070</u>	<u>0.76</u>	<u>-31</u>	<u>6.6</u>
0702	<u>↓</u>	<u>/</u>	<u>21.36</u>	<u>21.3</u>	<u>7.09</u>	<u>1070</u>	<u>0.74</u>	<u>-32</u>	<u>6.2</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
0725	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
0725	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton

name signature

WG-1620-FD04-20190117

# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UP RR HWPU Date: 1-20-2020  
 Sampling Location (well ID, etc.): MW-87C Starting Water Level (ft. BMP): 15.79  
 Sample Number: WG-1620 MW 87C 20200120 Casing Stickup (ft.): 0  
 Sampled by: J.M. McSpodden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 65.0 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe) Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H<sub>2</sub>O drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P2111 Heron Other: Geo Amp P1102  
 Multi Meter: HoriBa C1103  
 Field Calibration: Ph. 4.0  
 Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:20	5.79	.2	.2	23.49	1.52	1.05	9.13	-165	84.2	clear/no
13:30	6.22	.3	.2	23.65	0.21	1.03	9.41	-174	87.7	
13:35	6.12	.4	.2	23.67	0.05	.985	10.00	-189	51.3	
13:40	5.96	.5	.2	23.51	0.0	.968	10.24	-190	49.9	
13:45	5.81	.6	.2	23.23	0.0	.976	10.59	-204	39.4	
13:50	5.84	.7	.2	23.01	0.0	.975	10.69	-209	37.5	
13:55	5.88	.8	.2	22.64	0	.974	10.93	-208	36.9	
14:00										

WL (ft. BMP) at End of Purge: 15.88 Sample Intake Depth (ft. BMP): 60"

## SAMPLE INVENTORY

Time	Bottles Collected			No.	Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)					
14:00				6			

Comments:

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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>24.64</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>24.64</u>	ft. BGL
MW ID	<u>MW88C</u>	Location	Other _____	Total MW Depth	<u>-</u>	ft. BGL
Sample ID	<u>WG-162D-MW88C-20190114</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1117				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1127	<u>2</u>	<u>/</u>	<u>24.72</u>	<u>23.4</u>	<u>7.21</u>	<u>1210</u>	<u>0.67</u>	<u>-77</u>	<u>5.2</u>
1132	<u>↓</u>	<u>/</u>	<u>24.71</u>	<u>23.1</u>	<u>7.16</u>	<u>1240</u>	<u>0.42</u>	<u>-71</u>	<u>4.1</u>
1137	<u>↓</u>	<u>/</u>	<u>24.72</u>	<u>23.2</u>	<u>7.17</u>	<u>1230</u>	<u>0.43</u>	<u>-72</u>	<u>4.2</u>

Purging was completed based on:  stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1150</u>	<u>60ML</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1150</u>	<u>40ML / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON name John Brayton signature



**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-16-2020  
 Sampling Location (well ID, etc.): MW 89B Starting Water Level (ft. BMP): 6.19  
 Sample Number: WG-1620 MW 89B 20200116 Casing Stickup (ft.): 0  
 Sampled by: Jim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or (PVC) TD (ft. BMP): 40.80 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID) 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): low flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: D/Aleconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H2O drum 1-7-2020

**INSTRUMENTS** (Indicate make, model, I.d.)

Water Level: P2111 Heron Other: Geo Pump P1102  
 Multi Meter: Horiba C1103  
 Field Calibration: Ph. 4.0  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:10	6.19	.2	.2	23.35	0.98	258	7.81	-138	27.5	clear/no
12:20	6.30	.3	.2	23.26	0.29	260	7.69	-146	23.6	
12:25	6.47	.4	.2	22.98	0	281	7.57	-155	21.3	
12:30	6.55	.5	.2	22.90	0	283	7.55	-157	21.1	
12:35	6.71	.6	.2	22.95	0	287	7.45	-171	23.4	
12:40	6.82	.7	.2	22.98	0	289	7.36	-182	23.9	

WL (ft. BMP) at End of Purge: 6.82 Sample Intake Depth (ft. BMP): 36.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
12:45			6			

Comments: WG-1620 FB08 20200116 12:00

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**GROUNDWATER SAMPLING RECORD**

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPW Date: 1-20-2020  
 Sampling Location (well ID, etc.): MW 90B Starting Water Level (ft. BMP): 0.89  
 Sample Number: WG-1620 MW 90B 2020 01 20 Casing Stickup (ft.): 0  
 Sampled by: J.M. McSpurda WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 36.40 (ft. BGL): 2  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): \_\_\_\_\_  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe) Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Rinse  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: Purge H2O drum 1-7-2020

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: P2111 Horon Other: Geo Pump P1103  
 Multi Meter: Horiba C1103  
 Field Calibration: Ph. 4.0  
 Filter / Filter Size: 10 microns

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
14:40	0.89	.2	.2	21.76	1.52	.854	9.10	-16	32.2	Clear/wd
14:50	1.31	.3	.2	22.05	1.76	.861	8.03	-27	14.3	
14:55	1.47	.4	.2	22.00	1.48	.857	8.86	-37	14.0	
15:00	2.52	.5	.2	21.73	1.10	.858	8.32	-40	18.4	
15:05	2.68	.6	.2	21.59	1.33	.840	8.28	-42	19.4	
15:10	2.84	.7	.2	21.50	1.26	.844	8.05	-42	19.2	
15:15	2.96	.8	.2	21.71	1.12	.836	7.02	-44	14.1	

WL (ft. BMP) at End of Purge: 2.96 Sample Intake Depth (ft. BMP): 31.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
15:20			6			

Comments:



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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: **1911232** Project Name: **UPRR HWPW** Date: **1-13-2020**  
 Sampling Location (well ID, etc.): **P-10** Starting Water Level (ft. BMP): **3.18**  
 Sample Number: **WG-1620 MW P10 2020.0113** Casing Stickup (ft.): \_\_\_\_\_  
 Sampled by: **Tim Mespaddeu** WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or **PVC** TD (ft. BMP): **42.90** (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): **2"**  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal.): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): **Low Flow - Dedicated Tubing/Equipment**  
 Cleaning Equipment: **DI/Alconox Rinse**  
 Purge: **Peristaltic Pump / SS Pump / Bailer / Bladder** Sampling: **Peristaltic Pump / SS Pump / Bailer / Bladder**  
 Disposal of Discharged Water: **Purge 170 drum 1-7-2020**

INSTRUMENTS (Indicate make, model, I.d.)  
 Water Level: **P2111 Horiba** Other: \_\_\_\_\_  
 Multi Meter: **Horiba C1103**  
 Field Calibration: **PH-3.99 NTU+0 count 445**  
 Filter / Filter Size: **-**

**SAMPLING MEASUREMENTS** Begin Purge:


Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:35	3.18	.2	.2	19.79	2.81	1.981	6.95	114	13.8	clear/no
13:45	4.61	.3	.2	19.61	2.41	1.01	6.97	113	8.9	
13:50	4.64	.4	.2	19.70	1.81	1.07	7.08	110	6.5	
13:55	4.64	.5	.2	19.70	1.57	1.08	7.19	108	6.1	
14:00	4.65	.6	.2	19.63	1.52	1.08	7.17	107	4.9	
14:05	4.66	.7	.2	19.65	1.38	1.09	7.21	106	3.3	
14:10	4.68	.8	.2	19.61	1.36	1.09	7.21	106	1.4	
14:15	4.70	.9	.2	19.60	1.38	1.09	7.21	106	2.4	

WL (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): **32**

**SAMPLE INVENTORY**

Time	Bottles Collected			No.	Filtration (Y (N))	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)					
14:20				2		heat	

Comments: \_\_\_\_\_



**GOLDER**  
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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>4.26</u>	ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u>	ft.
Date	<u>1-14-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>4.26</u>	ft. BGL
MW ID	<u>P11</u>	Location	Other _____	Total MW Depth	<u>42.75</u>	ft. BGL
Sample ID	<u>WG-162D-P11 -20190114</u>	Water Quality		MW Diameter	<u>2.0</u>	inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u>	gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u>	ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1023				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1033	<u>2</u>	<u>/</u>	<u>4.42</u>	<u>21.7</u>	<u>6.91</u>	<u>1060</u>	<u>1.34</u>	<u>-99</u>	<u>7.9</u>
1038	<u>↓</u>	<u>/</u>	<u>4.43</u>	<u>21.6</u>	<u>6.91</u>	<u>1020</u>	<u>0.93</u>	<u>-96</u>	<u>7.1</u>
1043	<u>↓</u>	<u>/</u>	<u>4.43</u>	<u>21.6</u>	<u>6.92</u>	<u>1030</u>	<u>0.92</u>	<u>-96</u>	<u>7.4</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
1055	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO3</u>	<u>METALS</u>
1055	<u>40mL / 1L</u>	<u>G/G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCs / SVOCs</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON  
name

John Brayton  
signature

# GROUNDWATER SAMPLING RECORD

1

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UPRR HWPLW Date: 1-13-2020  
 Sampling Location (well ID, etc.): P-12 Starting Water Level (ft. BMP): 3.96  
 Sample Number: WG-1620 MW P12 20200113 Casing Stickup (ft.): 14'  
 Sampled by: Tim McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): 42.75 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment

Cleaning Equipment: DI/Alconox-Rinse

Purge: Peristaltic Pump / SS Pump / Bailer / Bladder

Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder

Disposal of Discharged Water: Purge 1/30 drum 1-7-2020

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: P2111 Herra

Other: Geo pump P1102

Multi Meter: Numba C1103

Field Calibration: PH-3.99 NTU-0 cond 4.49

Filter / Filter Size: 10

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
<u>9:05</u>	<u>9:45</u>	<u>3.96</u>	<u>.2</u>	<u>20.01</u>	<u>7.01</u>	<u>1.34</u>	<u>7.39</u>	<u>99</u>	<u>73.3</u>	<u>Clear/No</u>
<u>9:15</u>	<u>9:55</u>	<u>4.89</u>	<u>.2</u>	<u>20.09</u>	<u>1.82</u>	<u>1.33</u>	<u>7.30</u>	<u>100</u>	<u>76.0</u>	
<u>9:30</u>	<u>10:00</u>	<u>4.91</u>	<u>.2</u>	<u>20.11</u>	<u>1.44</u>	<u>1.31</u>	<u>7.19</u>	<u>103</u>	<u>72.8</u>	
<u>9:35</u>	<u>10:05</u>	<u>4.95</u>	<u>.2</u>	<u>20.14</u>	<u>1.30</u>	<u>1.30</u>	<u>7.14</u>	<u>104</u>	<u>70.8</u>	
<u>9:40</u>	<u>10:10</u>	<u>4.98</u>	<u>.2</u>	<u>20.17</u>	<u>1.20</u>	<u>1.30</u>	<u>7.09</u>	<u>106</u>	<u>68.5</u>	
<u>9:45</u>	<u>10:15</u>	<u>4.99</u>	<u>.2</u>	<u>20.26</u>	<u>1.00</u>	<u>1.28</u>	<u>6.99</u>	<u>111</u>	<u>59.9</u>	
<u>9:50</u>	<u>10:20</u>	<u>4.99</u>	<u>.2</u>	<u>20.33</u>	<u>0.67</u>	<u>1.28</u>	<u>6.90</u>	<u>112</u>	<u>59.5</u>	
<u>9:55</u>	<u>10:25</u>	<u>5.02</u>	<u>.2</u>	<u>20.33</u>	<u>0.49</u>	<u>1.28</u>	<u>6.89</u>	<u>112</u>	<u>53.4</u>	

WL (ft. BMP) at End of Purge: 5.02

Sample Intake Depth (ft. BMP): 37

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
<u>10:30</u>			<u>2</u>	<u>-</u>	<u>Heat</u>	

Comments:

Collected MS MSD  
Anal Fred blank @ P-12



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# Groundwater Sample Collection



Project/Phase	<u>19119232</u>	Equipment Decon	<input checked="" type="checkbox"/> Dedicated equipment	Depth to Water	<u>5.53</u> ft. BMP
Site Location	<u>UPRR-HWPW</u>		<input type="checkbox"/> Decon between locations	Casing Stickup	<u>-</u> ft.
Date	<u>1-13-20</u>	Reference Point	<input checked="" type="checkbox"/> Top of casing	<input type="checkbox"/> Depth to Water	<u>5.52</u> ft. BGL
MW ID	<u>TW41B</u>	Location	Other _____	Total MW Depth	<u>42.25</u> ft. BGL
Sample ID	<u>WG-1620-TW41B-2019</u>	Water Quality		MW Diameter	<u>2.0</u> inches
Pump	<input type="checkbox"/> Waterra <input type="checkbox"/> Submersible	Meter Model	<u>HORIBA</u>	MW Volume	<u>1</u> gallons
	<input checked="" type="checkbox"/> Peristaltic <input type="checkbox"/> Bladder	Unit Number	<u>U-50</u>	Pump Intake Depth	<u>1</u> ft. BGL

Standard volume capacity of monitoring wells: Schedule 40 PVC (1-inch OD, 0.04 gallons/foot; 2-inch OD, 0.16 gallons/foot; 4-inch OD, 0.65 gallons/foot)

Time	Purge Rate (L/min)	Cumulative Purge Volume (L)	Depth to Water (ft)	Temp (°C)	pH	Conductivity (µS/m or mS/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)	Turbidity (NTU)
1326				<input type="checkbox"/> NM	<input type="checkbox"/> NM	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> not measured	<input type="checkbox"/> NM
1336	<u>2</u>	<u>/</u>	<u>5.74</u>	<u>22.7</u>	<u>6.91</u>	<u>1460</u>	<u>0.79</u>	<u>-36</u>	<u>21</u>
1341	<u>↓</u>	<u>/</u>	<u>5.74</u>	<u>22.3</u>	<u>6.84</u>	<u>1490</u>	<u>0.61</u>	<u>-27</u>	<u>14</u>
1346	<u>↓</u>	<u>/</u>	<u>5.77</u>	<u>22.4</u>	<u>6.86</u>	<u>1490</u>	<u>0.62</u>	<u>-29</u>	<u>16</u>

Purging was completed based on:

stabilization of water quality parameters  removal of three well volumes  removal of at least one half well volume (low yield well)

Bottles Collected				Filtration (10µm / 45µm)	Preserved (type)	Notes and Observations (quality control sample, other)
Time	Volume	Composition (G / P)	No.			
<u>1400</u>	<u>60mL</u>	<u>P</u>	<u>1</u>	<input type="checkbox"/> filtered <input type="checkbox"/> unfiltered	<u>HNO<sub>3</sub></u>	<u>METALS</u>
<u>1400</u>	<u>40mL / 1L</u>	<u>G / G</u>	<u>1</u>	<input type="checkbox"/> filtered <input checked="" type="checkbox"/> unfiltered	<u>HCL / NONE</u>	<u>VOCS / SVOCS</u>

Notes: record time at which purging is started. For low flow sampling, recommended stabilization criteria: temp ± 0.5°C; pH ± 0.1 units; conductivity ± 3%; DO ± 1 mg/L; ORP ± 10 mV; for at least three successive measurements that are made every 3-5 minutes with <1 foot of stable draw down.

Unless otherwise noted, groundwater sample collection was completed in accordance with the applicable requirements of Golder's Quality Assurance Program and Standard Operating Procedure  9 Conventional Groundwater Sample Collection  10 Low Flow Groundwater Sample Collection.

Field Team Leader JOHN BRAYTON John Brayton  
 name signature

# GROUNDWATER SAMPLING RECORD

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Project Number: **19119232** Project Name: **HWPW** Date: **3-17-2021**  
 Sampling Location (well ID, etc.): **MW 60B** Starting Water Level (ft. BMP): **24.91**  
 Sample Number: **WG-1620 MW 60B 20200317** Casing Stickup (ft.): **0**  
 Sampled by: \_\_\_\_\_ WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC - Steel or PVC TD (ft. BMP): **39.50** (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): \_\_\_\_\_  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): **Low Flow - Dedicated Tubing/Equipment**

Cleaning Equipment: **DI/Alconex Rinse**

Purge: **Peristaltic Pump / SS Pump / Bailer / Bladder**

Sampling: **Peristaltic Pump / SS Pump / Bailer / Bladder**

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: **Heron D2106**

Other: **Geo pump 1122**

Multi Meter: **Horiba C1122**

Field Calibration: **PH buffer 4.0, Field Cal @ 4.0**

Filter / Filter Size: **10 micron**

## SAMPLING MEASUREMENTS

Begin Purge: \_\_\_\_\_

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
14:30	24.91	.2	.2	26.85	0.36	1.35	6.73	-82	306	Clear
14:40	26.11	.3	.2	27.24	1.00	1.36	6.71	-87	0.0	LT Brown
14:45	26.40	.4	.2	26.69	0.31	1.40	6.58	-105	0.0	LT Brown
14:50	27.58	.5	.2	26.20	0.19	1.41	6.55	-111	0.0	LT Brown / Clearing
14:55	27.71	.6	.2	26.08	0.11	1.42	6.54	-113	0.0	clear
15:00	27.84	.7	.2	25.96	0.14	1.42	6.53	-116	0.0	clear
15:05	27.91	.8	.2	25.78	0.10	1.42	6.51	-117	757	clear
15:10	28.04	.9	.2	25.93	0.01	1.41	6.50	-117	691	clear
				25.93						

WL (ft. BMP) at End of Purge: **28.04**

Sample Intake Depth (ft. BMP): **32.0**

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
15:15		Purge	9	10 micron		

Comments: \_\_\_\_\_



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# GROUNDWATER SAMPLING RECORD

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Project Number: 19119232 Project Name: HWPW Date: 3-17-2020  
 Sampling Location (well ID, etc.): MW-88B Starting Water Level (ft. BMP): 7.82  
 Sample Number: WG 1620 MW 88B 20200317 Casing Stickup (ft.): 3'  
 Sampled by: T.M. McSpadden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC Steel or PVC TD (ft. BMP): 43.0 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: \_\_\_\_\_ Casing Dia. (In ID): \_\_\_\_\_  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

## QUALITY ASSURANCE

Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment

Cleaning Equipment: DI/Alconox Rinse

Purge: Peristaltic Pump / SS Pump / Bailer / Bladder

Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder

Disposal of Discharged Water: \_\_\_\_\_

INSTRUMENTS (Indicate make, model, I.d.)

Water Level: Hervo D2106

Other: Geo Pump P1122

Multi Meter: Homba C1122

Field Calibration: PH buffer 4.0 Field cal @ 4.0

Filter / Filter Size: \_\_\_\_\_

## SAMPLING MEASUREMENTS

Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L / m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
13:10	7.82	.2	.2	25.11	4.38	1.15	6.65	37	18.8	Clear
13:20	8.44	.3	.2	25.27	2.33	1.16	6.67	36	1.4	
13:25	8.41	.4	.2	25.35	2.14	1.16	6.52	37	0.0	
13:30	8.52	.5	.2	25.25	2.08	1.16	6.46	50	0.0	
13:35	8.57	.6	.2	25.09	1.80	1.16	6.40	16	0.0	
13:40	8.47	.7	.2	25.09	1.61	1.16	6.38	9	0.1	
13:45	8.41	.8	.2	25.23	1.51	1.16	6.35	-10	0.9	
13:50	8.44	.9	.2	25.11	1.21	1.17	6.35	-17	0.8	

WL (ft. BMP) at End of Purge: \_\_\_\_\_

Sample Intake Depth (ft. BMP): 40.0

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
1350		Pump	9			

Comments: \_\_\_\_\_



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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: Harpw Date: 3-17-2021  
 Sampling Location (well ID, etc.): MW 88A Starting Water Level (ft. BMP): 8.00  
 Sample Number: WG-1020 MW 88A 20200317 Casing Stickup (ft.): 3'  
 Sampled by: Tim McSpedden WL (ft. BMP): \_\_\_\_\_ (ft. BGL): \_\_\_\_\_  
 Measuring Point (MP) of Well: TOC Steel or PVC TD (ft. BMP): 25.50 (ft. BGL): \_\_\_\_\_  
 Screened Interval (ft. BGL): \_\_\_\_\_ Ft. water: 2.8 Casing Dia. (In ID): 2'  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ 1X Casing Vol (gal.): \_\_\_\_\_ 3X (gal): \_\_\_\_\_

**QUALITY ASSURANCE** Gallons/Foot: 2": 0.16 4": 0.65 5.25": 1.12 6": 1.47 6.25": 1.59

METHODS (describe): Low Flow - Dedicated Tubing/Equipment  
 Cleaning Equipment: DI/Alconox Base  
 Purge: Peristaltic Pump / SS Pump / Bailer / Bladder Sampling: Peristaltic Pump / SS Pump / Bailer / Bladder  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS** (Indicate make, model, I.d.)  
 Water Level: Heron D 2106 Other: Geo pump P1122  
 Multi Meter: Hanba C1122  
 Field Calibration: Ph buffer 4.0 = Ph 4.0  
 Filter / Filter Size: 10 micron

**SAMPLING MEASUREMENTS** Begin Purge:

Time	DTW (ft BTOC)	Cum. Vol. (gal. or L)	Purge Rate (gal. or L /m)	Temp. (oC)	DO (mg/L)	Spec. Cond. (µs/cm)	pH	ORP (mV)	Turbidity (NTU)	Color & Sediment
12:20	8.00	.2	.2	24.93	0.72	0.954	6.47	-43	980	clear
12:30	8.64	.3	.2	25.09	0.55	0.957	6.41	-48	926	
12:35	8.69	.4	.2	25.19	0.31	0.957	6.37	-49	910	
12:40	8.57	.5	.2	25.64	0.40	0.948	6.27	-51	832	
12:45	8.51	.6	.2	25.33	0.33	0.957	6.23	-50	773	
12:50	8.47	.7	.2	26.01	0.29	0.938	6.29	-57	684	
12:55	8.41	.8	.2	26.13	0.27	0.935	6.30	-58	637	

WL (ft. BMP) at End of Purge: 8.41 Sample Intake Depth (ft. BMP): 23.

**SAMPLE INVENTORY**

Time	Bottles Collected				Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.				
12:55		Pump	9		10 micron		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_


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# GROUNDWATER SAMPLING RECORD

PAGE      of     

Project Number: 19119232 Project Name: HWPW Date: 3-17-2020

Sample Number: WG-1620 MW 95A 20200317 Starting Water Level (ft. BMP): 3.88

Sampling Location (well ID, etc.): MW-95A Casing Stickup (ft.): 0

Sampled by: T.M. McSpalden Starting Water Level (ft. BGL): 3.88

Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 24.90

Screened Interval (ft. BGL): Casing Diameter (In ID): 2"

Filter Pack Interval (ft. BGL): Casing Volume (gal.): 3.3

## QUALITY ASSURANCE

METHODS (describe): Low Flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: Heraeus D2106 Thermometer: Heraeus C1122

pH Meter: Heraeus C1122 Field Calibration: \_\_\_\_\_

Conductivity Meter: Heraeus C1122 Field Calibration: Ph buffer 4.0 - Field cal Ph-4.0

Filter / Filter Size: \_\_\_\_\_ Other: Geopump P1122

## SAMPLING MEASUREMENTS

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
				± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
8:30	3.88	.2	.2	22.56	5.71	0.832	2.49	235	20.7	Clear
8:40	4.83	.3	.2	22.49	6.05	0.819	2.37	217	16.7	
8:45	4.90	.4	.2	22.31	6.19	0.806	2.22	209	14.8	
8:50	5.99544	.5	.2	22.10	6.23	0.802	2.18	205	12.6	
8:55	5.48	.6	.2	22.25	6.30	0.799	2.07	199	11.4	
9:00	5.37	.7	.2	22.22	6.31	0.799	2.03	194	9.2	
9:05	5.41	.8	.2	22.36	6.36	0.797	1.91	188	9.7	
9:10	5.42	.9	.2	22.42	6.37	0.797	1.90	186	9.1	

Water Level (ft. BMP) at End of Purge: 5.42 Sample Intake Depth (ft. BMP): 22.1

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G.)	No.			
9:10		Pump	9			

Comments: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_



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# GROUNDWATER SAMPLING RECORD

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Project Number: 19119232 Project Name: HWPu Date: 3-17-2020

Sample Number: WG 1620 MW 613 20200317 Starting Water Level (ft. BMP): 8.31

Sampling Location (well ID, etc.): MW 613 Casing Stickup (ft.): 0

Sampled by: Tim McSpadden Starting Water Level (ft. BGL): 8.31

Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 33.25

Screened Interval (ft. BGL): Casing Diameter (In ID): 2"

Filter Pack Interval (ft. BGL): Casing Volume (gal.): 3.9

## QUALITY ASSURANCE

METHODS (describe): Low flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.D.)

Water Level: Heraeus D2106 Thermometer: Hanna C1122

pH Meter: Hanna C1122 Field Calibration: \_\_\_\_\_

Conductivity Meter: Hanna C1122 Field Calibration: PH buffer 4.0 Field Cal. = 4.0

Filter / Filter Size: \_\_\_\_\_ Other: Geo pump P1122

## SAMPLING MEASUREMENTS


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
				± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
11:10	8.31	.2	.2	23.76	7.01	1.18	3.80	129	0.0	Clear
11:20	7.99	.3	.2	23.58	6.72	1.21	3.41	140	0.0	
11:25	8.71	.4	.2	23.41	6.58	1.22	3.18	147	0.0	
11:30	8.59	.5	.2	23.53	6.52	1.22	3.06	149	0.0	
11:35	8.41	.6	.2	23.66	6.50	1.22	3.00	150	0.0	
11:40	8.37	.7	.2	23.63	6.46	1.22	2.91	152	0.0	
11:45	8.31	.8	.2	23.59	6.44	1.22	2.89	153	0.0	

Water Level (ft. BMP) at End of Purge: 8.31 Sample Intake Depth (ft. BMP): 30.

## SAMPLE INVENTORY

Time	Bottles Collected		Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G. #)			
11:45		Pump	9	NV	

Comments:

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# GROUNDWATER SAMPLING RECORD

PAGE \_\_\_ of \_\_\_

Project Number: 19119232 Project Name: UP CSA Date: 3-11-2020  
 Sample Number: WG-1620 MW 54B 20200311 Starting Water Level (ft. BMP): 13.42  
 Sampling Location (well ID, etc.): MW 54B Casing Stickup (ft.): 0  
 Sampled by: TIM M S parker Starting Water Level (ft. BGL): 13.42  
 Measuring Point (MP) of Well: \_\_\_\_\_ Total Depth (ft. BMP): 40.08  
 Screened Interval (ft. BGL): \_\_\_\_\_ Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): \_\_\_\_\_ Casing Volume (gal.): 4.2

## QUALITY ASSURANCE

METHODS (describe): Low Flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_

Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: Herau H-oil D2106

Thermometer: Hanba C1122

pH Meter: Hanba C1122

Field Calibration: \_\_\_\_\_

Conductivity Meter: Hanba C1122

Field Calibration: Buffer 4.0 = Field cal @ 4.0

Filter / Filter Size: 10 micron

Other: Geo pump P1122

## SAMPLING MEASUREMENTS

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L /m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
13:55	13.42	.2	.2	25.98	7.45	1.57	4.15	130	33.4	clear
14:05	15.05	.3	.2	25.87	7.19	1.56	3.71	113	24.9	clear
14:10	15.01	.4	.2	25.79	7.14	1.56	3.54	107	22.2	
14:15	14.97	.5	.2	25.83	7.11	1.57	3.65	102	21.2	
14:20	14.93	.6	.2	25.84	7.09	1.57	3.63	99	21.8	

Water Level (ft. BMP) at End of Purge: 14.93

Sample Intake Depth (ft. BMP): 37.0

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
14:20		Pump	9	Y		

Comments: \_\_\_\_\_



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UP CSA Date: 3-11-2020  
 Sample Number: WG 1620 MW 99C 2020034 Starting Water Level (ft. BMP): 14.99  
 Sampling Location (well ID, etc.): MW 99C Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden Starting Water Level (ft. BGL): 14.99  
 Measuring Point (MP) of Well: TOG Total Depth (ft. BMP): 68.25  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2 1/2  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.): 8.5

**QUALITY ASSURANCE**

METHODS (describe): Low flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: Heron H-OIL D2100 Thermometer: Homba C1122  
 pH Meter: Homba C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Homba C1122 Field Calibration: Buffer 4.0 = Field Cal @ 4.0  
 Filter / Filter Size: \_\_\_\_\_ Other: Geo pump P1122

**SAMPLING MEASUREMENTS**

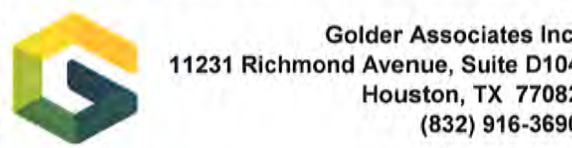
Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	- Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
15:00	14.99	.2	.2	28.20	11.48	1.62	1.13	-88	8.8	clear
15:10	15.25	.3	.2	28.51	11.74	1.62	0.57	-151	5.4	
15:15	15.19	.4	.2	28.87	11.78	1.63	0.40	-160	3.6	
15:20	15.10	.5	.2	28.27	11.80	1.63	0.31	-166	0.1	
15:25	15.02	.6	.2	28.23	11.80	1.63	0.30	-167	0.0	
15:30	14.99	.7	.2	28.26	11.81	1.63	0.28	-169	0.0	
15:35	14.92	.8	.2	28.08	11.78	1.63	0.22	-175	0.0	

Water Level (ft. BMP) at End of Purge: 14.99 Sample Intake Depth (ft. BMP): 210.0

**SAMPLE INVENTORY**

Bottles Collected				Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
Time	Volume	Composition (G, P)	No.			
15:35		Pump	9			

Comments: \_\_\_\_\_



**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UP CSA HWPW Date: 3-12-2020  
 Sample Number: UG-1620 MW 91A 2020 0312 Starting Water Level (ft. BMP): 5.21  
 Sampling Location (well ID, etc.): MW 91A Casing Stickup (ft.): 0  
 Sampled by: Tim McSpedde Starting Water Level (ft. BGL): 5.21  
 Measuring Point (MP) of Well: TOG Total Depth (ft. BMP): 24.10  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: Heraon H-016 D2106 Thermometer: Heraon C1122  
 pH Meter: Heraon C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Heraon C1122 Field Calibration: Ph buffer 4.0 - Field Cal = 4.0  
 Filter / Filter Size: 10 microns Other: Geo pump P1122

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L /m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
8:45	5.21	.2	.2	22.73	7.82	0.733	2.18	-53	332	clear
8:55	5.31	.3	.2	22.60	7.52	0.742	0.71	-69	391	
9:00	5.34	.4	.2	22.50	7.19	0.747	0.24	-81	272	
9:05	5.36	.5	.2	22.57	6.68	0.801	0.08	-89	208	
9:10	5.33	.6	.2	22.65	6.55	0.838	0.00	-91	131	
9:15	5.34	.7	.2	22.64	6.54	0.835	0.00	-92	108	
9:20	5.31	.8	.2	22.67	6.52	0.845	0.00	-92	102	

Water Level (ft. BMP) at End of Purge: 5.31 Sample Intake Depth (ft. BMP): 22.0

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
9:20		Pump	9	10 microns		

Comments:



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: UP Harp Date: 3-12-2020  
 Sample Number: WG-1620 MW 70C 20200312 Starting Water Level (ft. BMP): 16.04  
 Sampling Location (well ID, etc.): MW 70C Casing Stickup (ft.): 0  
 Sampled by: TIM MCS peadden Starting Water Level (ft. BGL): 16.04  
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 67.02  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.): 8.1

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: Horan H-Oil D2106 Thermometer: Horiba C1122  
 pH Meter: Horiba C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Horiba C1122 Field Calibration: PH buffer 4.0 - Field Cal = 4.0  
 Filter / Filter Size: 10 MICRON Other: Geo pump P1122

**SAMPLING MEASUREMENTS**


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
9:55	16.04	.2	.2	24.37	7.01	1.00	0.99	-7	70.3	Clear
10:05	16.25	.3	.2	24.58	7.04	0.997	0.61	-13	70.5	
10:10	16.19	.4	.2	24.63	7.13	0.999	0.35	-41	58.9	
10:15	16.18	.5	.2	24.72	7.10	1.01	0.22	-46	59.7	
10:20	16.20	.6	.2	24.87	7.06	1.11	0.00	-80	56.2	
10:25	16.18	.7	.2	24.76	6.99	1.18	0.02	-91	53.2	
10:30	16.19	.8	.2	24.83	6.98	1.20	0.01	-93	51.4	

Water Level (ft. BMP) at End of Purge: 16.19 Sample Intake Depth (ft. BMP): \_\_\_\_\_

**SAMPLE INVENTORY**

Bottles Collected				Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
Time	Volume	Composition (G/P)	No.			
10:35		Pump	18	10micron		

Comments: Duplicate  
WG-1620 FD 01 2020 0312



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119732 Project Name: UP HWPW Date: 3-12-2020  
 Sample Number: WC-1020 MW 84A 20200312 Starting Water Level (ft. BMP): 4.19  
 Sampling Location (well ID, etc.): MW 84A Casing Stickup (ft.): 0  
 Sampled by: T.M. McSpadden Starting Water Level (ft. BGL): 4.19  
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 23.80  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.): 3.1

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.D.)**

Water Level: Heron H-GIL D2100 Thermometer: Hanna-C1122  
 pH Meter: Hanna C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Hanna C1122 Field Calibration: PH buffer 4.0-Field Cal = 4.0  
 Filter / Filter Size: 0.1 micron Other: Geopung P1122

**SAMPLING MEASUREMENTS**


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L /m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
11:25	4.19	26.09	.2	26.04	6.96	0.970	1.52	53	78.8	clear
11:35	4.41	.3	.2	26.50	6.57	0.985	0.51	34	65.7	
11:40	4.29	.4	.2	26.59	6.54	0.984	0.45	33	71.4	
11:45	4.24	.5	.2	26.74	6.52	0.984	0.31	34	53.5	
11:50	4.22	.6	.2	26.85	6.57	0.984	0.26	36	55.6	
11:55	4.21	.7	.2	27.04	6.57	0.980	0.25	34	53.4	
12:00	4.22	.8	.2	27.19	6.50	0.981	0.24	31	44.8	
12:05	4.20	.9	.2	27.35	6.50	0.978	0.24	30	50.8	

Water Level (ft. BMP) at End of Purge: 4.20 Sample Intake Depth (ft. BMP): 21

**SAMPLE INVENTORY**

Bottles Collected				Filtration (Y / N)	Preservation	Remarks (quality control sample, other)
Time	Volume	Composition (G.P)	No.			
12:15		Pump	9	Whose		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: HWPW Date: 3-12-2020  
 Sample Number: WG-1620 MW 92B 20200312 Starting Water Level (ft. BMP): 4.19  
 Sampling Location (well ID, etc.): MW 92B Casing Stickup (ft.): 0  
 Sampled by: Jim McSpadden Starting Water Level (ft. BGL): 4.19  
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 34.60  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.): 4.8

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: Heron H-01C D2106 Thermometer: Hanna C1122  
 pH Meter: Hanna C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Hanna C1122 Field Calibration: PH buffer 4.0 Field Cal. PH 4.0  
 Filter / Filter Size: 20 Microns Other: Geo pump P1122

**SAMPLING MEASUREMENTS**


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
12:45	4.19	.2	.2	30.89	6.56	1.77	0.53	17	3.7	clear
12:55	4.90	.3	.2	29.99	6.36	1.83	0.26	-4	4.0	
13:00	4.72	.4	.2	29.92	6.34	1.82	0.21	-6	4.8	
13:05	4.64	.5	.2	29.91	6.32	1.83	0.19	-10	5.7	
13:10	4.51	.6	.2	29.90	6.32	1.83	0.17	-12	7.4	
13:15	4.49	.7	.2	29.93	6.32	1.83	0.15	-19	9.2	

Water Level (ft. BMP) at End of Purge: 4.49 Sample Intake Depth (ft. BMP): 32.0

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/P)	No.			
13:20		Pump	9	N		

Comments: \_\_\_\_\_



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: HWPW Date: 3-12-2020  
 Sample Number: WG-1020 MW 93B 20200312 Starting Water Level (ft. BMP): 4.06  
 Sampling Location (well ID, etc.): MW 93B Casing Stickup (ft.): 0  
 Sampled by: Tim McSpadden Starting Water Level (ft. BGL): 4.06  
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 34.01  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.): 4.7

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: Heraeus H-OIL D 2106 Thermometer: Hanba C1122  
 pH Meter: Hanba C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Hanba C1122 Field Calibration: Ph buffer 4.0 Solid Cal = 4.0  
 Filter / Filter Size: 10 MICRON Other: Geo pump P 1122

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L / m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
13:55	4.06	.2	.2	30.37	6.69	1.24	1.09	11	10.8	Clear
14:05	4.49	.3	.2	30.39	6.54	1.28	0.50	-26	12.4	
14:10	4.36	.4	.2	29.56	6.45	1.34	0.25	-50	12.8	
14:15	4.29	.5	.2	28.83	6.40	1.37	0.15	-59	16.0	
14:20	4.28	.6	.2	28.08	6.34	1.37	0.05	-66	21.0	
14:25	4.28	.7	.2	27.79	6.30	1.36	0.01	-70	25.1	
14:30	4.31	.8	.2	27.01	6.31	1.36	0.00	-72	22.4	
14:35	4.28	.9	.2	27.67	6.23	1.34	0.00	-72	28.6	

Water Level (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): 32.0

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G/C)	No.			
14:35		Pump	9	10micron		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: HWPu Date: 3-12-2020  
 Sample Number: WG1020MW94A20200312 Starting Water Level (ft. BMP): 3.99  
 Sampling Location (well ID, etc.): MW 94A Casing Stickup (ft.): 0  
 Sampled by: TIM McSpadden Starting Water Level (ft. BGL): 3.99  
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 21.60  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.):

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.D.)**

Water Level: Hanna H-016 D2106 Thermometer: Hanna C1122  
 pH Meter: Hanna C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Hanna C1122 Field Calibration: PH buffer 4.0, Field Cal = 4.0  
 Filter / Filter Size: 10 micron Other: Geo pump P1122

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
15:05	3.99	.2	.2	30.83	6.78	0.598	0.62	-72	138	Clean
15:15	4.21	.3	.2	30.37	6.82	0.605	0.44	-78	154	
15:20	4.01	.4	.2	29.71	6.77	0.612	0.33	-84	163	
15:25	4.00	.5	.2	29.16	6.74	0.617	0.26	-88	215	
15:30	4.00	.6	.2	27.99	6.65	0.626	0.09	-95	197	
15:35	4.01	.7	.2	27.87	6.63	0.627	0.07	-94	161	
15:40	4.01	.8	.2	27.41	6.59	0.630	0.00	-95	157	
15:45	4.00	.9	.2	26.70	6.34	0.633	0.00	-85	92.8	

Water Level (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): \_\_\_\_\_

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G.P)	No.			
15:45		Pump	9	10micron		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



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# GROUNDWATER SAMPLING RECORD

PAGE 1 of 1

Project Number: 19119232 Project Name: HWPW Date: 3.12.20

Sample Number: 76B W61-1620-MW76B-20200312 Starting Water Level (ft. BMP): ~~4.00~~ 5.40

Sampling Location (well ID, etc.): MW-76B Casing Stickup (ft.): -

Sampled by: *[Signature]* Starting Water Level (ft. BGL): -

Measuring Point (MP) of Well: TOC/PVC Total Depth (ft. BMP): 35.50

Screened Interval (ft. BGL): 31-36 Casing Diameter (in ID): 2

Filter Pack Interval (ft. BGL): 29-36 Casing Volume (gal.): 4.91 = 18.58 L

## QUALITY ASSURANCE

### METHODS (describe):

Cleaning Equipment: Dedicated Equipment

Purging: Peristaltic or bladder pump with dedicated tubing Sampling: Same

Disposal of Discharged Water: 55 gal Drum CSA

### INSTRUMENTS (Indicate make, model, I.D.)

Water Level: Keck

pH Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Conductivity Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Thermometer: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Turbidimeter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

ORP Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

DO Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Other: Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Time	Cum. Vol. (L)	Purge Rate (L/m)	Temp. (oC)	pH	Spec. Cond. (µS/cm)	Color	Turbidity & Sediment (NTU)	ORP (mV)	DO (mg/L)	Water Level (ft BMP)	Remarks
10:10	-	0.2	-	-	-	-	-	-	-	5.40	Pump is on
10:20	2	↓	24.67	6.87	1520	Neutral	30.7	204	2.17	6.20	
10:25	3	↓	24.88	6.88	1530	↓	15.8	194	1.89	6.55	
10:30	4	↓	25.24	6.88	1530	↓	15.6	179	1.75	6.90	

Water Level (ft. BMP) at End of Purge: 6.90 Sample Intake Depth (ft. BMP): 32.50

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
10:35	1 L	G1	2	N	Neat	SVOCs - site specific
10:35	120 mL	P	1	Y	HNO3	Metals - As
10:35	40 mL	G1	6	N	HCL	VOC site specific / TPH

Comments:



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# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: HWPW Date: 3-12-20  
 Sample Number: W61-1620-MW97A-20200312 Starting Water Level (ft. BMP): 5.20  
 Sampling Location (well ID, etc.): MW-97A Casing Stickup (ft.): -  
 Sampled by: *AS* Starting Water Level (ft. BGL): -  
 Measuring Point (MP) of Well: TOC/PVC Total Depth (ft. BMP): 19.85  
 Screened Interval (ft. BGL): 10.5 - 20.5 Casing Diameter (in ID): 2  
 Filter Pack Interval (ft. BGL): 8.5 - 21 Casing Volume (gal.): 2.39 = 9.05L

## QUALITY ASSURANCE

### METHODS (describe):

Cleaning Equipment: Dedicated Equipment  
 Purging: Peristaltic or bladder pump with dedicated tubing Sampling: Same  
 Disposal of Discharged Water: 55 gal Drum CSA

### INSTRUMENTS (Indicate make, model, I.D.)

Water Level: Keck  
 pH Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Conductivity Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Thermometer: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Turbidimeter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 ORP Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 DO Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Other: Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Time	Cum. Vol. (L)	Purge Rate (L/m)	Temp. (oC)	pH	Spec. Cond. (µS/cm)	Color	Turbidity & Sediment (NTU)	ORP (mV)	DO (mg/L)	Water Level (ft BMP)	Remarks
11:19	-	-	-	-	-	-	-	-	-	5.20	Pump is on
11:29	2	0.2	26.31	6.58	1910	Neutral	109	45	0.00	5.75	
11:34	3	↓	26.52	6.58	1890	↓	89.7	38	0.00	5.82	
11:39	4	↓	26.68	6.57	1890	↓	77.6	34	0.00	5.91	
11:42	4.5	↓	26.68	6.59	1880	↓	64.9	33	0.00	5.92	
11:45	5	↓	26.70	6.59	1870	↓	89.2	33	0.00	5.93	

Water Level (ft. BMP) at End of Purge: 5.93

Sample Intake Depth (ft. BMP): 16.85

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
11:47	1 L	G	2	N	Heat	SiOCs
11:47	20 mL	P	1	Y	HNO <sub>3</sub>	Metals - As
11:47	40 mL	G	6	N	HCL	VOCs / TPH

Comments:



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# GROUNDWATER SAMPLING RECORD

Project Number: 1919232 Project Name: HWPW Date: 3.12.20  
 Sample Number: W61-1620-MW98A-20200312 Starting Water Level (ft. BMP): 7.19  
 Sampling Location (well ID, etc.): MW-98A Casing Stickup (ft.): -  
 Sampled by: SAS Starting Water Level (ft. BGL): -  
 Measuring Point (MP) of Well: TOC/PVC Total Depth (ft. BMP): 20.40  
 Screened Interval (ft. BGL): 10.5-20.5 Casing Diameter (in ID): 2  
 Filter Pack Interval (ft. BGL): 8.5-21 Casing Volume (gal.): 2.16 = 8.18 L

## QUALITY ASSURANCE

### METHODS (describe):

Cleaning Equipment: Dedicated Equipment  
 Purging: Peristaltic or bladder pump with dedicated tubing Sampling: Same  
 Disposal of Discharged Water: 55 gal drum CSA

### INSTRUMENTS (Indicate make, model, I.D.)

Water Level: Keck  
 pH Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Conductivity Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Thermometer: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Turbidimeter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 ORP Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 DO Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Other: Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Time	Cum. Vol. (L)	Purge Rate (L/m)	Temp. (oC)	pH	Spec. Cond. (µS/cm)	Color	Turbidity & Sediment (NTU)	ORP (mV)	DO (mg/L)	Water Level (ft BMP)	Remarks
12:26	-	0.2	-	-	-	-	-	-	-	7.19	Pump is on
12:36	2	↓	30.18	6.89	1810	Neutral	76.7	-88	0.00	7.91	
12:41	3	↓	30.01	6.89	1820	↓	63.3	-92	0.00	7.95	
12:46	4	↓	29.76	6.87	1850	↓	48.7	-94	0.00	7.97	
12:49	4.5	↓	29.66	6.86	1860	↓	45.0	-95	0.00	7.98	

Water Level (ft. BMP) at End of Purge: 7.98 Sample Intake Depth (ft. BMP): 17.40

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
12:54	1 L	G	2	N	Neat	JVOCs
12:51	120 mL	P	1	Y	HNO3	Metals - As
12:54	40 mL	G	6	N	HCL	VOCs / TPH

Comments:



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# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: HWPW Date: 3.12.20  
 Sample Number: W6-1620-MW98B-20200312 Starting Water Level (ft. BMP): 8.30  
 Sampling Location (well ID, etc.): MW-98B Casing Stickup (ft.): -  
 Sampled by: *[Signature]* Starting Water Level (ft. BGL): 35 -  
 Measuring Point (MP) of Well: TOC/PVC Total Depth (ft. BMP): 39.80  
 Screened Interval (ft. BGL): 35-40 Casing Diameter (in ID): 2  
 Filter Pack Interval (ft. BGL): 31-40 Casing Volume (gal.): 5.14 = 19.46 L

## QUALITY ASSURANCE

### METHODS (describe):

Cleaning Equipment: Dedicated Equipment

Purging: Peristaltic or bladder pump with dedicated tubing Sampling: Same

Disposal of Discharged Water: 55 gal Drum CSB

### INSTRUMENTS (Indicate make, model, I.D.)

Water Level: Keck

pH Meter: Horiba U-52

Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Conductivity Meter: Horiba U-52

Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Thermometer: Horiba U-52

Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Turbidimeter: Horiba U-52

Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

ORP Meter: Horiba U-52

Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

DO Meter: Horiba U-52

Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution

Other:

Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Time	Cum. Vol. (L)	Purge Rate (L/m)	Temp. (oC)	pH	Spec. Cond. (µS/cm)	Color	Turbidity & Sediment (NTU)	ORP (mV)	DO (mg/L)	Water Level (ft BMP)	Remarks
13:25	-	0.2	-	-	-	-	-	-	-	8.30	Pump is on
13:35	2	↓	24.26	7.11	991	Neutral	30.7	74	0.42	11.30	
13:40	3	↓	24.78	7.12	981	↓	25.7	73	0.37	12.25	
13:45	4	↓	24.03	7.17	993	↓	4.2	74	0.37	13.81	
13:48	4.5	↓	24.17	7.12	991	↓	6.2	75	0.44	14.89	
13:51	5	↓	24.32	7.12	987	↓	6.8	75	0.47	15.23	

Water Level (ft. BMP) at End of Purge: 15.23

Sample Intake Depth (ft. BMP): 36.80

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y / N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
13:50	1 L	G	2	N	Neat	SVOcs
13:50	120 mL	P	1	N	HNO3	Metals-As
13:50	40 mL	G	6	N	HCL	VOCs / TPH

Comments:



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# GROUNDWATER SAMPLING RECORD

PAGE 1 of 1

Project Number: 19119237 Project Name: HWPW Date: 3.12.20  
 Sample Number: WGI-1620-MW50B-20200312 Starting Water Level (ft. BMP): 8.37  
 Sampling Location (well ID, etc.): 50B Casing Stickup (ft.): -  
 Sampled by: SB Starting Water Level (ft. BGL): -  
 Measuring Point (MP) of Well: TOC/PVC Total Depth (ft. BMP): 39.5 - Silty bottom  
 Screened Interval (ft. BGL): 34.5 - 39.5 Casing Diameter (in ID): 2  
 Filter Pack Interval (ft. BGL): 31 - 40 Casing Volume (gal.): 4.92 = 18.62 L

## QUALITY ASSURANCE

### METHODS (describe):

Cleaning Equipment: Dedicated Equipment  
 Purging: Peristaltic or bladder pump with dedicated tubing Sampling: Same  
 Disposal of Discharged Water: 55 gal Drum CSA

### INSTRUMENTS (Indicate make, model, I.D.)

Water Level: Keck  
 pH Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Conductivity Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Thermometer: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Turbidimeter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 ORP Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 DO Meter: Horiba U-52 Field Calibration: Auto Calibration - 100-4 Horiba Calibration Solution  
 Other: Filter / Filter Size: 10 micron

## SAMPLING MEASUREMENTS

Time	Cum. Vol. (L)	Purge Rate (L/m)	Temp. (oC)	pH	Spec. Cond. (µS/cm)	Color	Turbidity & Sediment (NTU)	ORP (mV)	DO (mg/L)	Water Level (ft BMP)	Remarks
14:49	-	0.2	-	-	-	-	-	-	-	8.37	Pump is on
14:59	2	↓	30.75	6.98	1620	cloudy	360	26	2.72	11.65	
15:04	3	↓	30.53	6.96	1620	Neutral	83.8	32	2.84	12.93	
15:09	4	↓	30.43	6.97	1620	↓	60.3	22	2.82	14.10	
15:12	4.5	↓	30.53	6.97	1620	↓	87.2	19	2.76	14.57	
15:15	5	↓	30.11	6.97	1630	↓	50.1	13	2.65	15.10	
15:18	5.5	↓	30.27	6.97	1630	↓	54.7	9	2.50	15.50	

Water Level (ft. BMP) at End of Purge: 15.50

Sample Intake Depth (ft. BMP): 35.5

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation (type)	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
15:20	1 L	G	2	N	Neut	SVOCs
15:20	120 mL	P	1	Y	HNO <sub>3</sub>	Metals-As
15:20	40 mL	G	6	N	HCL	VOCs/TPH

Comments: Very Silty Bottom



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**GROUNDWATER SAMPLING RECORD**

Project Number: 19119232 Project Name: HWPW Date: 3-20-2020  
 Sample Number: UG-1620 MW 60 AR 20200320 Starting Water Level (ft. BMP): 8.03  
 Sampling Location (well ID, etc.): MW 60 AR Casing Stickup (ft.): 0  
 Sampled by: T.M. McSpodden Starting Water Level (ft. BGL): 8.03  
 Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 29.60  
 Screened Interval (ft. BGL): Casing Diameter (In ID): 2"  
 Filter Pack Interval (ft. BGL): Casing Volume (gal.): 3.4

**QUALITY ASSURANCE**

METHODS (describe): Low Flow  
 Cleaning Equipment: \_\_\_\_\_  
 Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_  
 Disposal of Discharged Water: \_\_\_\_\_

**INSTRUMENTS (Indicate make, model, I.d.)**

Water Level: Horiba D2106 Thermometer: Horiba C1122  
 pH Meter: Horiba C1122 Field Calibration: \_\_\_\_\_  
 Conductivity Meter: Horiba C1122 Field Calibration: Ph buffer 4.0 - Field Cal = 4.0  
 Filter / Filter Size: 10 MICR Other: Geo Pump P1122

**SAMPLING MEASUREMENTS**

Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
8:00	8.03	.2	.2	23.96	5.90	1.08	1.48	190	249	
8:10	8.51	.3	.2	23.67	6.09	1.05	1.18	164	139	
8:15	8.33	.4	.2	23.52	6.17	1.04	0.92	145	142	
8:20										
8:25	Rain delay		20 minutes							
8:30										
8:35	8.51	.5	.2	23.63	6.82	0.999	1.02	103	147	
8:40	8.23	.6	.2	23.38	6.73	1.01	1.00	78	103	
8:45	8.24	.7	.2	23.26	6.72	1.01	1.14	78	74.8	
8:50	8.22	.8	.2	23.15	6.70	1.01	1.02	65	68.9	
8:55	8.23	.9	.2	23.07	6.68	1.01	0.80	56	67.1	

Water Level (ft. BMP) at End of Purge: 8.23 Sample Intake Depth (ft. BMP): 26.

**SAMPLE INVENTORY**

Time	Bottles Collected			Filtration (Y) (N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
8:55		Pump	9	10 MICR		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



# GROUNDWATER SAMPLING RECORD

Project Number: 19119232 Project Name: HWPW Date: 3-20-2020

Sample Number: WG-1620 MW 47A 20200320 Starting Water Level (ft. BMP): 9.26

Sampling Location (well ID, etc.): MW 47A Casing Stickup (ft.): 0

Sampled by: TIM McSpadden Starting Water Level (ft. BGL): 9.26

Measuring Point (MP) of Well: TOC Total Depth (ft. BMP): 2.5-03

Screened Interval (ft. BGL): Casing Diameter (In ID): 2'

Filter Pack Interval (ft. BGL): Casing Volume (gal.): 2.5

## QUALITY ASSURANCE

METHODS (describe): Low Flow

Cleaning Equipment: \_\_\_\_\_

Purging: \_\_\_\_\_ Sampling: \_\_\_\_\_

Disposal of Discharged Water: \_\_\_\_\_

## INSTRUMENTS (Indicate make, model, I.d.)

Water Level: Herow D2106 Thermometer: Horiba C1122

pH Meter: Horiba C1122 Field Calibration: \_\_\_\_\_

Conductivity Meter: Horiba C1122 Field Calibration: Ph buffer 4.0 Field cal: 4.0

Filter / Filter Size: \_\_\_\_\_ Other: Geo pump P1122

## SAMPLING MEASUREMENTS


Time	Water Depth (ft BMP)	Cum. Vol. (gal. or L)	Purge Rate (L/m)	Temp. (°C)	pH (S.U.)	Spec. Cond. (mS/cm)	D.O. (mg/L)	Redox (mV)	Turbidity (NTU)	Color
--	--	--	--	± 3%	± 0.1	± 3%	± 10% if >0.5	± 10	± 10%	--
9:50	9.26	.2	.2	21.69	6.99	0.851	9.62	109	169	
10:00	9.37	.3	.2	22.03	7.18	0.856	2.21	119	81.4	
10:05	9.29	.4	.2	22.14	6.89	0.826	0.80	112	69.7	
10:10	9.27	.5	.2	22.17	6.84	0.863	0.42	109	58.4	
10:15	9.28	.6	.2	22.19	6.80	0.863	0.33	105	59.6	
10:20	9.26	.7	.2	22.29	6.78	0.859	0.38	101	86.4	
10:25	9.31	.8	.2	22.38	6.79	0.859	0.30	99	117	
10:30	9.32	.9	.2	22.39	6.78	0.858	0.24	100	97.3	

Water Level (ft. BMP) at End of Purge: \_\_\_\_\_ Sample Intake Depth (ft. BMP): 22.

## SAMPLE INVENTORY

Time	Bottles Collected			Filtration (Y/N)	Preservation	Remarks (quality control sample, other)
	Volume	Composition (G, P)	No.			
10:30		Purge	9			

Comments: \_\_\_\_\_



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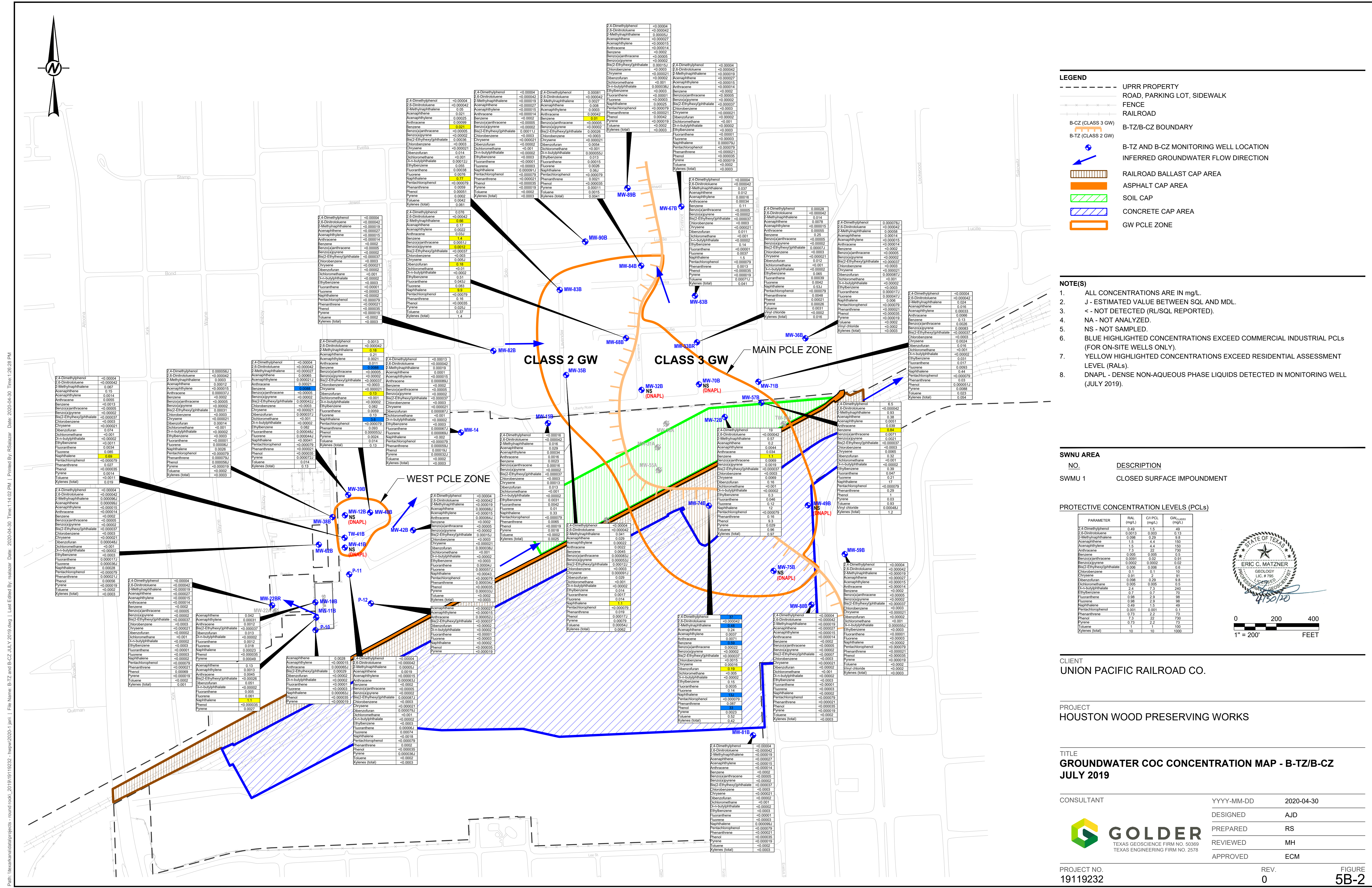
**ATTACHMENT D**

**July 2019 Sampling Event COC  
Concentration Maps**









Path: \\usar\ar\dat\reports\ground\cok\_2019\01\19232\_inwp\2020-1.jm... | File Name: B-TZ and B-CZ JULY 2019.dwg | L: User: E:\res\res\res | Date: 2020-04-30 | Time: 14:02 PM | Printed By: R.Saizab | Date: 2020-04-30 | Time: 2:28:25 PM

**LEGEND**

- UPRR PROPERTY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-CZ (CLASS 3 GW)
- B-TZ (CLASS 2 GW)
- B-TZ AND B-CZ MONITORING WELL LOCATION
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA
- GW PCLE ZONE

- NOTE(S)**
- ALL CONCENTRATIONS ARE IN mg/L.
  - J - ESTIMATED VALUE BETWEEN SQL AND MDL.
  - < - NOT DETECTED (RL/SQL REPORTED).
  - NA - NOT ANALYZED.
  - NS - NOT SAMPLED.
  - BLUE HIGHLIGHTED CONCENTRATIONS EXCEED COMMERCIAL INDUSTRIAL PCLs (FOR ON-SITE WELLS ONLY).
  - YELLOW HIGHLIGHTED CONCENTRATIONS EXCEED RESIDENTIAL ASSESSMENT LEVEL (RALs).
  - DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).

**SWMU AREA**

NO.	DESCRIPTION
SWMU 1	CLOSED SURFACE IMPONDEMENT

**PROTECTIVE CONCENTRATION LEVELS (PCLs)**

PARAMETER	RAL (mg/L)	CI PCL (mg/L)	GW <sub>USE</sub> (mg/L)
2,4-Dimethylphenol	0.49	1.5	49
2,6-Dinitrotoluene	0.00113	0.001	0.13
2-Methylnaphthalene	0.006	0.20	9.8
Acenaphthene	1.5	4.4	150
Acenaphthylene	1.5	4.4	150
Anthracene	7.3	22	730
Benzo(a)anthracene	0.0001	0.002	0.01
Benzo(a)pyrene	0.0001	0.002	0.01
Bis(2-Ethylhexyl)phthalate	0.006	0.006	0.6
Chlorobenzene	0.1	0.1	10
Chrysene	0.01	0.01	0.1
Dibenzofuran	0.006	0.20	9.8
Dibenzothiophene	0.005	0.005	0.5
Di-n-butylphthalate	2.4	7.3	240
Fluorene	0.08	2.9	98
Fluoranthene	0.08	2.9	98
Furan	0.001	0.001	0.1
Naphthalene	0.49	1.5	49
Phenanthrene	0.001	0.001	0.1
Phenol	7.3	22	730
Pyrene	0.73	2.2	73
Toluene	1.73	1	100
Xylenes (total)	10	10	1000

**CLIENT**  
UNION PACIFIC RAILROAD CO.

**PROJECT**  
HOUSTON WOOD PRESERVING WORKS

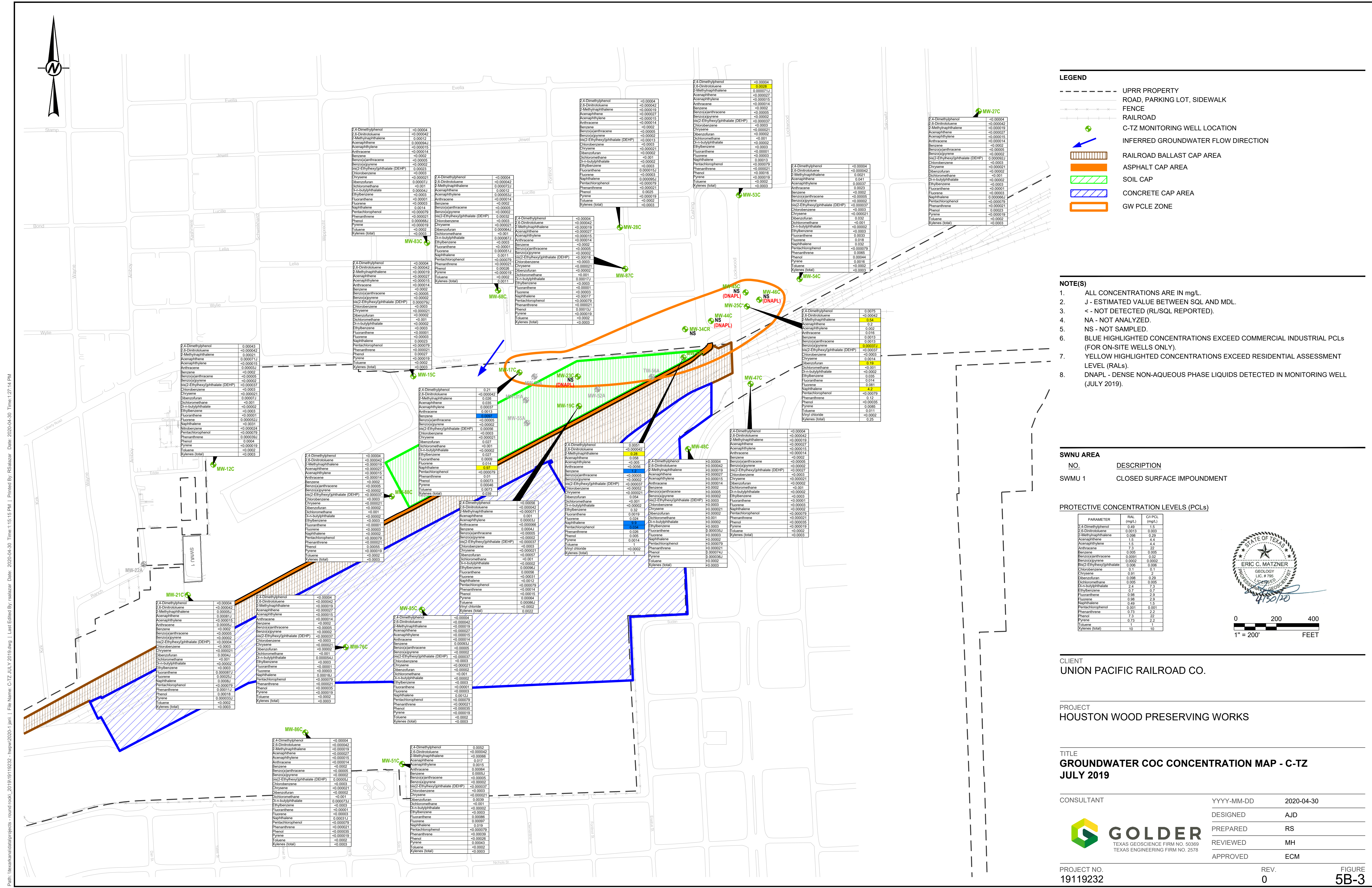
**TITLE**  
GROUNDWATER COC CONCENTRATION MAP - B-TZ/B-CZ  
JULY 2019

CONSULTANT	DATE
YYYY-MM-DD	2020-04-30
DESIGNED	AJD
PREPARED	RS
REVIEWED	MH
APPROVED	ECM

PROJECT NO. 19119232 REV. 0 FIGURE 5B-2

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM A NS/D





**LEGEND**

- UPRR PROPERTY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- INFERRED GROUNDWATER FLOW DIRECTION
- ▨ RAILROAD BALLAST CAP AREA
- ▨ ASPHALT CAP AREA
- ▨ SOIL CAP
- ▨ CONCRETE CAP AREA
- ▨ GW PCLE ZONE

- NOTE(S)**
- ALL CONCENTRATIONS ARE IN mg/L.
  - J - ESTIMATED VALUE BETWEEN SQL AND MDL.
  - < - NOT DETECTED (RL/SQL REPORTED).
  - NA - NOT ANALYZED.
  - NS - NOT SAMPLED.
  - BLUE HIGHLIGHTED CONCENTRATIONS EXCEED COMMERCIAL INDUSTRIAL PCLs (FOR ON-SITE WELLS ONLY).
  - YELLOW HIGHLIGHTED CONCENTRATIONS EXCEED RESIDENTIAL ASSESSMENT LEVEL (RALs).
  - DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).

**SWNU AREA**

NO.	DESCRIPTION
SWMU 1	CLOSED SURFACE IMPONDEMENT

**PROTECTIVE CONCENTRATION LEVELS (PCLs)**

PARAMETER	RAL (mg/L)	CI PCL (mg/L)
2,4-Dimethylphenol	0.49	1.5
2,6-Dinitrotoluene	0.013	0.03
2-Methylnaphthalene	0.988	0.29
Acenaphthylene	1.5	4.4
Benzo(a)anthracene	0.005	0.005
Benzo(a)pyrene	0.001	0.02
Bis(2-Ethylhexyl)phthalate (DEHP)	0.002	0.002
Chlorobenzene	0.1	0.1
Chrysene	0.01	0.1
Dibenzofuran	0.005	0.005
Dibenzothiazene	0.7	0.7
Dibenzotriazole	2.4	7.3
Ethylbenzene	0.98	2.9
Fluorene	0.88	2.8
Fluoranthene	0.98	2.9
Naphthalene	0.49	1.5
Pentachlorophenol	0.001	0.001
Phenanthrene	0.73	2.2
Phenol	7.3	22
Pyrene	0.73	2.2
Toluene	1	1
Xylenes (total)	10	10



CLIENT  
UNION PACIFIC RAILROAD CO.

PROJECT  
HOUSTON WOOD PRESERVING WORKS

TITLE  
GROUNDWATER COC CONCENTRATION MAP - C-TZ  
JULY 2019

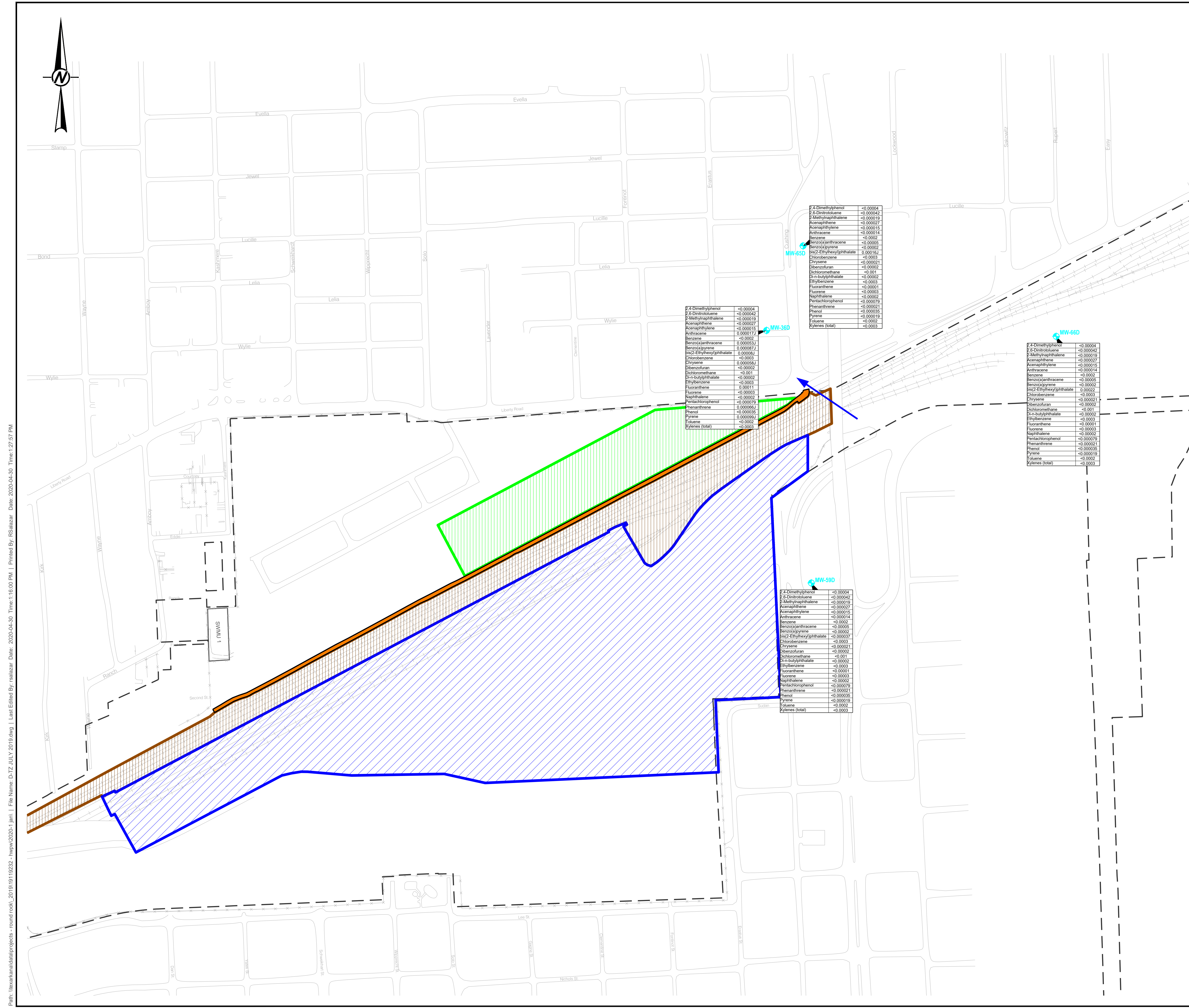
CONSULTANT	DATE
YYYY-MM-DD	2020-04-30
DESIGNED	AJD
PREPARED	RS
REVIEWED	MH
APPROVED	ECM

PROJECT NO. 19119232 REV. 0 FIGURE 5B-3

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IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM A NS/D





**LEGEND**

- UPRR PROPERTY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- D-TZ MONITORING WELL LOCATION
- INFERRED GROUNDWATER FLOW DIRECTION
- [Hatched] RAILROAD BALLAST CAP AREA
- [Orange] ASPHALT CAP AREA
- [Green] SOIL CAP
- [Blue] CONCRETE CAP AREA

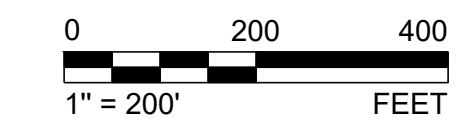
- NOTE(S)**
- ALL CONCENTRATIONS ARE IN mg/L.
  - J - ESTIMATED VALUE BETWEEN SQL AND MDL.
  - < - NOT DETECTED (RL/SQL REPORTED).
  - NA - NOT ANALYZED.
  - NS - NOT SAMPLED.
  - BLUE HIGHLIGHTED CONCENTRATIONS EXCEED COMMERCIAL INDUSTRIAL PCLs (FOR ON-SITE WELLS ONLY).
  - YELLOW HIGHLIGHTED CONCENTRATIONS EXCEED RESIDENTIAL ASSESSMENT LEVEL (RALs).
  - DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).

**SWNU AREA**

NO.	DESCRIPTION
SWMU 1	CLOSED SURFACE IMPOUNDMENT

**PROTECTIVE CONCENTRATION LEVELS (PCLs)**

PARAMETER	RAL (mg/L)	CI PCL (mg/L)
2,4-Dimethylphenol	0.49	1.5
2,6-Dinitrotoluene	0.0013	0.003
2-Methylnaphthalene	0.098	0.29
Acenaphthylene	1.5	4.4
Acenaphthylene	1.5	4.4
Anthracene	7.3	22
Benzene	0.005	0.005
Benzo(a)anthracene	0.0001	0.02
Benzo(a)pyrene	0.0002	0.0002
Benzo(b)fluoranthene	0.0001	0.0001
Benzo(k)fluoranthene	0.0001	0.0001
Chlorobenzene	0.1	0.1
Chrysene	0.01	0.01
Dibenzofuran	0.098	0.29
Dibenzomethane	0.095	0.095
Di-n-butylphthalate	2.4	7.3
Ethylbenzene	0.7	0.7
Fluoranthene	0.98	2.9
Fluorene	0.98	2.9
Naphthalene	0.49	1.5
Pentachlorophenol	0.001	0.001
Phenanthrene	0.73	2.2
Phenol	7.3	22
Pyrene	1.73	2.2
Toluene	1	1
Xylenes (total)	10	10



CLIENT  
UNION PACIFIC RAILROAD CO.

PROJECT  
HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - D-TZ**  
**JULY 2019**

CONSULTANT

YYYY-MM-DD	2020-04-30
DESIGNED	AJD
PREPARED	RS
REVIEWED	MH
APPROVED	ECM

PROJECT NO. 19119232      REV. 0      FIGURE 5B-4

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1" = 200'      IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM A3.









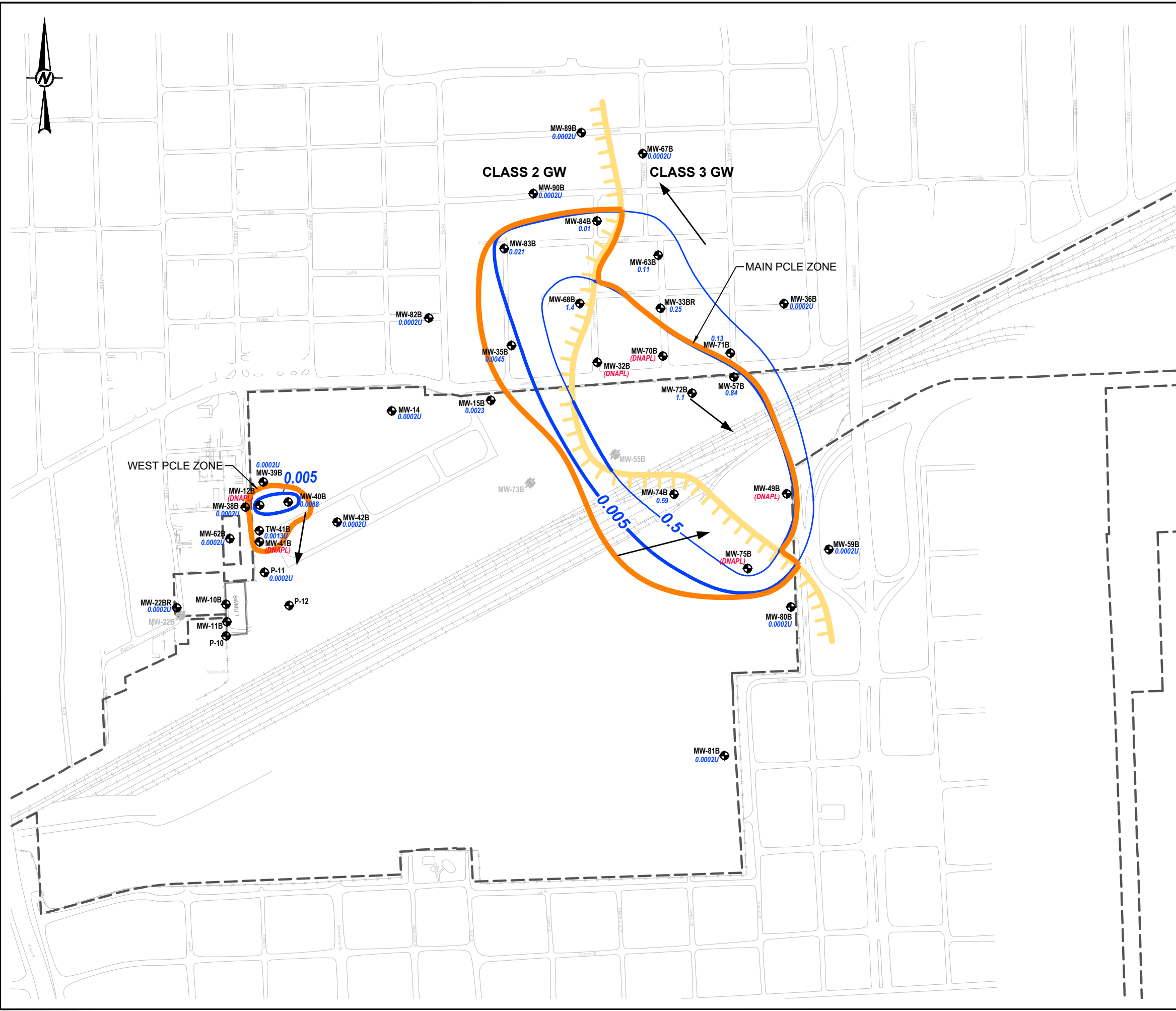








Path: \\uswest\erdm\dat\proj\houston\ground\_coc\_2019\01\19232 - houston\01\19232 - houston\01\19232.dwg | File Name: FIG 5B-10 - B-TZ Benzene Map | Last Edited By: erdm | Printed By: Release | Date: 2020-04-30 | Time: 14:13 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW) B-TZ/B-CZ BOUNDARY
- B-TZ (CLASS 2 GW) **0.346** BENZENE CONCENTRATION (mg/L)
- 0.005** BENZENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

**NOTE(S)**

1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
2. NS - NOT SAMPLED
3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
 CLASS 2 GW PCL: 0.005 mg/L  
 CLASS 3 GW PCL: 0.5 mg/L

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

**TITLE**  
 GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 BENZENE - JULY 2019

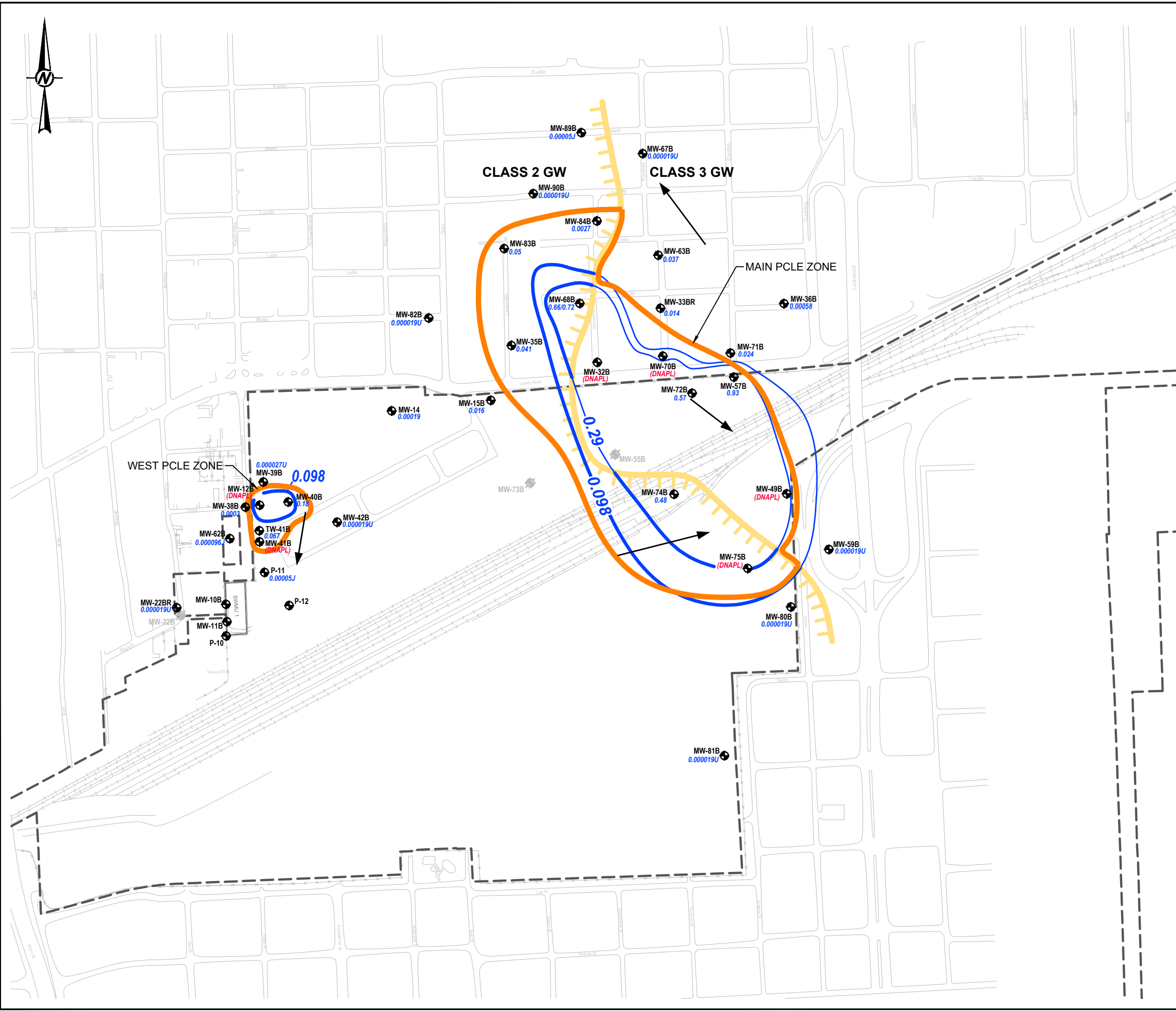
CONSULTANT	DATE
<b>GOLDER</b> <small>TEXAS GEOSCIENCE FIRM NO. 50369 TEXAS ENGINEERING FIRM NO. 2578</small>	YYYY-MM-DD 2020-04-30
	DESIGNED AJD
	PREPARED AJD
	REVIEWED MH
	APPROVED ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5B-10

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erdm\data\projects\houston\cok\_2019\01\19232 - houston\01\19232 - houston\01\19232.dwg | File Name: FIG 5B-12 - BTZ\_2-Methylnaphthalene.dwg | Last Edited By: malarik | Date: 2020-04-30 | Time: 1:33:17 PM | Printed By: RSebastian | Date: 2020-04-30 | Time: 1:43:19 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW) B-TZ/B-CZ BOUNDARY
- B-TZ (CLASS 2 GW) **0.276** 2-METHYLNAPHTHALENE CONCENTRATION (mg/L)
- 0.098** 2-METHYLNAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
 CLASS 2 GW PCL: 0.098 mg/L (RES.) & 0.29 mg/L (C/I)  
 CLASS 3 GW PCL: 9.8 mg/L (RES.) & 29 mg/L (C/I)

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 2-METHYLNAPHTHALENE - JULY 2019**

CONSULTANT	DATE	BY
GOLDER	2020-04-30	AJD
DESIGNED		AJD
PREPARED		AJD
REVIEWED		MH
APPROVED		ECM

PROJECT NO. 19119232      REV. 0      FIGURE 5B-12

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

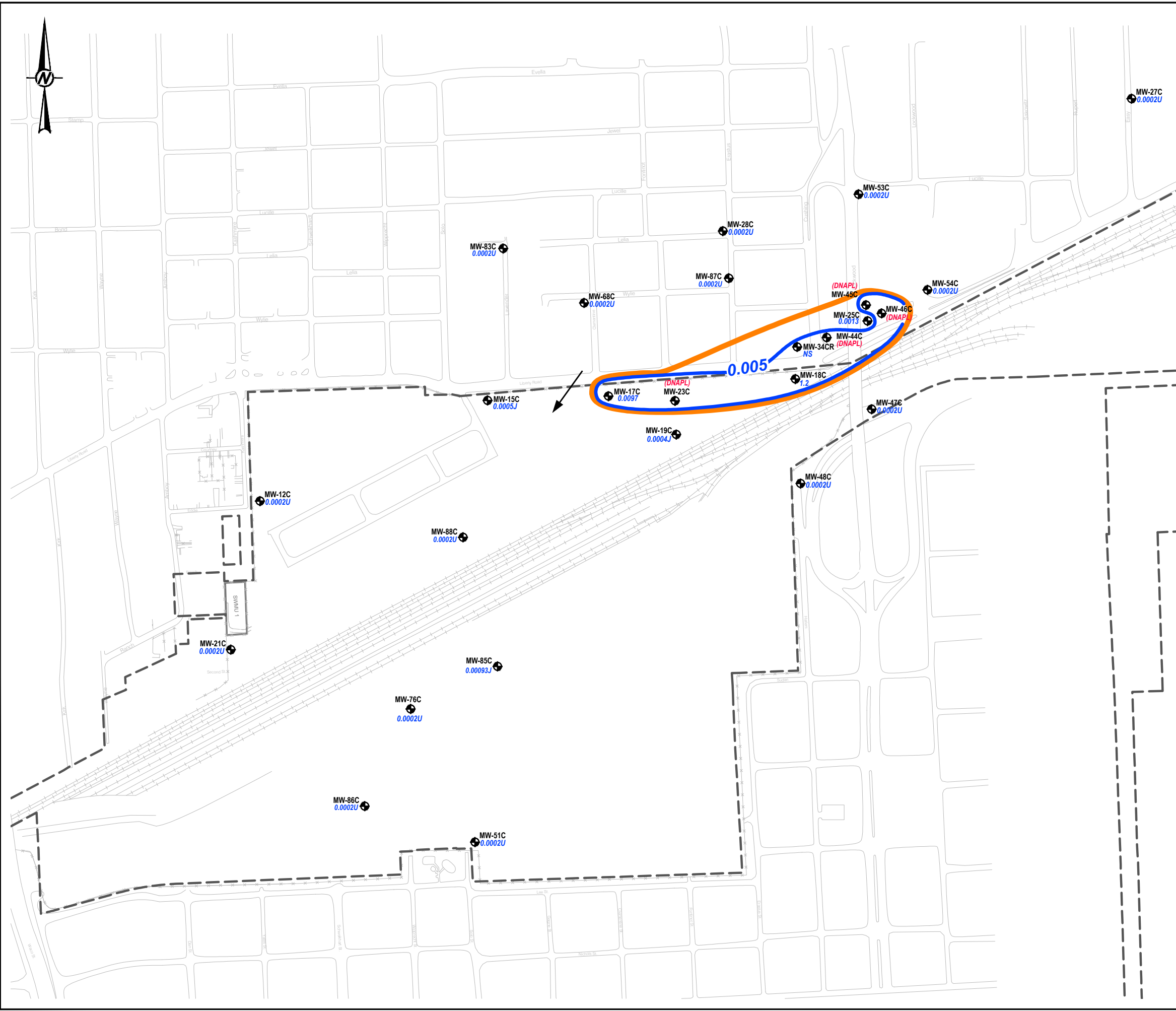








Path: \\uswest\erdm\data\projects\19119232 - houston\2019\01\19232 - houston\2019\04\12\Draw1 - Benzene.dwg | File Name: FIG 5B-15 - C-TZ Benzene Map | Last Edited By: rbarbar | Date: 2020-04-30 Time: 1:30:01 PM | Printed By: RBarbar | Date: 2020-04-30 Time: 1:46:26 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- - - - - FENCE
- - - - - RAILROAD
- ⊕ C-TZ MONITORING WELL LOCATION
- 8.01 BENZENE CONCENTRATION (mg/L)
- 0.005** BENZENE CONCENTRATION CONTOUR (mg/L)
- ↖ INFERRED GROUNDWATER FLOW DIRECTION
- ▭ GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.005 mg/L).

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

STATE OF TEXAS  
 ERIC C. MATZNER  
 GEOLOGY  
 LIC. # 795  
 LICENSED  
 PROFESSIONAL GEOSCIENTIST  
 4/30/20

0 200 400  
 1" = 400' FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - C-TZ  
 BENZENE - JULY 2019**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED	AJD	
PREPARED	AJD	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-15

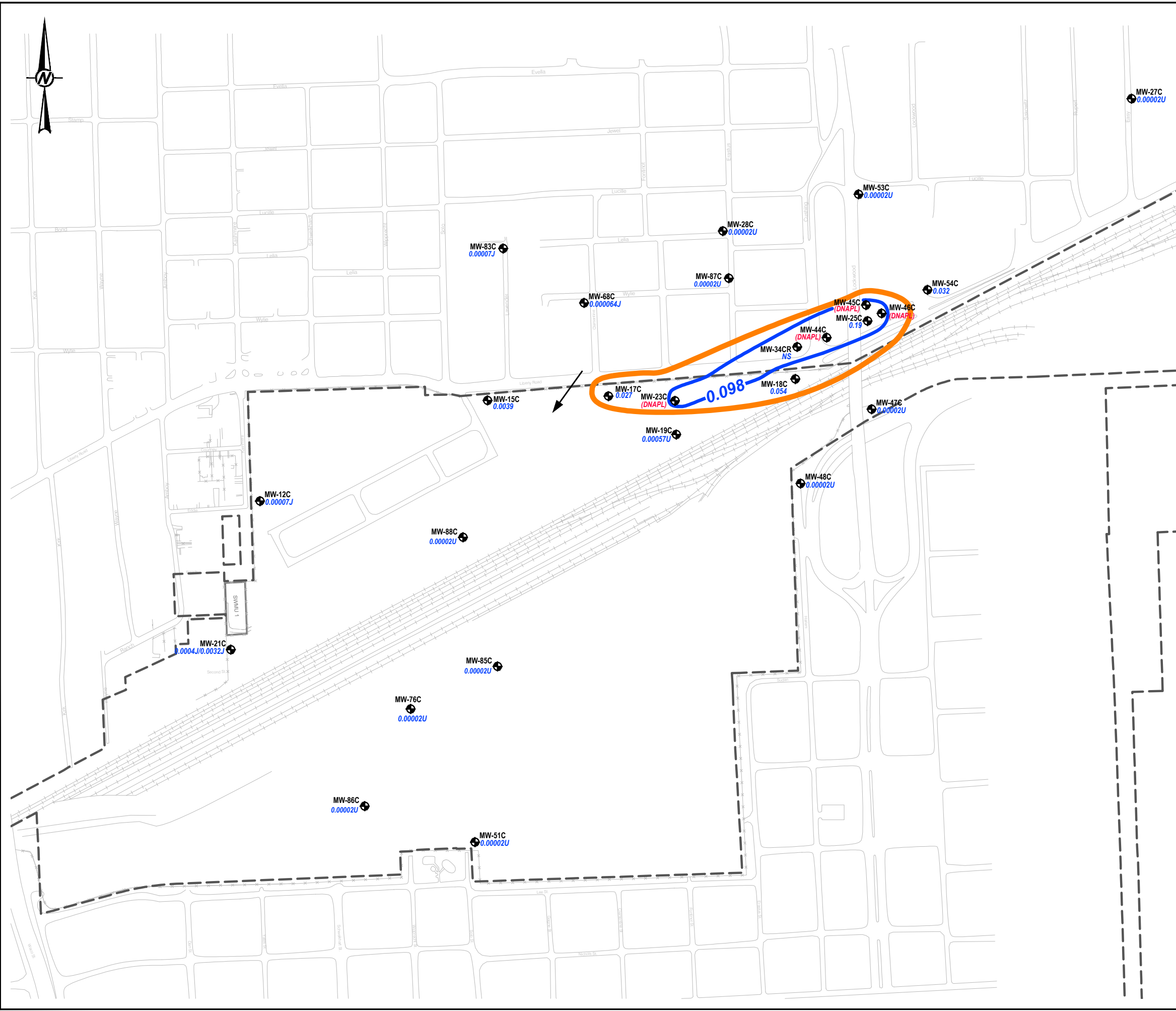
1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B







Path: \\uswest\erdm\dat\proj\19119232 - houston\2019-04-12\Draw1 - houston\2019-04-12\Draw1 - File Name: FIG 5B-18 - C-TZ - Dibenzofuran.cdw | Last Edited By: matzner Date: 2020-04-30 Time: 1:42:21 PM | Printed By: RSalazar Date: 2020-04-30 Time: 1:43:05 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- 0.207 DIBENZOFURAN CONCENTRATION (mg/L)
- DIBENZOFURAN CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDDED AT THE RAL AND C/I PCL (0.098 mg/L AND 0.29 mg/L, RESPECTIVELY).

**REFERENCE(S)**  
BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

0 200 400  
1" = 400' FEET

CLIENT  
UNION PACIFIC RAILROAD CO.

PROJECT  
HOUSTON WOOD PRESERVING WORKS

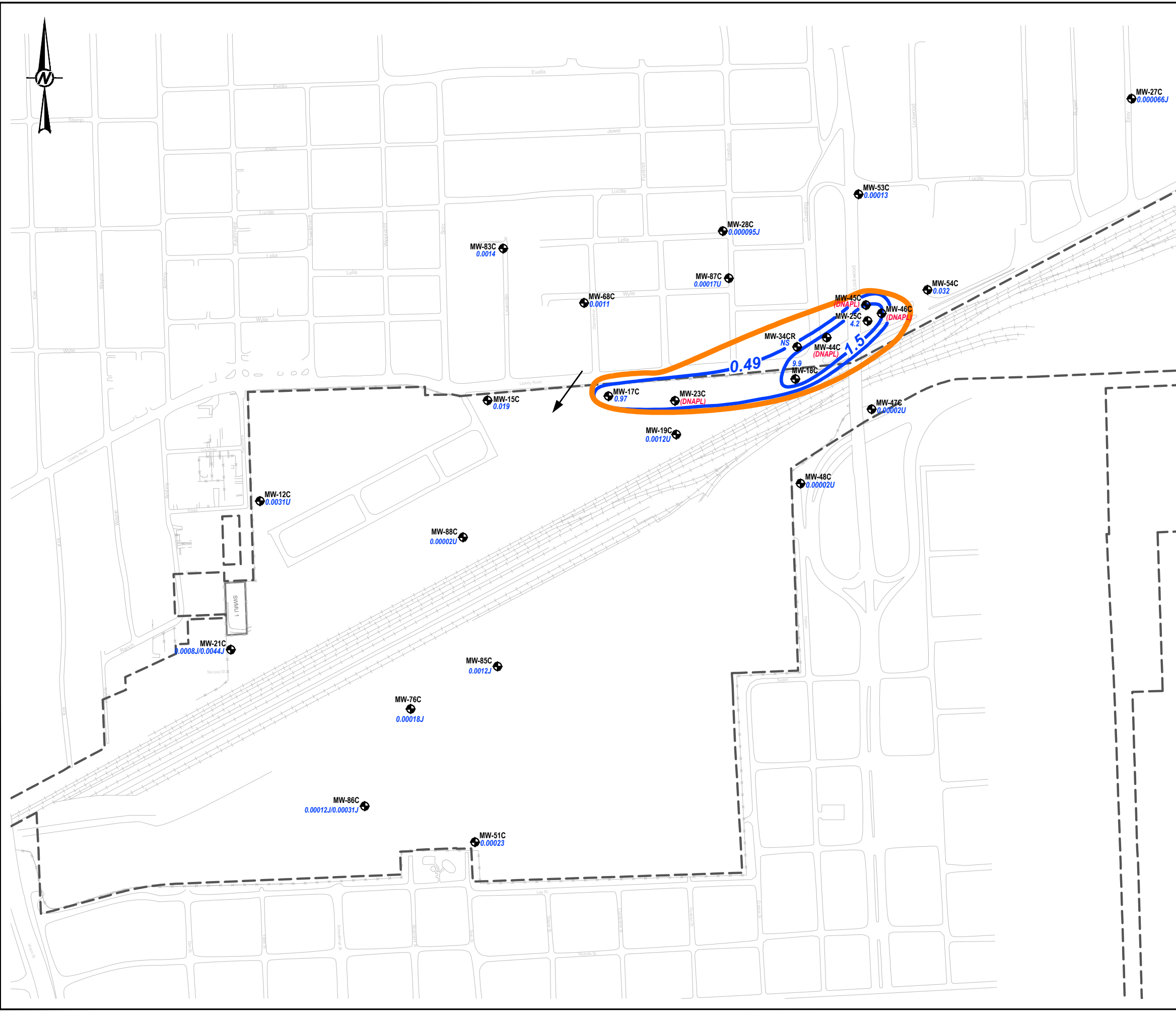
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - C-TZ  
DIBENZOFURAN - JULY 2019**

CONSULTANT	YYYY-MM-DD	2020-04-30
<b>GOLDER</b> <small>TEXAS GEOSCIENCE FIRM NO. 50369 TEXAS ENGINEERING FIRM NO. 2578</small>	DESIGNED	AJD
	PREPARED	AJD
	REVIEWED	MH
	APPROVED	ECM

PROJECT NO. 19119232      REV. 0      FIGURE 5B-18

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\ermsouthwest\projects\ground\cok\2019\01\19232 - hwp\2019\04\12\Desk1 - hwp\2019\04\12\Desk1 - File Name: FIG 5B-10 - C-TZ Naphthalene.docx | Last Edited By: makbar | Date: 2020-04-30 | Time: 1:43:57 PM | Printed By: Release | Date: 2020-04-30 | Time: 1:40:38 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- 14.7 NAPHTHALENE CONCENTRATION (mg/L)
- NAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.49 mg/L AND 1.5 mg/L, RESPECTIVELY).

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

STATE OF TEXAS  
 ERIC C. MATZNER  
 GEOLOGY  
 LIC. # 795  
 LICENSED  
 PROFESSIONAL GEOSCIENTIST  
 4/30/20

0      200      400  
 1" = 400'      FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

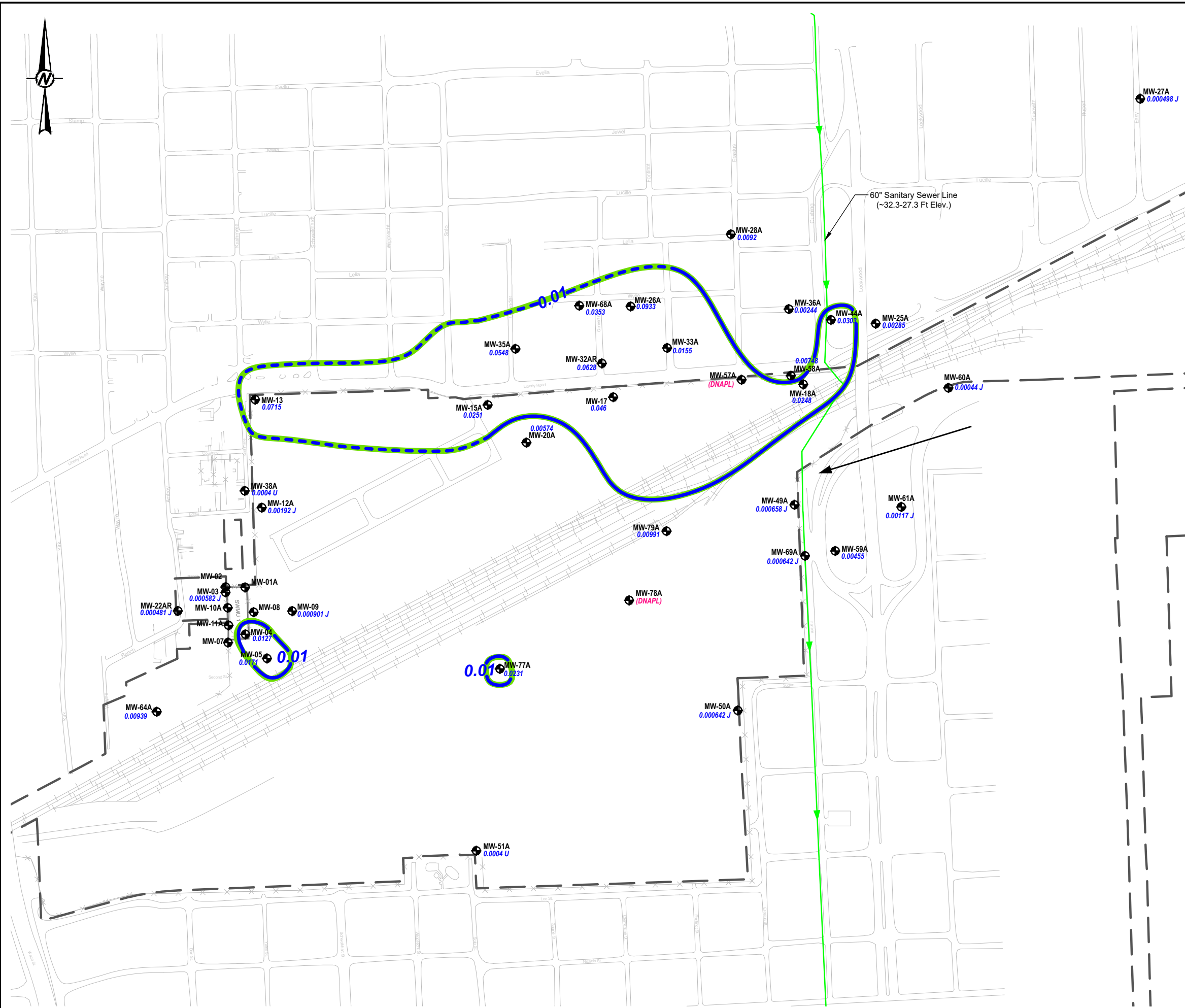
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - C-TZ  
 NAPHTHALENE - JULY 2019**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED	AJD	
PREPARED	AJD	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-19

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erdm\data\projects\houston\cok\_2019\01\10202 - Inpwr\2020-1\jmk | File Name: ARSENIC-A-TZ JULY 2019.dwg | Last Edited By: mshawn | Date: 2020-04-30 | Time: 1:46:14 PM | Printed By: RSealar | Date: 2020-04-30 | Time: 1:51:01 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARLING LOT, SIDEWALK
- FENCE
- RAILROAD
- A-TZ MONITORING WELL LOCATION
- 0.108** ARSENIC CONCENTRATION (mg/L)
- 0.01** ARSENIC CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- ARSENIC PCLE ZONE

**NOTE(S)**

1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
2. NS - NOT SAMPLED
3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.01 mg/L).

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1. DATED JUNE 2004.

STATE OF TEXAS  
 ERIC C. MATZNER  
 GEOLOGY  
 LIC. # 795  
 LICENSED  
 PROFESSIONAL GEOSCIENTIST  
 4/30/20

0 200 400  
 1" = 400'  
 FEET

**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

**TITLE**  
 GROUNDWATER COC CONCENTRATION MAP - A-TZ  
 ARSENIC - JULY 2019

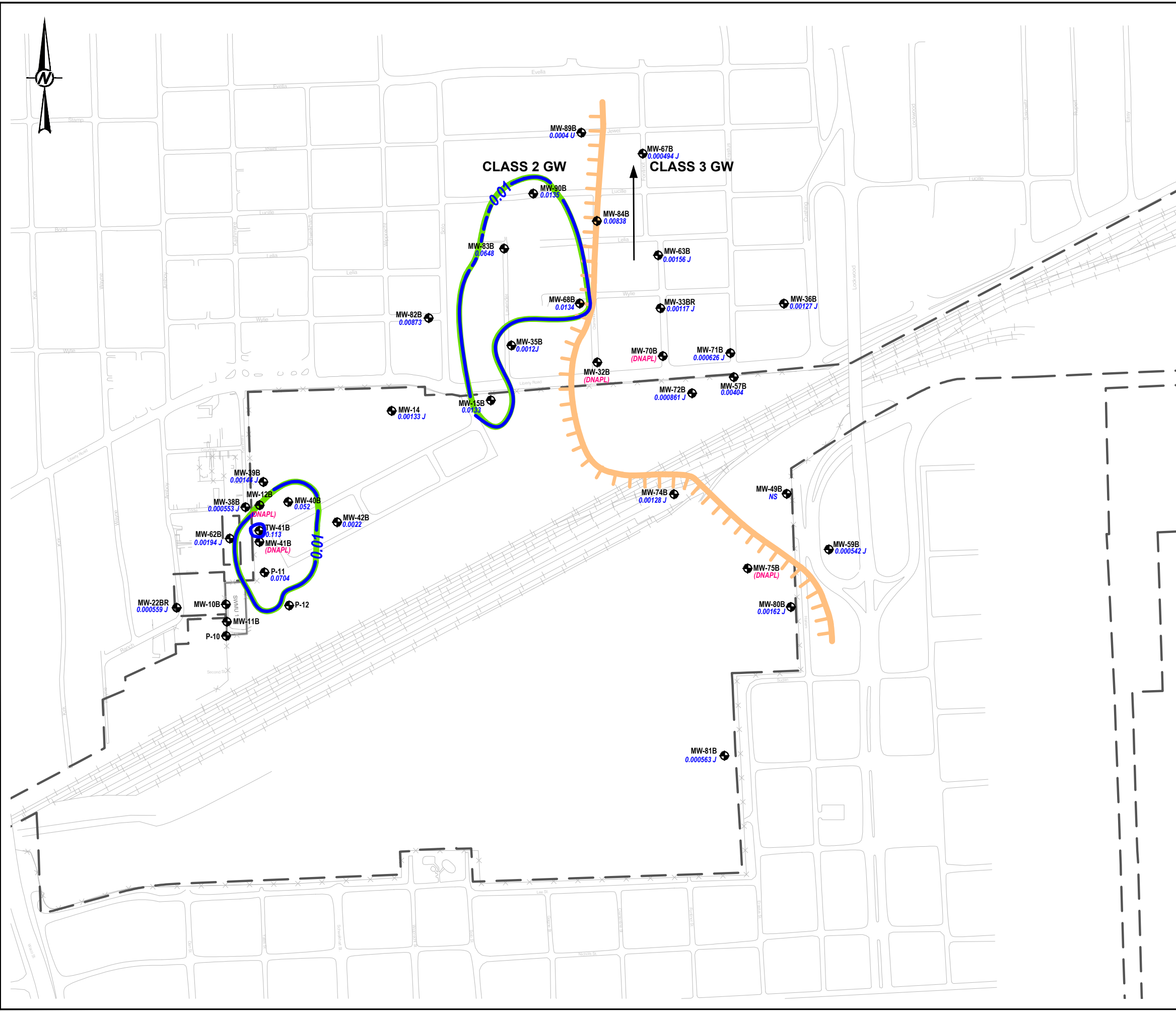
CONSULTANT	DATE	REVISION
	YYYY-MM-DD	2020-04-30
	DESIGNED	AJD
	PREPARED	RS
	REVIEWED	MH
	APPROVED	ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5B-20

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



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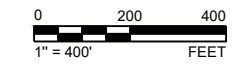
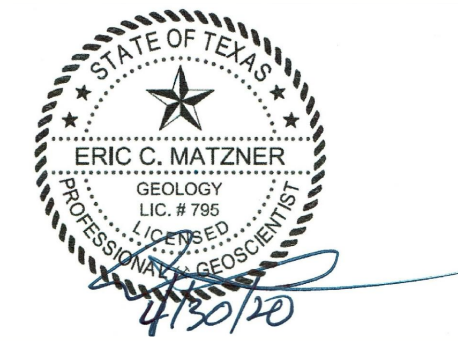
**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ B-TZ MONITORING WELL LOCATION
- B-CZ (CLASS 3 GW) B-TZ/B-CZ BOUNDARY
- B-TZ (CLASS 2 GW)
- 0.346 ARSENIC CONCENTRATION (mg/L)
- 0.01 ARSENIC CONCENTRATION CONTOUR (mg/L)
- ← INFERRED GROUNDWATER FLOW DIRECTION
- ARSENIC PCLE ZONE

**NOTE(S)**

1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
2. NS - NOT SAMPLED
3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
 CLASS 2 GW PCL: 0.01 mg/L  
 CLASS 3 GW PCL: 1 mg/L

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

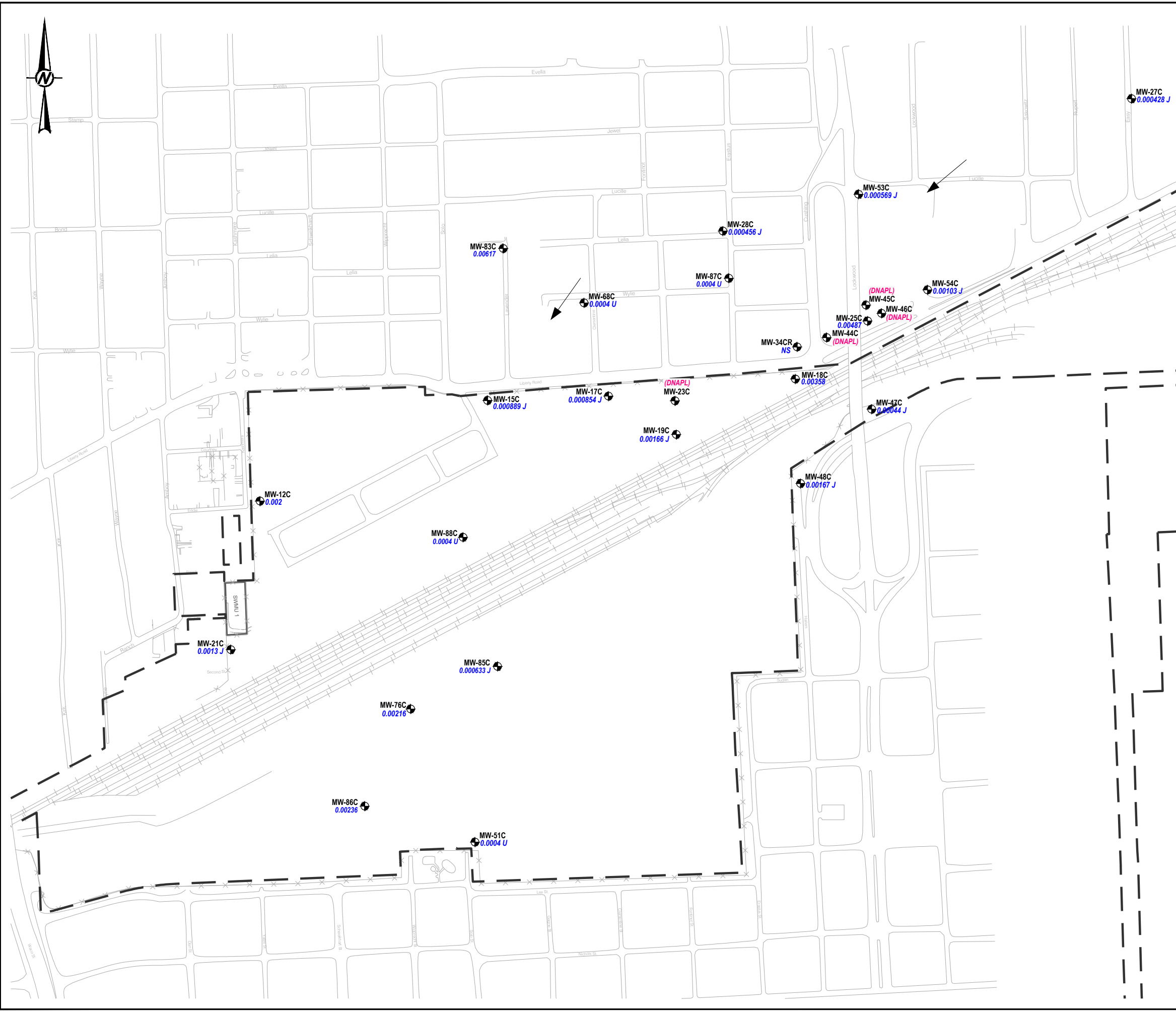
**TITLE**  
 GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 ARSENIC - JULY 2019

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED	AJD	
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5B-21

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erdm\data\project\5B-22\19119232 - Ground\cok\_2019\01\19232 - Inpvt\2020-1\jmh | File Name: ARSENIC C-TZ JULY 2019.dwg | Last Edited By: mshahar | Date: 2020-04-30 | Time: 1:48:44 PM | Printed By: RShahar | Date: 2020-04-30 | Time: 1:52:45 PM



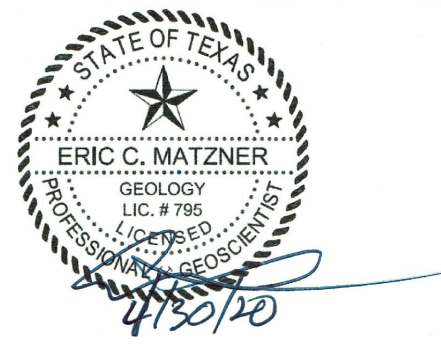
**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊙ C-TZ MONITORING WELL LOCATION
- 8.01** ARSENIC CONCENTRATION (mg/L)
- 0.01** ARSENIC CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION

**NOTE(S)**

1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JULY 2019).
2. NS - NOT SAMPLED
3. CONTOURS ARE BOLD ED AT THE RAL AND C/I PCL (0.01 mg/L).

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - C-TZ**  
**ARSENIC - JULY 2019**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED	AJD	
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-22

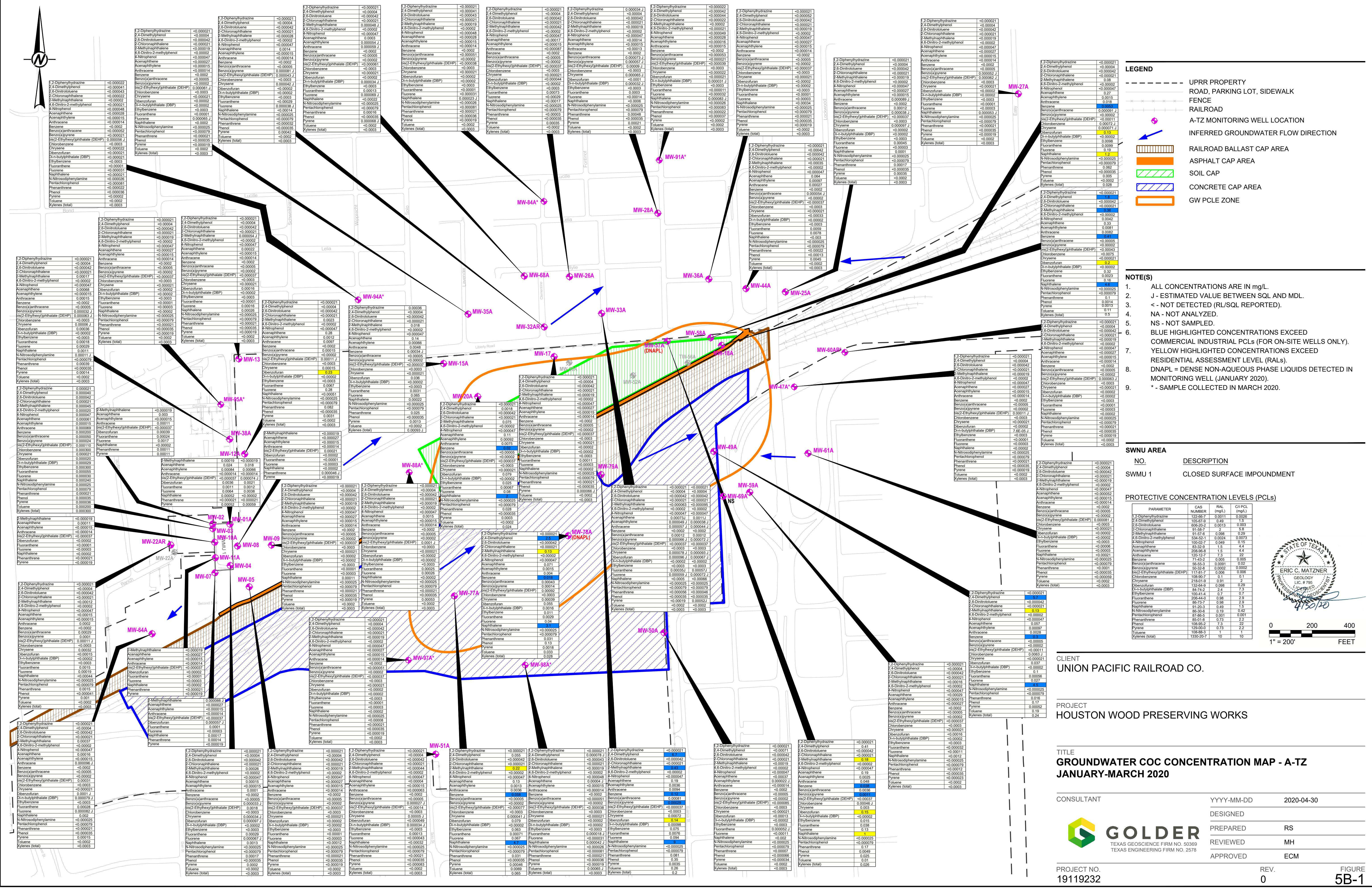
1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



**ATTACHMENT E**

# January-March 2020 Sampling Event COC Concentration Maps





**LEGEND**

- UPRR PROPERTY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- A-TZ MONITORING WELL LOCATION
- INFERRED GROUNDWATER FLOW DIRECTION
- RAILROAD BALLAST CAP AREA
- ASPHALT CAP AREA
- SOIL CAP
- CONCRETE CAP AREA
- GW PCLE ZONE

- NOTE(S)**
- ALL CONCENTRATIONS ARE IN mg/L.
  - J - ESTIMATED VALUE BETWEEN SQL AND MDL.
  - < - NOT DETECTED (RL/SQL REPORTED).
  - NA - NOT ANALYZED.
  - NS - NOT SAMPLED.
  - BLUE HIGHLIGHTED CONCENTRATIONS EXCEED COMMERCIAL INDUSTRIAL PCLs (FOR ON-SITE WELLS ONLY).
  - YELLOW HIGHLIGHTED CONCENTRATIONS EXCEED RESIDENTIAL ASSESSMENT LEVEL (RALs).
  - DNAPL = DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  - \* - SAMPLE COLLECTED IN MARCH 2020.

**SWNU AREA**

NO.	DESCRIPTION
SWNU 1	CLOSED SURFACE IMPONDMENT

**PROTECTIVE CONCENTRATION LEVELS (PCLs)**

PARAMETER	CAS NUMBER	RAL (mg/L)	CI PCL (mg/L)
2,2-Diphenylhydrazine	122-66-7	0.0011	0.0028
2,4-Dimethylphenol	105-67-6	0.63	1.61
2,6-Dinitrotoluene	86-22-3	0.0013	0.0033
2-Chloronaphthalene	91-58-7	0.0001	0.0001
2-Methylnaphthalene	91-58-7	0.0001	0.0001
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4
2,4-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
2,6-Dinitro-2-methylphenol	109-92-7	0.969	1.5
2-Methylphenol	83-32-1	1.5	4.4



CLIENT  
**UNION PACIFIC RAILROAD CO.**

PROJECT  
**HOUSTON WOOD PRESERVING WORKS**

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - A-TZ  
JANUARY-MARCH 2020**

CONSULTANT  
**GOLDER**  
TEXAS GEOSCIENCE FIRM NO. 55368  
TEXAS ENGINEERING FIRM NO. 25879

DATE  
2020-04-30

DESIGNED  
RS

PREPARED  
MH

REVIEWED  
ECM

APPROVED  
ECM

PROJECT NO.  
19119232

REV.  
0

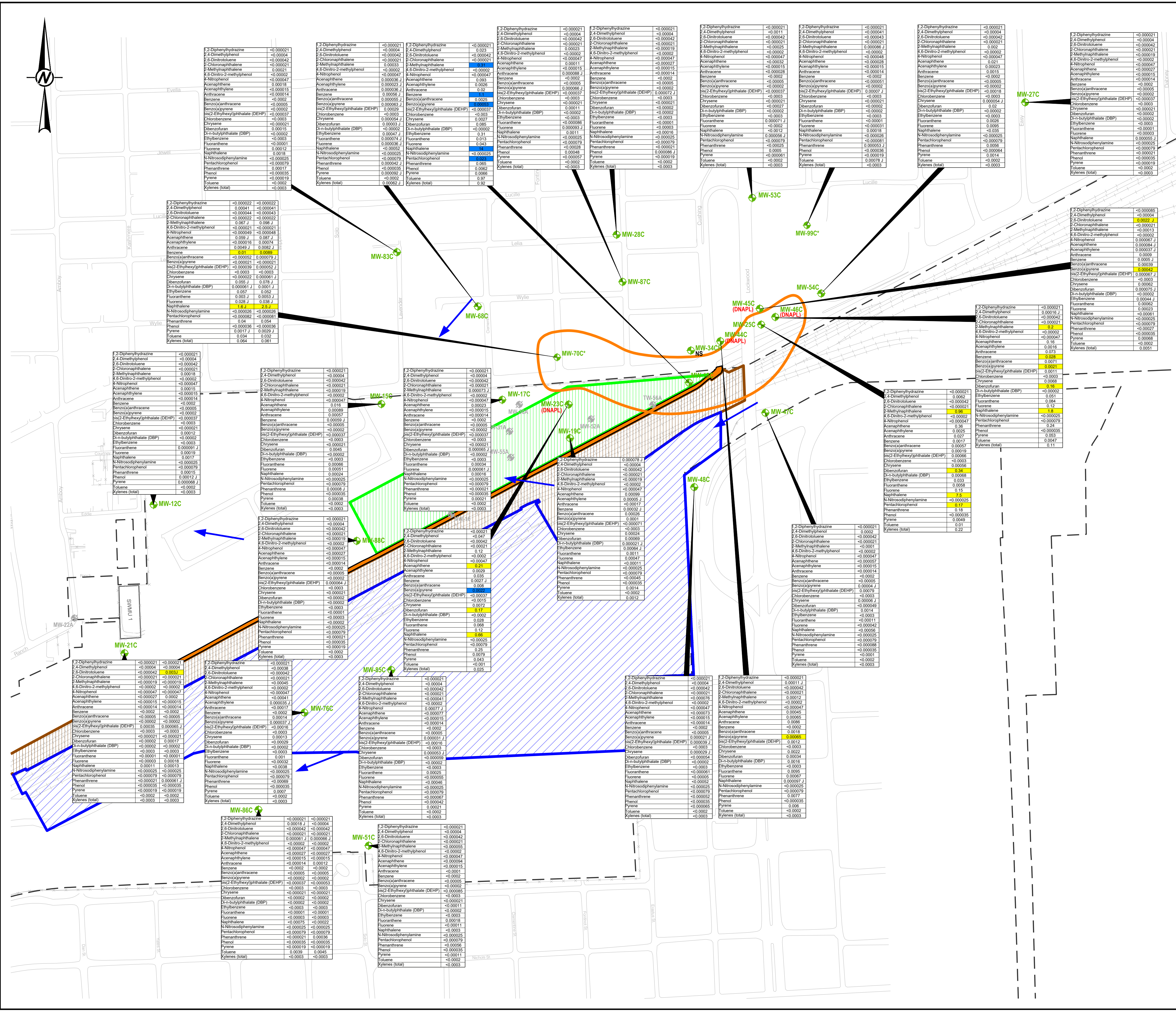
FIGURE  
5B-1

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN THE SHEET SIZE HAS BEEN MODIFIED FROM A35D









**LEGEND**

- UPRR PROPERTY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- INFERRED GROUNDWATER FLOW DIRECTION
- ▨ RAILROAD BALLAST CAP AREA
- ▨ ASPHALT CAP AREA
- ▨ SOIL CAP
- ▨ CONCRETE CAP AREA
- ▨ GW PCLE ZONE

- NOTE(S)**
- ALL CONCENTRATIONS ARE IN mg/L.
  - J - ESTIMATED VALUE BETWEEN SQL AND MDL.
  - < - NOT DETECTED (RL/SQL REPORTED).
  - NA - NOT ANALYZED.
  - NS - NOT SAMPLED.
  - BLUE HIGHLIGHTED CONCENTRATIONS EXCEED COMMERCIAL INDUSTRIAL PCLs (FOR ON-SITE WELLS ONLY).
  - YELLOW HIGHLIGHTED CONCENTRATIONS EXCEED RESIDENTIAL ASSESSMENT LEVEL (RALs).
  - DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  - \*\* - SAMPLE COLLECTED IN MARCH 2020.

**SWNU AREA**

NO.	DESCRIPTION
SWMU 1	CLOSED SURFACE IMPOUNDMENT

**PROTECTIVE CONCENTRATION LEVELS (PCLs)**

PARAMETER	CAS NUMBER	RIAL (mg/L)	CI PCL (mg/L)
2,2-Diphenylhydrazine	122-66-7	0.0011	0.0026
2,4-Dimethylphenol	105-67-0	0.49	0.15
2,6-Dinitrotoluene	606-20-2	0.013	0.003
2-Chloronaphthalene	91-57-6	0.086	0.8
2-Methylnaphthalene	91-58-7	0.086	0.8
4,6-Dinitro-2-methylphenol	534-52-1	0.0024	0.0073
4-Nitrophenol	100-02-7	0.049	0.15
Acenaphthene	83-32-9	1.5	4.4
Acenaphthylene	208-96-3	1.5	4.4
Anthracene	120-12-7	7.3	22
Benzo(a)anthracene	71-43-2	0.005	0.005
Benzo(a)pyrene	50-32-8	0.0002	0.0002
benz(b)fluoranthene	117-81-7	0.006	0.006
Chlorobenzene	108-90-7	0.1	0.1
Chrysene	123-44-9	0.008	0.29
Dibenzofuran	84-74-2	2.4	7.3
Dibenzylidene malonitrile (DBP)	100-41-4	0.17	0.7
Ethylbenzene	100-41-4	0.17	0.7
Fluoranthene	206-44-0	0.98	2.9
Fluorene	86-73-7	0.98	2.9
Naphthalene	91-20-3	0.49	1.5
Nitrofluorene	86-34-6	0.19	0.43
Nitroindole	87-86-5	0.001	0.001
Pentachlorophenol	85-01-8	0.73	2.2
Phenanthrene	85-01-8	0.73	2.2
Phenol	108-95-2	7.3	22
Pyrene	129-00-0	0.73	2.2
Toluene	108-88-3	1	1
Xylenes (total)	1330-20-7	10	10

**CLIENT**  
UNION PACIFIC RAILROAD CO.

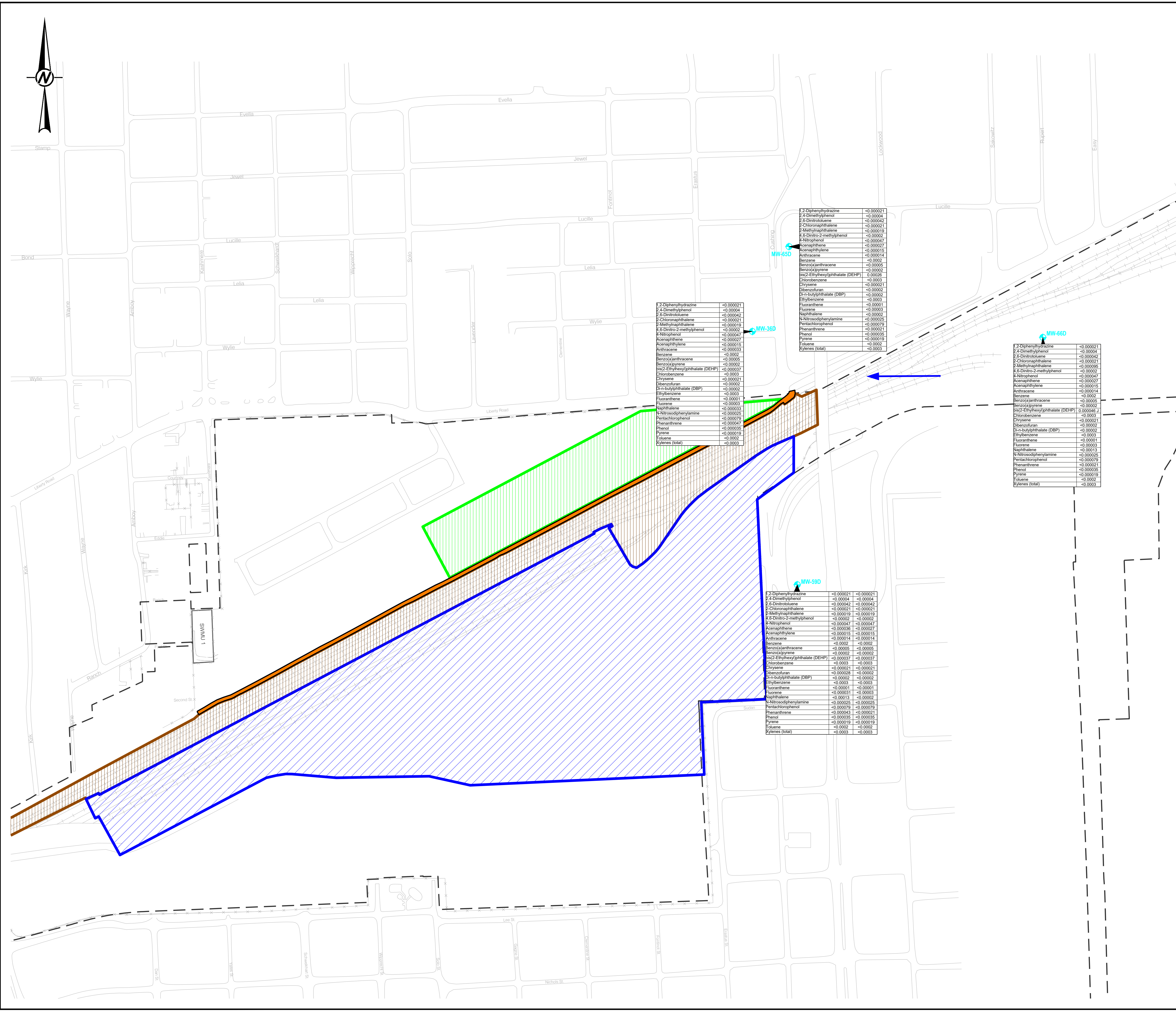
**PROJECT**  
HOUSTON WOOD PRESERVING WORKS

**TITLE**  
GROUNDWATER COC CONCENTRATION MAP - C-TZ  
JANUARY-MARCH 2020

CONSULTANT	DATE	REVISION
YYYY-MM-DD	2020-04-30	
DESIGNED		RS
PREPARED		MH
REVIEWED		ECM
APPROVED		

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM A3/D





2,2-Diphenylhydrazine	<0.000021
2,4-Dinitrophenol	<0.000041
2,6-Dinitrotoluene	<0.000042
2-Chloronaphthalene	<0.000021
2-Methylnaphthalene	<0.000019
4-Nitrophenol	<0.000041
4-Nitrophenol	<0.000041
Acenaphthene	<0.000027
Acenaphthylene	<0.000015
Anthracene	<0.000014
Benzo(a)anthracene	<0.000005
Benzo(a)pyrene	<0.000002
Benzo(b)fluoranthene	<0.000002
Benzo(k)fluoranthene	<0.000002
Chlorobenzene	<0.00003
Chrysene	<0.000021
Dibenzofuran	<0.000002
Dibenz(p,h)anthracene	<0.000002
Ethylbenzene	<0.00001
Fluoranthene	<0.00001
Fluorene	<0.000025
Naphthalene	<0.00002
Nitrofluorene	<0.000025
Phenanthrene	<0.000019
Phenol	<0.000035
Pyrene	<0.000019
Toluene	<0.00002
Xylenes (total)	<0.0003

2,2-Diphenylhydrazine	<0.000021
2,4-Dinitrophenol	<0.000041
2,6-Dinitrotoluene	<0.000042
2-Chloronaphthalene	<0.000021
2-Methylnaphthalene	<0.000019
4-Nitrophenol	<0.000041
4-Nitrophenol	<0.000041
Acenaphthene	<0.000027
Acenaphthylene	<0.000015
Anthracene	<0.000014
Benzo(a)anthracene	<0.000005
Benzo(a)pyrene	<0.000002
Benzo(b)fluoranthene	<0.000002
Benzo(k)fluoranthene	<0.000002
Chlorobenzene	<0.00003
Chrysene	<0.000021
Dibenzofuran	<0.000002
Dibenz(p,h)anthracene	<0.000002
Ethylbenzene	<0.00001
Fluoranthene	<0.00001
Fluorene	<0.000025
Naphthalene	<0.00002
Nitrofluorene	<0.000025
Phenanthrene	<0.000019
Phenol	<0.000035
Pyrene	<0.000019
Toluene	<0.00002
Xylenes (total)	<0.0003

2,2-Diphenylhydrazine	<0.000021
2,4-Dinitrophenol	<0.000041
2,6-Dinitrotoluene	<0.000042
2-Chloronaphthalene	<0.000021
2-Methylnaphthalene	<0.000019
4-Nitrophenol	<0.000041
4-Nitrophenol	<0.000041
Acenaphthene	<0.000027
Acenaphthylene	<0.000015
Anthracene	<0.000014
Benzo(a)anthracene	<0.000005
Benzo(a)pyrene	<0.000002
Benzo(b)fluoranthene	<0.000002
Benzo(k)fluoranthene	<0.000002
Chlorobenzene	<0.00003
Chrysene	<0.000021
Dibenzofuran	<0.000002
Dibenz(p,h)anthracene	<0.000002
Ethylbenzene	<0.00001
Fluoranthene	<0.00001
Fluorene	<0.000025
Naphthalene	<0.00002
Nitrofluorene	<0.000025
Phenanthrene	<0.000019
Phenol	<0.000035
Pyrene	<0.000019
Toluene	<0.00002
Xylenes (total)	<0.0003

2,2-Diphenylhydrazine	<0.000021
2,4-Dinitrophenol	<0.000041
2,6-Dinitrotoluene	<0.000042
2-Chloronaphthalene	<0.000021
2-Methylnaphthalene	<0.000019
4-Nitrophenol	<0.000041
4-Nitrophenol	<0.000041
Acenaphthene	<0.000027
Acenaphthylene	<0.000015
Anthracene	<0.000014
Benzo(a)anthracene	<0.000005
Benzo(a)pyrene	<0.000002
Benzo(b)fluoranthene	<0.000002
Benzo(k)fluoranthene	<0.000002
Chlorobenzene	<0.00003
Chrysene	<0.000021
Dibenzofuran	<0.000002
Dibenz(p,h)anthracene	<0.000002
Ethylbenzene	<0.00001
Fluoranthene	<0.00001
Fluorene	<0.000025
Naphthalene	<0.00002
Nitrofluorene	<0.000025
Phenanthrene	<0.000019
Phenol	<0.000035
Pyrene	<0.000019
Toluene	<0.00002
Xylenes (total)	<0.0003

**LEGEND**

- UPRR PROPERTY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- D-TZ MONITORING WELL LOCATION
- INFERRED GROUNDWATER FLOW DIRECTION
- ▨ RAILROAD BALLAST CAP AREA
- ▨ ASPHALT CAP AREA
- ▨ SOIL CAP
- ▨ CONCRETE CAP AREA

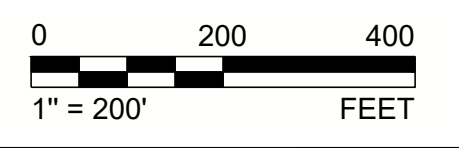
- NOTE(S)**
- ALL CONCENTRATIONS ARE IN mg/L.
  - J - ESTIMATED VALUE BETWEEN SQL AND MDL.
  - < - NOT DETECTED (RL/SQL REPORTED).
  - NS - NOT ANALYZED.
  - NA - NOT SAMPLED.
  - BLUE HIGHLIGHTED CONCENTRATIONS EXCEED COMMERCIAL INDUSTRIAL PCLs (FOR ON-SITE WELLS ONLY).
  - YELLOW HIGHLIGHTED CONCENTRATIONS EXCEED RESIDENTIAL ASSESSMENT LEVEL (RALs).
  - DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  - \* - SAMPLE COLLECTED IN MARCH 2020.

**SNWU AREA**

NO.	DESCRIPTION
SNWU 1	CLOSED SURFACE IMPOUNDMENT

**PROTECTIVE CONCENTRATION LEVELS (PCLs)**

PARAMETER	CAS NUMBER	RAL (mg/L)	C1 PCL (mg/L)
2,2-Diphenylhydrazine	122-66-7	0.0011	0.0026
2,4-Dinitrophenol	155-47-9	0.49	1.5
2,6-Dinitrotoluene	606-20-2	0.0013	0.003
2-Chloronaphthalene	91-58-7	0.008	0.29
2-Methylnaphthalene	91-57-5	0.049	0.15
4-Nitrophenol	534-52-1	0.0024	0.0073
4-Nitrophenol	100-02-7	0.049	0.15
Acenaphthene	83-32-9	1.5	4.4
Acenaphthylene	208-96-9	1.5	4.4
Anthracene	120-12-7	7.3	22
Benzo(a)anthracene	71-43-2	0.005	0.015
Benzo(a)pyrene	56-56-3	0.0011	0.002
Benzo(b)fluoranthene	50-32-8	0.0002	0.0007
Benzo(k)fluoranthene	117-81-7	0.006	0.006
Chlorobenzene	108-90-7	0.1	0.1
Chrysene	218-01-9	0.01	0.01
Dibenzofuran	132-64-9	0.008	0.29
Dibenz(p,h)anthracene	84-74-2	2.4	7.3
Ethylbenzene	100-41-4	0.17	0.7
Fluoranthene	206-44-0	0.98	2.9
Fluorene	86-73-7	0.98	2.9
Naphthalene	91-20-3	0.49	1.5
Nitrofluorene	86-30-4	0.19	0.42
Pentachlorophenol	87-86-5	0.001	0.001
Phenanthrene	85-01-8	0.73	2.2
Phenol	108-95-2	7.3	22
Pyrene	129-00-0	0.73	2.2
Toluene	108-88-3	1	1
Xylenes (total)	1330-20-7	10	10



CLIENT  
UNION PACIFIC RAILROAD CO.

PROJECT  
HOUSTON WOOD PRESERVING WORKS

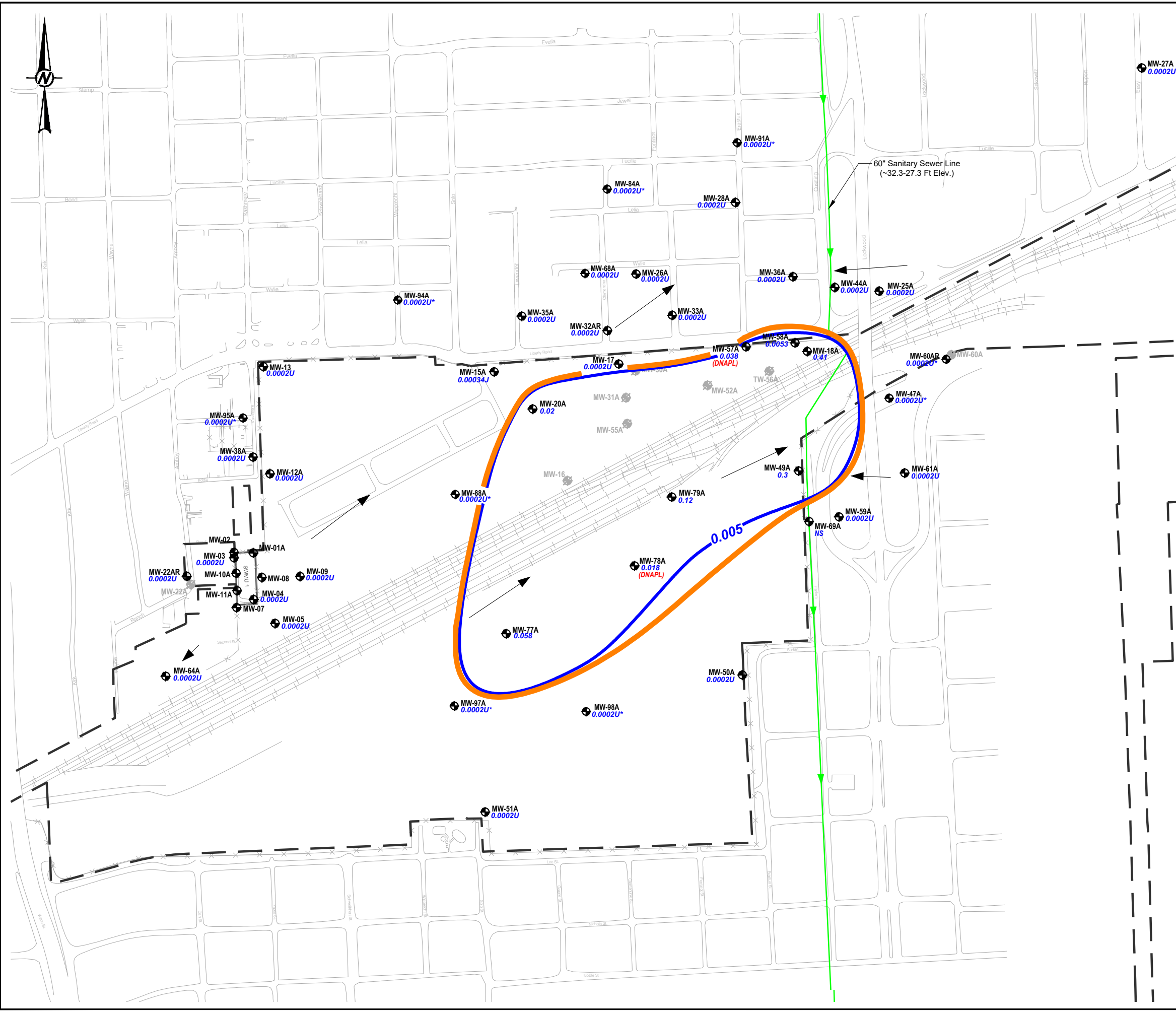
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**GROUNDWATER COC CONCENTRATION MAP - D-TZ**  
**JANUARY-MARCH 2020**

CONSULTANT	DATE
<p><b>GOLDER</b> TEXAS GEOSCIENCE FIRM NO. 50369 TEXAS ENGINEERING FIRM NO. 2578</p>	YYYY-MM-DD 2020-04-30
	DESIGNED
	PREPARED RS
	REVIEWED MH
	APPROVED ECM

PROJECT NO. 19119232 REV. 0 FIGURE 5B-4



Path: \\uswest\erand\data\projects - Round Rock\2019\19119232 - HWPW\2020-31 March\1 - File Name: 5B-5 - 5B-5 Groundwater COC Concentration Map A-TZ - January 2020.dwg | Last Edited By: rbsalar | Date: 2020-04-30 | Printed By: rbsalar | Date: 2020-04-30 | Time: 12:26:47 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- A-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- 0.038** BENZENE CONCENTRATION (mg/L)
- 0.005** BENZENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.005 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

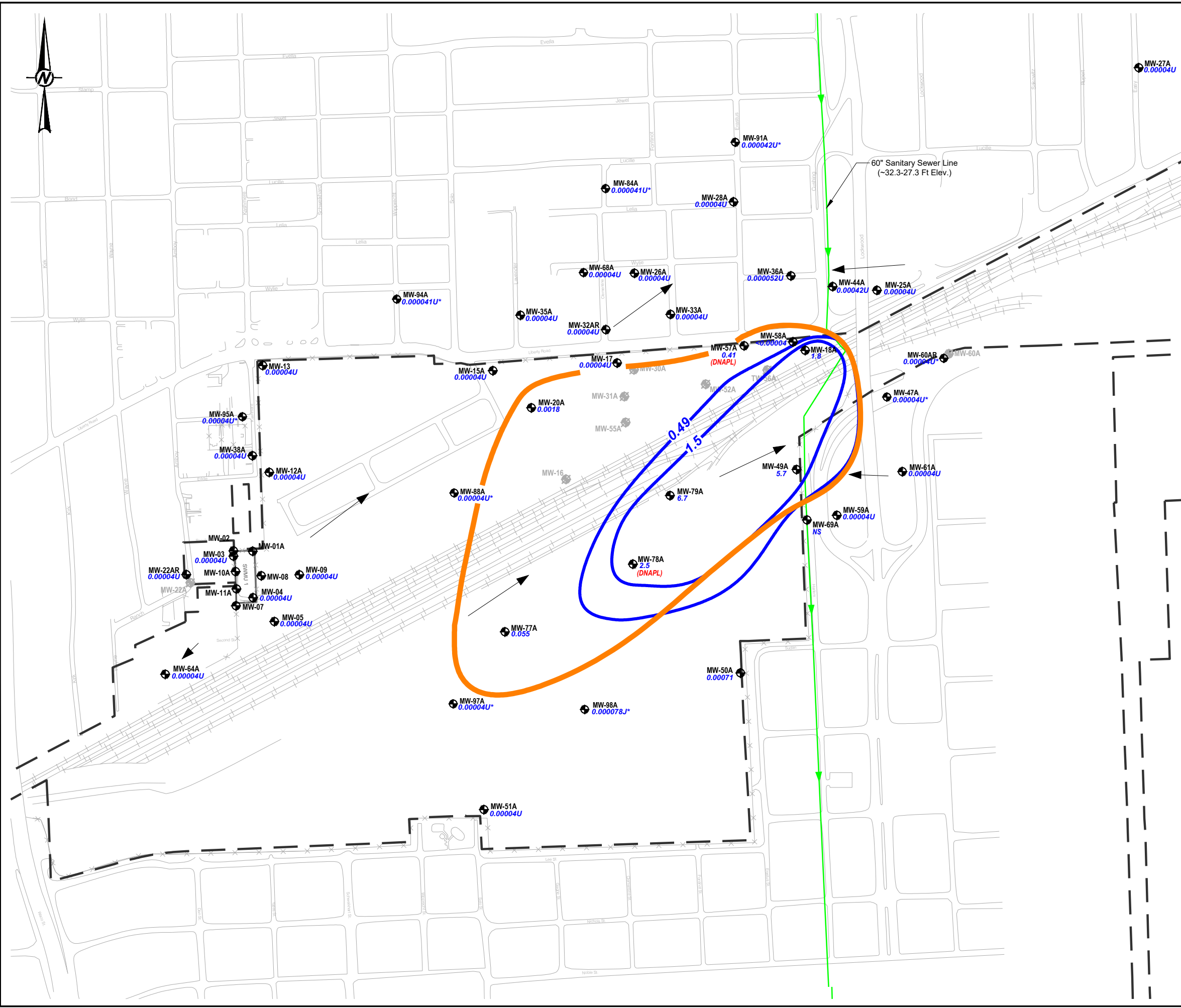
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - A-TZ  
 BENZENE - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
 GOLDER <small>TEXAS GEOSCIENCE FIRM NO. 50969          TEXAS ENGINEERING FIRM NO. 2578</small>	DESIGNED	
	PREPARED	RS
	REVIEWED	MH
	APPROVED	ECM

PROJECT NO. 19119232      REV. 0      FIGURE 5B-5

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\projects - Round Rock\2019\19119232 - HWPW\2020-31 March\1 - File Name: 5B-6 Groundwater COC Concentration Map A-TZ - January 2020.dwg | Last Edited By: rscholar | Date: 2020-04-30 | Printed By: rscholar | Date: 2020-04-30 | Time: 12:26:58 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- A-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- 0.41** 2,4 DIMETHYLPHENOL CONCENTRATION (mg/L)
- 0.49** 2,4 DIMETHYLPHENOL CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.49 mg/L AND 1.5 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

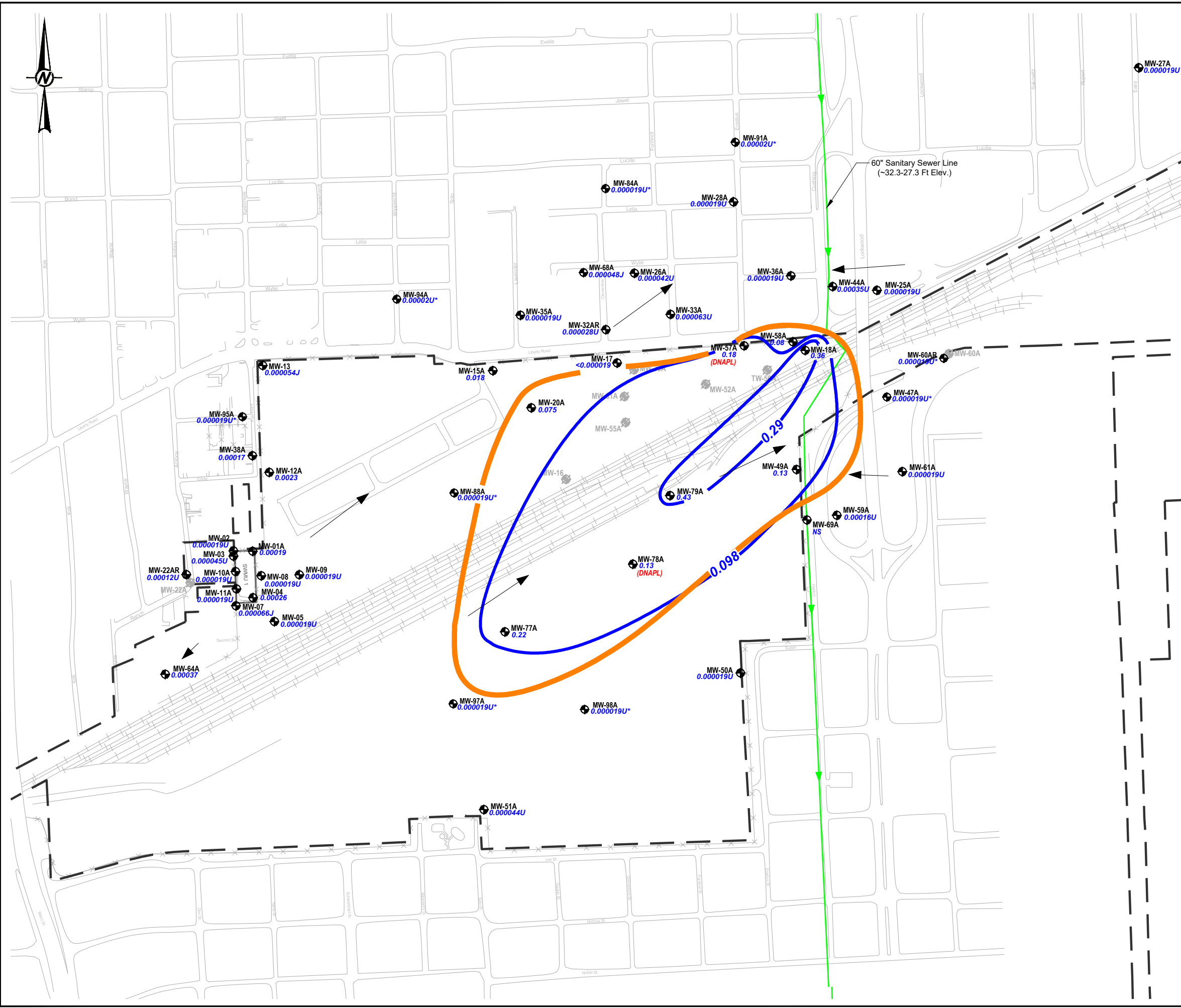
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**GROUNDWATER COC CONCENTRATION MAP - A-TZ**  
**2,4-DIMETHYLPHENOL - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-6

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\projects - Round Rock\19119232 - HWPW\2020-31\month 1 - File Name: 5B-5 - 5B-5 Groundwater COC Concentration Map A-TZ - January 2020.dwg | Last Edited By: rbsalar | Date: 2020-04-30 | Printed By: rbsalar | Date: 2020-04-30 | Time: 12:28:08 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ A-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- 0.18 2-METHYLNAPHTHALENE CONCENTRATION (mg/L)
- 0.098 2-METHYLNAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.098 mg/L AND 0.29 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - A-TZ  
 2-METHYLNAPHTHALENE - JANUARY-MARCH 2020**

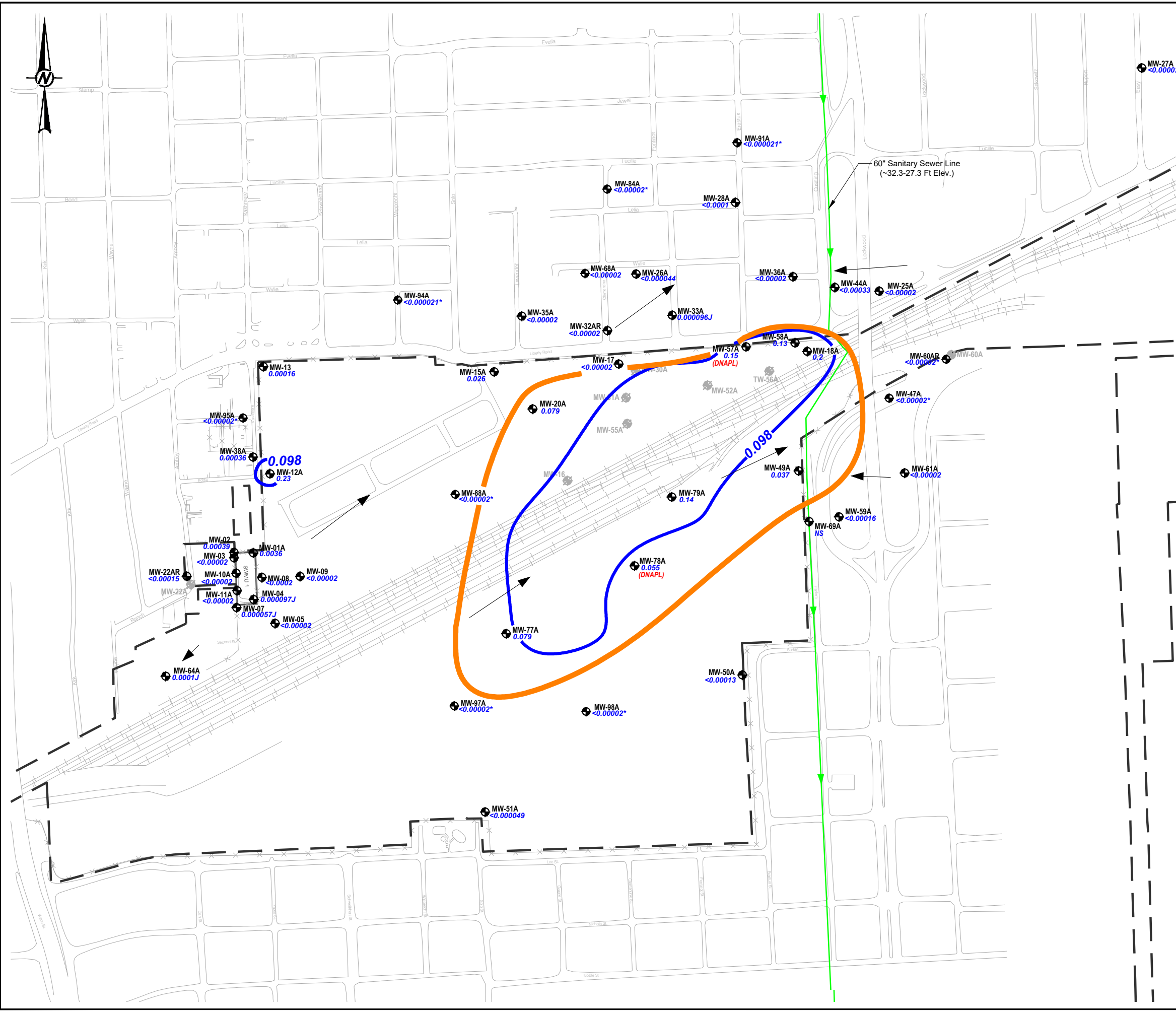
CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-7

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erand\data\projects - Round Rock\19119232 - HWPW\2020-31\Month 1 - File Name: 5B-5 - 5B-5 Groundwater COC Concentration Map A-TZ - January 2020.dwg | Last Edited By: rscholar | Date: 2020-04-30 | Printed By: rscholar | Date: 2020-04-30 | Time: 12:28:18 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ A-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- 0.15 DIBENZOFURAN CONCENTRATION (mg/L)
- 0.098** DIBENZOFURAN CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.098 mg/L AND 0.29 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

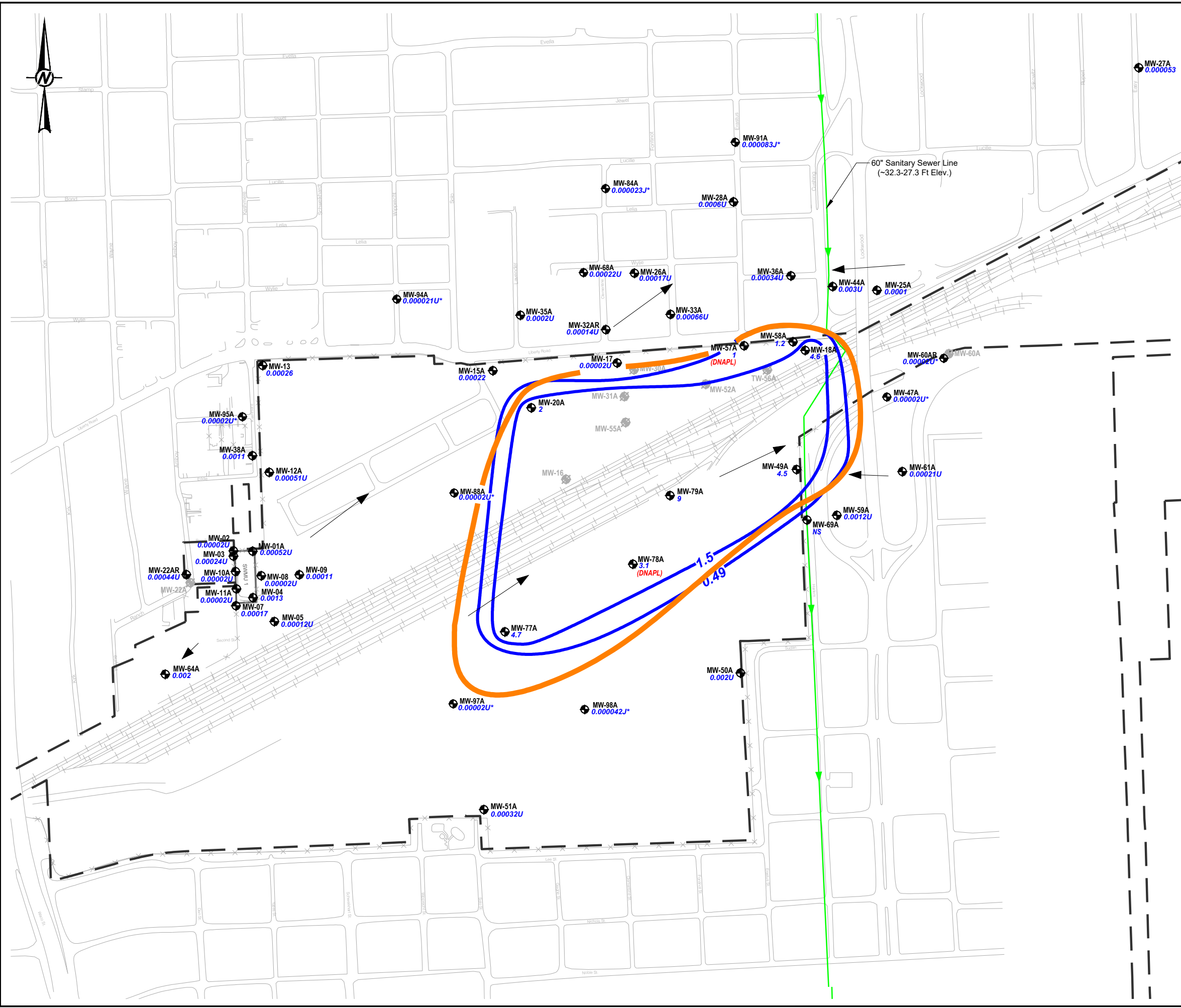
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - A-TZ  
 DIBENZOFURAN - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-8

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\2019\19119232 - HWPW\2020-31 March\1 - File Name: 5B-9 Groundwater COC Concentration Map A-TZ - January 2020.dwg | Last Edited By: rbsalar | Date: 2020-04-30 | Printed By: rbsalar | Date: 2020-04-30 | Time: 12:28:28 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- ⊕ A-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- <0.003 NAPHTHALENE CONCENTRATION (mg/L)
- 0.49** NAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- ➔ INFERRED GROUNDWATER FLOW DIRECTION
- ▭ GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.49 mg/L AND 1.5 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

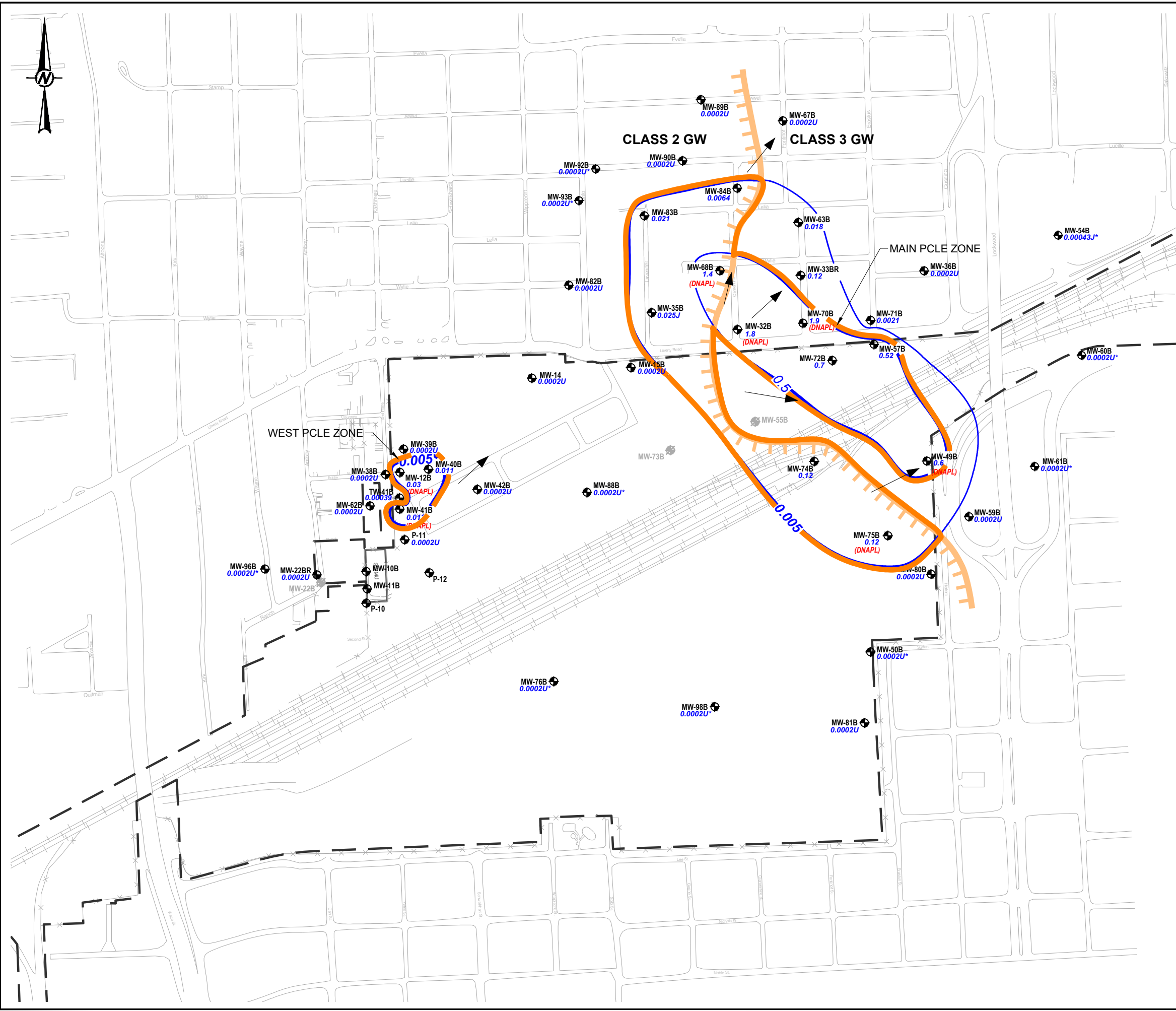
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - A-TZ  
 NAPHTHALENE - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-9

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\2019\01\10232 - HWPP\2020-31\March - File Name: 5B10 - 5B14 Groundwater COC Concentration Map B-TZ-CZ - January 2020.dwg | Last Edited By: rbsalar | Date: 2020-04-30 | Time: 12:27:01 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- X — FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW)
- B-TZ (CLASS 2 GW)
- <math>0.002</math> BENZENE CONCENTRATION (mg/L)
- 0.005** BENZENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
CLASS 2 GW PCL: 0.005 mg/L  
CLASS 3 GW PCL: 0.5 mg/L
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

0 200 400  
1" = 400' FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 BENZENE - JANUARY-MARCH 2020**

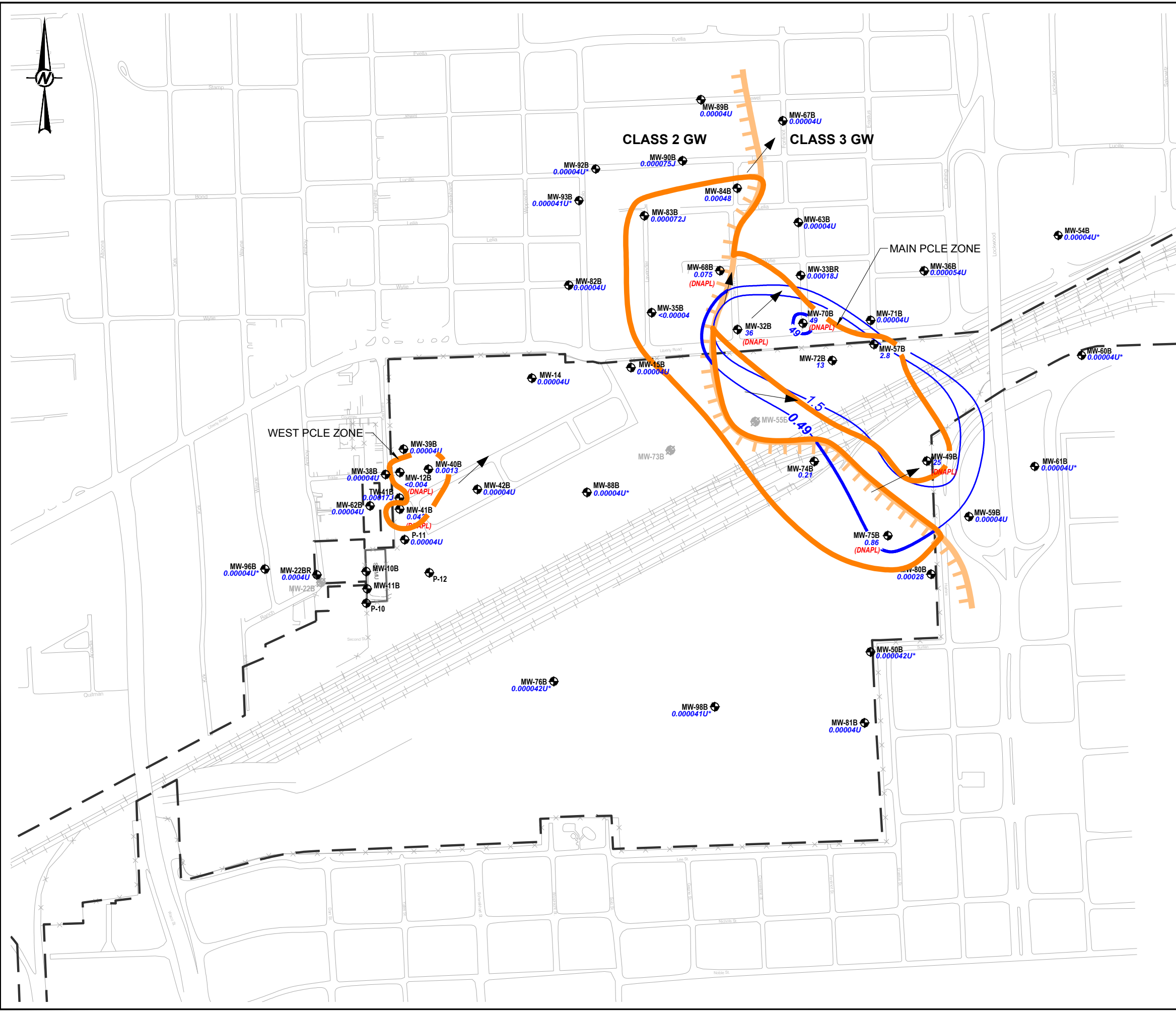
CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-10

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erand\data\Projects - Round Rock\2019\19119232 - HWPP\2020-31 March - File Name: 5B10 - 5B14 Groundwater COC Concentration Map B-TZ-CZ - January 2020.dwg | Last Edited By: rslabzar | Date: 2020-04-30 | Time: 12:27:14 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- X — FENCE
- RAILROAD
- ⊕ B-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW)
- B-TZ/B-CZ BOUNDARY
- 0.21 B-TZ (CLASS 2 GW) 2,4-DIMETHYLPHENOL CONCENTRATION (mg/L)
- 0.49 2,4-DIMETHYLPHENOL CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
CLASS 2 GW PCL: 0.49 mg/L (Res.) & 1.5 mg/L (C/I)  
CLASS 3 BW PCL: 49 mg/L (Res.) & 150 mg/L (C/I)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

STATE OF TEXAS  
 ERIC C. MATZNER  
 GEOLOGY  
 LIC. # 795  
 PROFESSIONAL GEOSCIENTIST  
 4/30/20

0 200 400  
 1" = 400' FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

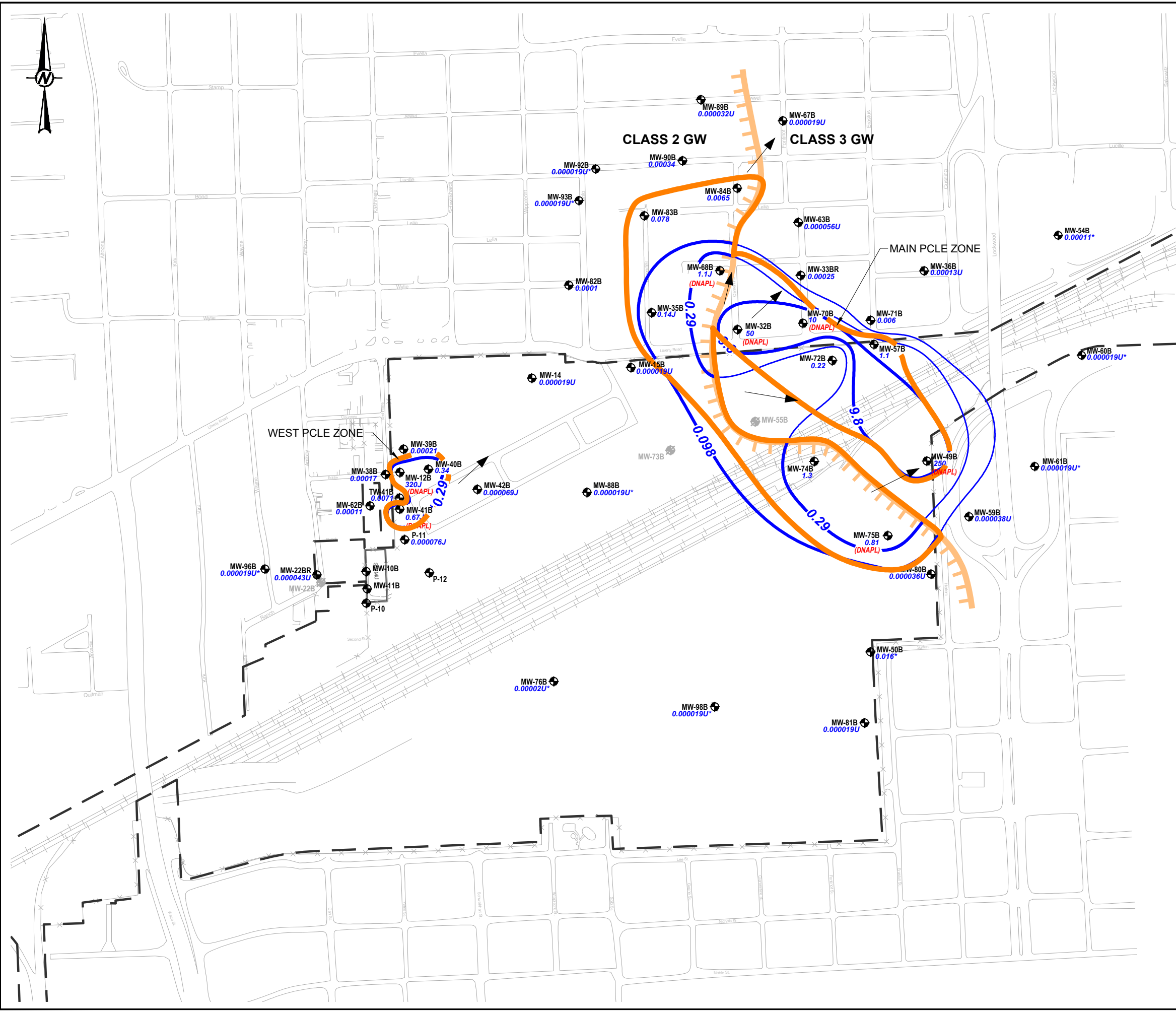
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 2,4-DIMETHYLPHENOL - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-11

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\2019\01\10232 - HWPP\2020-31\March - File Name: 5B10 - 5B14 Groundwater COC Concentration Map B-TZ-CZ - January 2020.dwg | Last Edited By: rbsalar | Date: 2020-04-30 | Time: 12:27:30 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW)
- B-TZ (CLASS 2 GW)
- 2-METHYLNAPHTHALENE CONCENTRATION (mg/L)
- 2-METHYLNAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
 CLASS 2 GW PCL: 0.098 mg/L (Res.) & 0.29 mg/L (C/I)  
 CLASS 3 GW PCL: 9.8 mg/L (Res.) & 29 mg/L (C/I)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

0 200 400  
1" = 400' FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

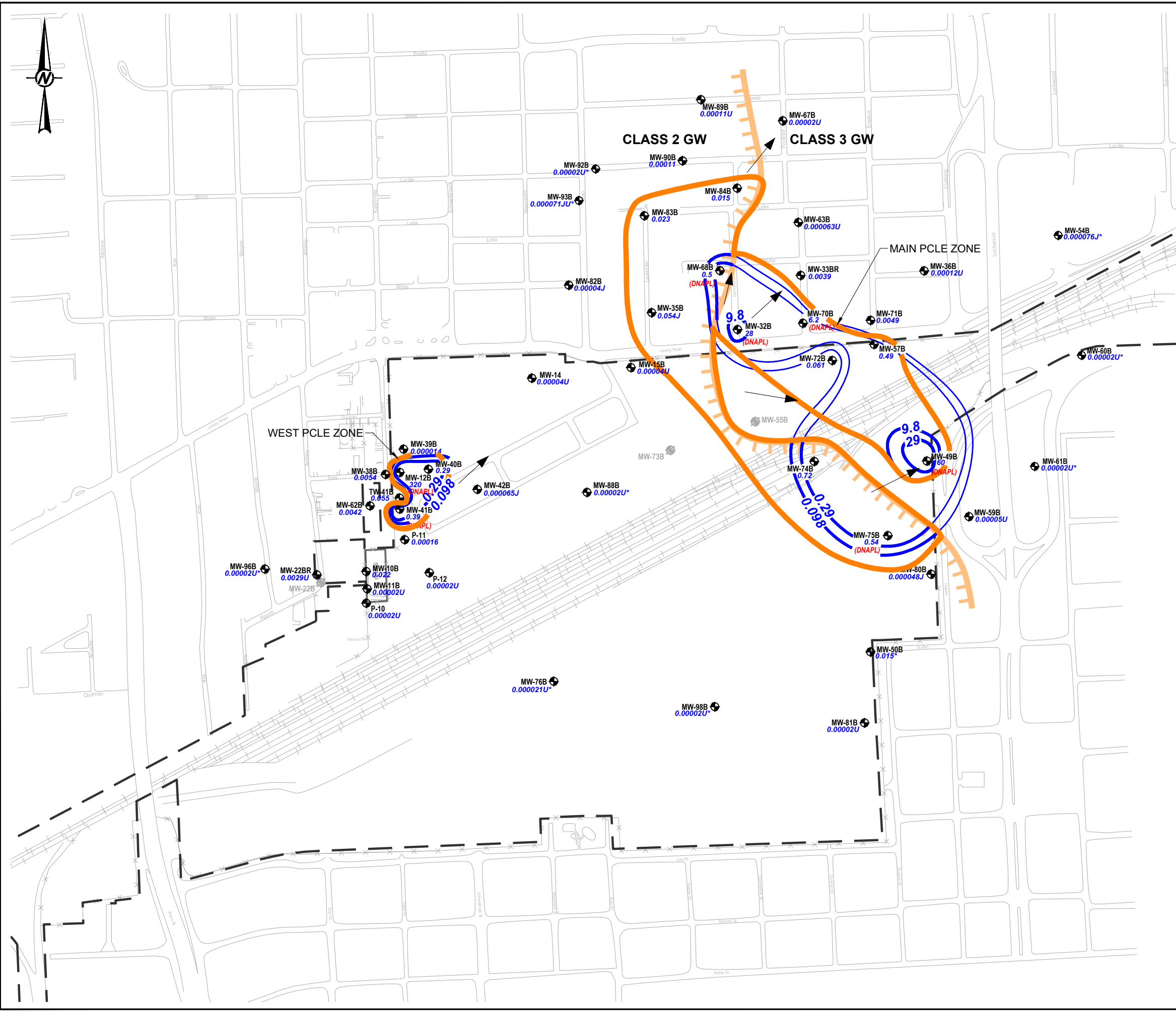
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**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 2-METHYLNAPHTHALENE - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-12

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\dat\Projects - Round Rock\2019\01\10232 - HWPW\2020-31\March - File Name: 5B10 - 5B14 Groundwater COC Concentration Map - B-TZ-CZ - January 2020.dwg | Last Edited By: rbsalar | Date: 2020-04-30 | Time: 12:27:44 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW)
- B-TZ (CLASS 2 GW)
- 0.0049 DIBENZOFURAN CONCENTRATION (mg/L)
- 0.098 DIBENZOFURAN CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
 CLASS 2 GW PCL: 0.098 mg/L (Res.) & 0.29 mg/L (C/I)  
 CLASS 3 GW PCL: 9.8 mg/L (Res.) & 29 mg/L (C/I)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

0 200 400  
 1" = 400' FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 DIBENZOFURAN - JANUARY-MARCH 2020**

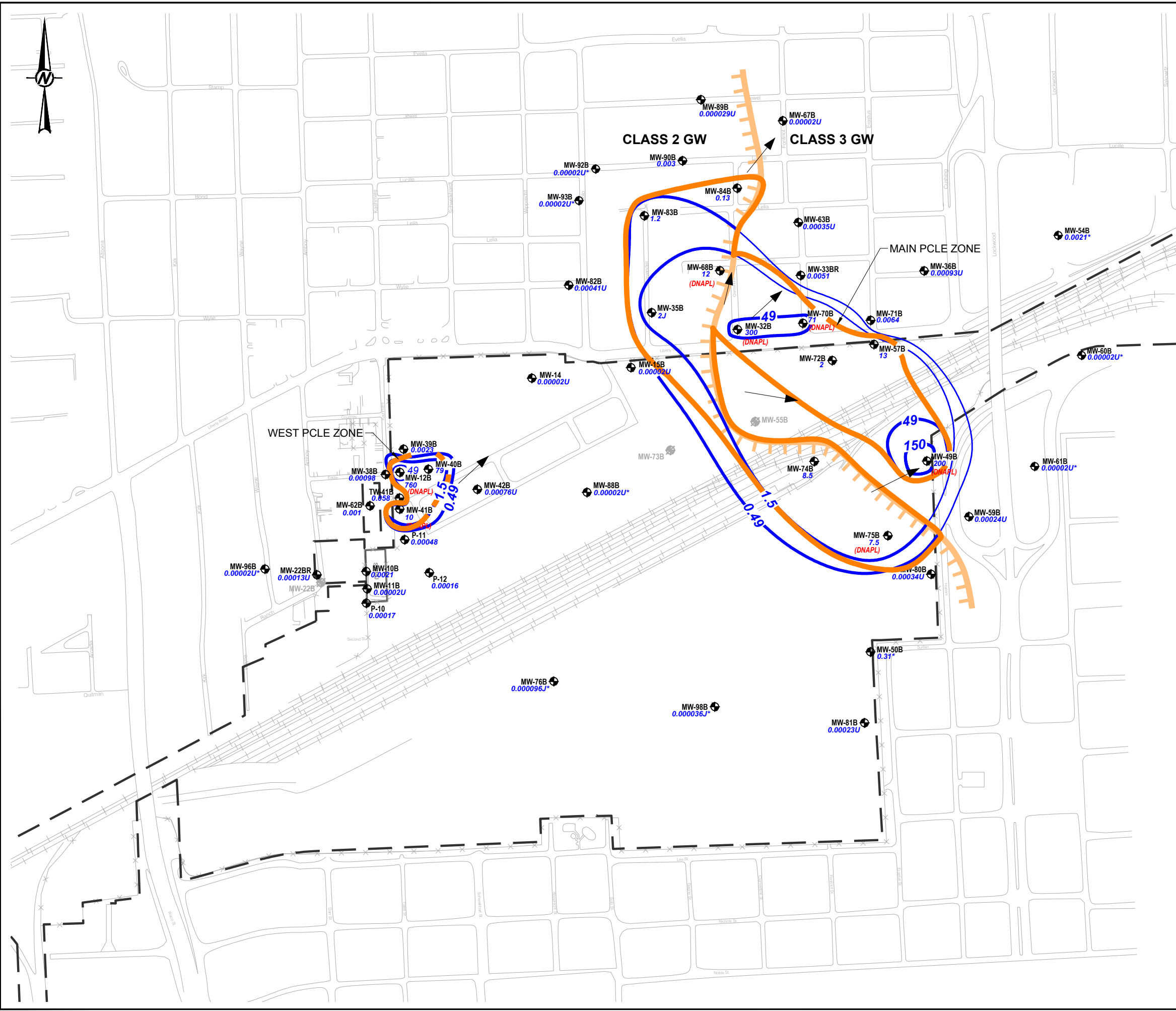
CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-13

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erand\data\Projects - Round Rock\2019\10110232 - HWPP\2020-31 March - File Name: 5B10 - 5B14 Groundwater COC Concentration Map - B-TZ/CZ - January 2020.dwg | Last Edited By: rslabzar | Date: 2020-04-30 | Time: 12:19:34 PM | Printed By: rslabzar | Date: 2020-04-30 | Time: 12:28:00 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- B-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW)
- B-TZ (CLASS 2 GW)
- 0.0021 NAPHTHALENE CONCENTRATION (mg/L)
- 0.49 NAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
CLASS 2 GW PCL: 0.49 mg/L (Res.) & 1.5 mg/L (C/I)  
CLASS 3 GW PCL: 49 mg/L (Res.) & 150 mg/L (C/I)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

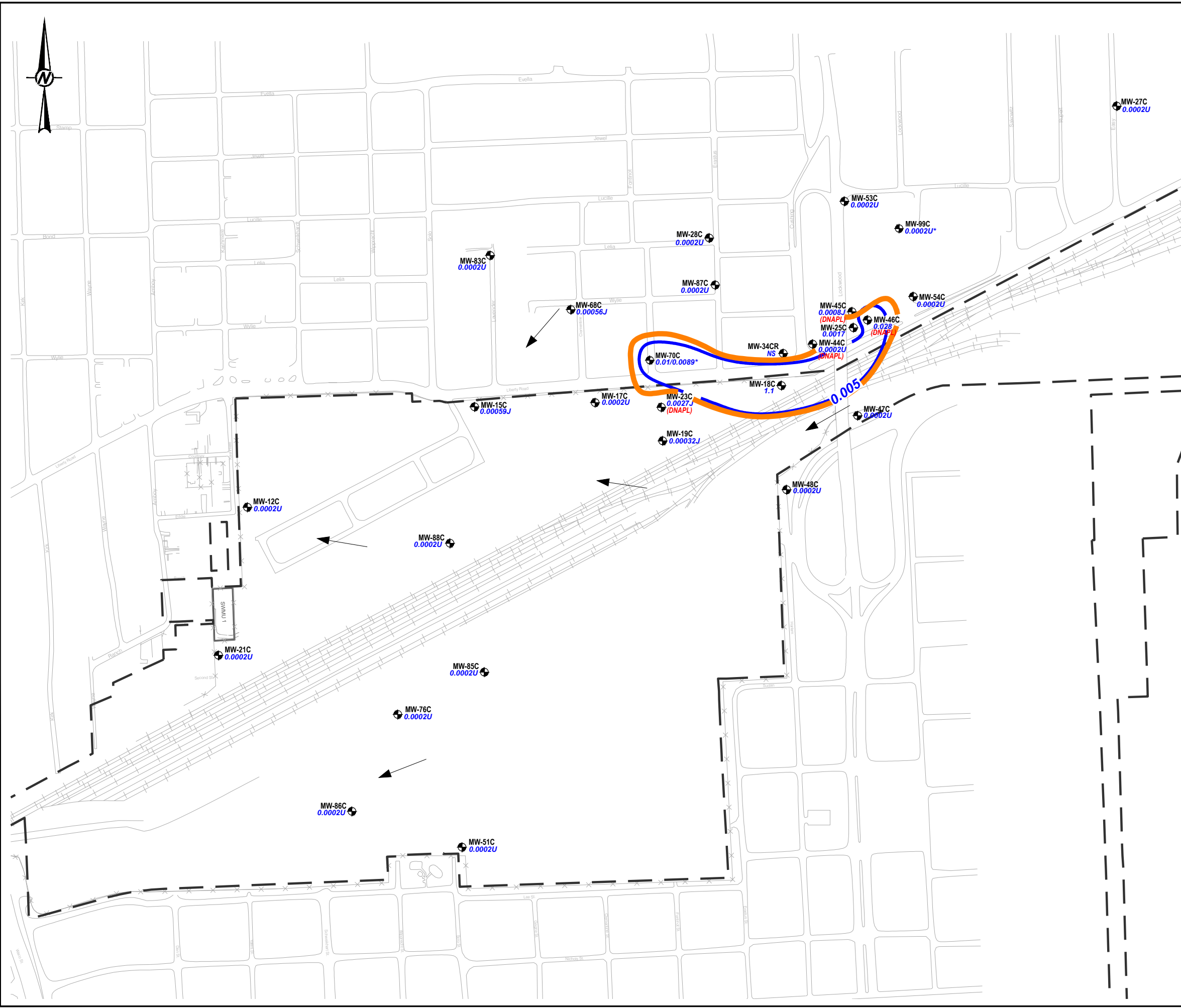
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ  
 NAPHTHALENE - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-14

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\2019\01\19232 - HWPW\2020-31\March - 5B15 - 5B19 Groundwater COC Concentration Map - C-TZ - January 2020.dwg | Last Edited By: rsalazar | Date: 2020-04-30 | Time: 12:21:14 PM | Printed By: RSalazar | Date: 2020-04-30 | Time: 12:28:38 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ C-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- <0.0002 BENZENE CONCENTRATION (mg/L)
- 0.005** BENZENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- ▭ GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.005 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

**TITLE**  
 GROUNDWATER COC CONCENTRATION MAP - C-TZ  
 BENZENE - JANUARY-MARCH 2020

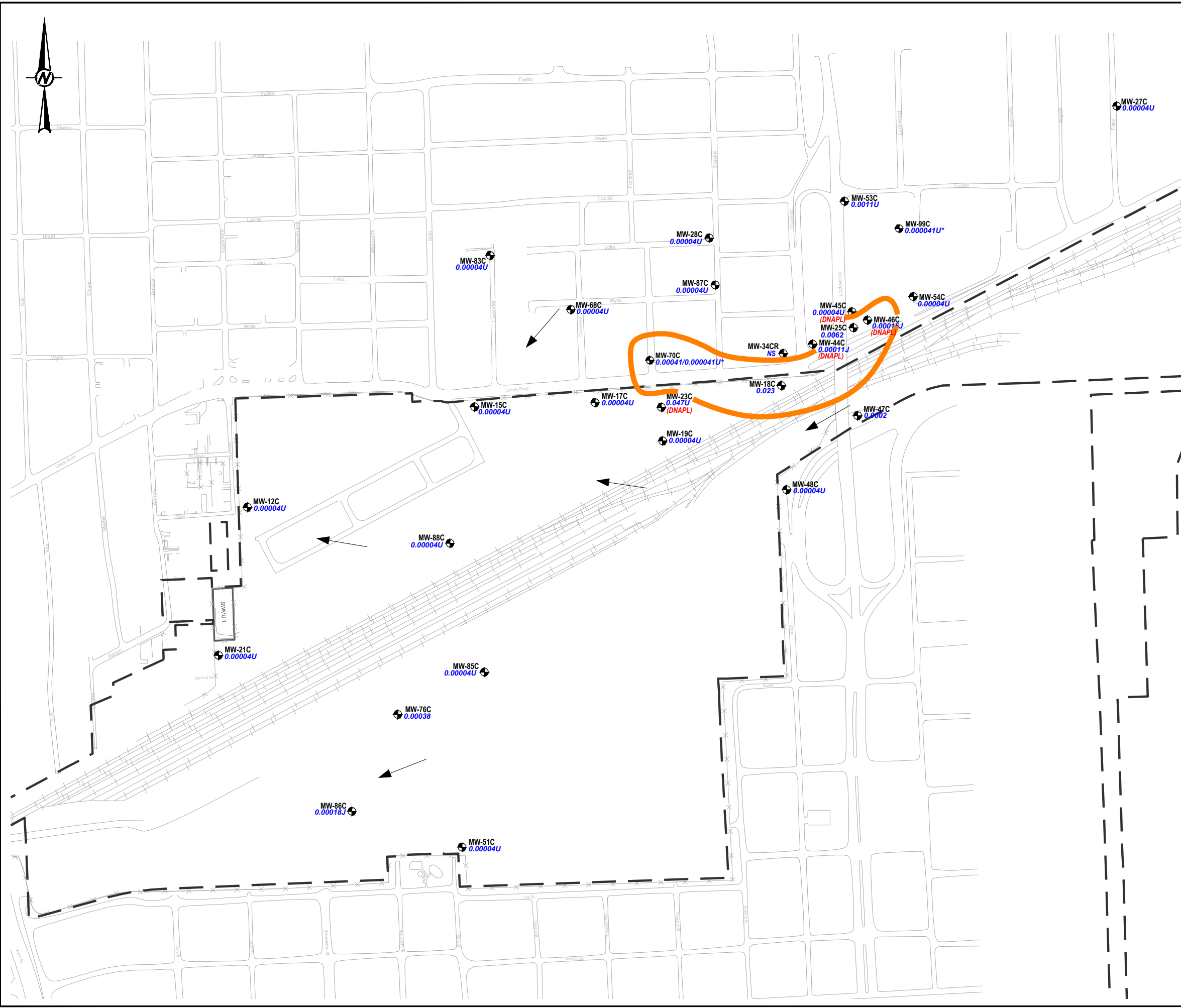
<b>CONSULTANT</b>	YYYY-MM-DD	2020-04-30
<b>DESIGNED</b>		
<b>PREPARED</b>	RS	
<b>REVIEWED</b>	MH	
<b>APPROVED</b>	ECM	

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5B-15

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erand\data\Projects - Round Rock\2019\01\19232 - HWPW\2020-3\March - File Name: 5B15 - 5B19 Groundwater COC Concentration Map C-TZ - January 2020.dwg | Last Edited By: rsalazar | Date: 2020-04-30 | Time: 12:21:14 PM | Printed By: RSalazar | Date: 2020-04-30 | Time: 12:28:50 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- <0.002 2,4-DIMETHYLPHENOL CONCENTRATION (mg/L)
- 2,4-DIMETHYLPHENOL CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.49 mg/L AND 1.5 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

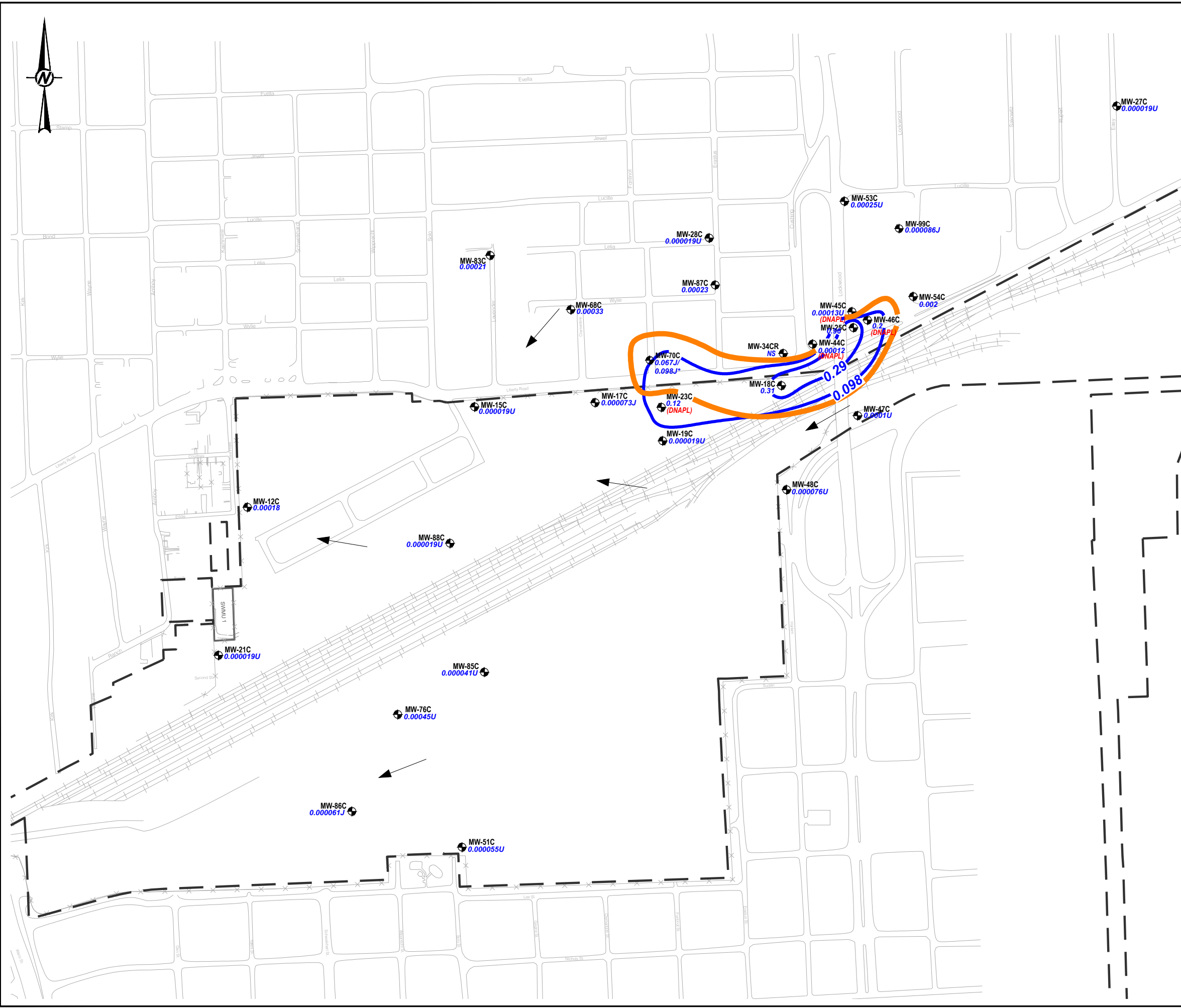
**TITLE**  
 GROUNDWATER COC CONCENTRATION MAP - C-TZ  
 2,4-DIMETHYLPHENOL - JANUARY-MARCH 2020

CONSULTANT	YYYY-MM-DD	2020-04-30
<b>GOLDER</b> <small>TEXAS GEOSCIENCE FIRM NO. 50969 TEXAS ENGINEERING FIRM NO. 2578</small>	DESIGNED	
	PREPARED	RS
	REVIEWED	MH
	APPROVED	ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5B-16

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\Projects - Round Rock\2019\01\19232 - HWPW\2020-3\March - 5B15 - 5B19 Groundwater COC Concentration Map C-TZ - January 2020.dwg | Last Edited By: rsalazar | Date: 2020-04-30 | Time: 12:21:14 PM | Printed By: RSalazar | Date: 2020-04-30 | Time: 12:20:08 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ C-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- <0.002 2-METHYLNAPHTHALENE CONCENTRATION (mg/L)
- 0.098** 2-METHYLNAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- ▭ GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.098 mg/L AND 0.29 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - C-TZ  
 2-METHYLNAPHTHALENE - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

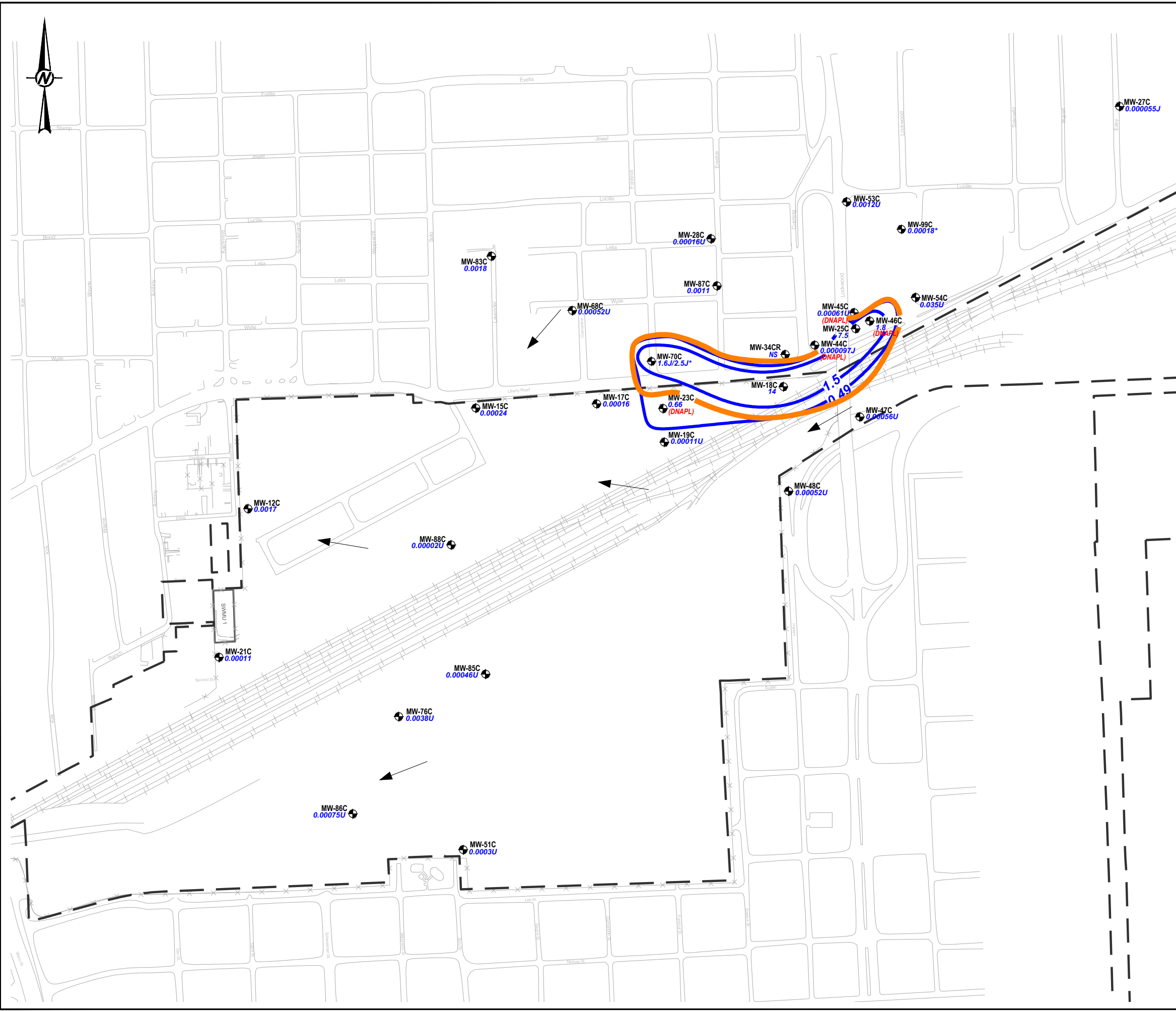
PROJECT NO. 19119232      REV. 0      FIGURE 5B-17

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B





Path: \\uswest\erand\data\Projects - Round Rock\2019\01\19232 - HWPW\2020-3\March - 5B19 Groundwater COC Concentration Map C-TZ - January 2020.dwg | Last Edited By: rsalazar | Date: 2020-04-30 | Time: 12:21:14 PM | Printed By: RSalazar | Date: 2020-04-30 | Time: 12:29:41 PM

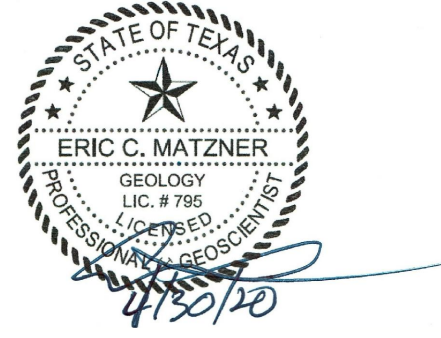


**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ C-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- <0.002 NAPHTHALENE CONCENTRATION (mg/L)
- 0.49** NAPHTHALENE CONCENTRATION CONTOUR (mg/L)
- ➔ INFERRED GROUNDWATER FLOW DIRECTION
- ▭ GW PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL (0.49 mg/L AND 1.5 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

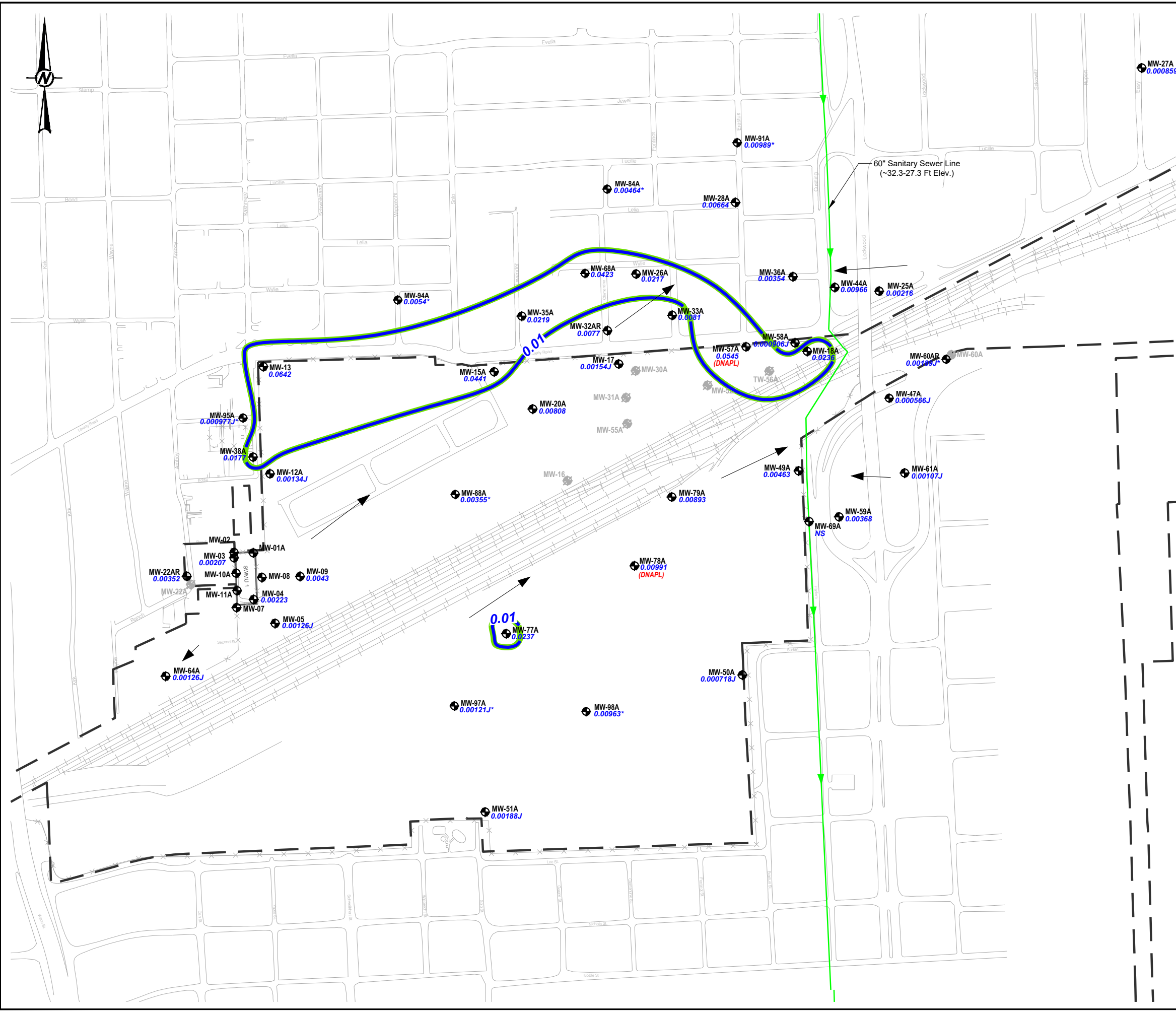
TITLE  
**GROUNDWATER COC CONCENTRATION MAP - C-TZ  
 NAPHTHALENE - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232      REV. 0      FIGURE 5B-19

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\data\projects - Round Rock\2019\01\19232 - HWPW\2020-3\March - Arsenic C/C Concentration Map A-TZ - January 2020.dwg | Last Edited By: realcar Date: 2020-04-30 Time: 12:27:02 PM | Printed By: RSalazar Date: 2020-04-30 Time: 12:31:33 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ A-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- 0.0077 ARSENIC CONCENTRATION IN (mg/L)
- 0.01** ARSENIC CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- █ ARSENIC PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLD AT THE RAL AND C/I PCL (0.01 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

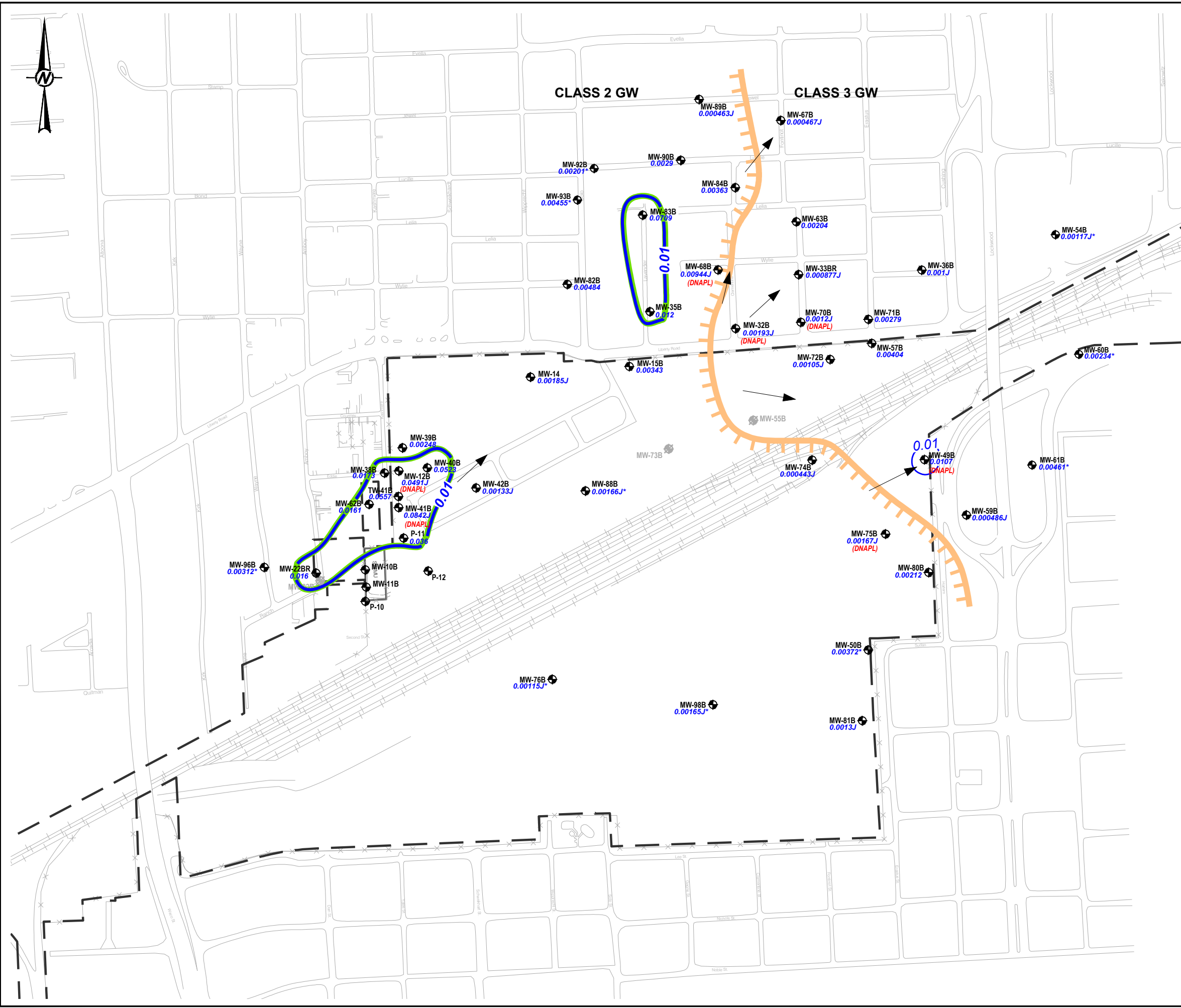
TITLE  
**GROUNDWATER C/C CONCENTRATION MAP - A-TZ  
 ARSENIC - JANUARY-MARCH 2020**

CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232  
 REV. 0  
 FIGURE 5B-20

1 in. IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

Path: \\uswest\erand\proj\19119232 - HWPP\2020-31 March - Arsenic CCG Concentration Map B-TZ B-CZ - January 2020.dwg | Last Edited By: malar | File Name: 5821 - Arsenic CCG Concentration Map B-TZ B-CZ - January 2020.dwg | Printed By: malar | Date: 2020-04-30 | Time: 12:28:22 PM |



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- × × × FENCE
- RAILROAD
- ⊕ B-TZ MONITORING WELL LOCATION
- ⊕ PLUGGED AND ABANDONED MONITORING WELL
- B-CZ (CLASS 3 GW)
- B-TZ (CLASS 2 GW)
- 0.001J ARSENIC CONCENTRATION (mg/L)
- 0.01** ARSENIC CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION
- ▭ ARSENIC PCLE ZONE

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLDED AT THE RAL AND C/I PCL:  
 CLASS 2 GW PCL: 0.01 mg/L  
 CLASS 3 GW PCL: 1 mg/L
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.

0 200 400  
 1" = 400' FEET

CLIENT  
 UNION PACIFIC RAILROAD CO.

PROJECT  
 HOUSTON WOOD PRESERVING WORKS

TITLE  
**GROUNDWATER COC CONCENTRATION MAP - B-CZ/B-TZ ARSENIC - JANUARY-MARCH 2020**

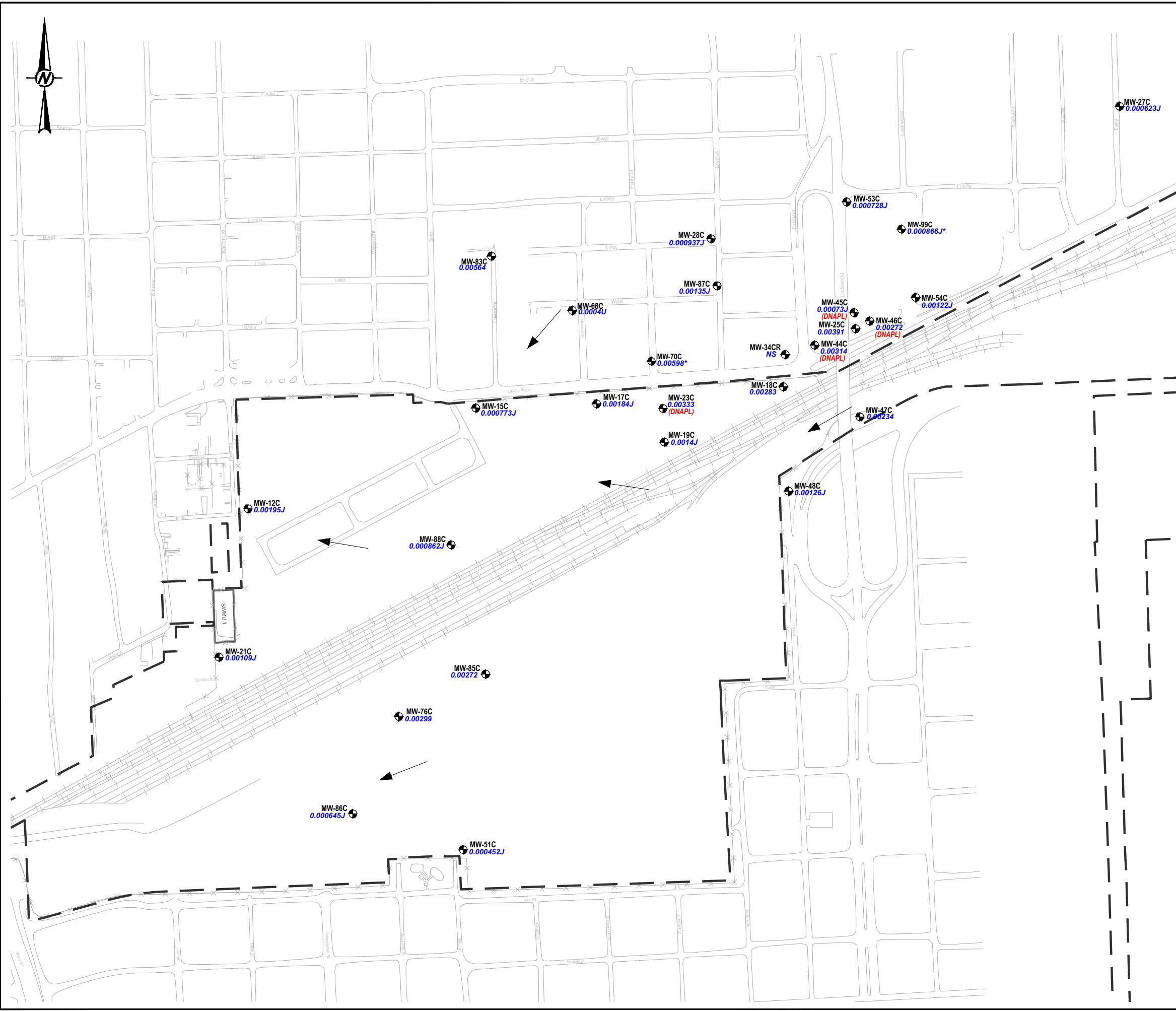
CONSULTANT	YYYY-MM-DD	2020-04-30
DESIGNED		
PREPARED	RS	
REVIEWED	MH	
APPROVED	ECM	

PROJECT NO. 19119232 REV. 0 FIGURE 5B-21

1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



Path: \\uswest\erand\data\Projects - Round Rock\2019\01\19232 - HWPM\2020-3\March - Arsenic C/C Concentration Map C-TZ - January 2020.dwg | File Name: 5B22 - Arsenic C/C Concentration Map C-TZ - January 2020.dwg | Last Edited By: reidmiller | Date: 2020-04-30 | Time: 12:28:46 PM | Printed By: reidmiller | Date: 2020-04-30 | Time: 12:32:33 PM



**LEGEND**

- UPRR PROPERTY BOUNDARY
- ROAD, PARKING LOT, SIDEWALK
- FENCE
- RAILROAD
- C-TZ MONITORING WELL LOCATION
- PLUGGED AND ABANDONED MONITORING WELL
- 0.00272 ARSENIC CONCENTRATION (mg/L)
- ARSENIC CONCENTRATION CONTOUR (mg/L)
- INFERRED GROUNDWATER FLOW DIRECTION

- NOTE(S)**
1. DNAPL - DENSE NON-AQUEOUS PHASE LIQUIDS DETECTED IN MONITORING WELL (JANUARY 2020).
  2. NS - NOT SAMPLED
  3. CONTOURS ARE BOLD AT THE RAL AND C/I PCL (0.01 mg/L)
  4. \* - SAMPLE COLLECTED IN MARCH 2020.

**REFERENCE(S)**  
 BASE MAP FROM ERM-SOUTHWEST, INC APAR ADDENDUM, FIG 3-1, DATED JUNE 2004.



**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

**TITLE**  
 GROUNDWATER C/C CONCENTRATION MAP - C-TZ  
 ARSENIC - JANUARY-MARCH 2020

CONSULTANT	DATE	BY
<b>GOLDER</b> <small>TEXAS GEOSCIENCE FIRM NO. 50969          TEXAS ENGINEERING FIRM NO. 2578</small>	YYYY-MM-DD	2020-04-30
	DESIGNED	
	PREPARED	RS
	REVIEWED	MH
	APPROVED	ECM

**PROJECT NO.** 19119232      **REV.** 0      **FIGURE** 5B-22

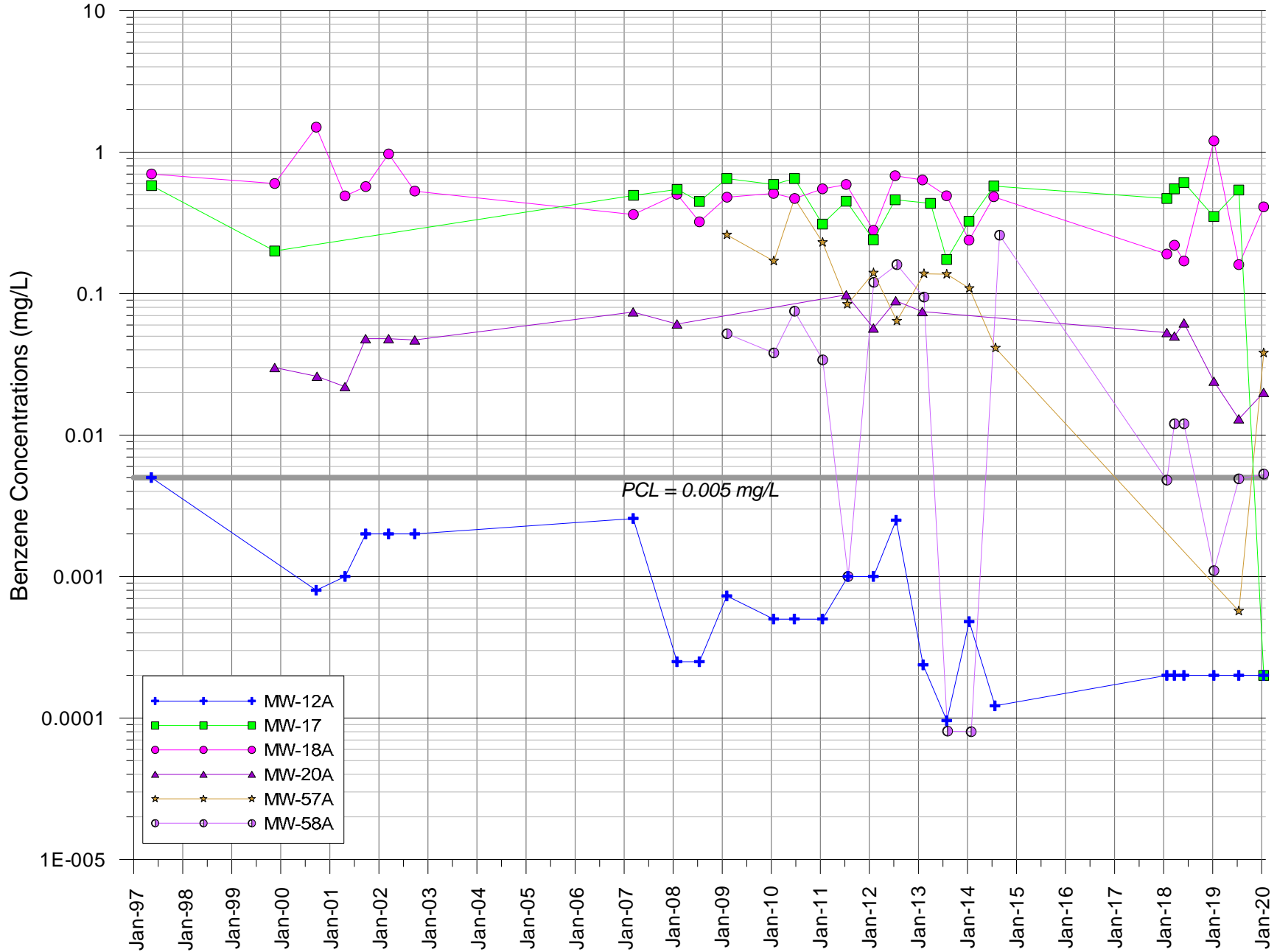
1 in IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B

**ATTACHMENT F**

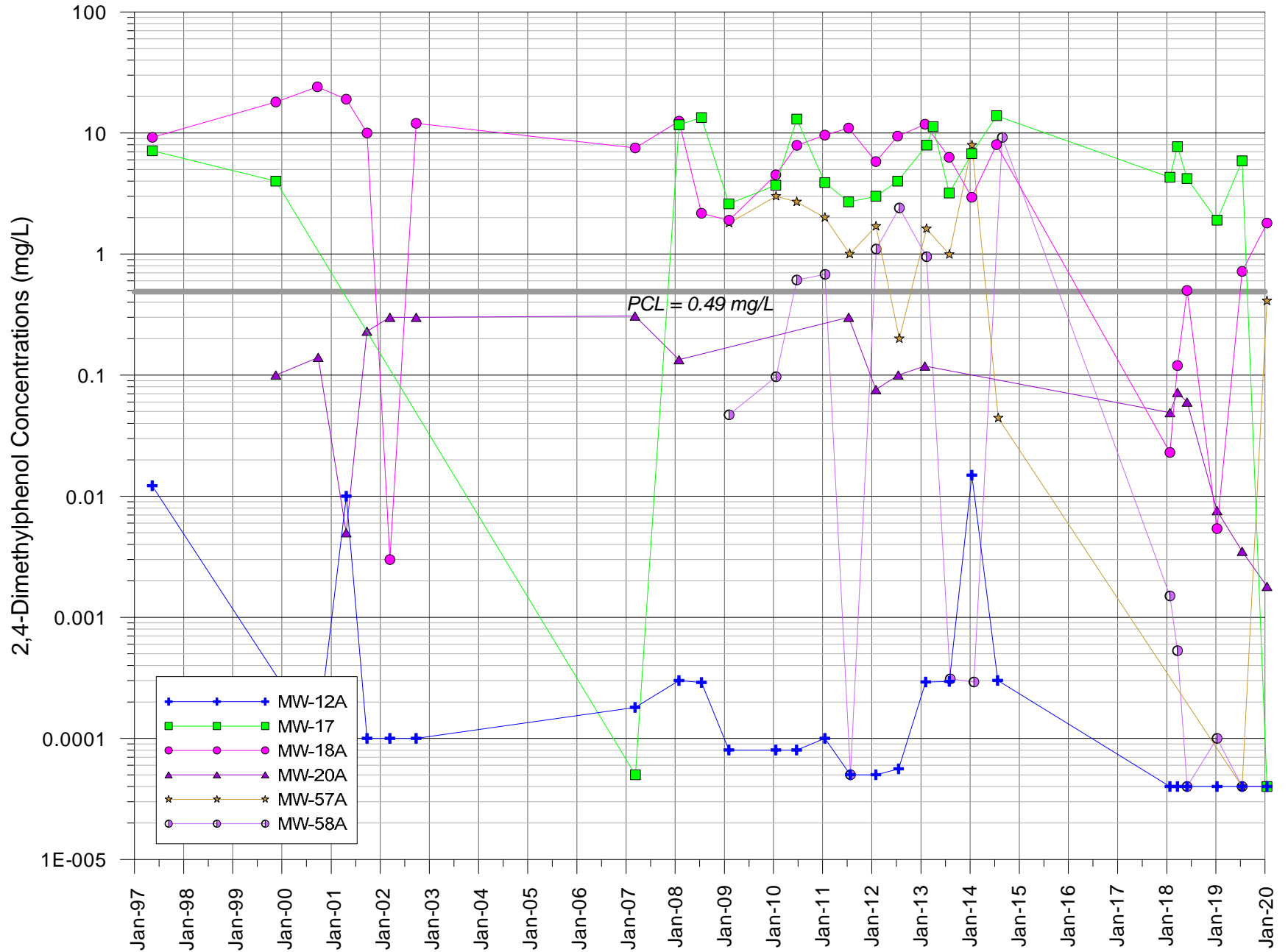
## **COC Concentration Graphs**



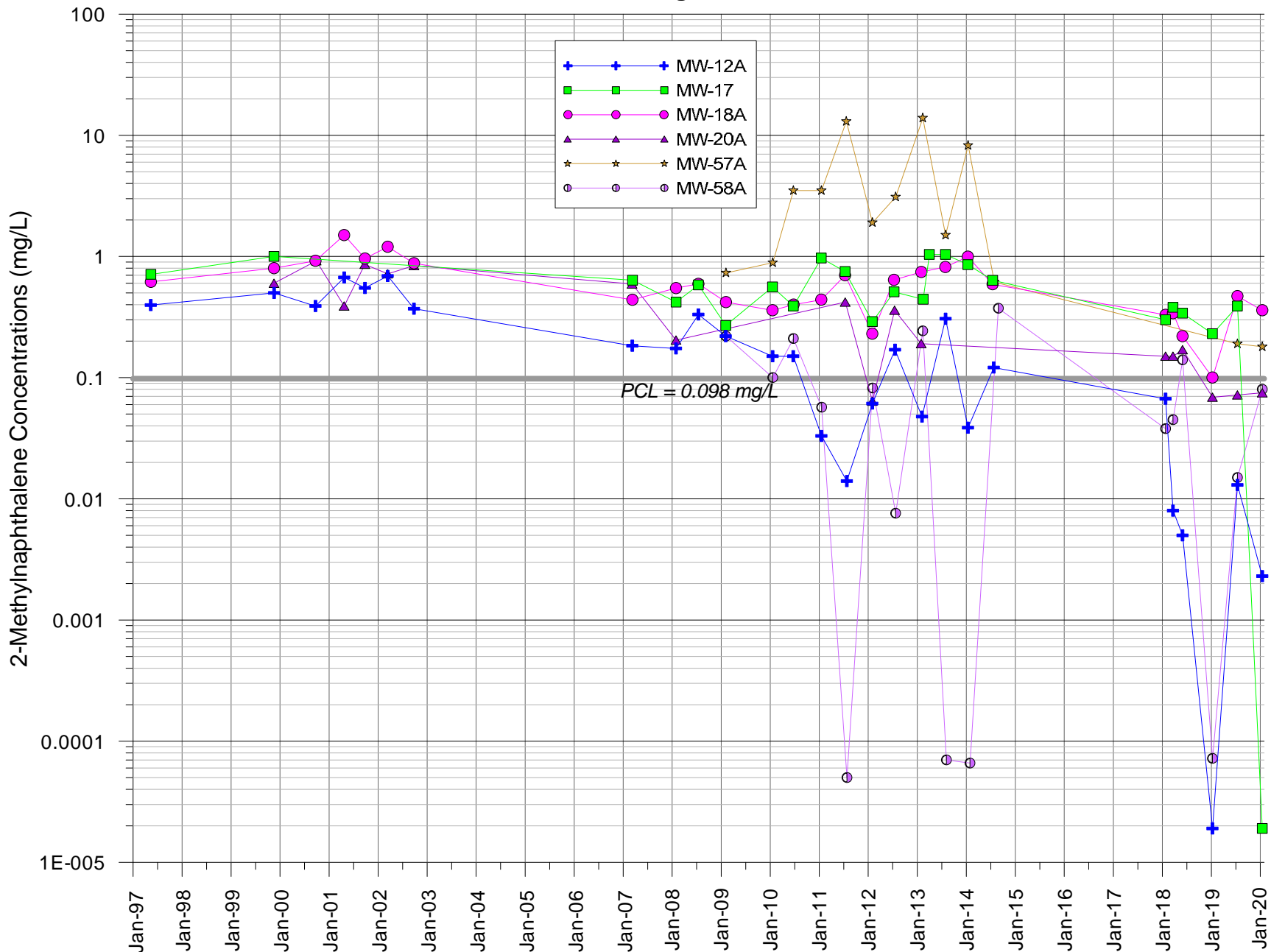
**ATTACHMENT 1B-1**  
**Benzene Concentrations at Source Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



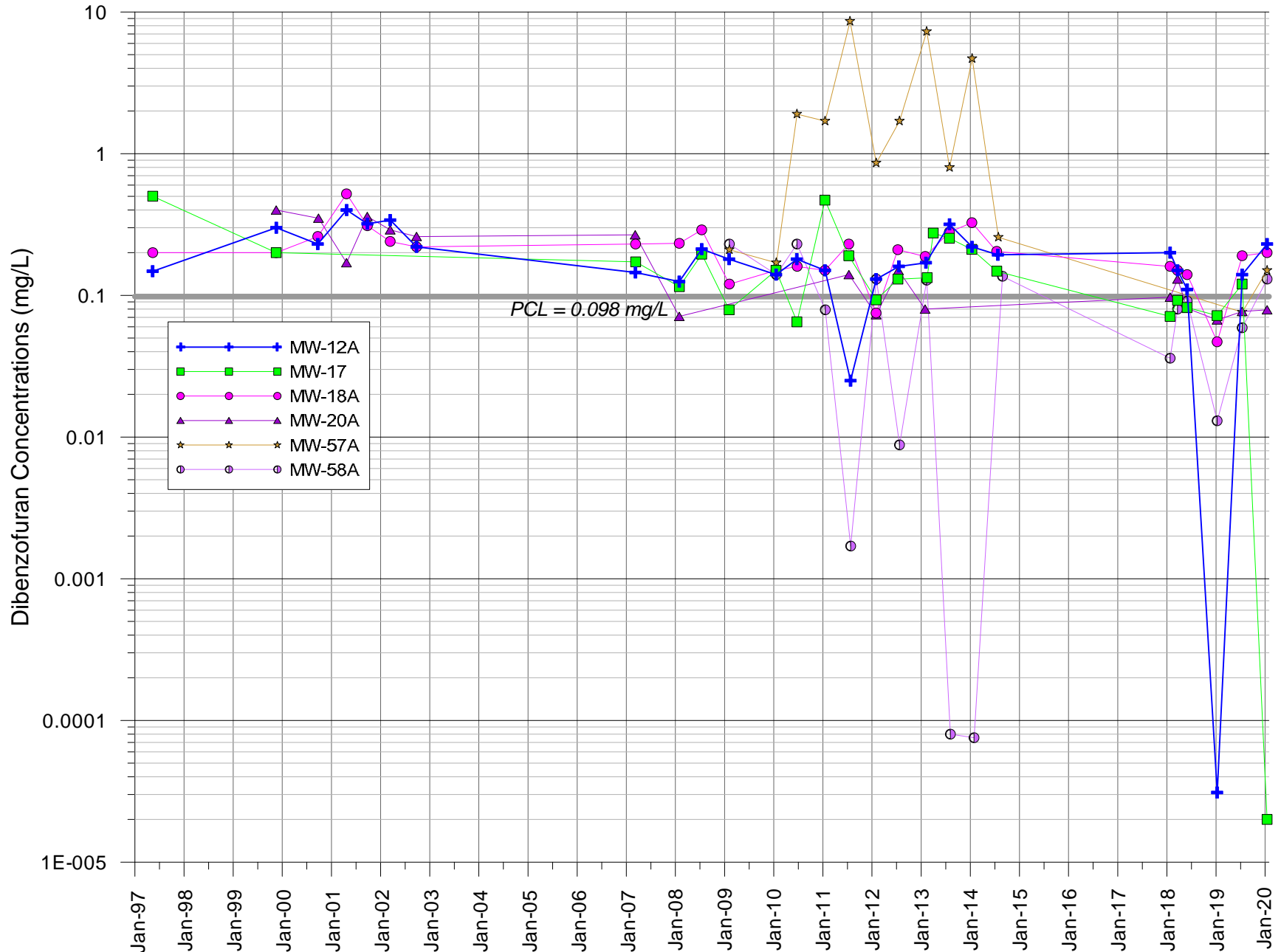
**ATTACHMENT 1B-2**  
**2,4-Dimethylphenol Concentrations at Source Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



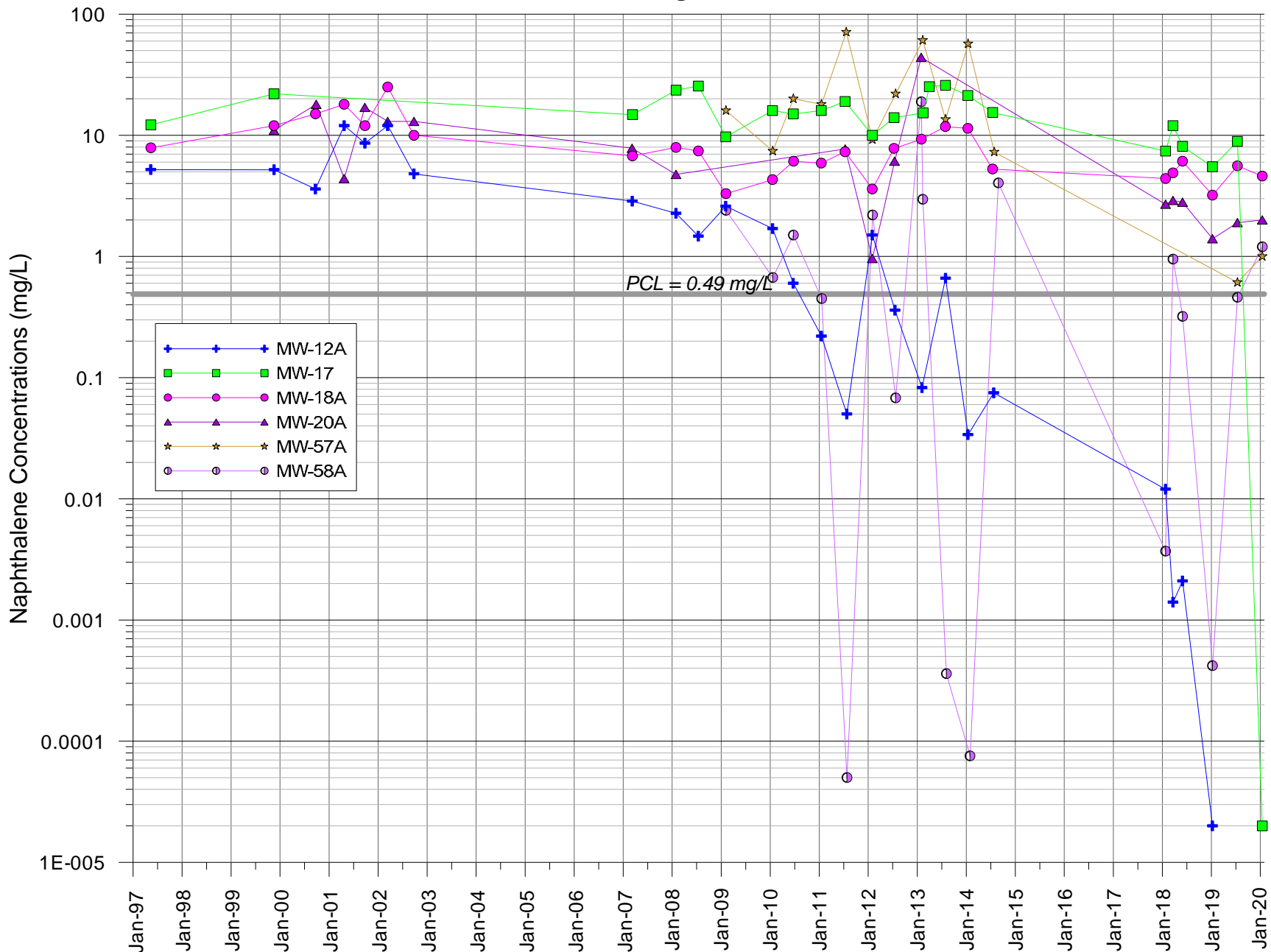
**ATTACHMENT 1B-3**  
**2-Methylnaphthalene Concentrations at Source Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-4**  
**Dibenzofuran Concentrations at Source Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



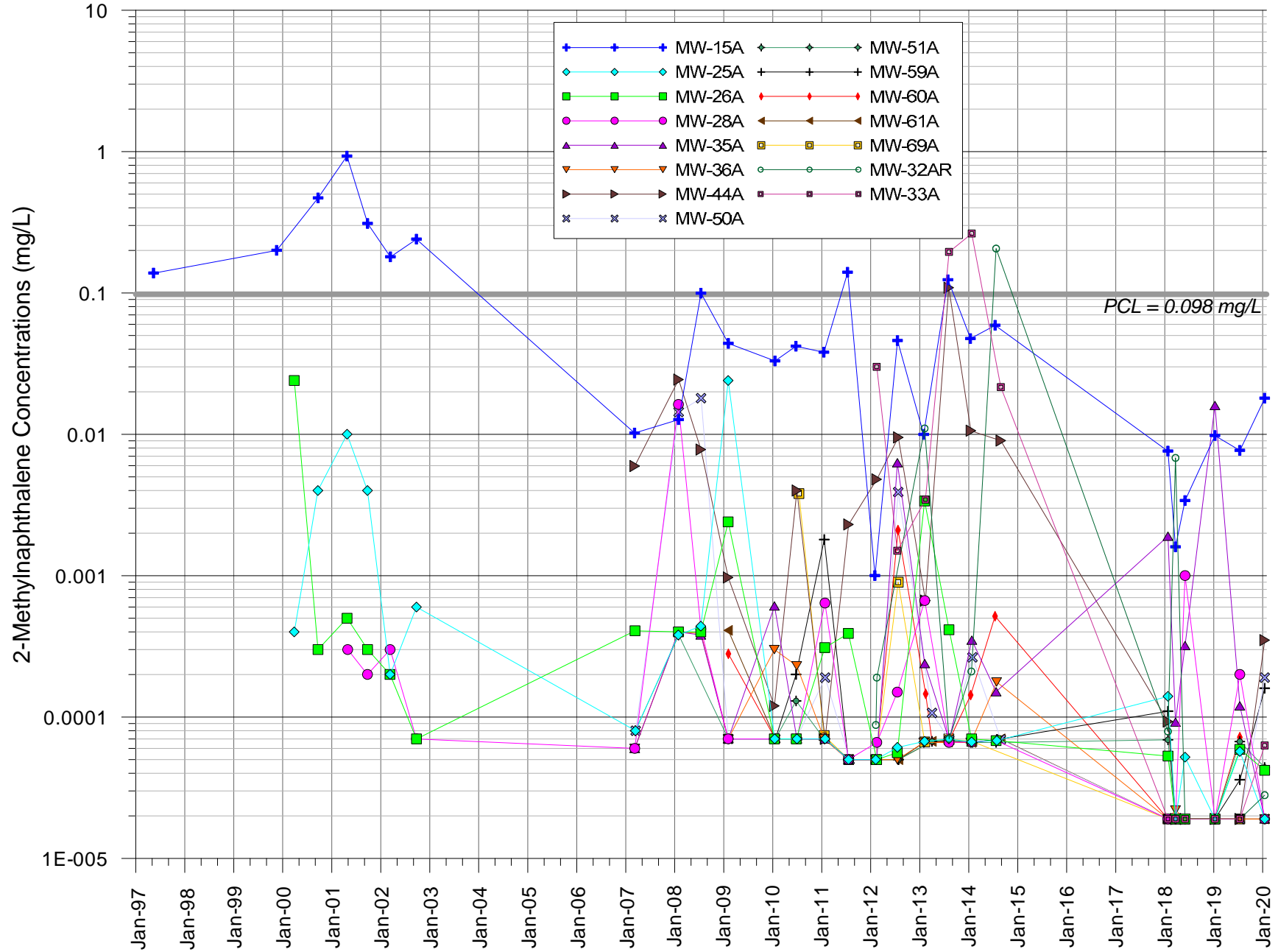
**ATTACHMENT 1B-5**  
**Naphthalene Concentrations at Source Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**







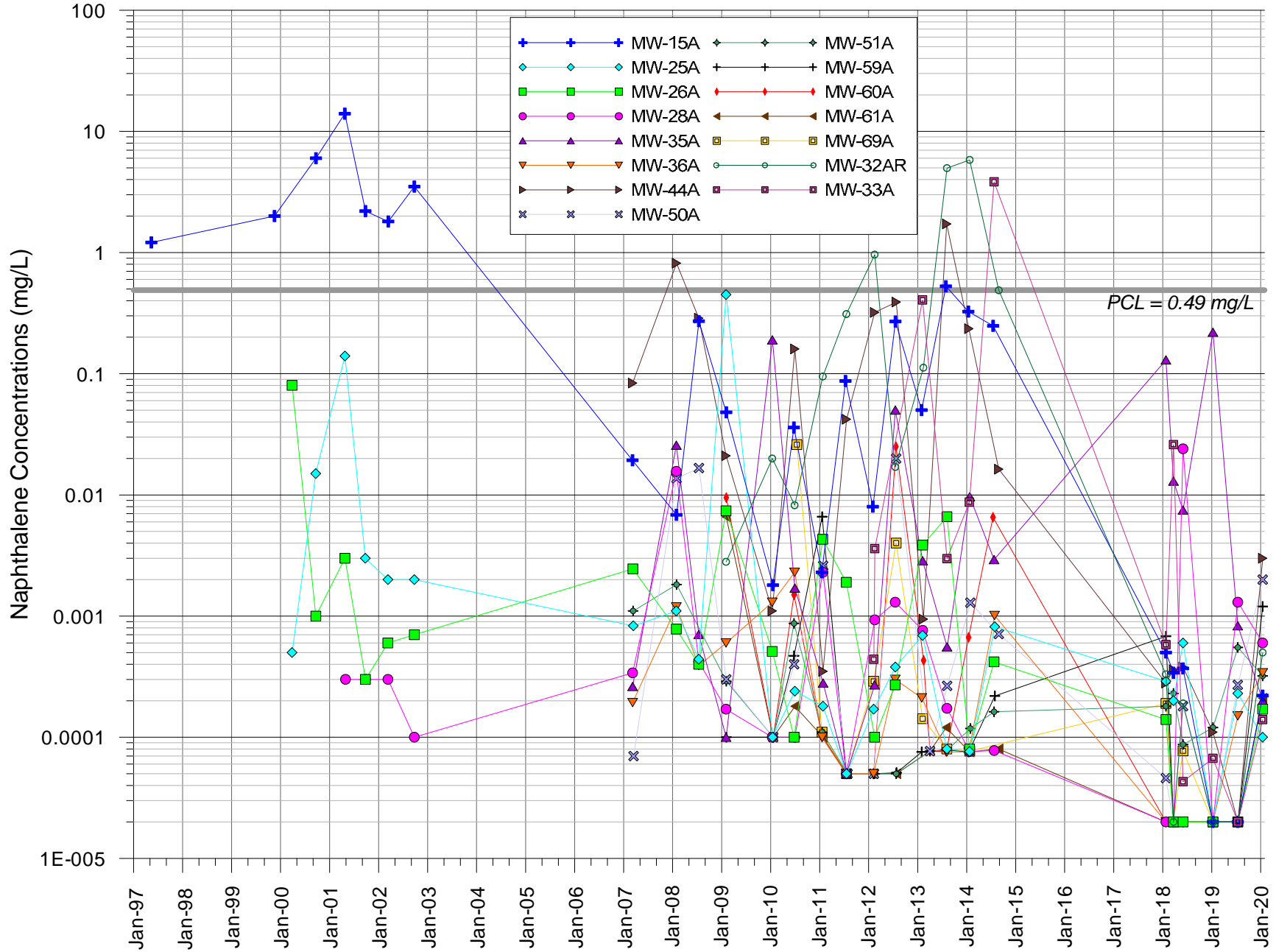
**ATTACHMENT 1B-8**  
**2-Methylnaphthalene Concentrations at Perimeter Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



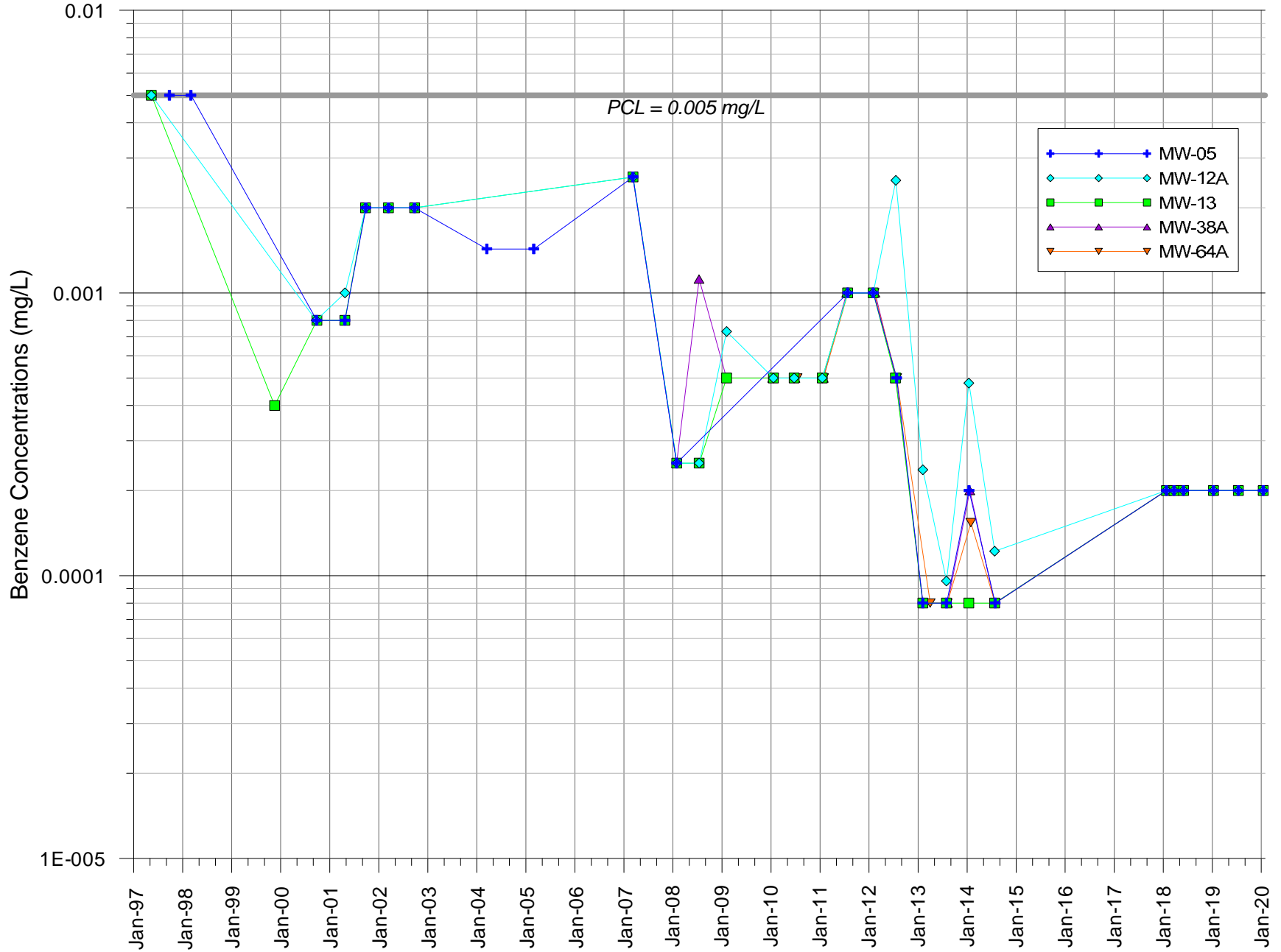




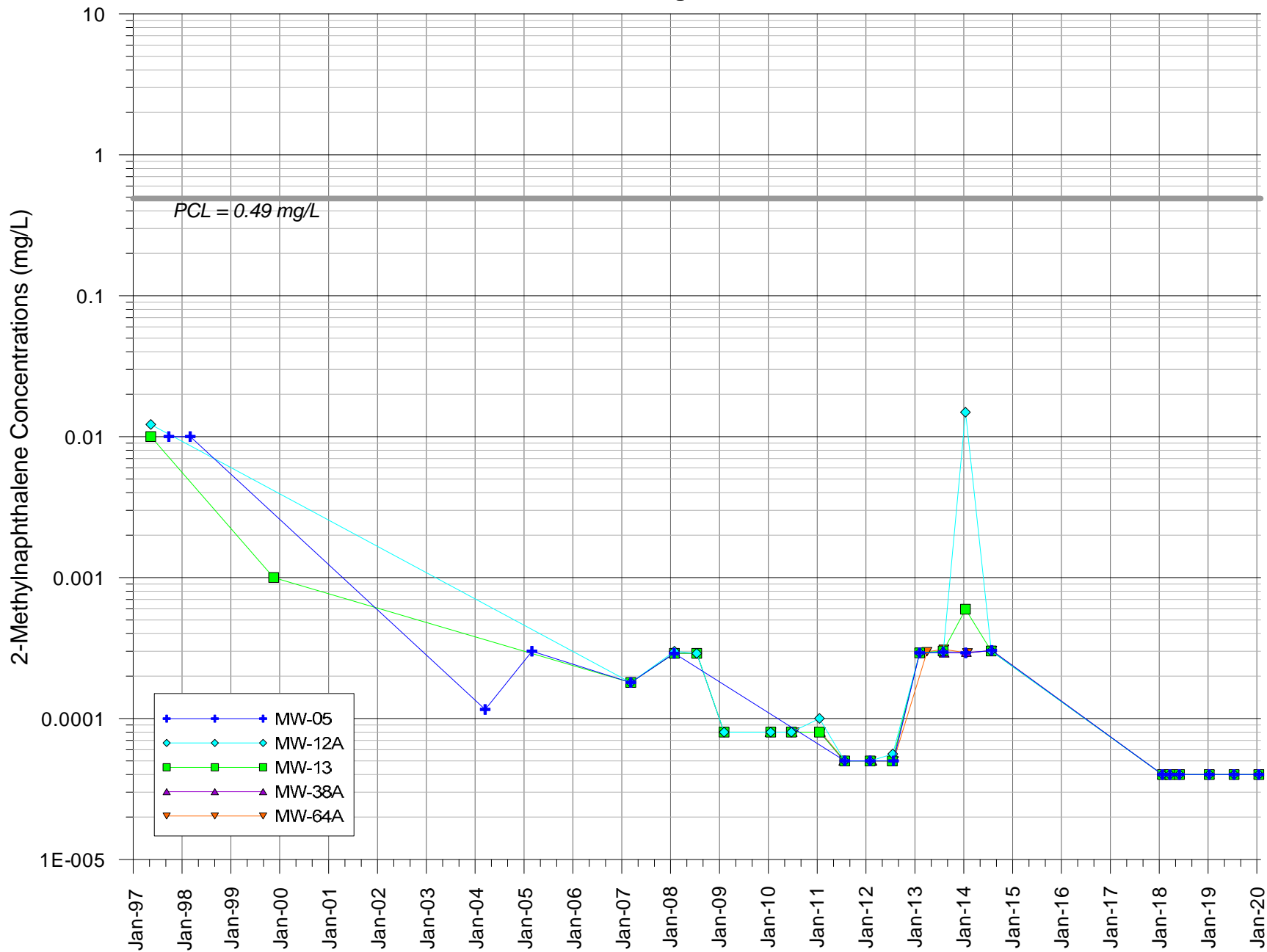
**ATTACHMENT 1B-10**  
**Naphthalene Concentrations at Perimeter Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



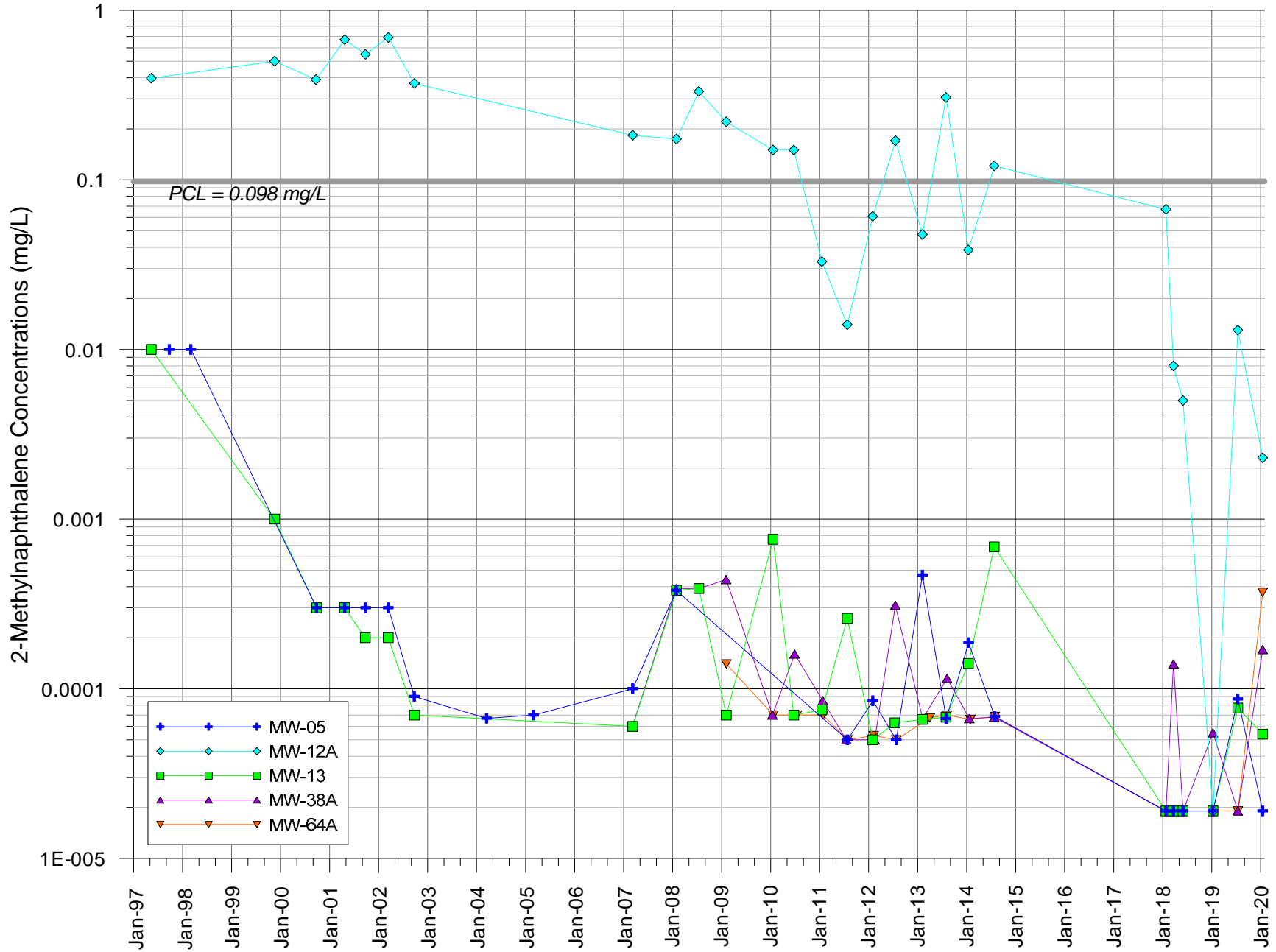
**ATTACHMENT 1B-11**  
**Benzene Concentrations at West End Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



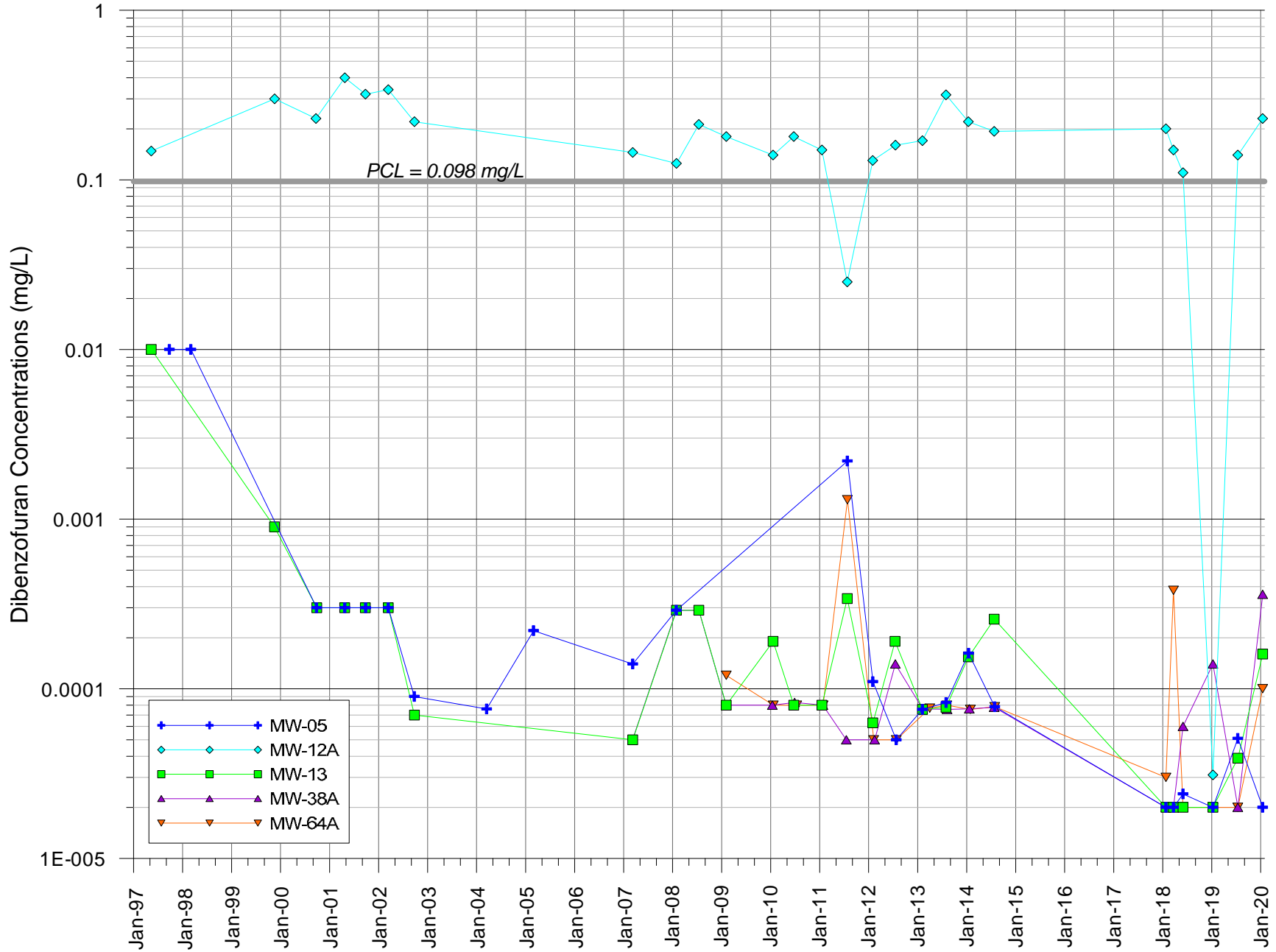
**ATTACHMENT 1B-12**  
**2,4-Dimethylphenol Concentrations at West End Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



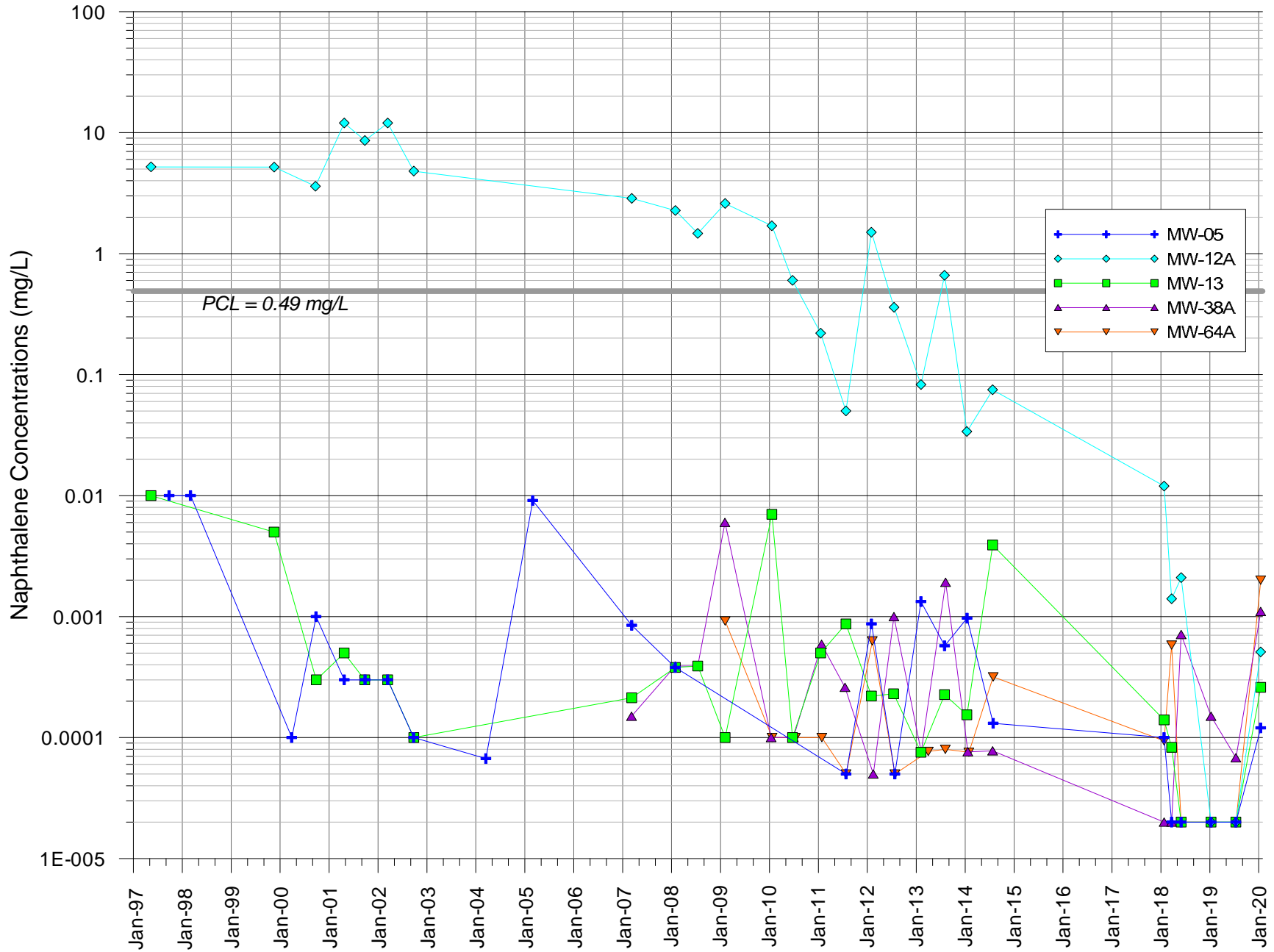
**ATTACHMENT 1B-13**  
**2-Methylnaphthalene Concentrations at West End Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



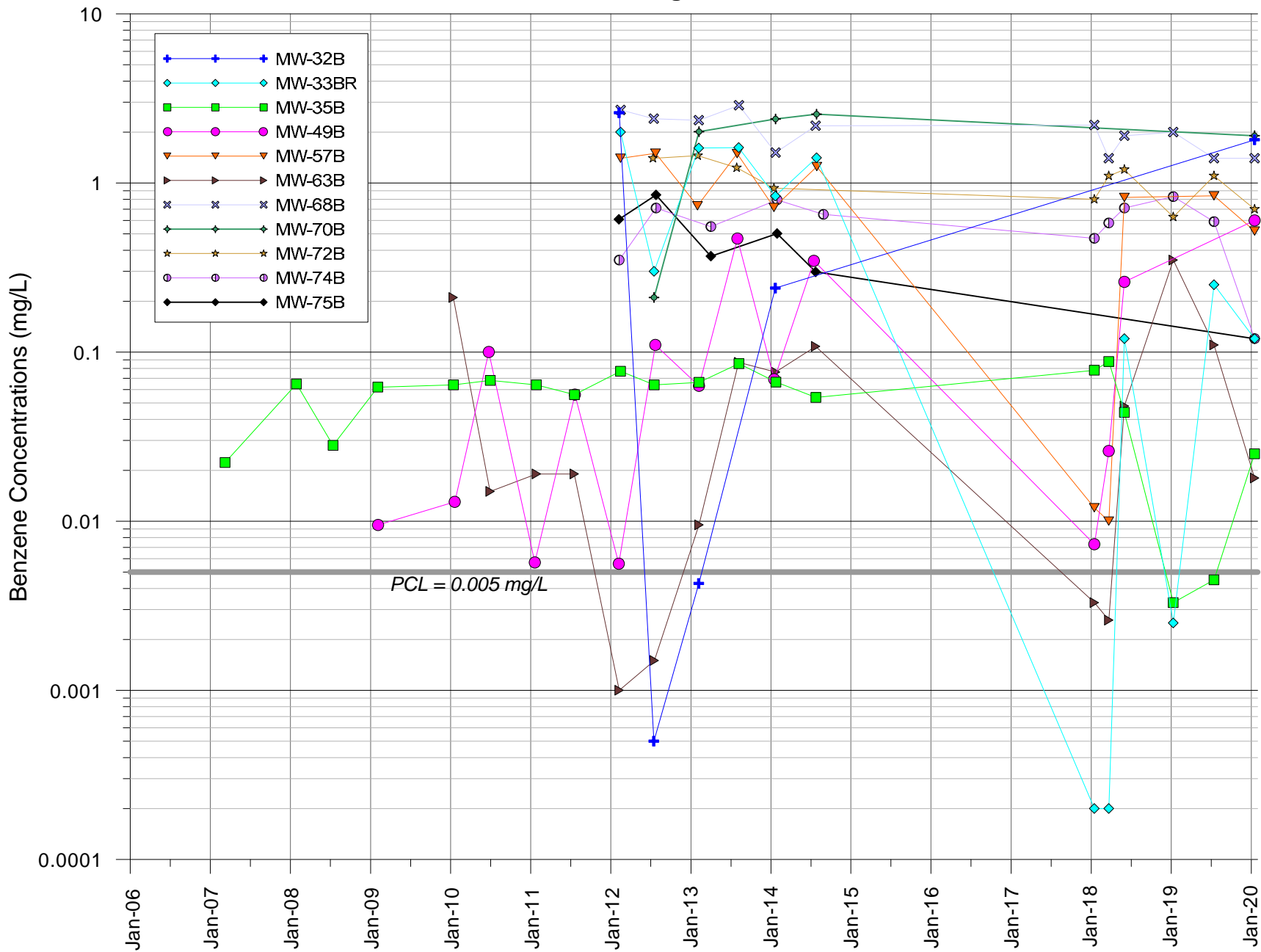
**ATTACHMENT 1B-14**  
**Dibenzofuran Concentrations at West End Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-15**  
**Naphthalene Concentrations at West End Area Wells - A-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

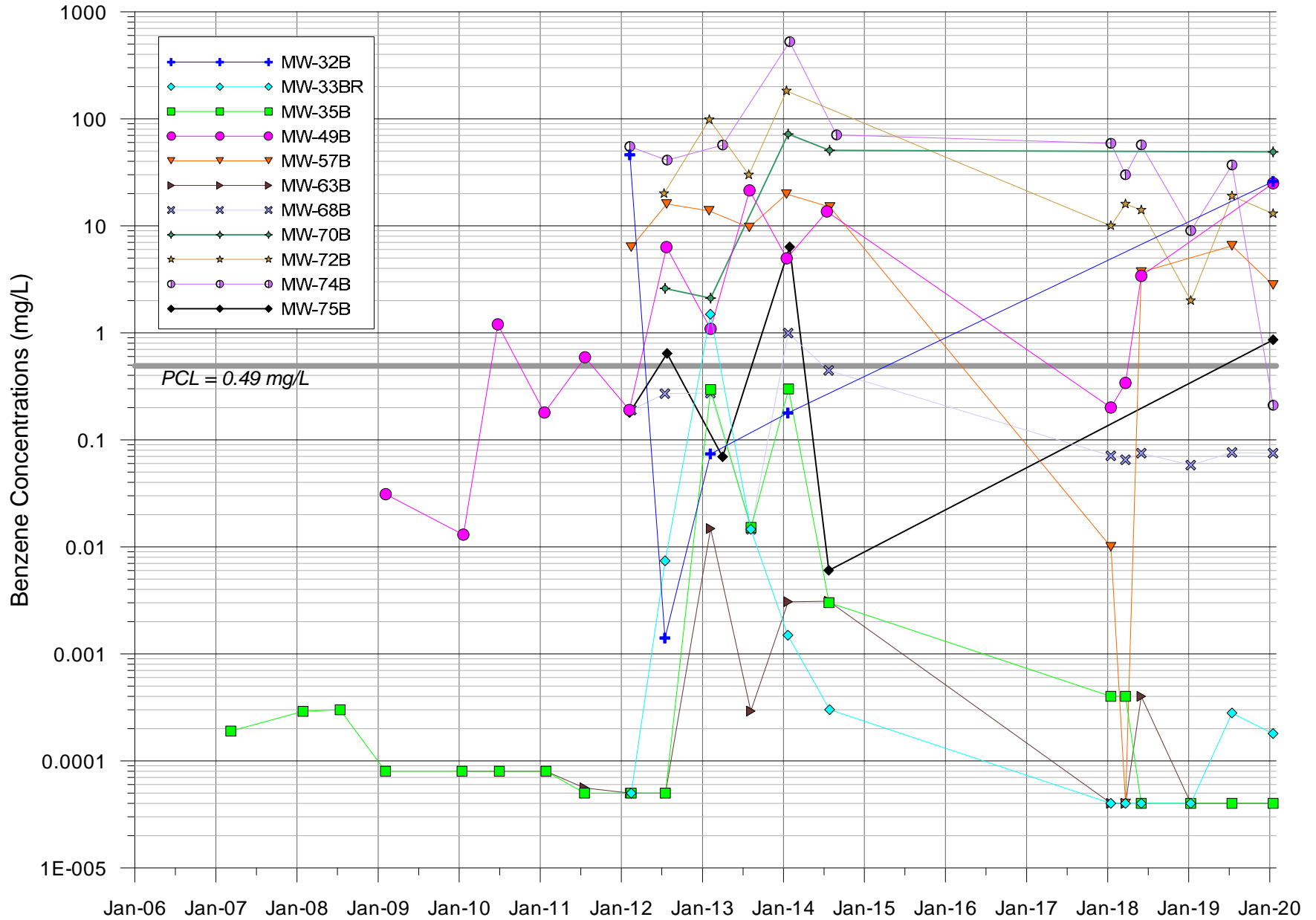


**ATTACHMENT 1B-16**  
**Benzene Concentrations at Source Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



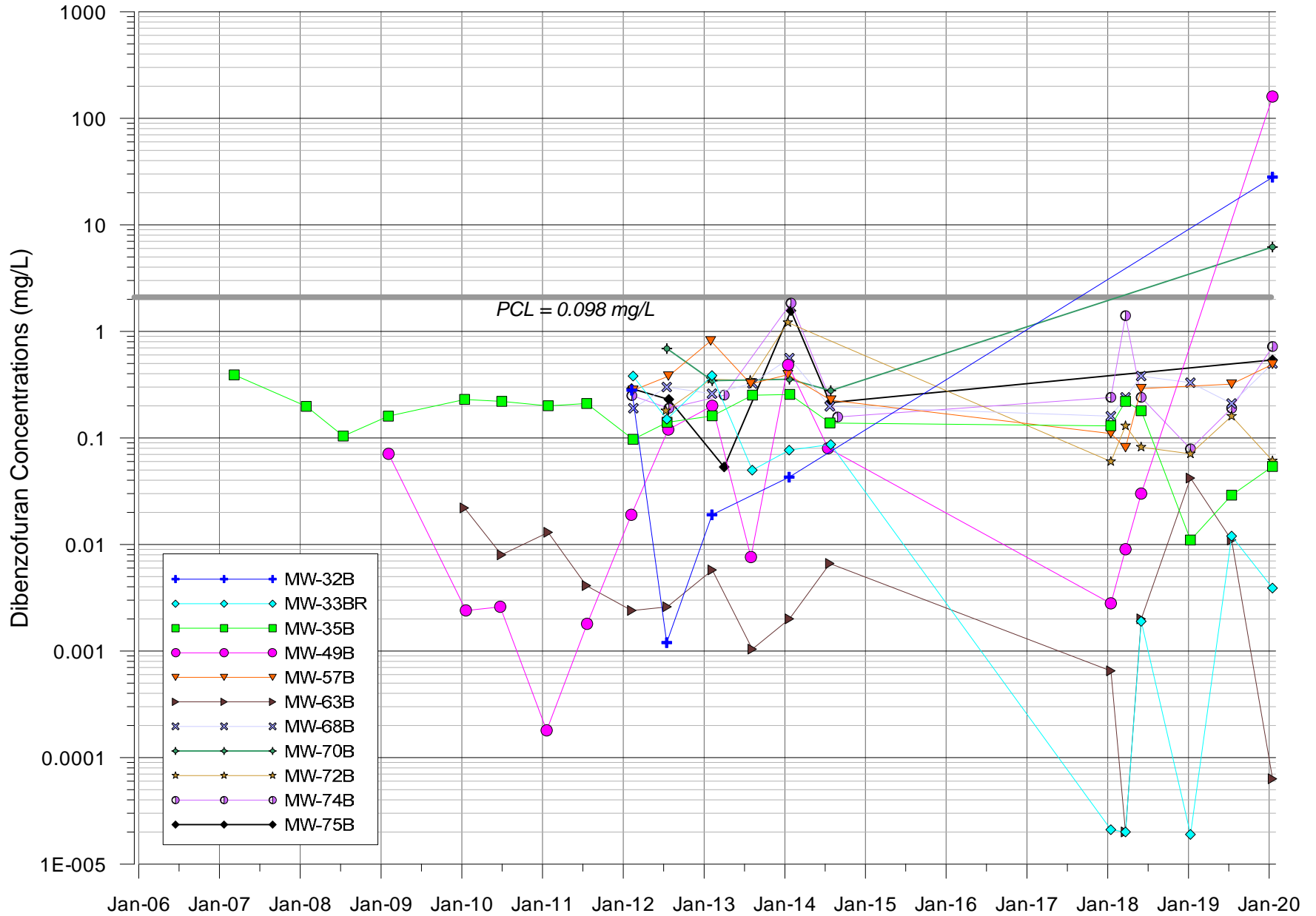


**ATTACHMENT 1B-17**  
**2,4-Dimethylphenol Concentrations at Source Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

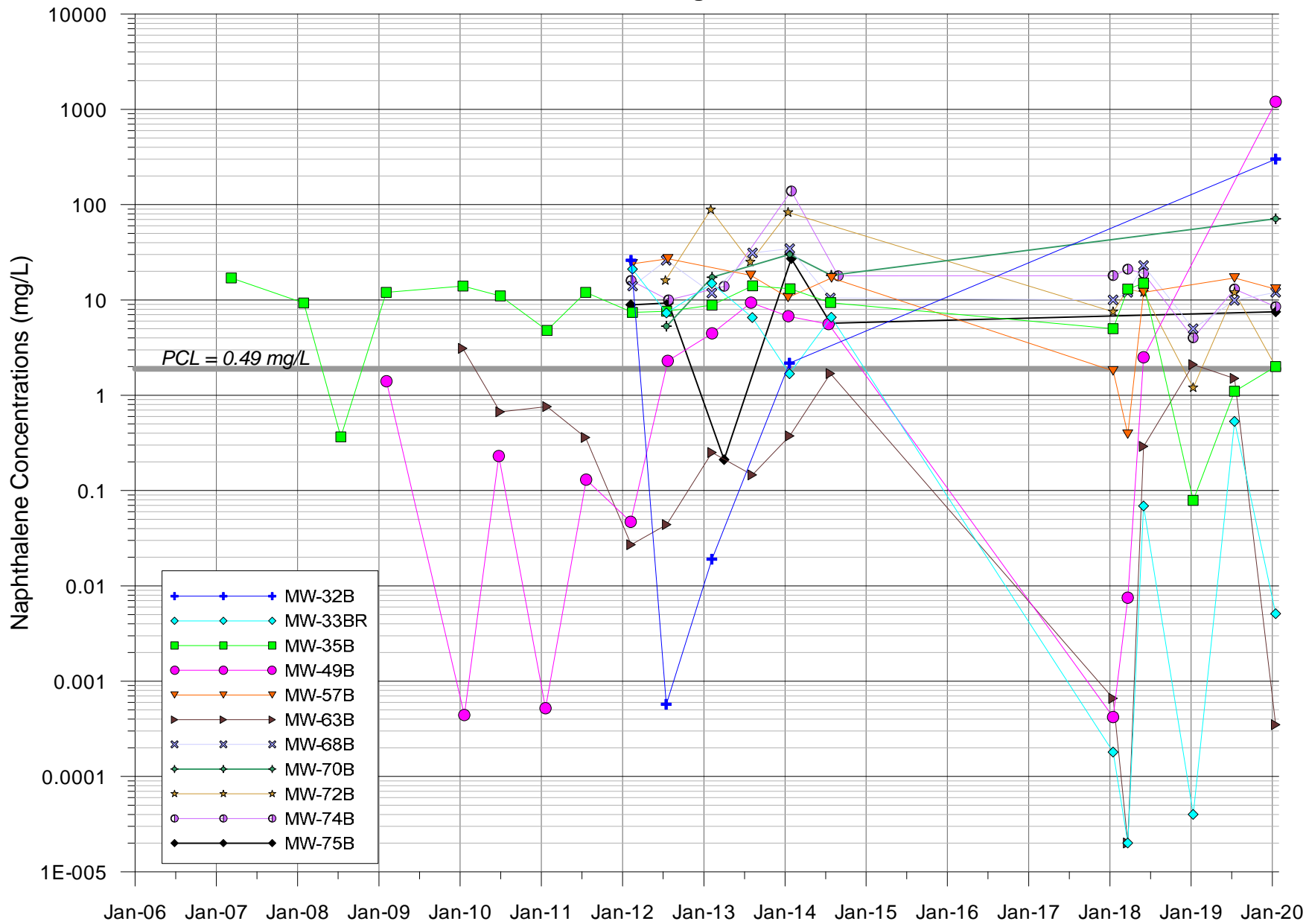




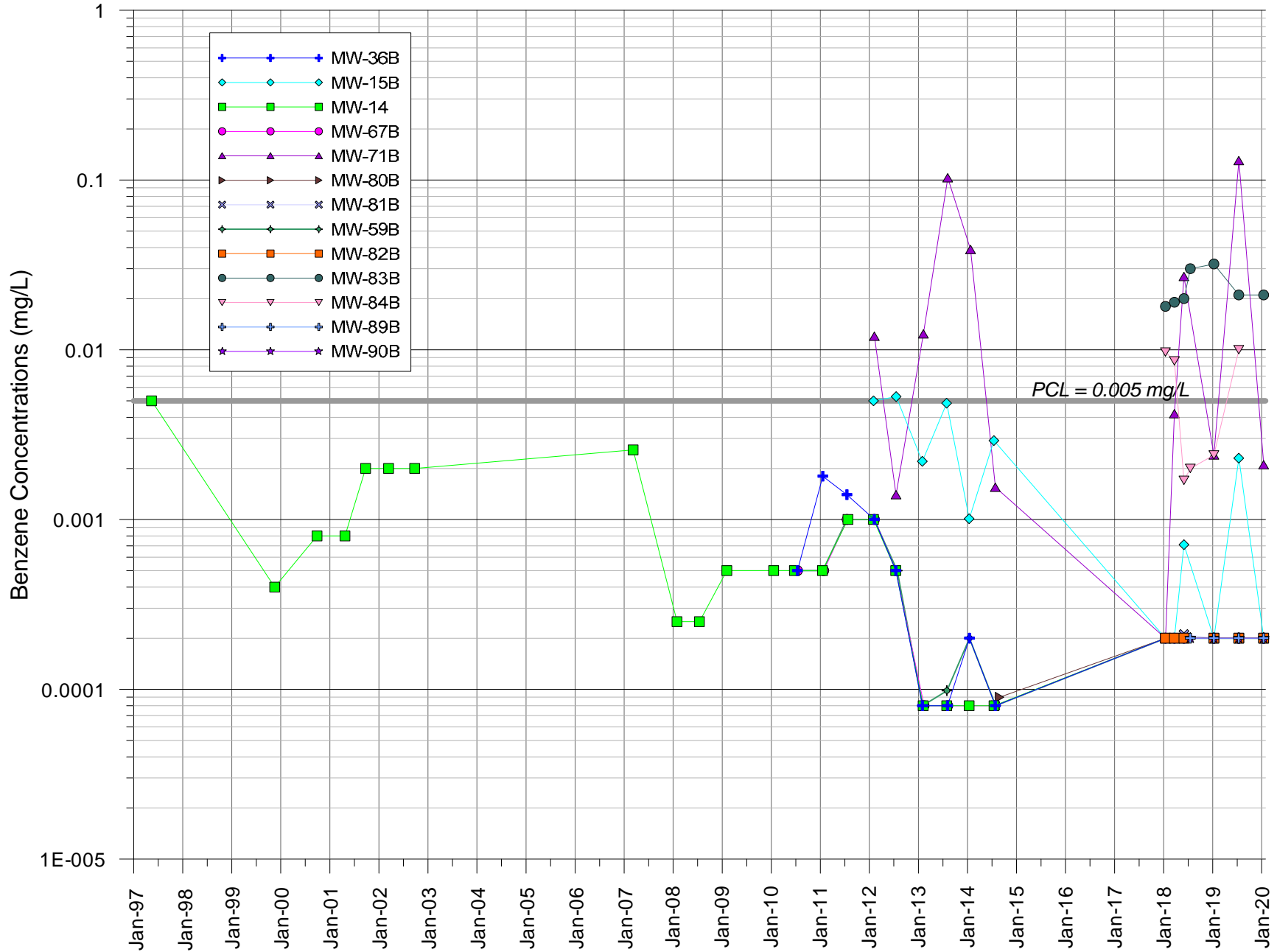
**ATTACHMENT 1B-19**  
**Dibenzofuran Concentrations at Source Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



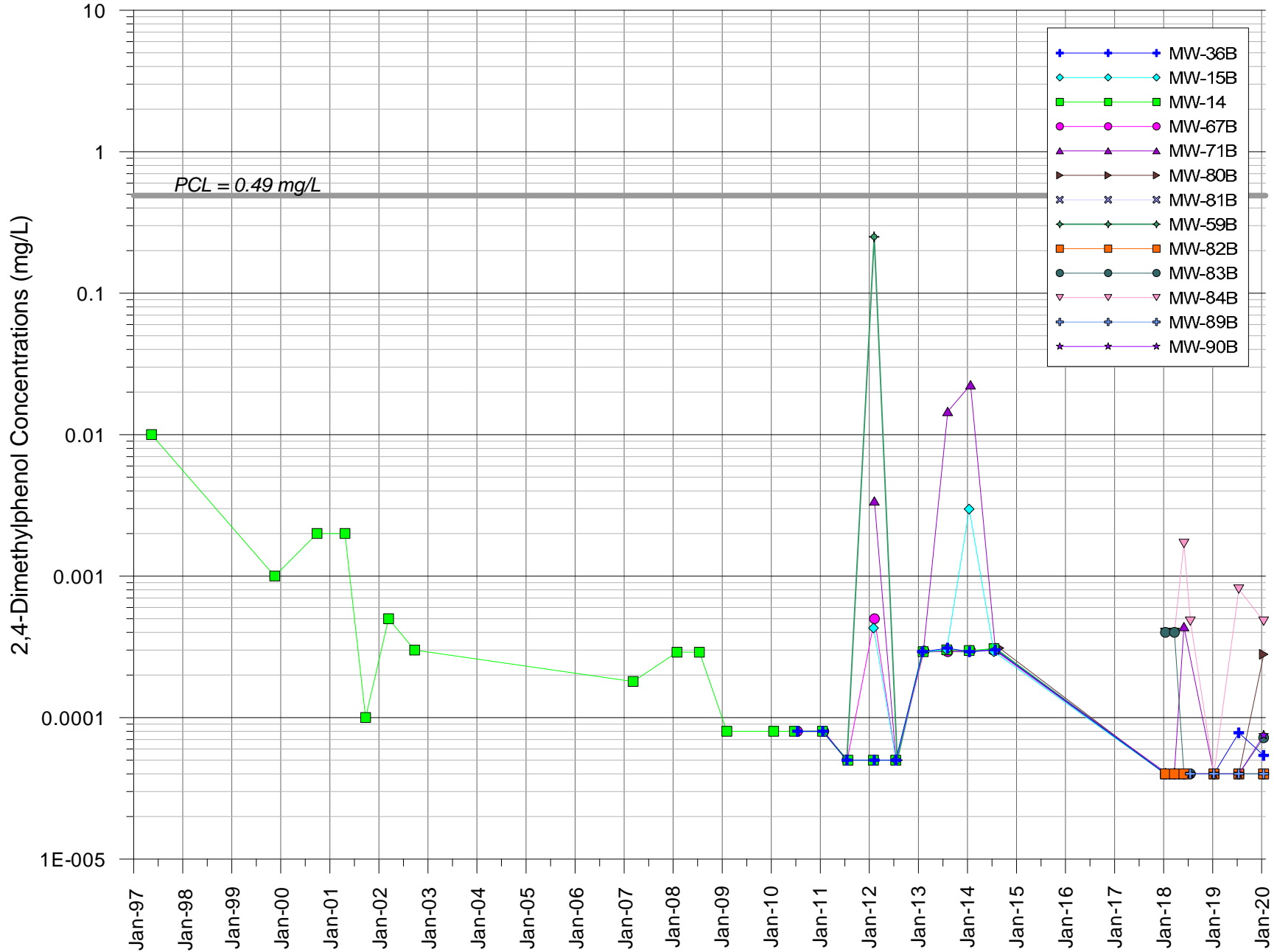
**ATTACHMENT 1B-20**  
**Naphthalene Concentrations at Source Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



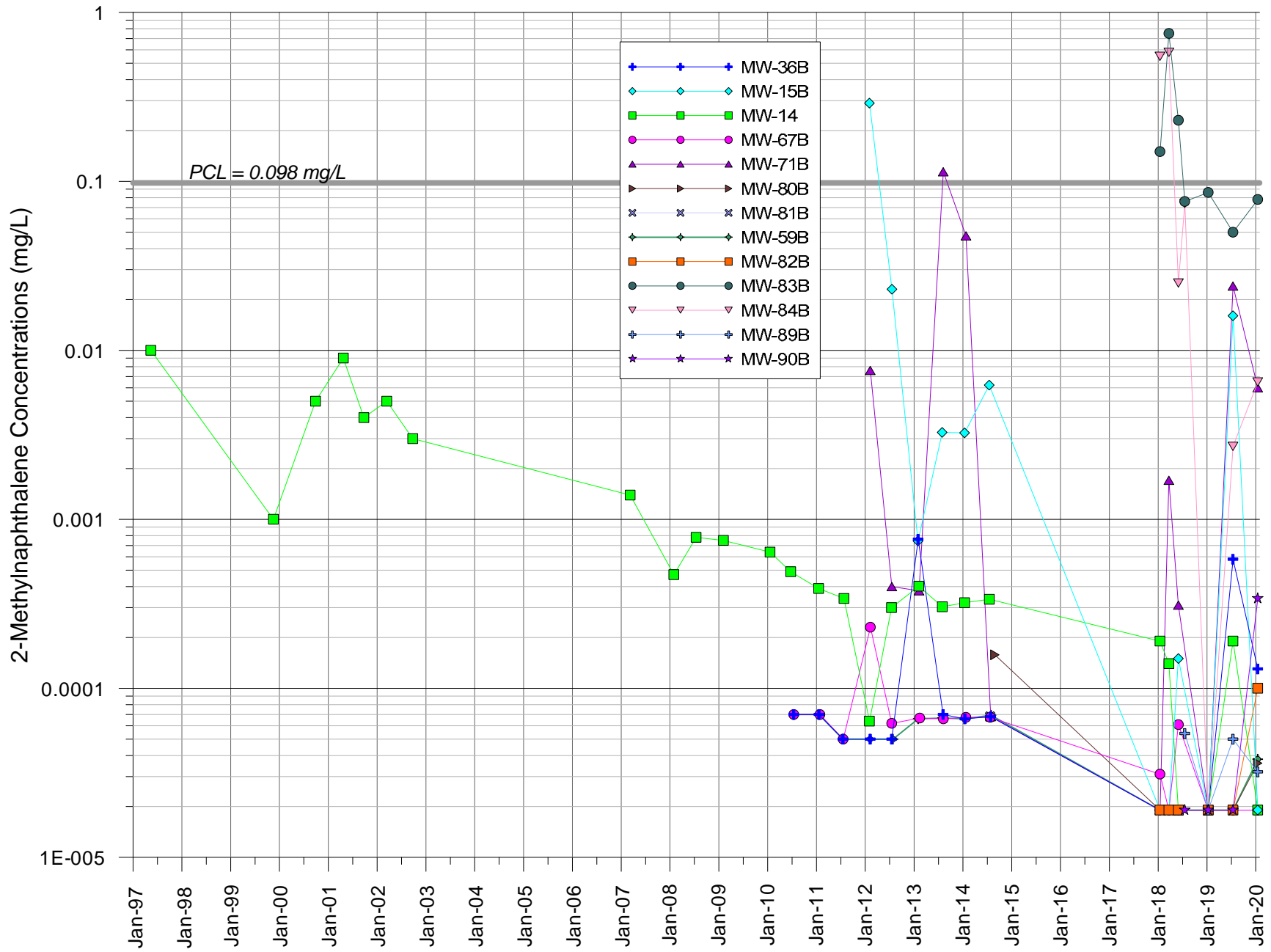
**ATTACHMENT 1B-21**  
**Benzene Concentrations at Perimeter Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



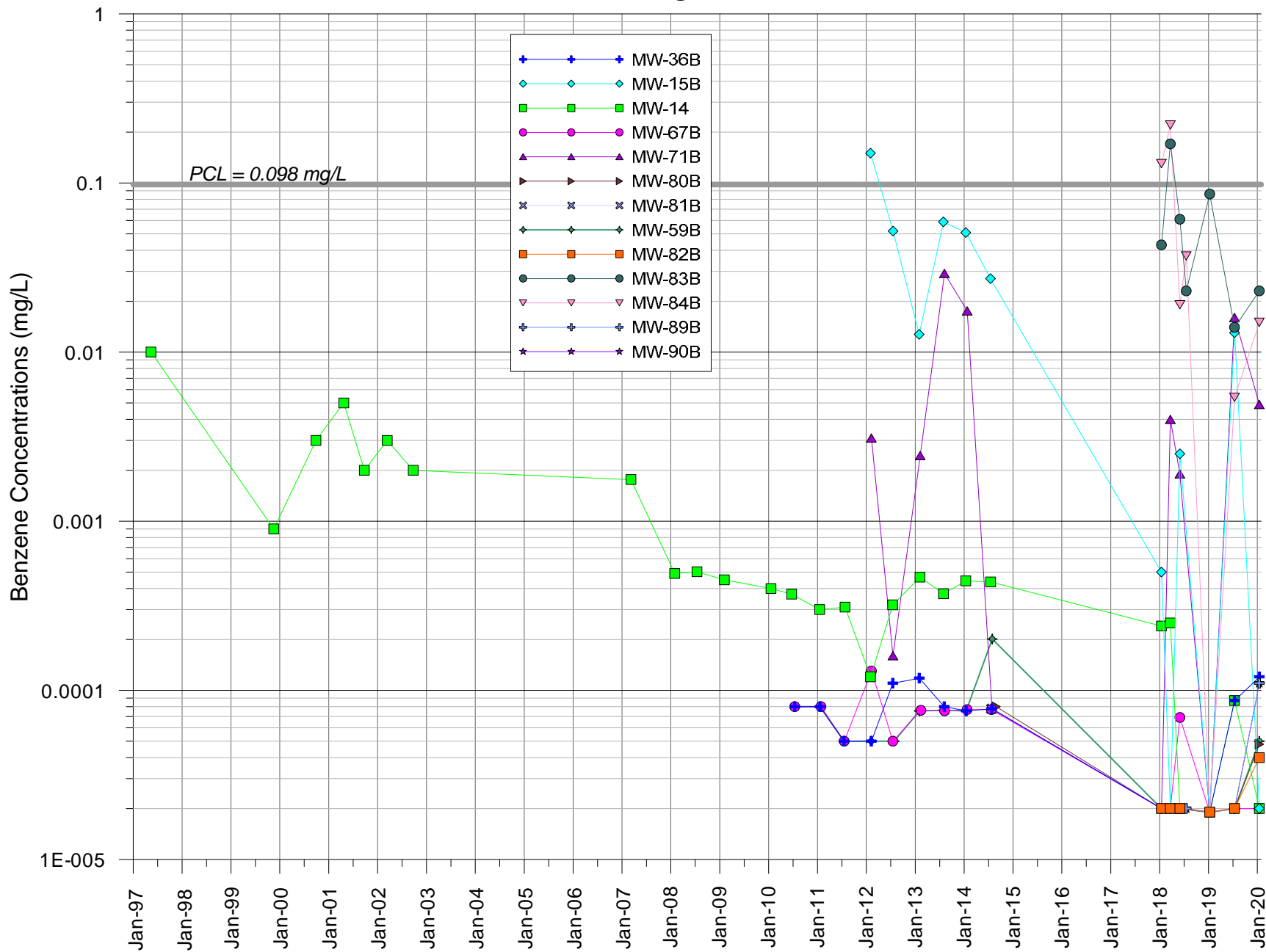
**ATTACHMENT 1B-22**  
**2,4-Dimethylphenol Concentrations at Perimeter Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-23**  
**2-Methylnaphthalene Concentrations at Perimeter Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

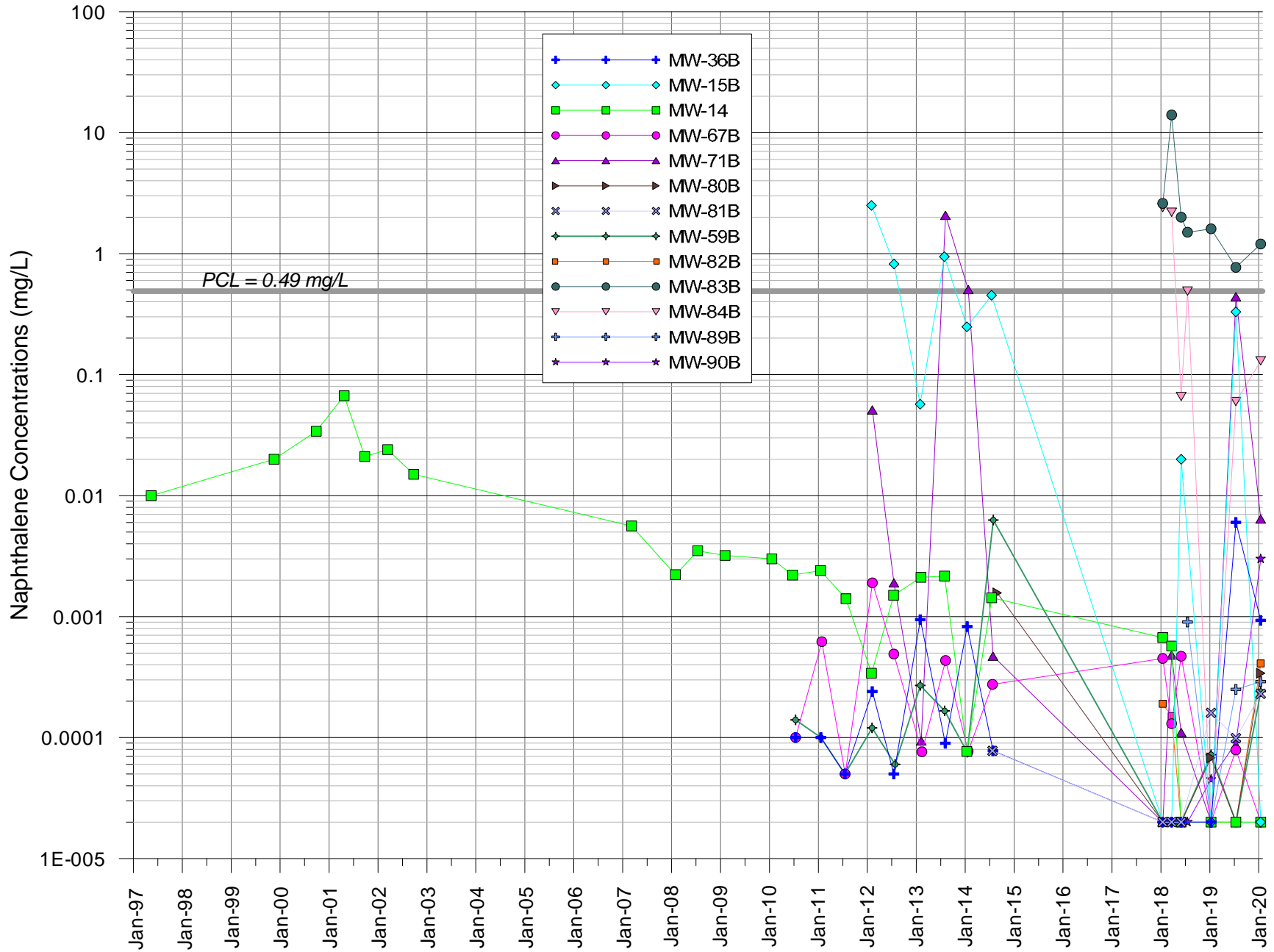


**ATTACHMENT 1B-24**  
**Dibenzofuran Concentrations at Perimeter Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

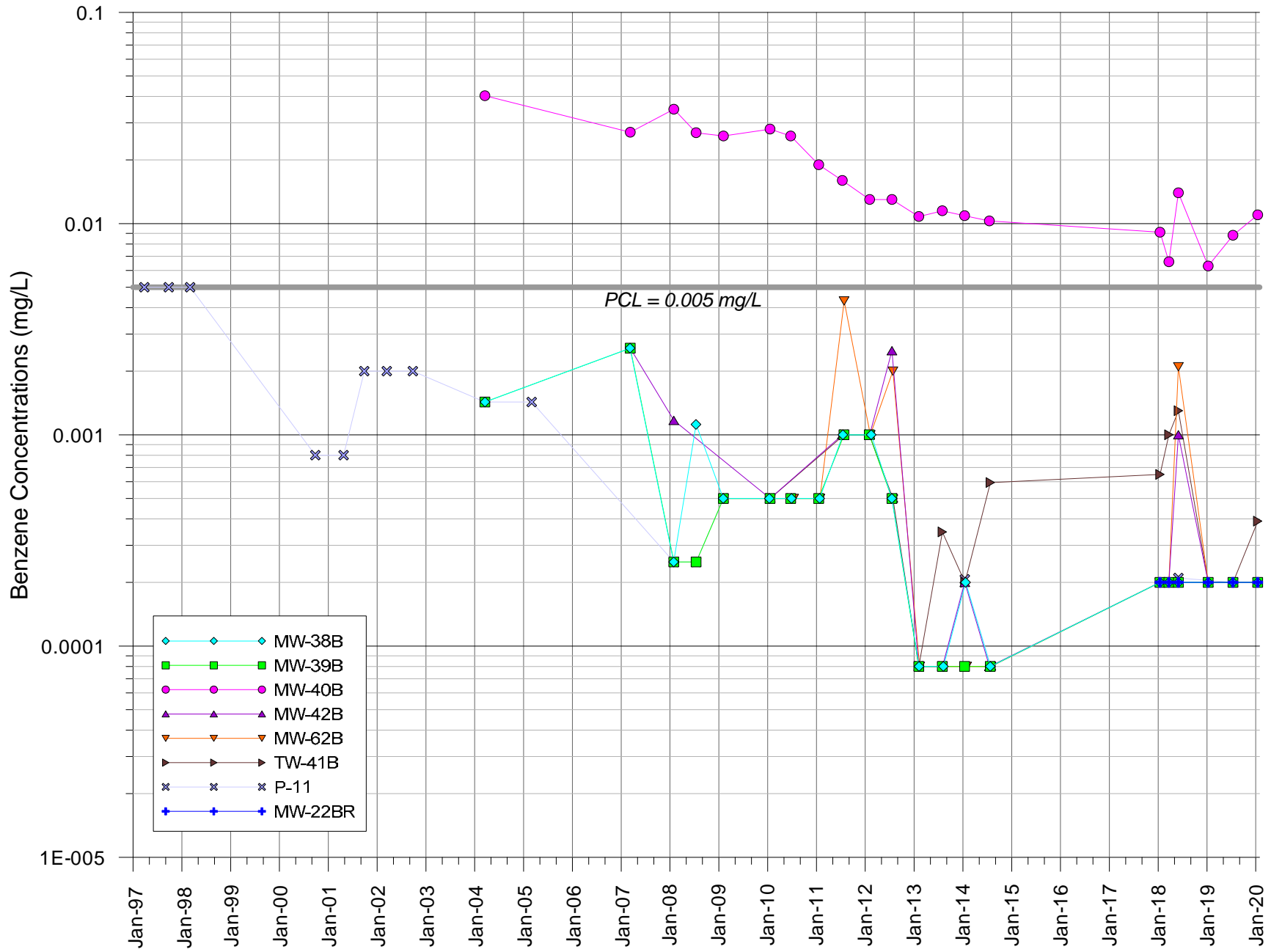




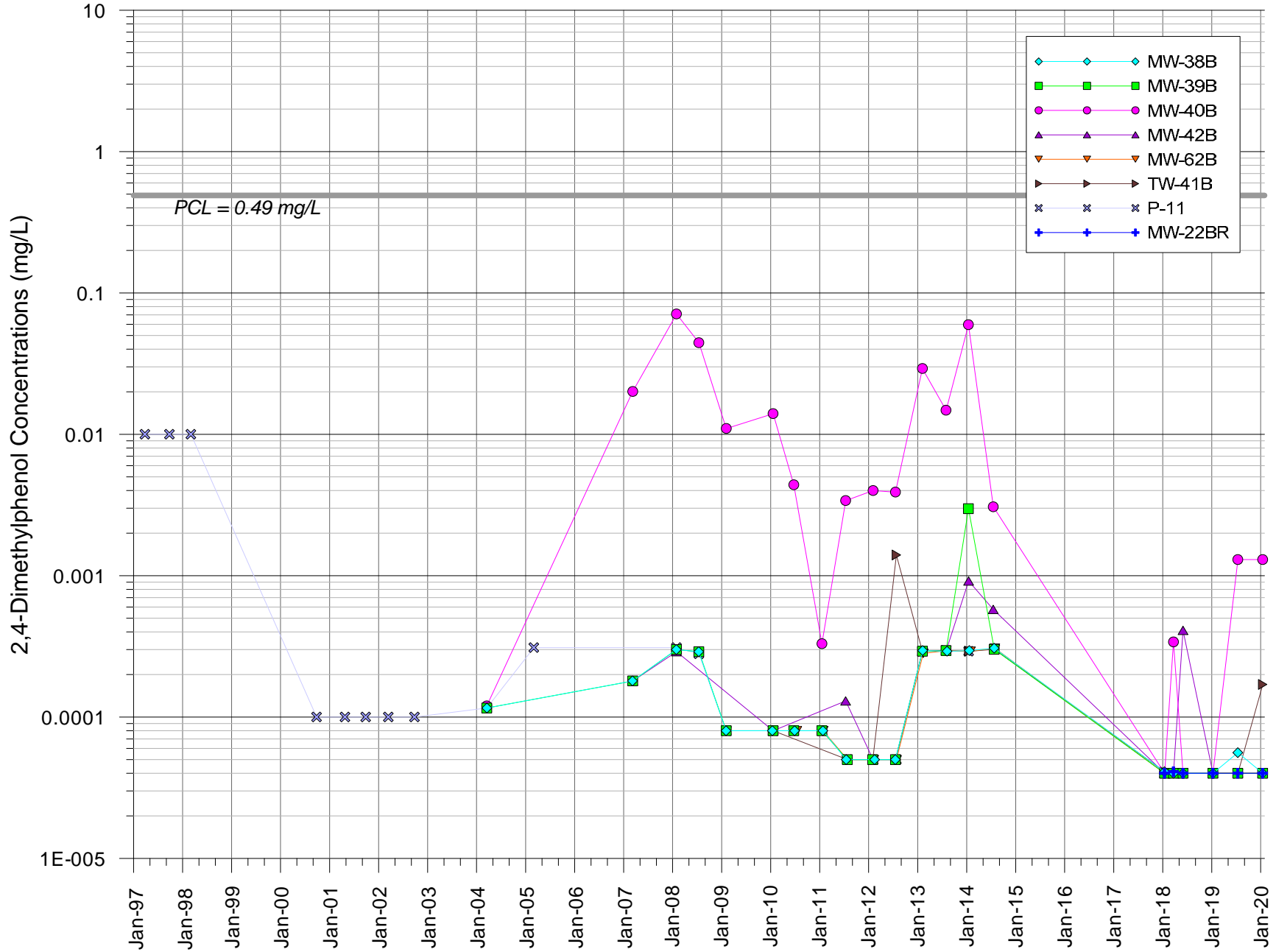
**ATTACHMENT 1B-25**  
**Naphthalene Concentrations at Perimeter Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-26**  
**Benzene Concentrations at West End Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

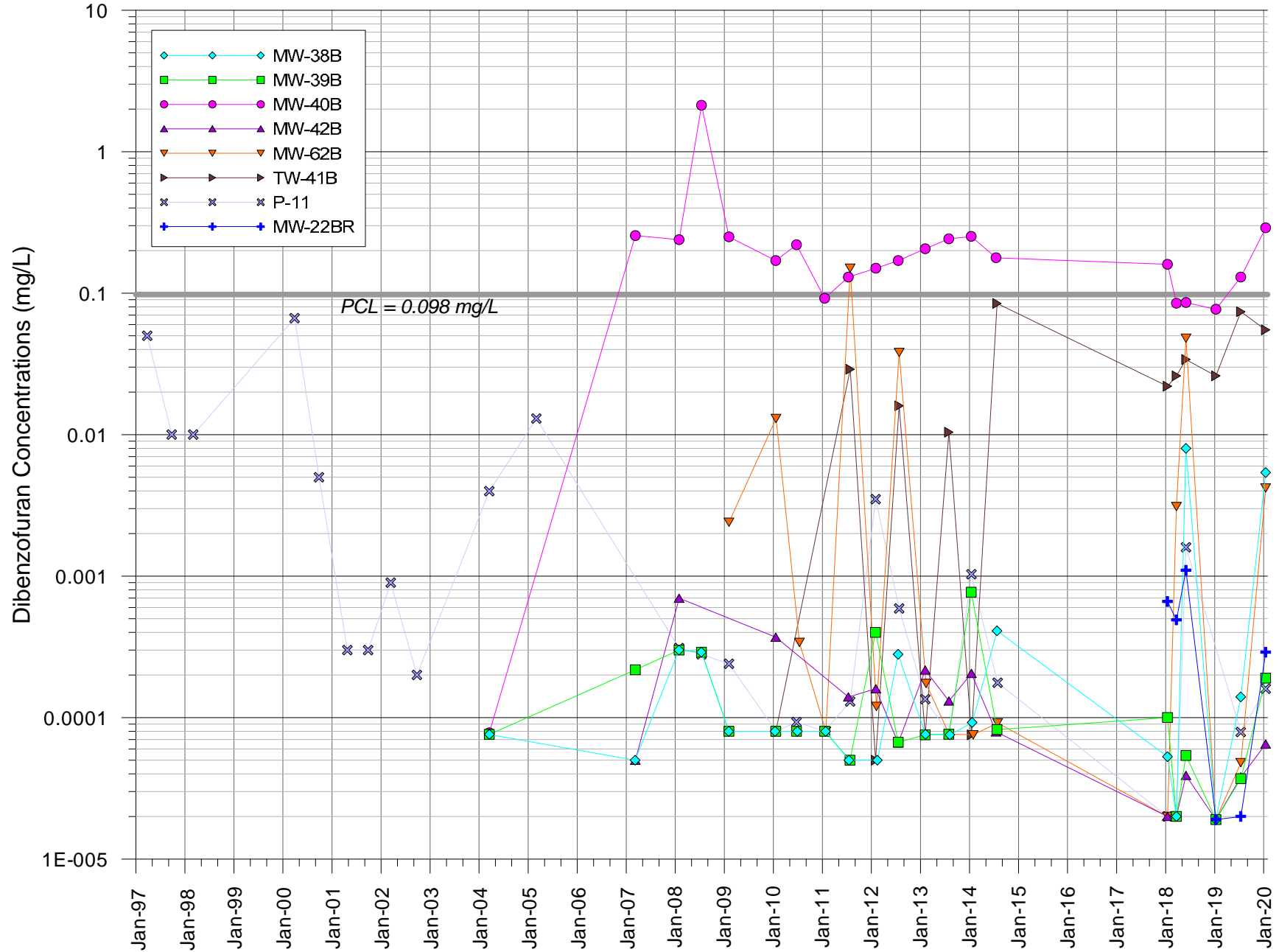


**ATTACHMENT 1B-27**  
**2,4-Dimethylphenol Concentrations at West End Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

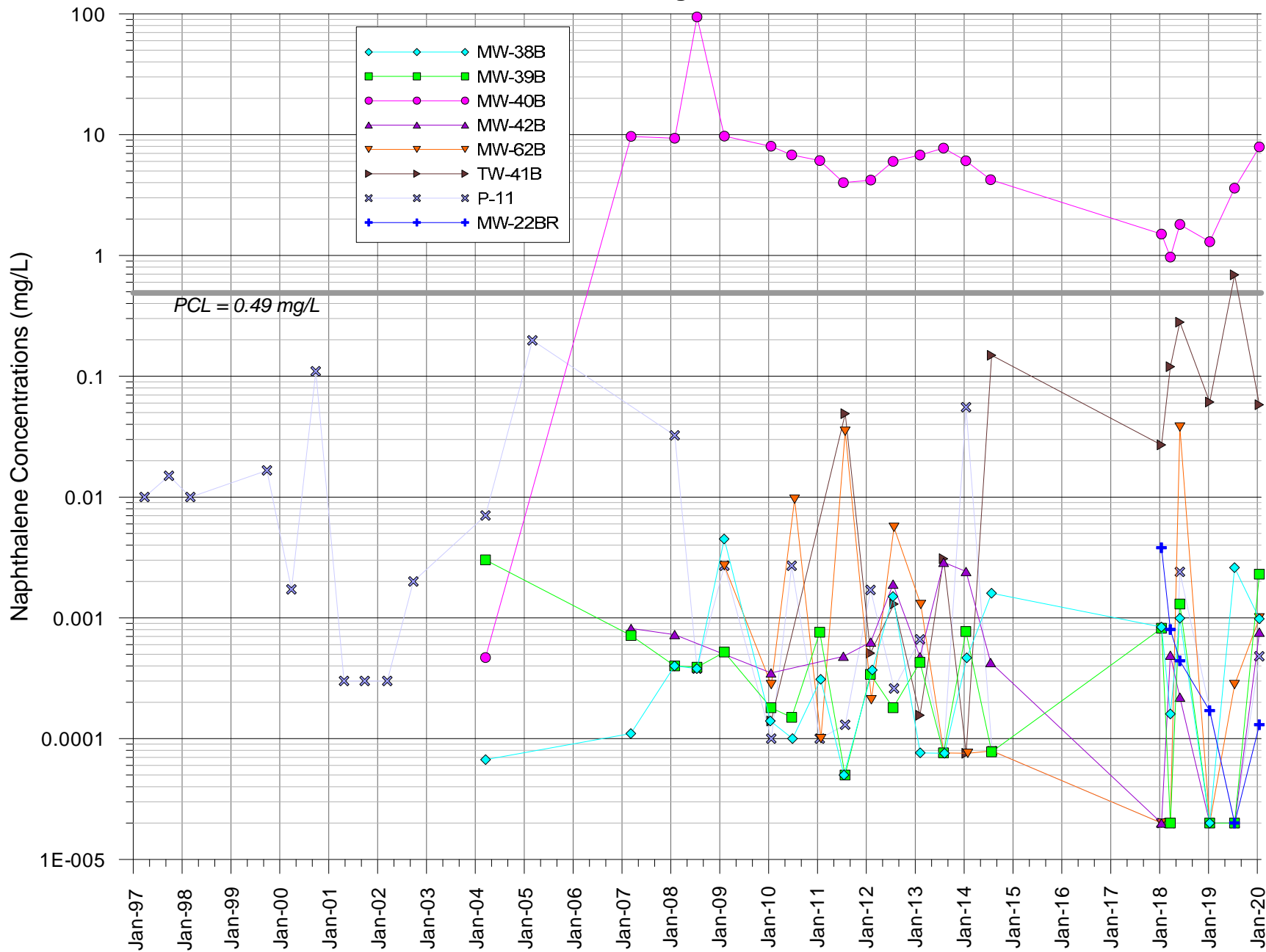




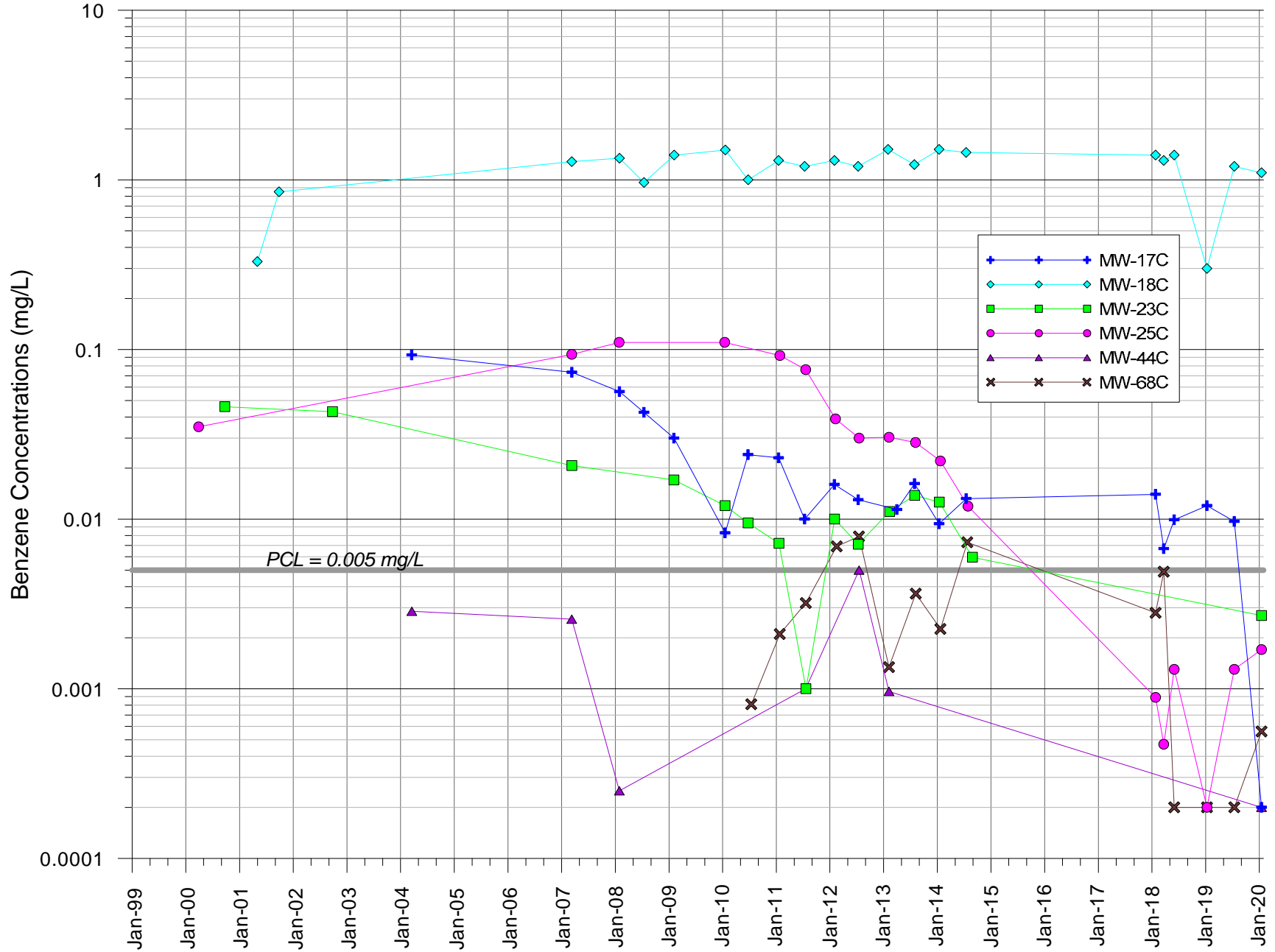
**ATTACHMENT 1B-29**  
**Dibenzofuran Concentrations at West End Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



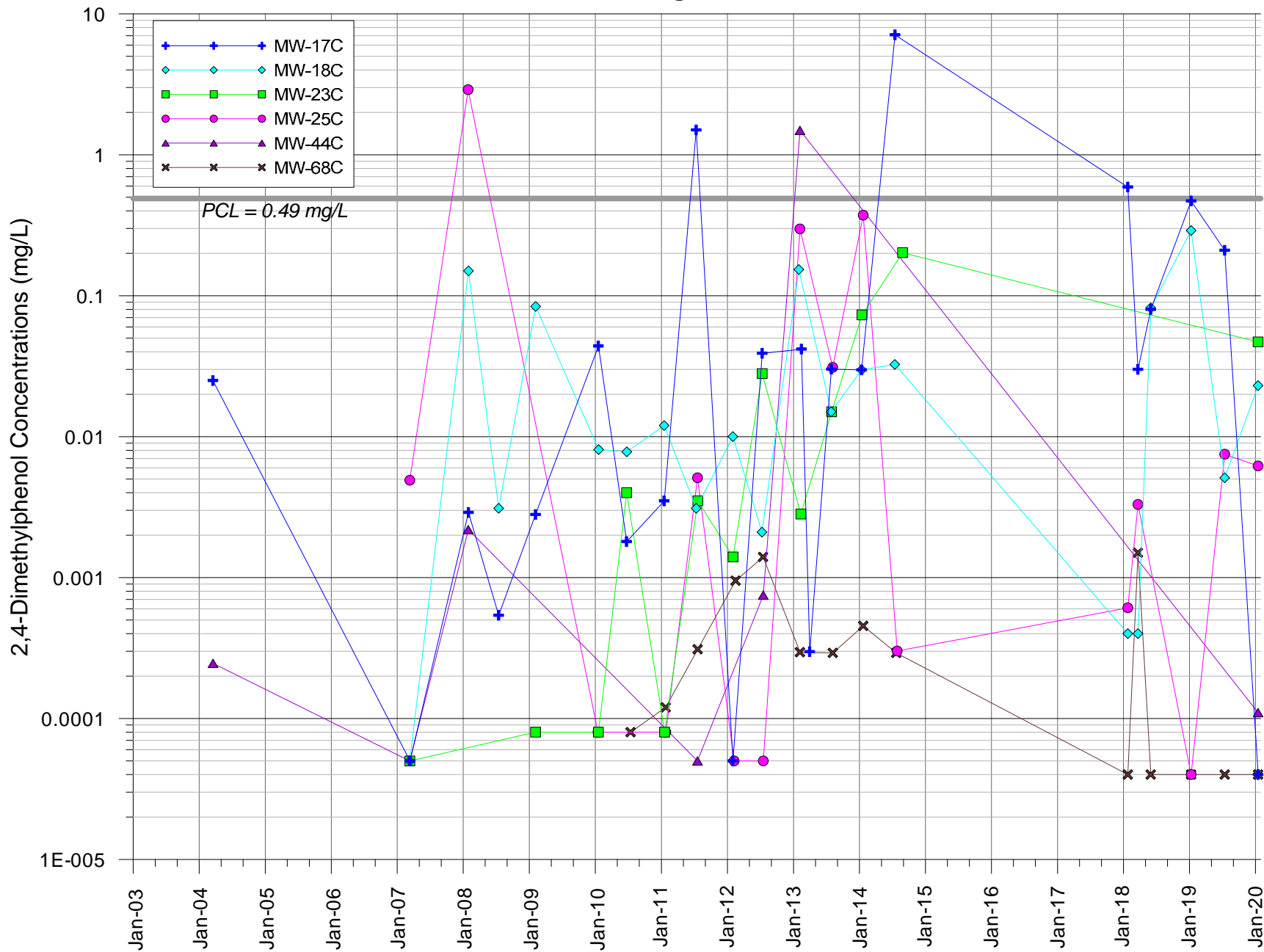
**ATTACHMENT 1B-30**  
**Naphthalene Concentrations at West End Area Wells - B-CZ/B-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-31**  
**Benzene Concentrations at Source Area Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

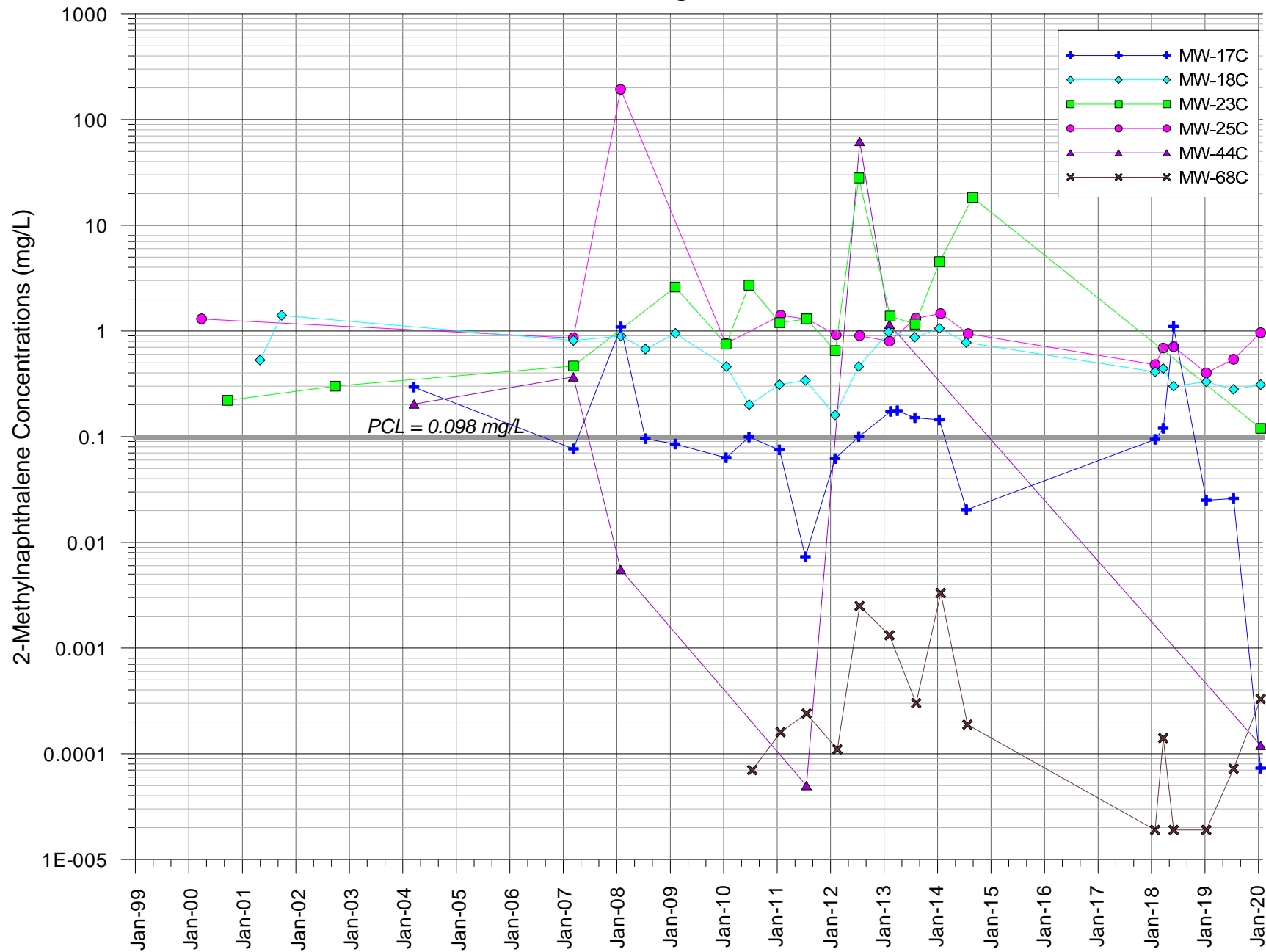


**ATTACHMENT 1B-32**  
**2,4-Dimethylphenol Concentrations at Source Area Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

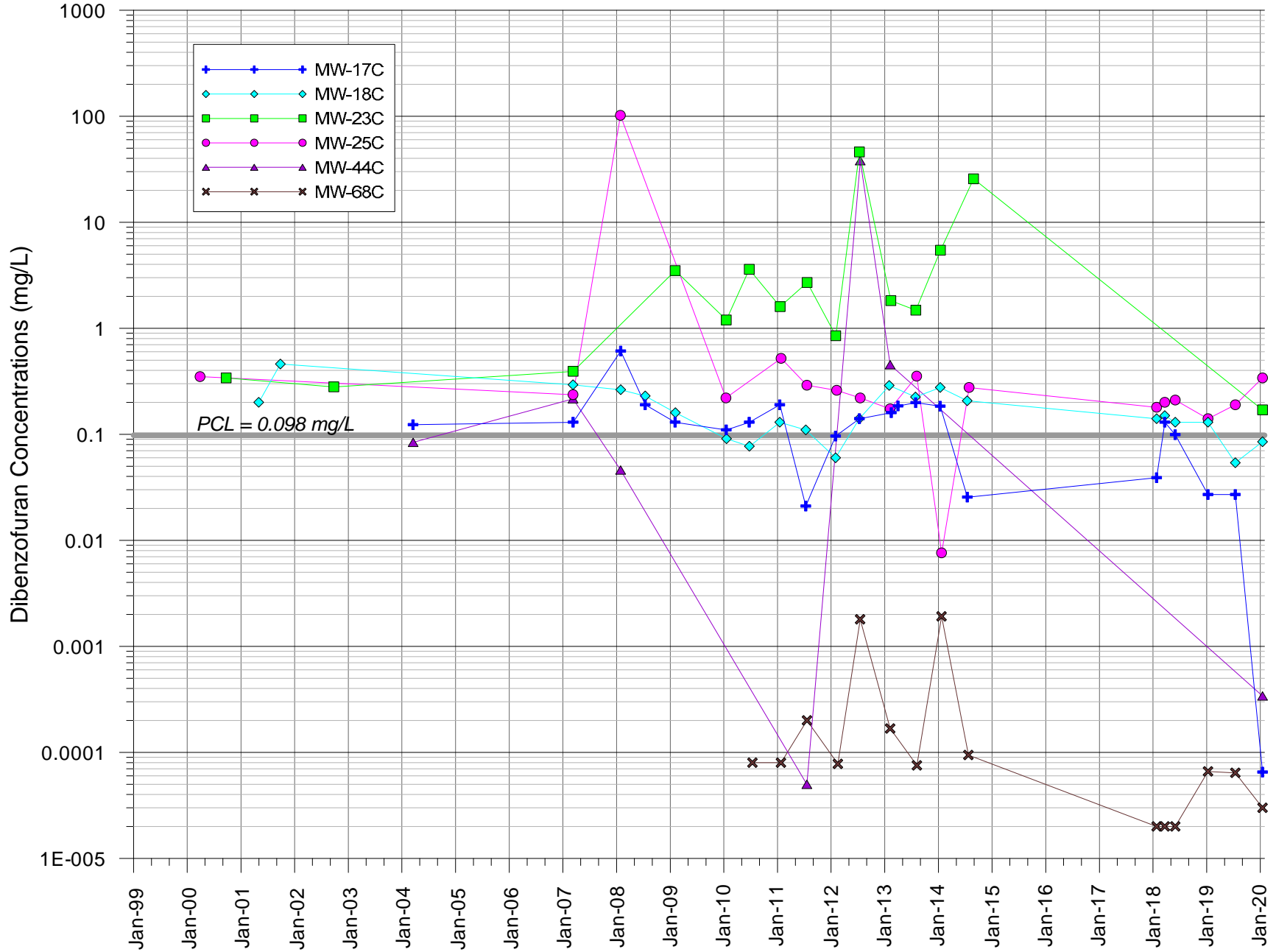




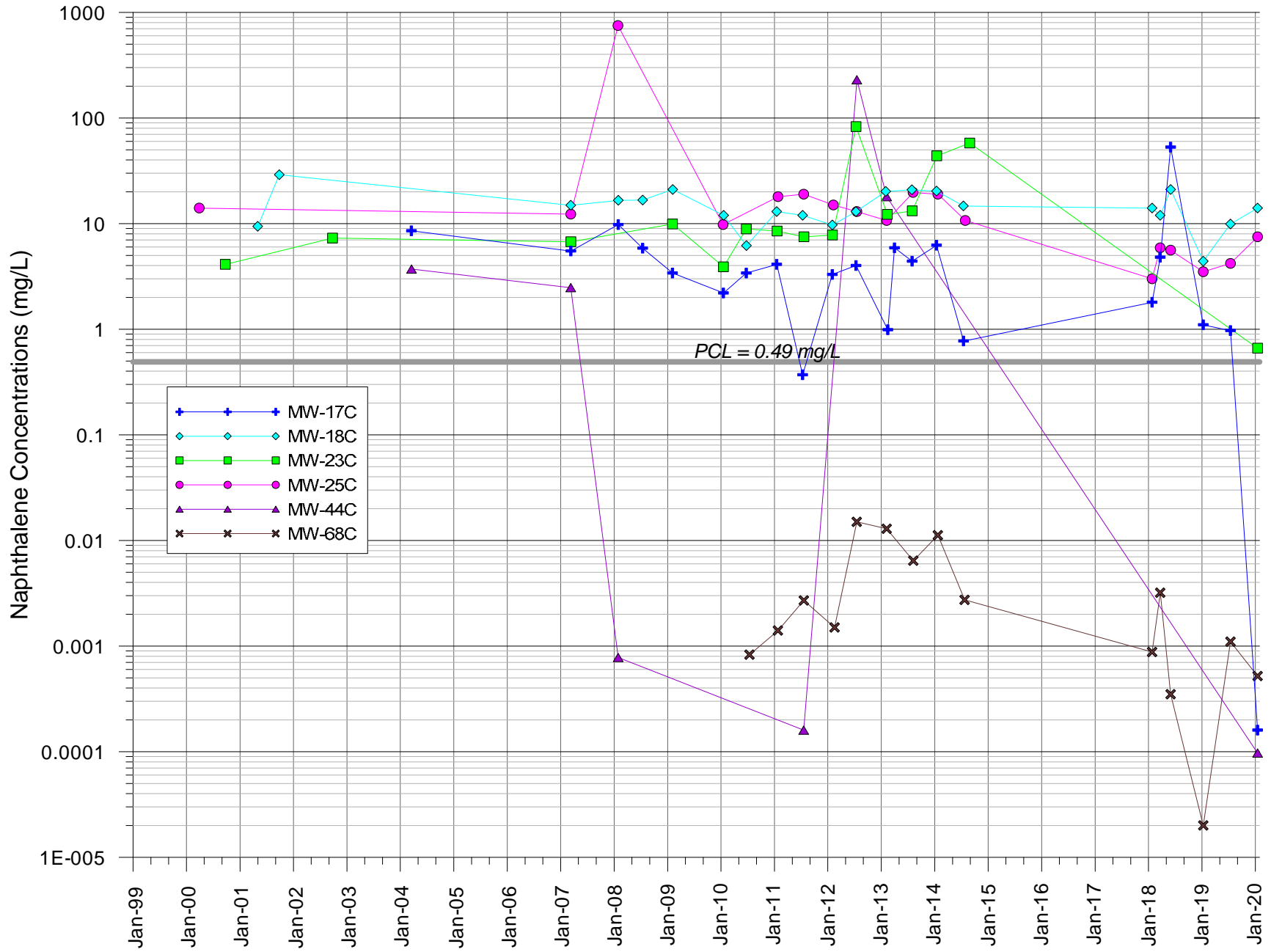
**ATTACHMENT 1B-33**  
**2-Methylnaphthalene Concentrations at Source Area Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



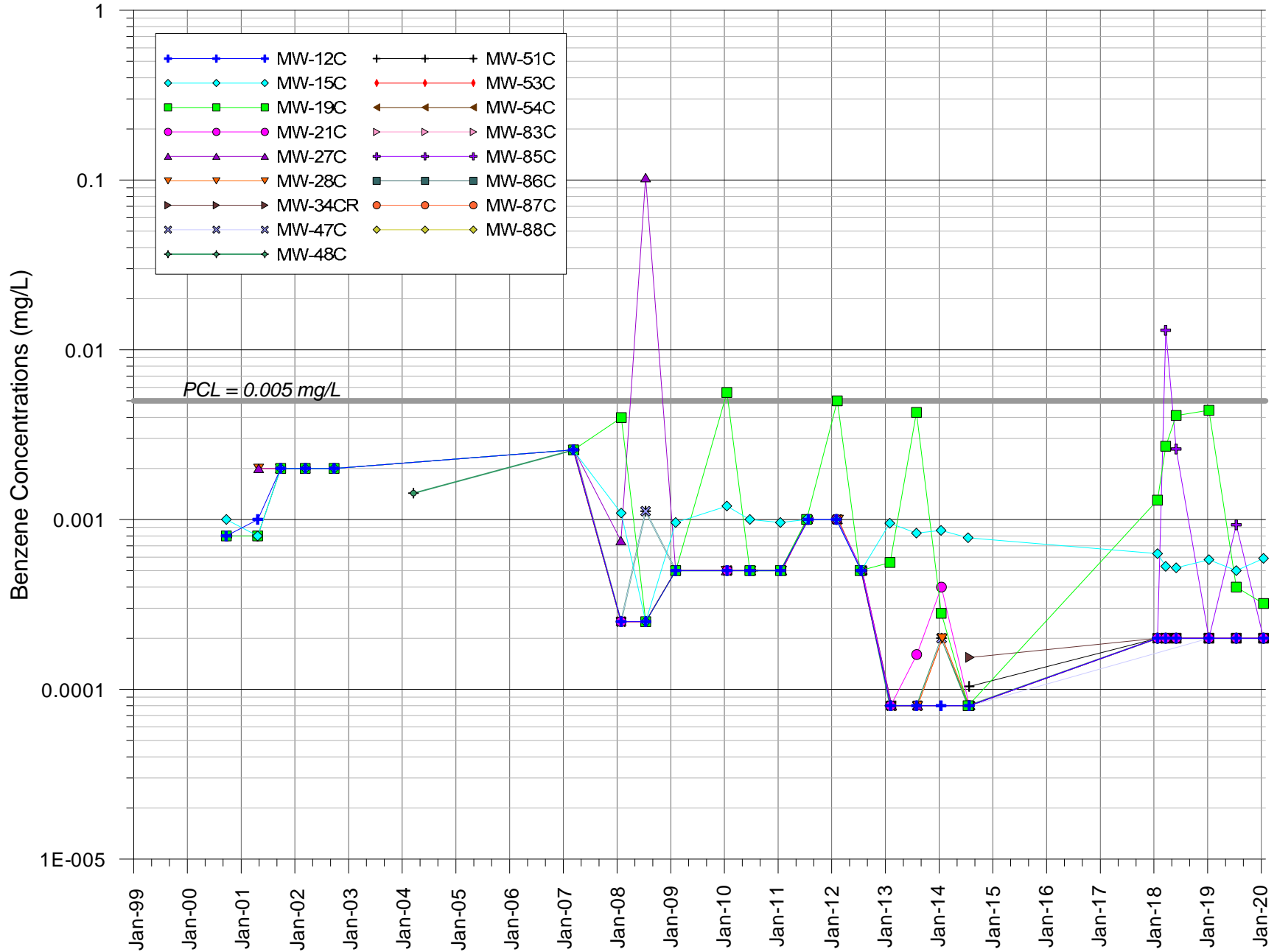
**ATTACHMENT 1B-34**  
**Dibenzofuran Concentrations at Source Area Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



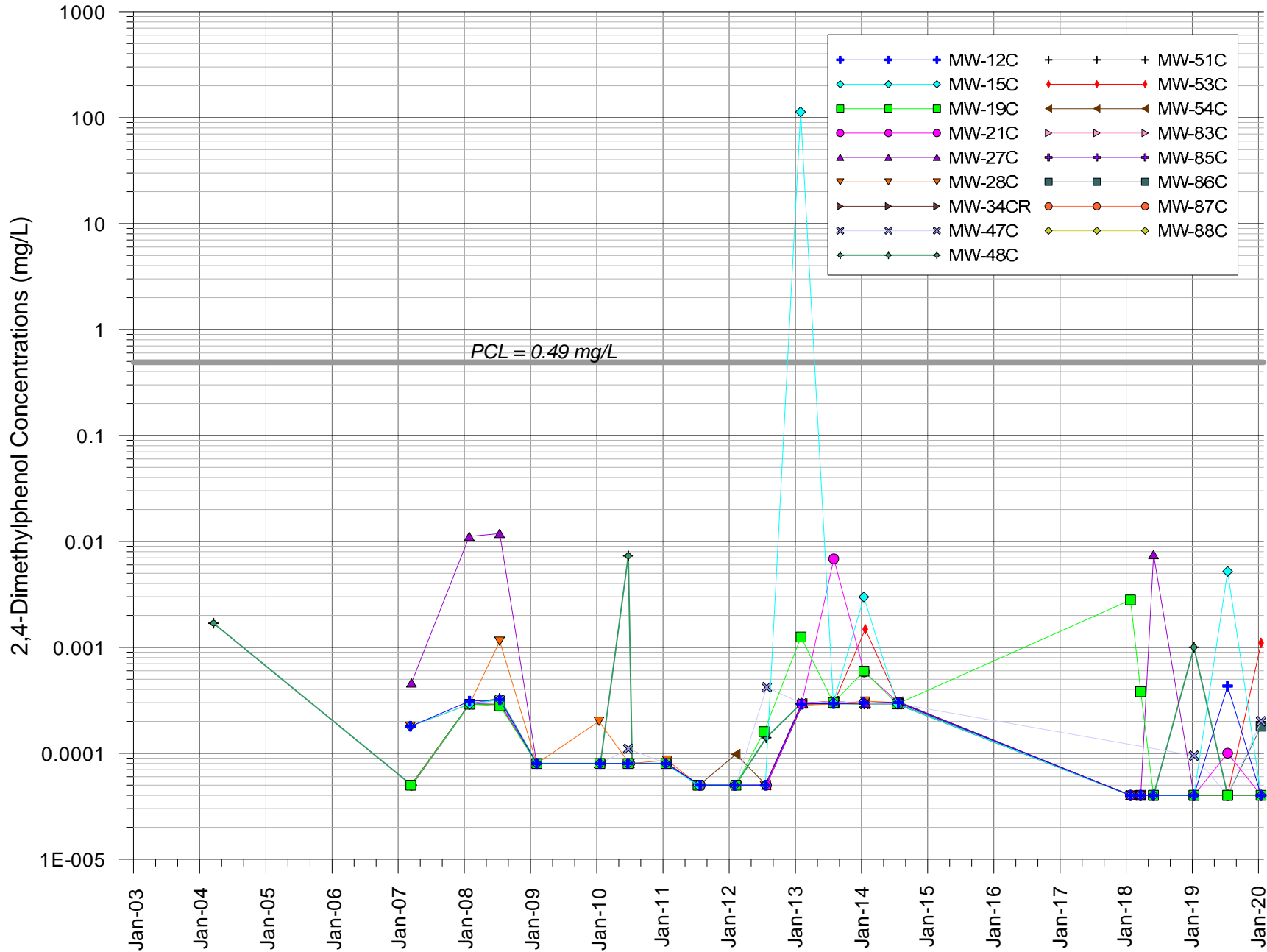
**ATTACHMENT 1B-35**  
**Naphthalene Concentrations at Source Area Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-36**  
**Benzene Concentrations at Perimeter Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



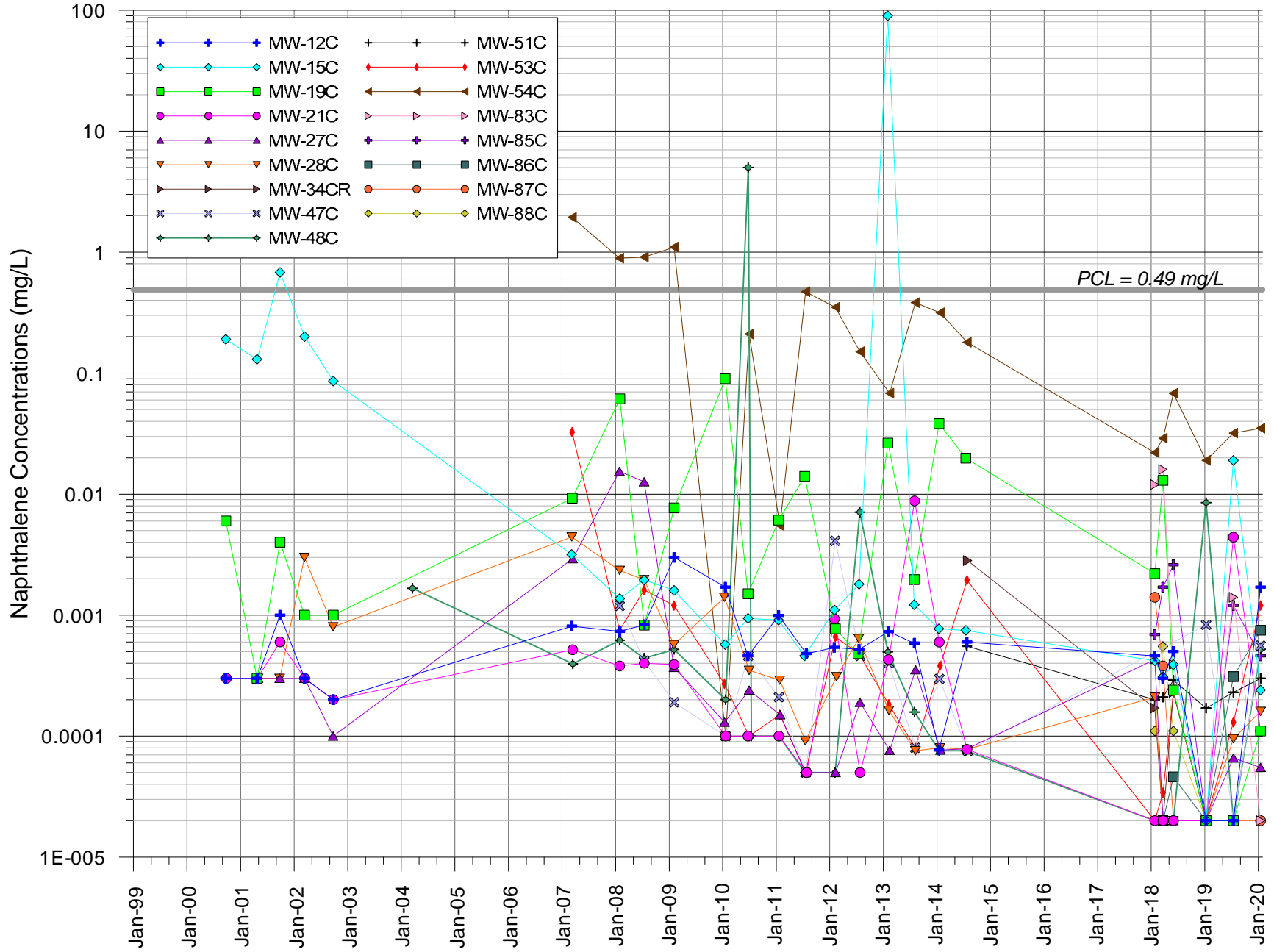
**ATTACHMENT 1B-37**  
**2,4-Dimethylphenol Concentrations at Perimeter Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**





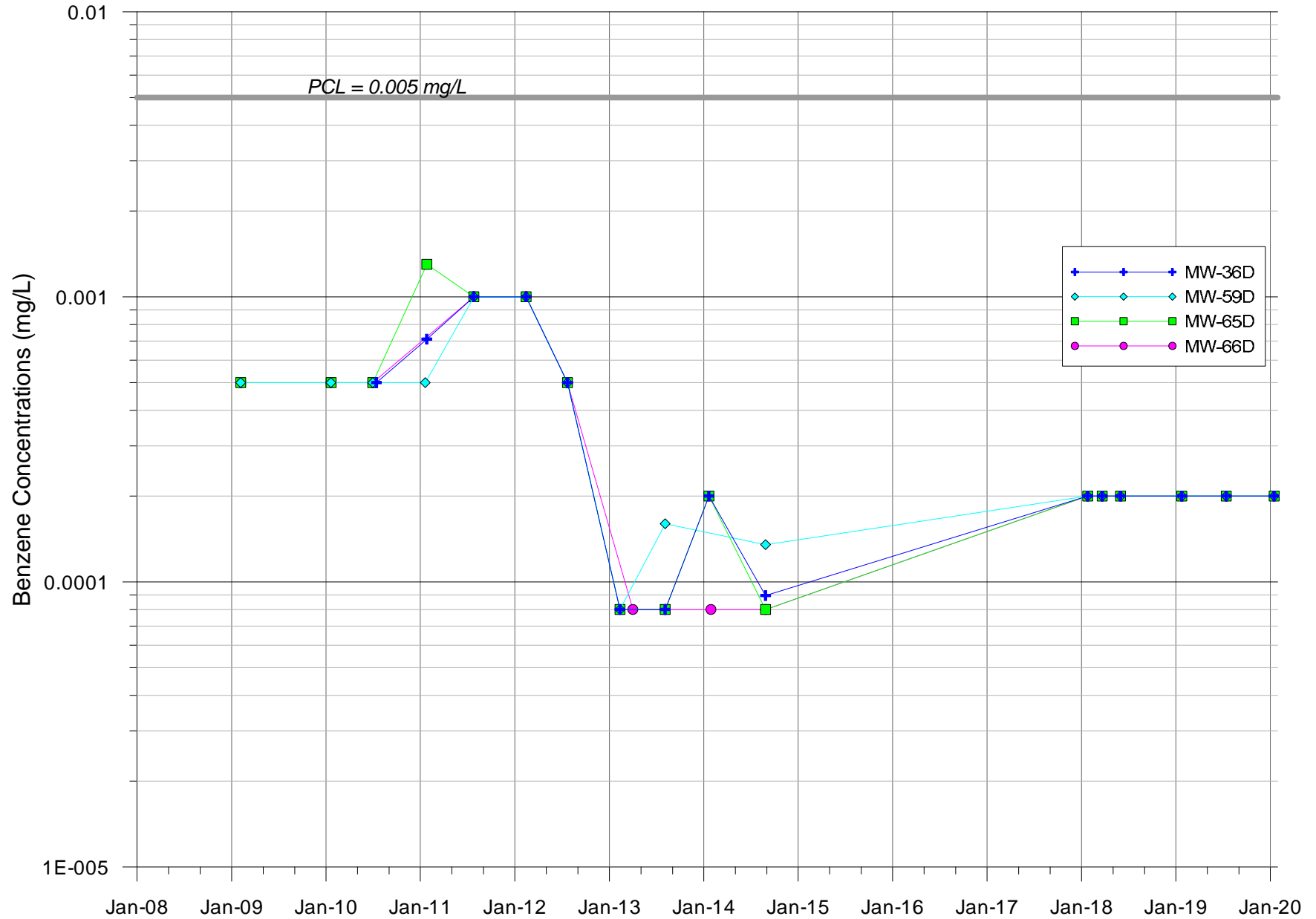


**ATTACHMENT 1B-40**  
**Naphthalene Concentrations at Perimeter Wells - C-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**

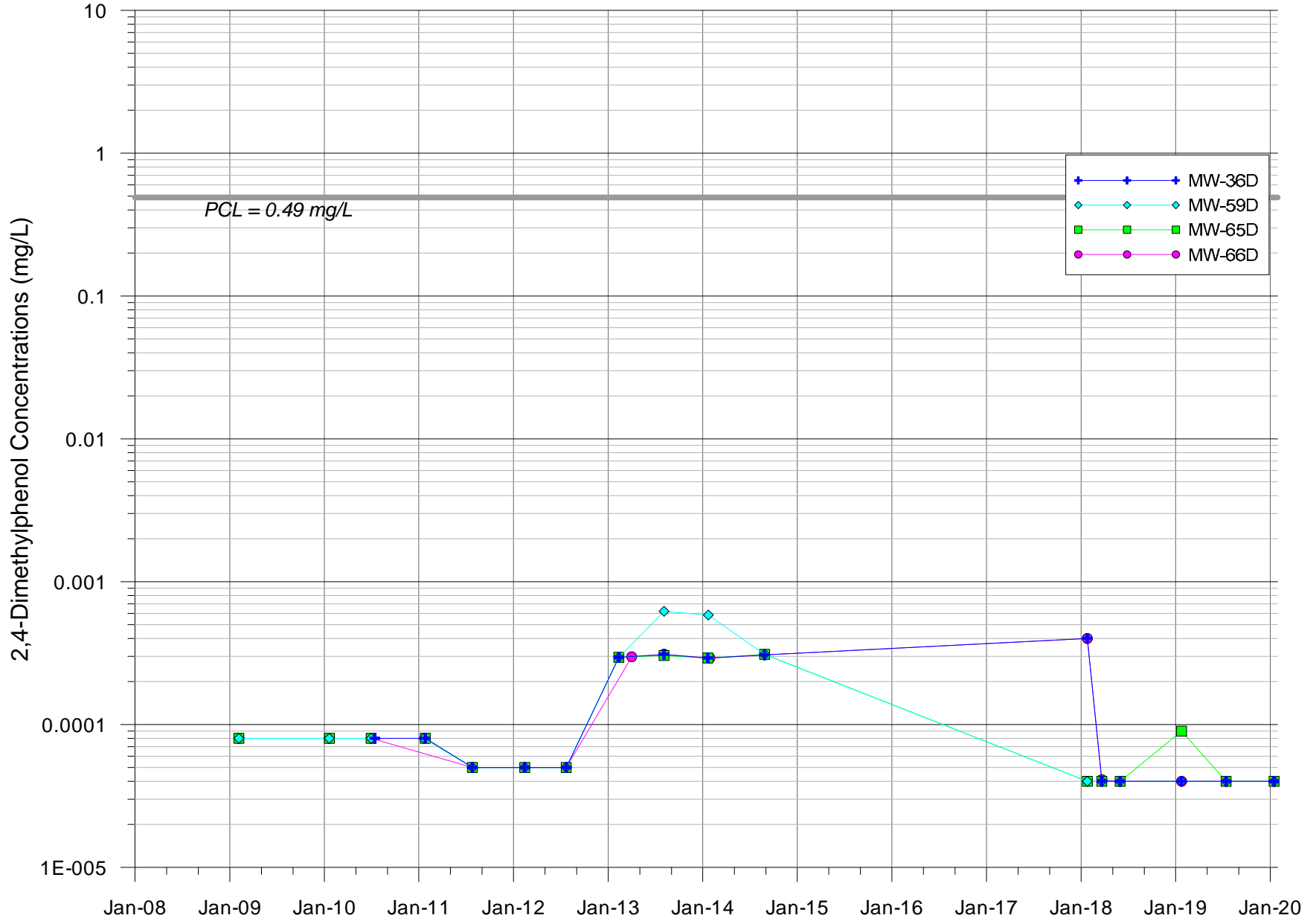




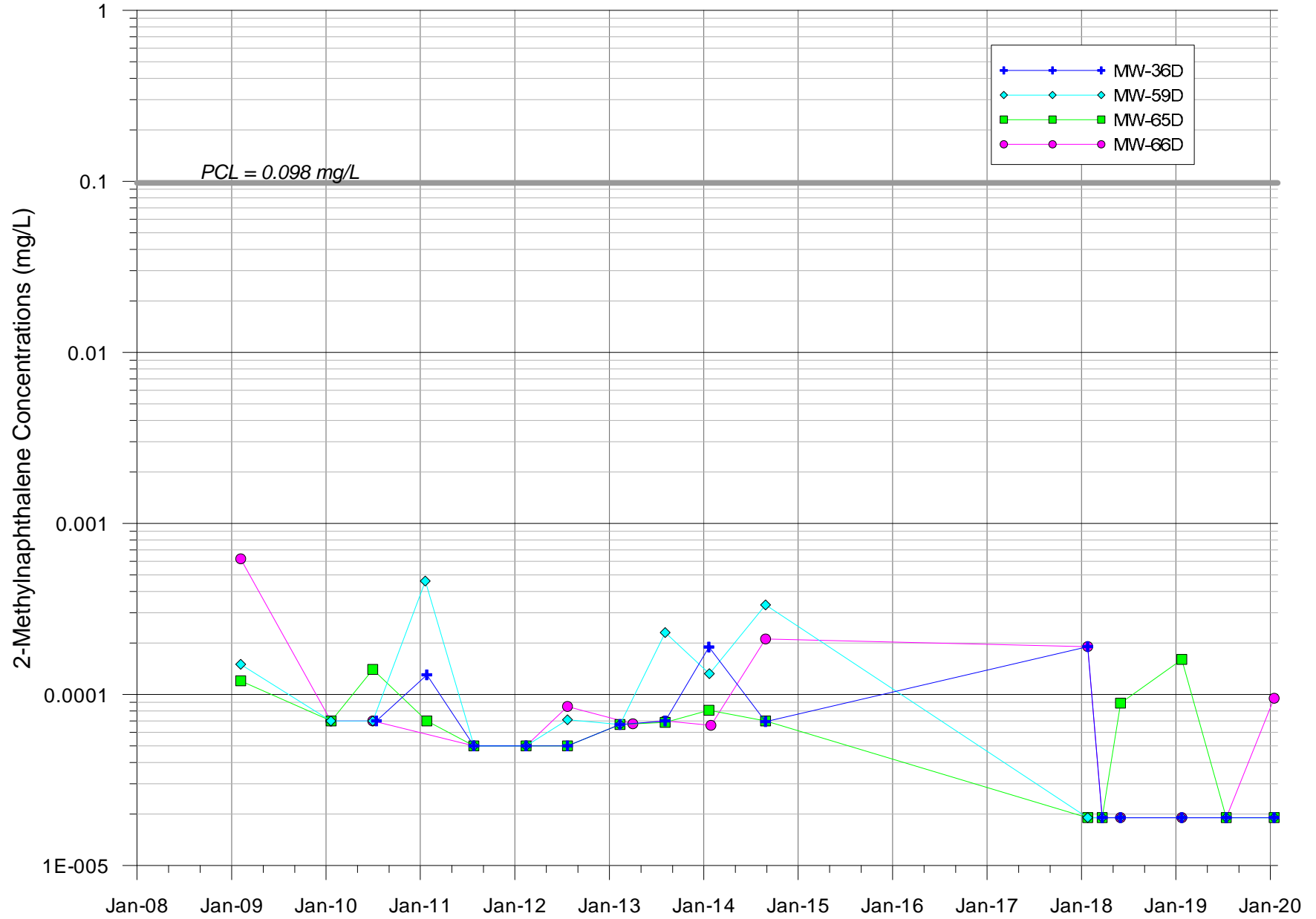
**ATTACHMENT 1B-41**  
**Benzene Concentrations at Source Area Wells - D-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



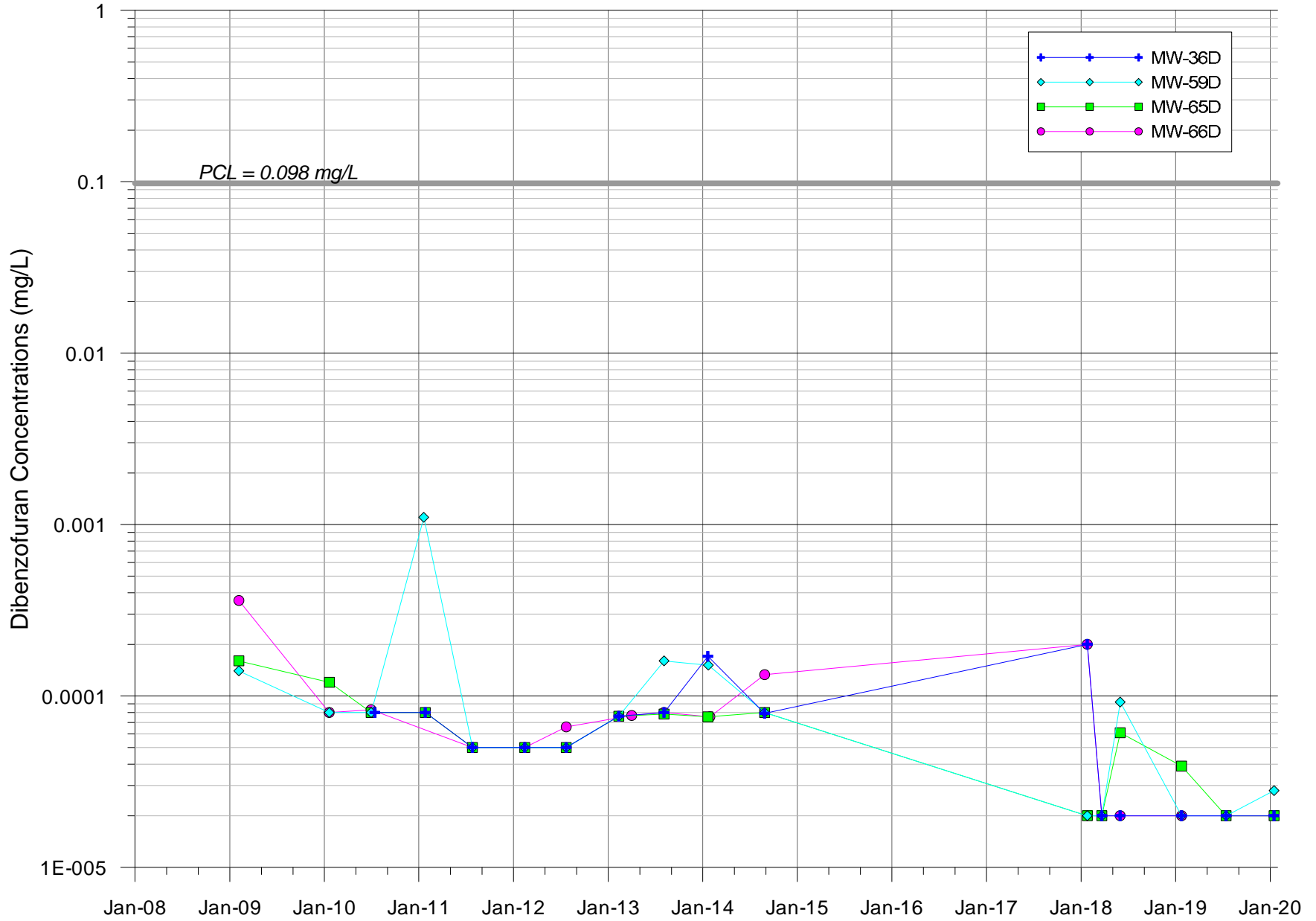
**ATTACHMENT 1B-42**  
**2,4-Dimethylphenol Concentrations at Source Area Wells - D-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



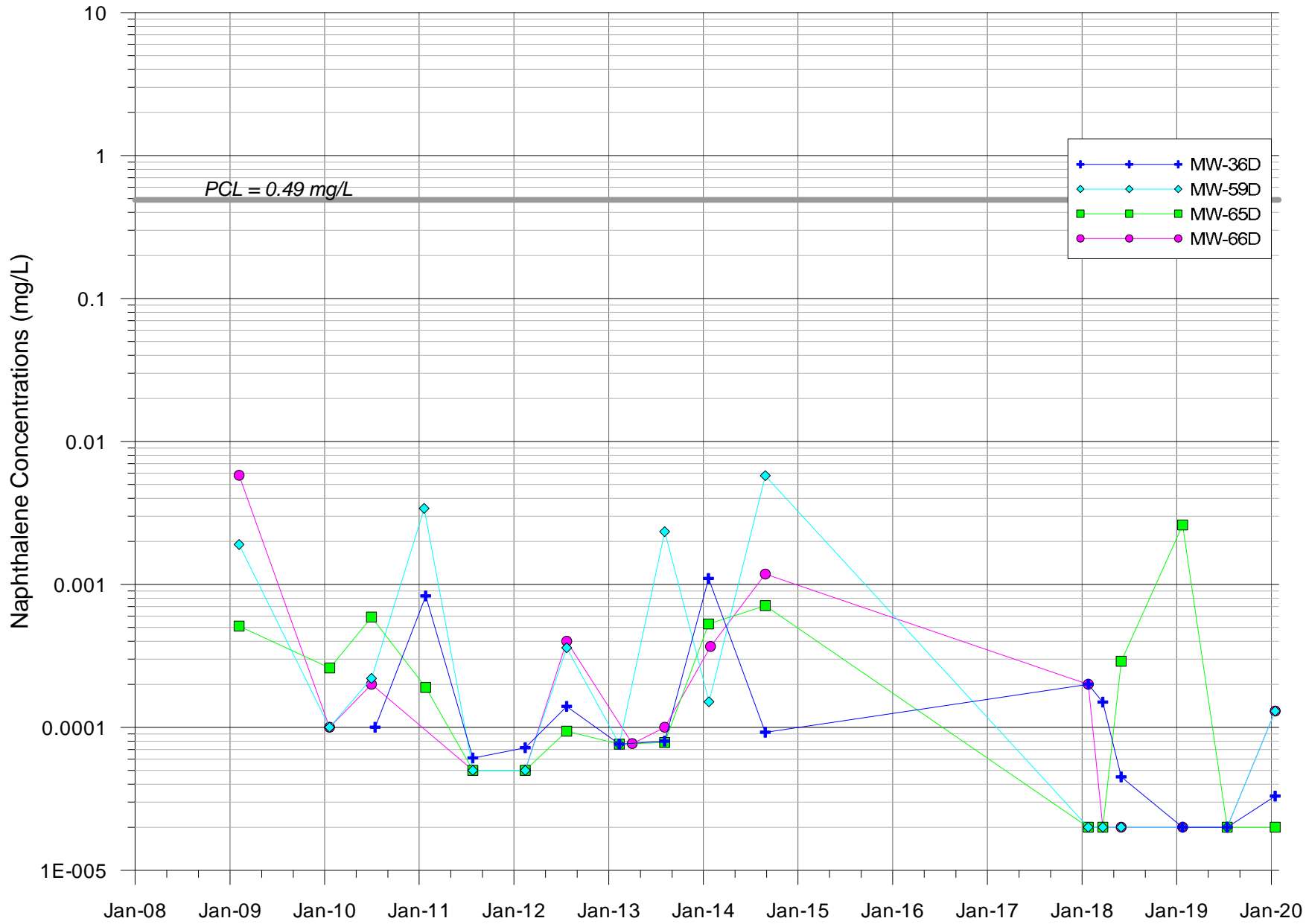
**ATTACHMENT 1B-43**  
**2-Methylnaphthalene Concentrations at Source Area Wells - D-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-44**  
**Dibenzofuran Concentrations at Source Area Wells - D-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT 1B-45**  
**Naphthalene Concentrations at Source Area Wells - D-TZ**  
**UPRR Houston Wood Preserving Works, Houston, Texas**



**ATTACHMENT G**

## Mann-Kendall Trend Test Results

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

Time Period: 1/22/2009 to 1/7/2020

Consolidation Period: No Time Consolidation

Consolidation Type: Median

Duplicate Consolidation: Average

ND Values: Detection Limit

J Flag Values : Actual Value

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
2,4-DIMETHYLPHENOL								
MW-03	T	6	0	0.00	0	42.3%	Yes	ND
MW-04	T	6	0	0.55	-5	76.5%	Yes	ND
MW-05	T	13	0	0.98	-18	84.7%	Yes	ND
MW-09	T	10	0	1.17	-21	96.4%	Yes	ND
MW-12A	T	17	2	3.69	-41	95.0%	No	D
MW-12C	T	17	1	0.97	-19	76.8%	No	S
MW-13	T	17	1	1.17	-38	93.6%	No	PD
MW-14	T	17	1	1.11	-31	89.0%	No	NT
MW-15A	T	17	3	3.13	-30	88.2%	No	NT
MW-15B	T	12	0	2.11	-29	97.4%	Yes	ND
MW-15C	T	17	2	4.12	-31	89.0%	No	NT
MW-17	S	18	17	0.69	9	61.7%	No	NT
MW-17C	S	18	13	2.96	25	81.6%	No	NT
MW-18A	S	17	17	0.86	-46	96.8%	No	D
MW-18C	S	17	11	1.69	9	62.7%	No	NT
MW-19C	S	17	3	1.63	26	84.6%	No	NT
MW-20A	S	10	10	1.11	-35	100.0%	No	D
MW-21C	T	17	2	2.73	-40	94.6%	No	PD
MW-22AR	T	6	0	0.01	-3	64.0%	Yes	ND
MW-22BR	T	6	0	0.01	-3	64.0%	Yes	ND
MW-23C	S	12	4	1.86	43	99.9%	No	I
MW-25A	T	17	1	2.56	-31	89.0%	No	NT
MW-25C	S	15	6	2.42	6	59.6%	No	NT
MW-26A	T	17	2	3.59	-64	99.6%	No	D
MW-27A	T	5	0	0.00	0	40.8%	Yes	ND
MW-27C	T	17	0	0.96	-46	96.8%	Yes	ND

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

2,4-DIMETHYLPHENOL

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-28A	T	17	1	3.01	-29	87.4%	No	NT
MW-28C	T	17	2	0.91	-45	96.5%	No	D
MW-32AR	T	12	5	2.74	-29	97.4%	No	D
MW-32B	S	5	4	1.45	2	59.2%	No	NT
MW-33A	T	17	7	2.83	-29	87.4%	No	NT
MW-33BR	T	10	2	3.12	-23	97.7%	No	D
MW-34CR	T	5	0	1.27	-4	75.8%	Yes	ND
MW-35A	T	17	1	2.12	-27	85.6%	No	NT
MW-35B	T	17	1	2.72	-15	71.5%	No	NT
MW-36A	T	17	0	0.94	-40	94.6%	Yes	ND
MW-36B	T	15	1	0.94	-20	82.3%	No	S
MW-36D	T	15	0	0.95	-27	89.9%	Yes	ND
MW-38A	T	17	0	0.96	-47	97.1%	Yes	ND
MW-38B	T	17	1	0.95	-42	95.4%	No	D
MW-39B	T	17	0	2.62	-46	96.8%	Yes	ND
MW-40B	T	17	10	1.71	-48	97.4%	No	D
MW-42B	T	14	3	1.23	-20	84.8%	No	NT
MW-44A	T	17	1	1.95	-34	91.2%	No	PD
MW-44C	S	4	1	2.00	2	62.5%	No	NT
MW-47C	T	14	6	0.74	6	60.6%	No	NT
MW-48C	T	17	3	2.30	-29	87.4%	No	NT
MW-49A	T	17	14	1.37	-46	96.8%	No	D
MW-49B	S	15	15	1.56	49	99.2%	No	I
MW-50A	T	17	1	1.16	-11	65.7%	No	NT
MW-51A	T	17	0	0.96	-53	98.5%	Yes	ND
MW-51C	T	7	0	1.25	-6	76.4%	Yes	ND
MW-53C	T	17	0	1.68	-23	81.5%	Yes	ND
MW-54C	T	17	1	0.93	-42	95.4%	No	D
MW-57A	S	13	12	1.15	-40	99.3%	No	D
MW-57B	S	11	10	0.80	-19	91.8%	No	PD
MW-58A	T	17	11	2.52	-49	97.7%	No	D
MW-59A	T	17	1	0.95	-52	98.3%	No	D
MW-59B	T	15	2	3.85	-41	97.7%	No	D



# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

2,4-DIMETHYLPHENOL

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-59D	T	17	0	0.97	-49	97.7%	Yes	ND
MW-60A	T	17	1	1.35	-40	94.6%	No	PD
MW-61A	T	17	0	0.96	-50	97.9%	Yes	ND
MW-62B	T	17	0	0.95	-46	96.8%	Yes	ND
MW-63B	T	16	2	2.08	-12	68.7%	No	NT
MW-64A	T	17	0	0.97	-50	97.9%	Yes	ND
MW-65D	T	17	0	0.93	-33	90.5%	Yes	ND
MW-66D	T	17	0	0.90	-32	89.8%	Yes	ND
MW-67B	T	15	1	1.02	-41	97.7%	No	D
MW-68A	T	3	0	0.00	0	0.0%	Yes	ND
MW-68B	S	12	11	0.88	-14	81.0%	No	S
MW-68C	T	15	7	1.25	-32	93.7%	No	PD
MW-69A	T	14	2	2.49	-40	98.5%	No	D
MW-70B	S	5	4	0.89	2	59.2%	No	NT
MW-71B	T	12	3	2.10	-32	98.4%	No	D
MW-72B	S	10	10	1.40	-19	94.6%	No	PD
MW-74B	S	11	11	1.72	-21	94.0%	No	PD
MW-75B	S	6	5	1.83	1	50.0%	No	NT
MW-76C	T	8	2	1.63	1	50.0%	No	NT
MW-77A	S	7	5	1.01	3	61.4%	No	NT
MW-79A	S	7	7	0.79	-4	66.7%	No	S
MW-80B	T	7	1	1.11	-1	50.0%	No	NT
MW-81B	T	7	0	1.28	-6	76.4%	Yes	ND
MW-82B	T	6	0	0.00	0	42.3%	Yes	ND
MW-83B	T	7	1	1.17	-6	76.4%	No	NT
MW-83C	T	6	0	0.00	0	42.3%	Yes	ND
MW-84B	T	7	4	0.87	0	43.7%	No	S
MW-85C	T	6	0	0.01	1	50.0%	Yes	ND
MW-86C	T	6	1	0.55	5	76.5%	No	NT
MW-87C	T	6	0	0.00	0	42.3%	Yes	ND
MW-88C	T	6	0	0.00	0	42.3%	Yes	ND
MW-89B	T	4	0	0.00	0	37.5%	Yes	ND
MW-90B	T	4	1	0.36	3	72.9%	No	NT

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## 2,4-DIMETHYLPHENOL

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
P-11	T	17	1	0.96	-47	97.1%	No	D
TW-41B	T	14	2	1.60	-29	93.7%	No	PD

## 2-METHYLNAPHTHALENE

MW-01A	T	23	19	1.69	-92	99.3%	No	D
MW-02	T	23	12	1.39	-105	99.7%	No	D
MW-03	T	6	1	0.43	9	93.2%	No	PI
MW-04	T	6	2	1.23	4	70.3%	No	NT
MW-05	T	13	4	1.38	-23	90.8%	No	PD
MW-07	T	23	3	1.28	-159	100.0%	No	D
MW-08	T	23	0	1.29	-178	100.0%	Yes	ND
MW-09	T	10	1	0.81	-21	96.4%	No	D
MW-10A	T	23	2	1.43	-147	100.0%	No	D
MW-11A	T	23	0	1.29	-182	100.0%	Yes	ND
MW-11B	T	1	1	0.00	0	0.0%	No	N/A
MW-12A	T	17	16	1.08	-69	99.8%	No	D
MW-12C	T	17	14	0.66	1	50.0%	No	NT
MW-13	T	17	7	1.52	-41	95.0%	No	D
MW-14	T	17	13	0.71	-94	100.0%	No	D
MW-15A	T	17	17	1.09	-30	88.2%	No	NT
MW-15B	T	12	7	2.89	-30	97.8%	No	D
MW-15C	T	17	4	4.10	-41	95.0%	No	D
MW-17	S	18	17	0.59	-37	91.2%	No	PD
MW-17C	S	18	18	1.78	-1	50.0%	No	NT
MW-18A	S	17	17	0.49	-18	75.5%	No	S
MW-18C	S	17	17	0.58	-24	82.6%	No	S
MW-19C	S	17	11	1.28	-47	97.1%	No	D
MW-20A	S	10	10	0.72	-18	93.4%	No	PD
MW-21C	T	17	5	2.99	-29	87.4%	No	NT
MW-22AR	T	6	0	1.15	5	76.5%	Yes	ND
MW-22BR	T	6	3	1.63	-8	89.8%	No	NT
MW-23C	S	12	12	1.67	2	52.7%	No	NT
MW-25A	T	17	5	3.97	-38	93.6%	No	PD
MW-25C	S	15	14	0.47	-31	93.0%	No	PD

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## 2-METHYLNAPHTHALENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-26A	T	17	7	2.12	-54	98.6%	No	D
MW-27A	T	5	1	1.73	2	59.2%	No	NT
MW-27C	T	17	1	1.72	-59	99.2%	No	D
MW-28A	T	17	4	1.48	-25	83.6%	No	NT
MW-28C	T	17	6	0.85	-80	100.0%	No	D
MW-32AR	T	12	7	3.16	-23	93.3%	No	PD
MW-32B	S	5	4	2.20	4	75.8%	No	NT
MW-33A	T	17	11	2.27	-22	80.4%	No	NT
MW-33BR	T	10	7	2.02	-20	95.5%	No	D
MW-34CR	T	5	1	1.31	-1	50.0%	No	NT
MW-35A	T	17	10	2.58	14	70.1%	No	NT
MW-35B	T	17	17	0.55	-62	99.5%	No	D
MW-36A	T	17	3	1.05	-79	100.0%	No	D
MW-36B	T	15	1	1.49	-15	75.2%	No	NT
MW-36D	T	15	2	0.84	-36	95.9%	No	D
MW-38A	T	17	7	1.00	-37	93.0%	No	PD
MW-38B	T	17	5	0.93	-16	72.9%	No	S
MW-39B	T	17	4	1.46	-5	56.4%	No	NT
MW-40B	T	17	17	0.53	-60	99.3%	No	D
MW-42B	T	14	6	0.97	-31	95.0%	No	D
MW-44A	T	17	9	2.92	-32	89.8%	No	NT
MW-44C	S	4	3	1.95	0	37.5%	No	NT
MW-47C	T	14	3	3.11	-24	89.4%	No	NT
MW-48C	T	17	3	4.01	-41	95.0%	No	D
MW-49A	T	17	13	1.06	-64	99.6%	No	D
MW-49B	S	15	13	3.84	38	96.7%	No	I
MW-50A	T	17	4	3.04	-42	95.4%	No	D
MW-51A	T	17	3	0.45	-60	99.3%	No	D
MW-51C	T	7	0	0.66	-1	50.0%	Yes	ND
MW-53C	T	17	4	1.01	-13	68.7%	No	NT
MW-54C	T	17	16	2.23	-28	86.5%	No	NT
MW-57A	S	13	13	1.20	-17	83.2%	No	NT
MW-57B	S	11	11	0.57	-13	82.1%	No	S

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## 2-METHYLNAPHTHALENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-58A	T	17	14	1.14	-20	78.0%	No	NT
MW-59A	T	17	4	2.43	-29	87.4%	No	NT
MW-59B	T	15	1	0.47	-48	99.1%	No	D
MW-59D	T	17	6	1.19	-66	99.7%	No	D
MW-60A	T	17	6	2.13	-38	93.6%	No	PD
MW-61A	T	17	1	1.34	-85	100.0%	No	D
MW-62B	T	17	6	1.98	-31	89.0%	No	NT
MW-63B	T	16	14	1.42	-40	96.1%	No	D
MW-64A	T	17	3	1.13	-58	99.1%	No	D
MW-65D	T	17	5	0.61	-34	91.2%	No	PD
MW-66D	T	17	3	1.36	-38	93.6%	No	PD
MW-67B	T	15	3	0.84	-59	99.9%	No	D
MW-68A	T	3	3	0.00	0	0.0%	No	N/A
MW-68B	S	12	12	0.37	-10	72.7%	No	S
MW-68C	T	15	10	1.70	-12	70.4%	No	NT
MW-69A	T	14	3	2.70	-54	99.9%	No	D
MW-70B	S	5	5	1.32	8	95.8%	No	I
MW-71B	T	12	8	2.00	-11	74.9%	No	NT
MW-72B	S	10	10	1.15	-23	97.7%	No	D
MW-74B	S	11	11	1.25	7	67.6%	No	NT
MW-75B	S	6	6	1.13	1	50.0%	No	NT
MW-76C	T	8	6	0.89	0	45.2%	No	S
MW-77A	S	7	6	0.68	4	66.7%	No	NT
MW-79A	S	7	7	0.68	-7	80.9%	No	S
MW-80B	T	7	1	1.26	-1	50.0%	No	NT
MW-81B	T	7	0	0.71	-6	76.4%	Yes	ND
MW-82B	T	6	1	1.02	5	76.5%	No	NT
MW-83B	T	7	7	1.23	-11	93.2%	No	PD
MW-83C	T	6	5	1.26	-4	70.3%	No	NT
MW-84B	T	7	6	1.42	-11	93.2%	No	PD
MW-85C	T	6	3	0.63	-8	89.8%	No	S
MW-86C	T	6	1	0.69	5	76.5%	No	NT
MW-87C	T	6	3	1.15	-4	70.3%	No	NT

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## 2-METHYLNAPHTHALENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-88C	T	6	1	0.55	-3	64.0%	No	S
MW-89B	T	4	2	0.42	-2	62.5%	No	S
MW-90B	T	4	1	1.62	3	72.9%	No	NT
P-10	T	1	1	0.00	0	0.0%	No	N/A
P-11	T	17	6	0.79	-22	80.4%	No	S
P-12	T	2	0	0.00	0	0.0%	Yes	ND
TW-41B	T	14	9	1.69	41	98.7%	No	I

## BENZENE

MW-03	T	6	0	0.00	0	42.3%	Yes	ND
MW-04	T	6	0	0.00	0	42.3%	Yes	ND
MW-05	T	13	0	1.01	-13	76.4%	Yes	ND
MW-09	T	10	0	0.91	-17	92.2%	Yes	ND
MW-12A	T	17	5	1.12	-59	99.2%	No	D
MW-12C	T	17	0	0.83	-44	96.2%	Yes	ND
MW-13	T	17	0	0.83	-44	96.2%	Yes	ND
MW-14	T	17	0	0.83	-44	96.2%	Yes	ND
MW-15A	T	17	15	0.53	-78	100.0%	No	D
MW-15B	T	12	7	0.96	-36	99.3%	No	D
MW-15C	T	17	14	0.29	-89	100.0%	No	D
MW-17	S	17	16	0.42	-23	81.5%	No	S
MW-17C	S	17	16	0.53	-60	99.3%	No	D
MW-18A	S	17	17	0.56	-40	94.6%	No	PD
MW-18C	S	17	17	0.23	-26	84.6%	No	S
MW-19C	S	17	12	1.05	-11	65.7%	No	NT
MW-20A	S	10	10	0.53	-33	99.9%	No	D
MW-21C	T	17	0	0.80	-49	97.7%	Yes	ND
MW-22AR	T	6	0	0.00	0	42.3%	Yes	ND
MW-22BR	T	6	0	0.00	0	42.3%	Yes	ND
MW-23C	S	12	10	0.39	-19	88.9%	No	S
MW-25A	T	17	1	0.76	-51	98.1%	No	D
MW-25C	S	15	13	1.20	-81	100.0%	No	D
MW-26A	T	20	9	2.30	-78	99.4%	No	D
MW-27A	T	5	0	0.00	0	40.8%	Yes	ND

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## BENZENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-27C	T	17	0	0.80	-49	97.7%	Yes	ND
MW-28A	T	17	0	0.80	-49	97.7%	Yes	ND
MW-28C	T	17	0	0.80	-49	97.7%	Yes	ND
MW-32AR	T	12	3	2.29	-31	98.1%	No	D
MW-32B	S	5	4	1.29	2	59.2%	No	NT
MW-33A	T	17	11	2.31	-37	93.0%	No	PD
MW-33BR	T	10	8	1.18	-17	92.2%	No	PD
MW-34CR	T	5	1	0.11	4	75.8%	No	NT
MW-35A	T	17	2	0.74	-65	99.7%	No	D
MW-35B	T	17	17	0.45	-25	83.6%	No	S
MW-36A	T	17	0	0.80	-49	97.7%	Yes	ND
MW-36B	T	15	2	1.15	-35	95.4%	No	D
MW-36D	T	15	2	0.88	-28	90.8%	No	PD
MW-38A	T	17	0	0.80	-49	97.7%	Yes	ND
MW-38B	T	17	0	0.80	-49	97.7%	Yes	ND
MW-39B	T	17	0	0.83	-44	96.2%	Yes	ND
MW-40B	T	17	17	0.48	-94	100.0%	No	D
MW-42B	T	14	0	1.26	-12	72.3%	Yes	ND
MW-44A	T	17	9	1.09	-41	95.0%	No	D
MW-44C	S	4	1	1.21	-4	83.3%	No	NT
MW-47C	T	14	0	0.78	-37	97.6%	Yes	ND
MW-48C	T	17	0	0.80	-49	97.7%	Yes	ND
MW-49A	T	17	15	0.97	-70	99.8%	No	D
MW-49B	S	15	15	1.32	39	97.1%	No	I
MW-50A	T	17	0	0.83	-44	96.2%	Yes	ND
MW-51A	T	17	0	0.83	-44	96.2%	Yes	ND
MW-51C	T	7	1	0.19	6	76.4%	No	NT
MW-53C	T	17	0	0.80	-49	97.7%	Yes	ND
MW-54C	T	17	0	0.80	-49	97.7%	Yes	ND
MW-57A	S	13	13	0.85	-56	100.0%	No	D
MW-57B	S	11	11	0.63	-23	95.7%	No	D
MW-58A	T	17	15	1.39	-35	91.8%	No	PD
MW-59A	T	17	0	0.80	-49	97.7%	Yes	ND

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

BENZENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-59B	T	15	1	0.88	-31	93.0%	No	PD
MW-59D	T	17	1	0.79	-47	97.1%	No	D
MW-60A	T	16	0	0.79	-45	97.7%	Yes	ND
MW-61A	T	17	0	0.80	-49	97.7%	Yes	ND
MW-62B	T	17	3	1.48	-24	82.6%	No	NT
MW-63B	T	16	15	1.40	18	77.5%	No	NT
MW-64A	T	17	1	0.81	-43	95.8%	No	D
MW-65D	T	17	1	0.90	-50	97.9%	No	D
MW-66D	T	16	0	0.87	-35	93.6%	Yes	ND
MW-67B	T	15	0	0.89	-31	93.0%	Yes	ND
MW-68A	T	3	0	0.00	0	0.0%	Yes	ND
MW-68B	S	12	12	0.25	-42	99.8%	No	D
MW-68C	T	16	13	0.92	-33	92.4%	No	PD
MW-69A	T	14	0	0.88	-29	93.7%	Yes	ND
MW-70B	S	5	5	0.52	4	75.8%	No	NT
MW-71B	T	12	11	1.55	2	52.7%	No	NT
MW-72B	S	10	10	0.27	-24	98.2%	No	D
MW-74B	S	11	11	0.36	2	53.0%	No	NT
MW-75B	S	6	6	0.56	-11	97.2%	No	D
MW-76C	T	7	2	0.10	5	71.9%	No	NT
MW-77A	S	7	5	0.69	6	76.4%	No	NT
MW-79A	S	7	7	1.04	-3	61.4%	No	NT
MW-80B	T	7	1	0.23	6	76.4%	No	NT
MW-81B	T	7	1	0.25	5	71.9%	No	NT
MW-82B	T	6	0	0.00	0	42.3%	Yes	ND
MW-83B	T	7	7	0.24	12	94.9%	No	PI
MW-83C	T	6	0	0.00	0	42.3%	Yes	ND
MW-84B	T	7	7	0.64	1	50.0%	No	NT
MW-85C	T	6	3	1.77	-4	70.3%	No	NT
MW-86C	T	6	0	0.00	0	42.3%	Yes	ND
MW-87C	T	6	0	0.00	0	42.3%	Yes	ND
MW-88C	T	6	0	0.00	0	42.3%	Yes	ND
MW-89B	T	4	0	0.00	0	37.5%	Yes	ND

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## BENZENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-90B	T	4	0	0.00	0	37.5%	Yes	ND
P-11	T	17	2	0.79	-54	98.6%	No	D
TW-41B	T	14	6	0.67	25	90.4%	No	PI

## DIBENZOFURAN

MW-01A	T	23	20	1.10	4	53.2%	No	NT
MW-02	T	23	16	1.67	-26	74.4%	No	NT
MW-03	T	6	0	0.00	0	42.3%	Yes	ND
MW-04	T	6	2	0.69	4	70.3%	No	NT
MW-05	T	13	6	2.66	-40	99.3%	No	D
MW-07	T	23	2	1.19	-175	100.0%	No	D
MW-08	T	23	1	1.17	-169	100.0%	No	D
MW-09	T	10	1	0.85	-21	96.4%	No	D
MW-10A	T	23	3	2.43	-153	100.0%	No	D
MW-10B	T	25	24	1.40	-4	52.8%	No	NT
MW-11A	T	23	3	1.04	-168	100.0%	No	D
MW-11B	T	23	17	1.04	-91	99.2%	No	D
MW-12A	T	17	17	0.46	1	50.0%	No	NT
MW-12C	T	17	11	4.02	-3	53.2%	No	NT
MW-13	T	17	6	0.84	-41	95.0%	No	D
MW-14	T	17	13	0.59	-65	99.7%	No	D
MW-15A	T	17	17	0.48	-42	95.4%	No	D
MW-15B	T	12	10	1.43	-41	99.8%	No	D
MW-15C	T	17	17	1.44	-53	98.5%	No	D
MW-17	S	18	17	0.73	-37	91.2%	No	PD
MW-17C	S	18	18	0.63	-41	93.4%	No	PD
MW-18A	S	17	17	0.38	5	56.4%	No	NT
MW-18C	S	17	17	0.48	-12	67.2%	No	S
MW-19C	S	17	11	0.77	22	80.4%	No	NT
MW-20A	S	10	10	0.32	-15	89.2%	No	S
MW-21C	T	17	5	2.41	-14	70.1%	No	NT
MW-22AR	T	6	1	1.22	7	86.4%	No	NT
MW-22BR	T	6	4	0.79	-8	89.8%	No	S
MW-23C	S	12	12	1.77	0	47.3%	No	NT



# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

DIBENZOFURAN

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-25A	T	17	7	3.72	-66	99.7%	No	D
MW-25C	S	15	14	0.58	-29	91.6%	No	PD
MW-26A	T	17	14	2.03	-48	97.4%	No	D
MW-27A	T	5	1	1.22	2	59.2%	No	NT
MW-27C	T	17	1	1.12	-61	99.4%	No	D
MW-28A	T	17	3	1.33	5	56.4%	No	NT
MW-28C	T	17	2	0.77	-81	100.0%	No	D
MW-32AR	T	12	8	2.60	-33	98.7%	No	D
MW-32B	S	5	4	2.20	4	75.8%	No	NT
MW-33A	T	17	15	2.10	-29	87.4%	No	NT
MW-33BR	T	10	7	1.92	-15	89.2%	No	NT
MW-34CR	T	5	1	0.77	-1	50.0%	No	S
MW-35A	T	17	16	1.08	-16	72.9%	No	NT
MW-35B	T	17	17	0.48	-46	96.8%	No	D
MW-36A	T	17	2	1.46	-61	99.4%	No	D
MW-36B	T	15	2	1.07	-11	68.7%	No	NT
MW-36D	T	15	1	0.80	-37	96.3%	No	D
MW-38A	T	18	5	0.89	-20	76.2%	No	S
MW-38B	T	17	5	2.54	12	67.2%	No	NT
MW-39B	T	17	5	1.40	-19	76.8%	No	NT
MW-40B	T	17	17	0.39	-21	79.2%	No	S
MW-42B	T	14	8	0.90	-48	99.6%	No	D
MW-44A	T	17	12	2.08	-63	99.6%	No	D
MW-44C	S	4	3	1.97	0	37.5%	No	NT
MW-47C	T	14	3	0.35	-45	99.3%	No	D
MW-48C	T	17	3	3.88	-75	99.9%	No	D
MW-49A	T	17	14	1.01	-65	99.7%	No	D
MW-49B	S	15	14	3.85	37	96.3%	No	I
MW-50A	T	17	4	2.66	-50	97.9%	No	D
MW-51A	T	17	2	0.51	-75	99.9%	No	D
MW-51C	T	7	0	0.90	1	50.0%	Yes	ND
MW-53C	T	17	0	1.12	-47	97.1%	Yes	ND
MW-54C	T	17	16	0.65	-5	56.4%	No	S

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

DIBENZOFURAN

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-57A	S	13	13	1.31	-16	81.6%	No	NT
MW-57B	S	11	11	0.59	-5	61.9%	No	S
MW-58A	T	17	15	0.85	-30	88.2%	No	S
MW-59A	T	17	3	1.62	-34	91.2%	No	PD
MW-59B	T	15	0	0.50	-45	98.6%	Yes	ND
MW-59D	T	17	3	2.06	-57	99.0%	No	D
MW-60A	T	17	5	1.70	-62	99.5%	No	D
MW-61A	T	17	0	0.51	-82	100.0%	Yes	ND
MW-62B	T	18	13	2.29	-40	92.9%	No	PD
MW-63B	T	16	14	1.06	-41	96.5%	No	D
MW-64A	T	17	5	1.97	-51	98.1%	No	D
MW-65D	T	17	4	0.59	-80	100.0%	No	D
MW-66D	T	17	4	1.01	-53	98.5%	No	D
MW-67B	T	15	2	0.57	-54	99.7%	No	D
MW-68A	T	3	1	0.00	0	0.0%	No	N/A
MW-68B	S	12	12	0.41	0	47.3%	No	S
MW-68C	T	15	8	2.00	-43	98.2%	No	D
MW-69A	T	14	3	2.50	-54	99.9%	No	D
MW-70B	S	5	5	1.65	0	40.8%	No	NT
MW-71B	T	12	8	1.41	-3	55.4%	No	NT
MW-72B	S	10	10	1.36	-19	94.6%	No	PD
MW-74B	S	11	10	1.15	-5	61.9%	No	NT
MW-75B	S	6	6	1.15	1	50.0%	No	NT
MW-76C	T	8	6	0.91	-13	92.9%	No	PD
MW-77A	S	7	6	0.64	3	61.4%	No	NT
MW-79A	S	7	7	0.85	-10	90.7%	No	PD
MW-80B	T	7	1	0.72	-1	50.0%	No	S
MW-81B	T	7	0	0.77	-6	76.4%	Yes	ND
MW-82B	T	6	1	0.35	5	76.5%	No	NT
MW-83B	T	7	7	1.09	-12	94.9%	No	PD
MW-83C	T	6	5	1.10	-5	76.5%	No	NT
MW-84B	T	7	6	1.29	-11	93.2%	No	PD
MW-85C	T	6	2	0.68	-2	57.0%	No	S

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

## DIBENZOFURAN

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-86C	T	6	0	0.00	0	42.3%	Yes	ND
MW-87C	T	6	3	1.27	-4	70.3%	No	NT
MW-88C	T	6	2	0.55	-7	86.4%	No	S
MW-89B	T	4	0	1.06	3	72.9%	Yes	ND
MW-90B	T	4	1	1.06	3	72.9%	No	NT
P-10	T	23	6	1.59	-107	99.8%	No	D
P-11	T	17	11	1.88	-16	72.9%	No	NT
P-12	T	23	0	1.20	-184	100.0%	Yes	ND
TW-41B	T	14	10	1.01	43	99.0%	No	I

## NAPHTHALENE

MW-01A	T	23	13	3.84	-113	99.9%	No	D
MW-02	T	23	14	1.92	-131	100.0%	No	D
MW-03	T	6	0	1.25	7	86.4%	Yes	ND
MW-04	T	6	3	0.93	-1	50.0%	No	S
MW-05	T	13	5	1.18	-18	84.7%	No	NT
MW-07	T	23	4	1.10	-151	100.0%	No	D
MW-08	T	23	0	1.19	-186	100.0%	Yes	ND
MW-09	T	10	3	2.55	5	63.6%	No	NT
MW-10A	T	23	4	3.26	-149	100.0%	No	D
MW-10B	T	24	24	1.98	-64	94.1%	No	PD
MW-11A	T	23	3	1.09	-170	100.0%	No	D
MW-11B	T	24	13	2.66	-29	75.4%	No	NT
MW-12A	T	17	13	1.63	-104	100.0%	No	D
MW-12C	T	17	10	0.89	-14	70.1%	No	S
MW-13	T	17	6	1.89	-6	58.0%	No	NT
MW-14	T	17	10	0.67	-85	100.0%	No	D
MW-15A	T	17	12	1.62	-38	93.6%	No	PD
MW-15B	T	12	8	1.62	-34	99.0%	No	D
MW-15C	T	17	8	4.12	-36	92.4%	No	PD
MW-17	S	18	17	0.50	-52	97.4%	No	D
MW-17C	S	18	18	2.10	-10	63.2%	No	NT
MW-18A	S	17	17	0.42	0	48.4%	No	S
MW-18C	S	17	17	0.37	-2	51.6%	No	S

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

NAPHTHALENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-19C	S	17	10	1.70	-46	96.8%	No	D
MW-20A	S	10	10	1.80	-15	89.2%	No	NT
MW-21C	T	17	8	1.96	-7	59.6%	No	NT
MW-22AR	T	6	1	1.24	-2	57.0%	No	NT
MW-22BR	T	6	4	1.62	-13	99.2%	No	D
MW-23C	S	12	12	1.21	16	84.5%	No	NT
MW-25A	T	17	8	3.82	9	62.7%	No	NT
MW-25C	S	15	15	0.62	-43	98.2%	No	D
MW-26A	T	17	7	1.48	-45	96.5%	No	D
MW-27A	T	5	2	2.12	5	82.1%	No	NT
MW-27C	T	17	7	1.23	-38	93.6%	No	PD
MW-28A	T	17	5	2.80	11	65.7%	No	NT
MW-28C	T	17	7	1.17	-65	99.7%	No	D
MW-32AR	T	12	8	3.09	-22	92.4%	No	PD
MW-32B	S	5	3	2.00	4	75.8%	No	NT
MW-33A	T	17	12	2.19	-28	86.5%	No	NT
MW-33BR	T	10	8	1.63	-21	96.4%	No	D
MW-34CR	T	5	1	1.48	0	40.8%	No	NT
MW-35A	T	17	14	1.92	26	84.6%	No	NT
MW-35B	T	17	17	0.56	-30	88.2%	No	S
MW-36A	T	17	5	1.56	-37	93.0%	No	PD
MW-36B	T	15	3	2.41	-11	68.7%	No	NT
MW-36D	T	15	6	1.58	-30	92.3%	No	PD
MW-38A	T	17	7	2.00	-18	75.5%	No	NT
MW-38B	T	17	9	1.35	18	75.5%	No	NT
MW-39B	T	17	7	1.17	20	78.0%	No	NT
MW-40B	T	17	17	0.52	-58	99.1%	No	D
MW-42B	T	14	7	1.04	-4	56.4%	No	NT
MW-44A	T	17	12	2.44	-40	94.6%	No	PD
MW-44C	S	4	3	1.81	-2	62.5%	No	NT
MW-47C	T	14	7	1.87	-1	50.0%	No	NT
MW-48C	T	17	5	4.09	-42	95.4%	No	D
MW-49A	T	17	13	1.06	-47	97.1%	No	D

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

NAPHTHALENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-49B	S	15	14	3.77	33	94.3%	No	PI
MW-50A	T	17	8	2.87	-14	70.1%	No	NT
MW-51A	T	17	6	1.06	39	94.1%	No	PI
MW-51C	T	7	4	0.46	1	50.0%	No	NT
MW-53C	T	17	5	1.26	-6	58.0%	No	NT
MW-54C	T	17	15	1.36	-38	93.6%	No	PD
MW-57A	S	13	13	1.01	-18	84.7%	No	NT
MW-57B	S	11	11	0.57	-28	98.4%	No	D
MW-58A	T	17	15	1.20	-12	67.2%	No	NT
MW-59A	T	17	5	2.62	-5	56.4%	No	NT
MW-59B	T	15	4	0.69	-4	55.8%	No	S
MW-59D	T	17	5	1.62	-42	95.4%	No	D
MW-60A	T	17	6	2.14	-48	97.4%	No	D
MW-61A	T	17	1	3.21	-50	97.9%	No	D
MW-62B	T	17	8	2.15	-39	94.1%	No	PD
MW-63B	T	16	13	1.29	-24	84.7%	No	NT
MW-64A	T	17	3	1.54	-7	59.6%	No	NT
MW-65D	T	17	6	1.73	-29	87.4%	No	NT
MW-66D	T	17	4	2.68	-43	95.8%	No	D
MW-67B	T	15	8	1.23	-30	92.3%	No	PD
MW-68A	T	3	2	0.00	0	0.0%	No	N/A
MW-68B	S	12	12	0.51	-22	92.4%	No	PD
MW-68C	T	15	12	1.20	-27	89.9%	No	NT
MW-69A	T	14	2	3.03	-35	96.9%	No	D
MW-70B	S	5	5	0.90	8	95.8%	No	I
MW-71B	T	12	8	2.34	-11	74.9%	No	NT
MW-72B	S	10	10	1.24	-26	98.9%	No	D
MW-74B	S	11	11	1.49	-7	67.6%	No	NT
MW-75B	S	6	6	0.93	-1	50.0%	No	S
MW-76C	T	8	6	1.00	2	54.8%	No	NT
MW-77A	S	7	6	0.79	-3	61.4%	No	S
MW-79A	S	7	7	0.67	1	50.0%	No	NT
MW-80B	T	7	2	1.95	1	50.0%	No	NT

# MAROS Mann-Kendall Statistics Summary

Project: HWPW

User Name: MHH

Location: Houston

State: Texas

NAPHTHALENE

Well	Source/ Tail	Number of Samples	Number of Detects	Coefficient of Variation	Mann- Kendall Statistic	Confidence in Trend	All Samples "ND" ?	Concentration Trend
MW-81B	T	7	2	0.91	10	90.7%	No	PI
MW-82B	T	6	1	1.14	-2	57.0%	No	NT
MW-83B	T	7	7	1.40	-15	98.5%	No	D
MW-83C	T	6	5	1.29	-3	64.0%	No	NT
MW-84B	T	7	6	1.33	-11	93.2%	No	PD
MW-85C	T	6	4	0.84	-3	64.0%	No	S
MW-86C	T	6	3	1.28	12	98.2%	No	I
MW-87C	T	6	3	1.15	-2	57.0%	No	NT
MW-88C	T	6	3	1.40	-11	97.2%	No	D
MW-89B	T	4	2	1.03	0	37.5%	No	NT
MW-90B	T	4	3	1.87	6	95.8%	No	I
P-10	T	23	6	3.35	-108	99.8%	No	D
P-11	T	17	8	3.29	-25	83.6%	No	NT
P-12	T	23	1	1.15	-169	100.0%	No	D
TW-41B	T	14	11	1.82	49	99.7%	No	I

Note: Increasing (I); Probably Increasing (PI); Stable (S); Probably Decreasing (PD); Decreasing (D); No Trend (NT); Not Applicable (N/A)-Due to insufficient Data (< 4 sampling events); Source/Tail (S/T)

The Number of Samples and Number of Detects shown above are post-consolidation values.

**ATTACHMENT H**

**Naphthalene Concentrations in  
MW-11B**

**Table 1**  
**Historical Naphthalene Concentrations at MW-11B**  
**Houston Wood Preserving Works**

Monitoring Well ID	Collection Date	Naphthalene (mg/L)	Qualifier
GWPS		<b>0.49</b>	
MW-11B	9/28/1994	<b>1.1</b>	
MW-11B	10/13/1994	<b>0.5</b>	
MW-11B	1/25/1995	<b>1.6</b>	
MW-11B	4/13/1995	<b>1.5</b>	
MW-11B	7/12/1995	<b>1</b>	
MW-11B	7/12/1995	<b>1.7</b>	
MW-11B	1/23/1996	<b>0.75</b>	
MW-11B	9/17/1996	<b>2.44</b>	
MW-11B	3/26/1997	0.1	U
MW-11B	9/25/1997	0.32	
MW-11B	3/3/1998	0.01	U
MW-11B	3/17/1999	<b>0.608</b>	
MW-11B	9/28/1999	0.409	D
MW-11B	3/29/2000	0.079	
MW-11B	9/22/2000	<b>1.5</b>	
MW-11B	4/25/2001	0.47	
MW-11B	9/27/2001	<b>0.5</b>	
MW-11B	3/14/2002	0.22	
MW-11B	9/24/2002	<b>0.7</b>	
MW-11B	3/16/2004	0.01168	
MW-11B	3/1/2005	0.00006	U
MW-11B	7/19/2005	0.186	
MW-11B	1/5/2006	0.0025	
MW-11B	7/31/2006	0.1	
MW-11B	1/23/2007	0.00013	
MW-11B	7/17/2007	0.0901	
MW-11B	1/28/2008	0.0354	
MW-11B	7/16/2008	0.0772	
MW-11B	1/22/2009	0.0008	U
MW-11B	7/22/2009	0.048	
MW-11B	1/21/2010	0.0006	U
MW-11B	7/13/2010	0.0068	
MW-11B	1/11/2011	0.0006	U
MW-11B	7/12/2011	0.06	
MW-11B	1/30/2012	0.0005	U
MW-11B	7/10/2012	0.004	J
MW-11B	1/13/2013	0.00008	U
MW-11B	7/11/2013	0.00535	
MW-11B	1/8/2014	0.000382	J
MW-11B	7/2/2014	0.0135	
MW-11B	1/15/2015	0.00008	U
MW-11B	7/8/2015	0.0021	
MW-11B	1/12/2016	0.00002	U
MW-11B	7/7/2016	0.0082	
MW-11B	1/11/2017	0.00002	U
MW-11B	7/12/2017	0.0019	
MW-11B	1/3/2018	0.00002	U
MW-11B	7/18/2018	0.34	
MW-11B	1/7/2019	0.00002	U
MW-11B	7/1/2019	<b>0.7</b>	
MW-11B	7/30/2019	<b>1.1</b>	
MW-11B	10/17/2019	<b>0.6</b>	
MW-11B	1/14/2020	<b>0.0002</b>	U

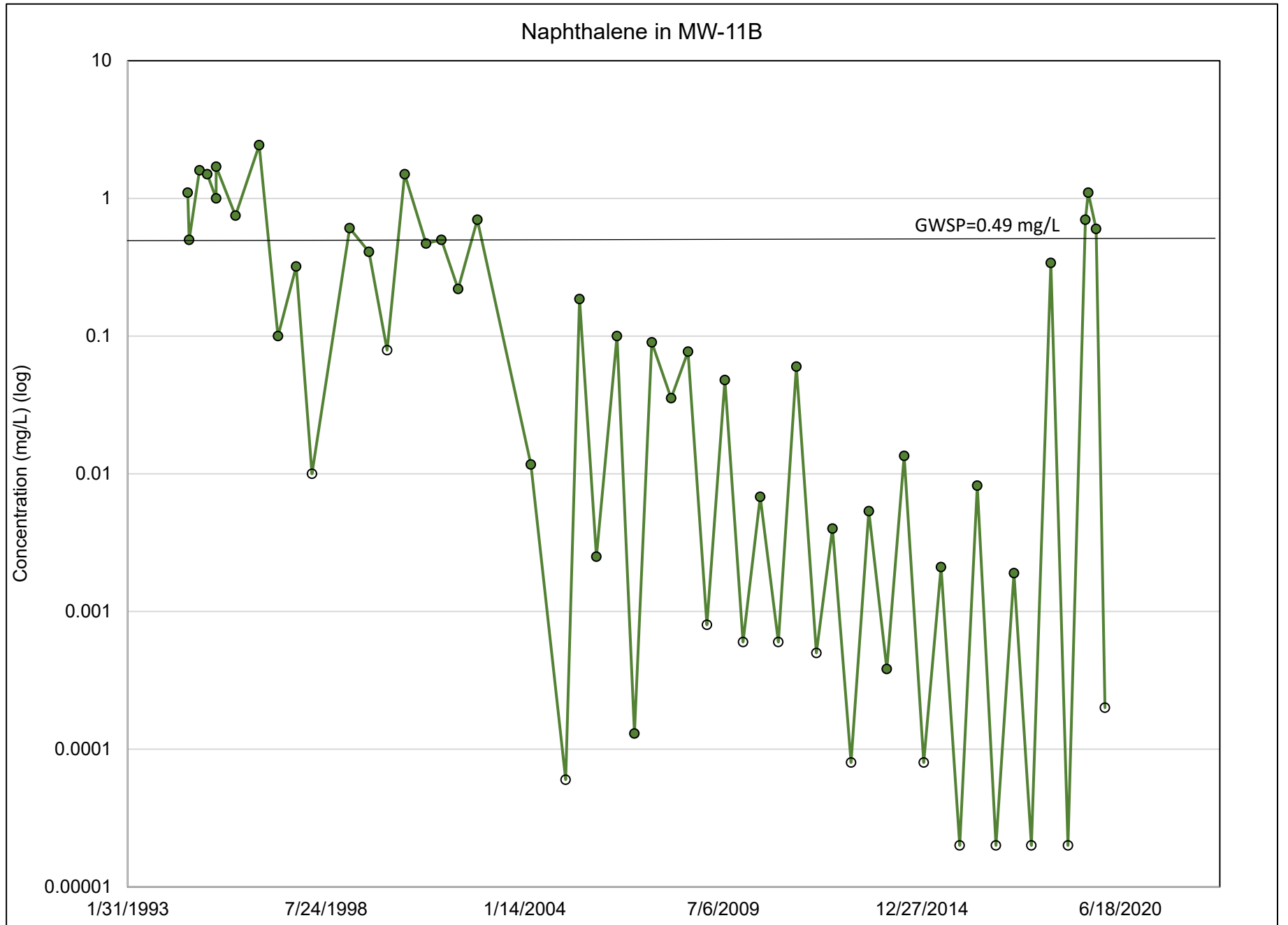
Notes:

1. Highlighted and bolded values exceed TCEQ Tier 1 Residential PCLs (November 2019)/ Groundwater Protection Standard (GWSP)

U - concentration below the method detection limit

J - estimate; concentration above the method detection limit & below the method quantitation limit





Note: Data points with no fill indicate that concentrations were below the method detection limit (MDL). The MDL is used as the result.



**ATTACHMENT I**

# TRRP Notifications

**Table I-1  
TRRP NOTIFICATIONS - PROPERTY OWNERS  
HOUSTON, TX - WOOD PRESERVING WORKS**

MAP ID	PARCEL ADDRESS	PARCEL OWNER
1	4705 Liberty Rd	Hall, Joannetta
2	4713 Liberty Rd	Howell, Evan
3	2902 Wipprecht St # 18	2902 Wipprecht LLC % T Goffney
4	2909 Lavender St	2013 Cottage LLC
5	2922 Solo St	Chaves, Armando R et Al
6	2925 Lavender St	Ellis, Johnnie Estate Of
7	2930 Solo St	Howell, Evan S
8	2934 Solo St	Green, Billie Joe
9	2938 Solo St	Ramos, Jose Rodolfo
10	2937 Lavender St	Limbrick, Mikell
11	2941 Lavender St	Brooks, Antonia Est Of % Patricia J Arthur
12	4910 Lucille St	Whitehead, Shirley A
13	5002 Lucille St	Toledano, Aristeo
14	5004 Lucille St	Fuentes, Maria I & Saul
15	2943 Lavender St	Apogee Real Estate Partners LLC
16	5006 Lucille St	Barrientos, Eine
17	5008 Lucille St	Norton Memorial Temple Cogic
18	2942 Lavender St	Hernandez, Roberto G
19	2938 Lavender St	Hutchins, Essie Lee
20	2934 Lavender St	Osborn, Zearlene
21	2926 Lavender St	Clark Investment Co
22	2924 Lavender St	Long, Lucille Estate Of
23	2922 Lavender St	Greater Mt Nebo Baptist Church
24	2910 Lavender St	Carrington, Ray
25	2906 Lavender St	Beal, Barbara
26	2904 Lavender St	Greater Mount Nebo Missionary Baptist
27	5005 Liberty	Greater Mount Nebo Baptist Church
28	0 Wylie St	Greater Mt Nebo Baptist Church
29	5006 Wylie St	Greater Mount Nebo Baptist Church
30	5010 Wylie St	Smith, Alberta
31	5005 Wylie St	Banda, Monico Duque & Martha Z
32	5007 Wylie St	Banda-Zuniga, Nancy G; Macia-Aranda, Guillermo
33	5011 Wylie St	Carr, Carrie Mae Estate Of
34	2901 Clementine St	Ross, Mary Bass
35	5002 Lelia St	Holmes Emmitt & Laura Estate % Mildred Wilder
36	0 Lelia St	Heptullabhai, Mustafa
37	5014 Lelia St	Rivera, Guadalupe Jr; Rivera Flor Esthela
38	2921 1/2 Clementine	Rivera, Guadalupe
39	2921 Clementine St	Pineda, Rogelio R & Olivia
40	0 Lelia St	Lelia Street Trust
41	5009 Lelia St	Current Owner
42	0 Lelia	Beach Property LLC
43	5015 Lelia St	Castillo, Manuel M
44	5012 Lucille	Norton'S Temple Church Of God In Christ Inc
45	5014 Lucille St	Blue Gavel Inc
46	5016 Lucille St	Current Owner
47	0 Lucille	Wahsh, Belal
48	5005 Lucille St	Correa, Neftali; Correa, Eliseo Sr
49	5005 Lucille St	Bryant, Michael A; White, Karen Y; White, David A
50	5007 Lucille St	White, Karen Y

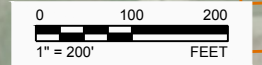
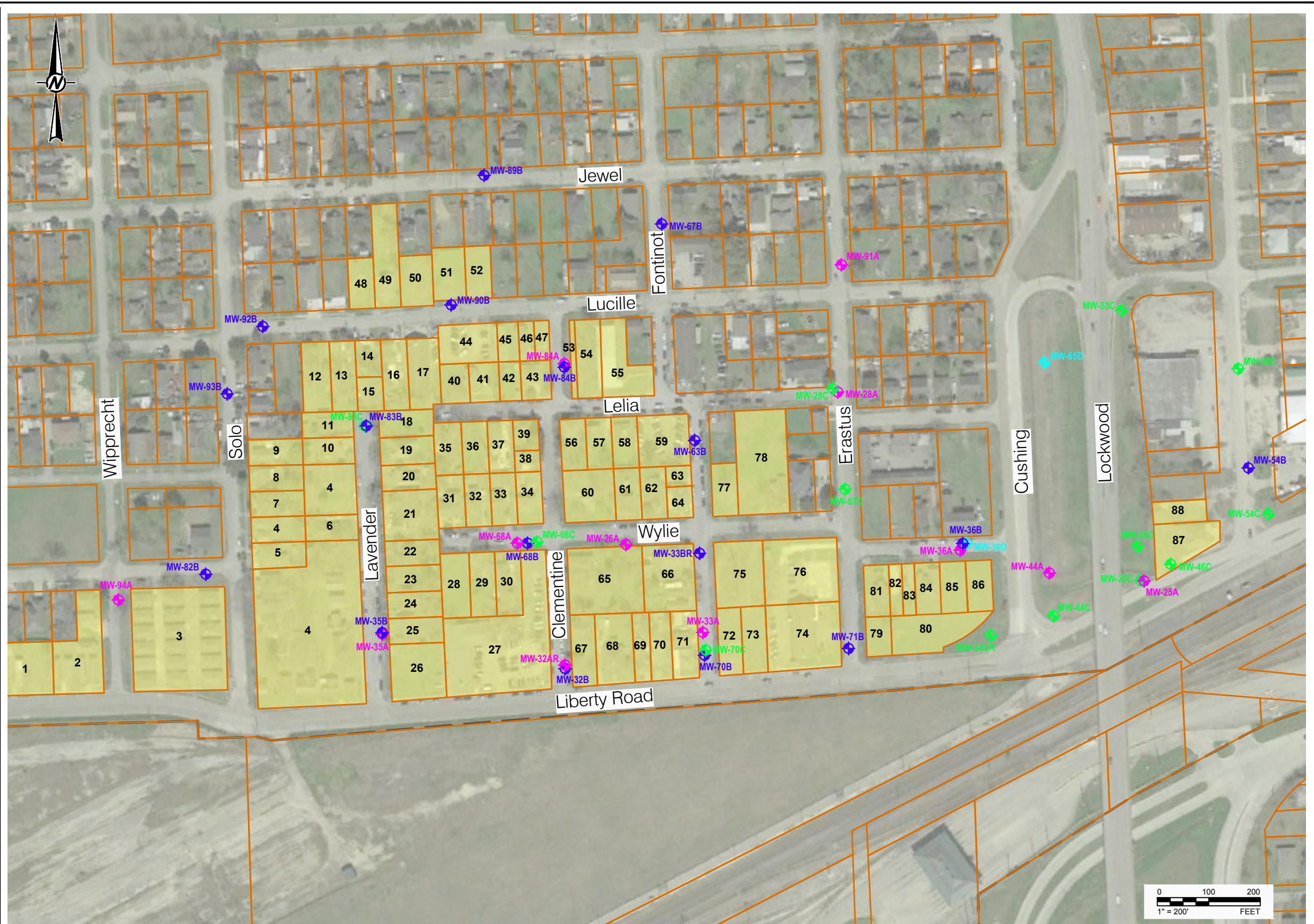
**Table I-1  
TRRP NOTIFICATIONS - PROPERTY OWNERS  
HOUSTON, TX - WOOD PRESERVING WORKS**

MAP ID	PARCEL ADDRESS	PARCEL OWNER
51	5009 Lucille St	White, David A
52	0 Lucille St	Vital, Percy L
53	0 Lelia	Schrinsky, Samuel J
54	705 Lucille	Pssw Construction Inc
55	0 Leila St	Greater True Vine Missionary Baptist Church
56	2918 Clementine St	City Of Houston
57	5116 Lelia St	Empire Holdings Real Estate Group LLC; CJO Investments LLC
58	5112 Lelia St	Potts, Avie
59	5118 Lelia St	Greater True Vine Baptist Church
60	5107 Wylie St	Potts, Maryland Estate
61	5107 Wylie St	Tolbert, Reginald & Leticia
62	5111 Wylie St	Perez, Aquilina Estate Of
63	2913 Fontinot St	Huerta, Noe; Rivera, Marely
64	5119 Wylie St	Gilliam, Martha
65	2820 Clementine St	Benson, Tillie Potts Estate Of
66	2813 Fontinot St	Coto, Jose A & Reina I; Coto, Christian A & Robin H
67	5101 Liberty Rd	Longoria, Wallace R & Janie
68	5105 Liberty Rd	Gonzalez, Alejandro
69	5109 Liberty Rd	Martinez, Joe H
70	5113 Liberty Rd	Medina, Claudia Elizabeth Delaportil; Leal, Jorge Daniel Rivera
71	5117 Liberty Rd	Rivera, Jorge D
72	5201 Liberty Rd	Full Gospel Christian Assn
73	5201 Liberty Rd	Full Gospel Christian Assn
74	2809 Erastus St # 1	Charity Baptist Church % Rev F W Mcilveen
75	5201 Wylie St	Charity Baptist Church
76	2809 Erastus St	Charity Baptist Church % Rev F W Mcilveen
77	5201 Wylie St	Barahona, Melvin
78	5202 Lelia St	Greater True Vine Mbc
79	5301 Liberty Rd	Jefferson, Sean
80	5311 Liberty Rd	Crawford Properties & States
81	2806 Erastus St	Roberts, Margaret et al
82	5304 Wylie St	Juarez, Irene Perez
83	5304 1/2 Wylie St	Delgado, Susie I Estate Of
84	5308 Wylie St	Brown, Audrey M
85	0 Liberty Rd	Crawford Properties & Estates
86	5311 Liberty Rd	Cls Enterprise Llc
87	3300 E Lockwood Dr	Damian, Robert
88	3300 E Lockwood	Damian, Robert
89	4518 Courtney St	Rosignon, Anthony Tyrone
90	2803 Kashmere St	Whitehead, Shirley A
91	4508 Eddie St	Henderson, Hester
92	0 Ranch St	Prince, James A
93	2604 Amboy St	Daniels, Joeed
94	0 Quitman St	Proler, L Michael
95	4425 Quitman St	Allende-Molina, Juan; Rodriguez-Campos, Micaela
96	0 1St St	Centerpoint Energy Intrastate Pipeline Inc



Path: \\nasar\land\land\Projects - Round Rock\19119232 - HWPW\2019\7 July - 1 File Name: FIG 1 - Off-Site Notification Properties.dwg | Last Edited By: nasar | Date: 2020-04-30 Time: 2:23:57 PM | Printed By: RSalazar | Date: 2020-04-30 Time: 3:37:21 PM

MAP ID	PARCEL ADDRESS	PARCEL OWNER
1	4705 Liberty Rd	Hall, Joannetta
2	4713 Liberty Rd	Howell, Evan
3	2902 Wipprecht St # 18	2902 Wipprecht LLC % T Goffney
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33	5011 Wylie St	Macia-Aranda, Guillermo
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37	5014 Lelia St	Heptullahai, Mustafa
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50	5007 Lucille St	Bryant, Michael A; White, Karen Y; White, David A
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53	0 Lelia St	Vital, Percy L
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87	3300 E Lockwood Dr	Cis Enterprise Llc
88	3300 E Lockwood	Damian, Robert



**REFERENCE(S)**  
 AERIAL PHOTO FROM GOOGLE EARTH, IMAGERY DATED 2/23/19, AND PROPERTY INFORMATION FROM HARRIS COUNTY APPRAISAL DISTRICT (HCAD), 2018.

**CLIENT**  
 UNION PACIFIC RAILROAD CO.

**PROJECT**  
 HOUSTON WOOD PRESERVING WORKS

**CONSULTANT**

YYYY-MM-DD	2020-04-30
DESIGNED	AJD
PREPARED	RS
REVIEWED	MH
APPROVED	ECM

**TITLE**  
 OFF-SITE NOTIFICATION PROPERTIES

**PROJECT NO.**  
 19119232

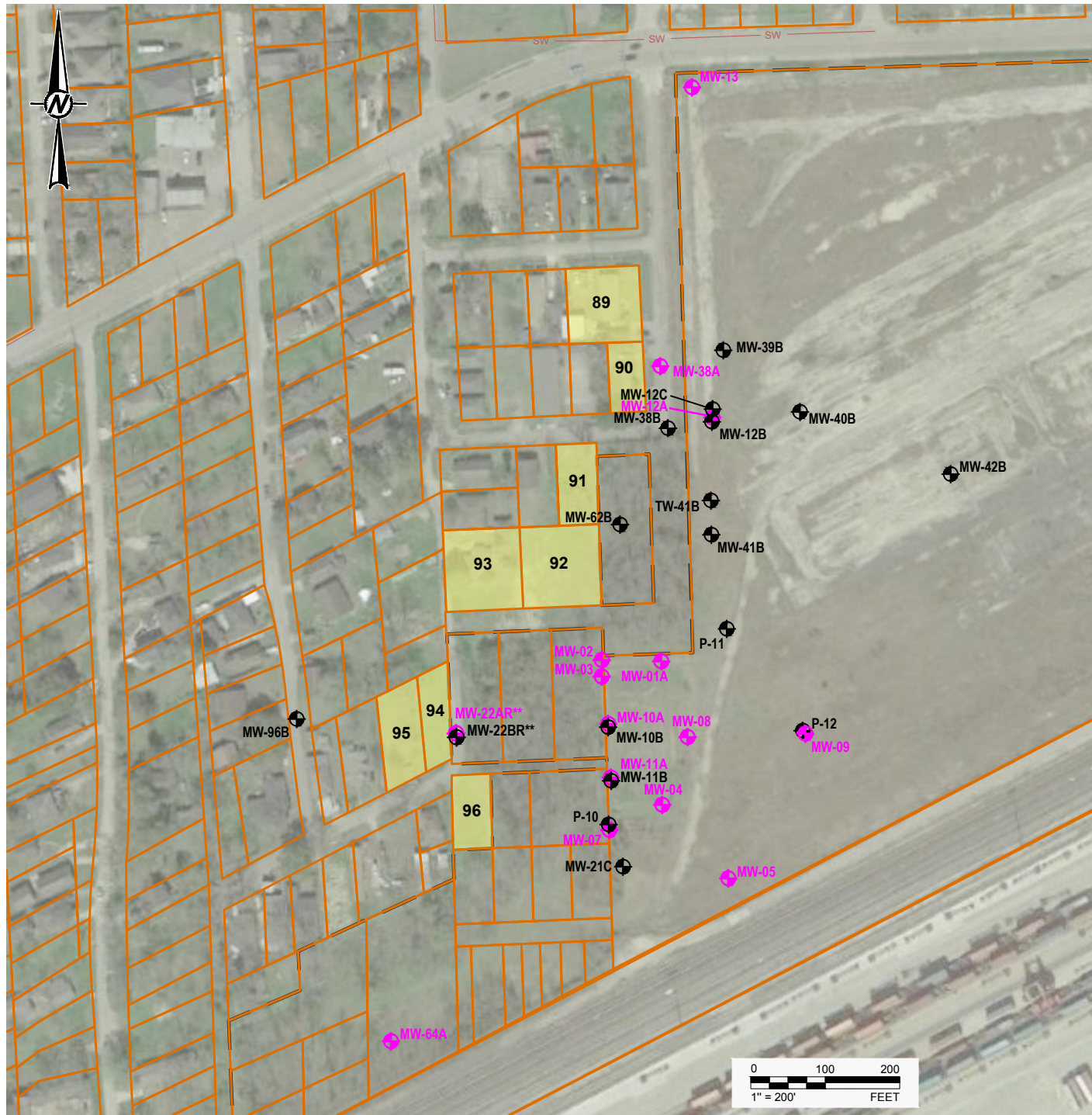
**REV.**  
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**FIGURE**  
 1

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM ANSI B



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 Path: \\solar\hna\data\Projects - Round Rock\_2019\19119232 - HWPPV\2019-7 July | File Name: FIG 2 - Off-Site Notification Properties.dwg



CLIENT  
UNION PACIFIC RAILROAD CO.

PROJECT  
HOUSTON WOOD PRESERVING WORKS

TITLE  
**OFF-SITE NOTIFICATION PROPERTIES**

CONSULTANT	YYYY-MM-DD	2020-04-30
	DESIGNED	AJD
	PREPARED	RS
	REVIEWED	MH
	APPROVED	ECM

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91	4508 Eddie St	Henderson, Hester
92	0 Ranch St	Prince, James A
93	2604 Amboy St	Daniels, Joeed
94	0 Quitman St	Proler, L Michael
95	4425 Quitman St	Allende-Molina, Juan; Rodriguez-Campos, Micaela
96	0 1St St	Centerpoint Energy Intrastate Pipeline Inc

**REFERENCE(S)**  
 AERIAL PHOTO FROM GOOGLE EARTH, IMAGERY DATED 2/23/19, AND PROPERTY INFORMATION FROM HARRIS COUNTY APPRAISAL DISTRICT (HCAD), 2018.

PROJECT NO. 19119232      REV. 0      FIGURE 1

IF THIS MEASUREMENT DOES NOT MATCH WHAT IS SHOWN, THE SHEET SIZE HAS BEEN MODIFIED FROM: ANSI A