

MCP PHASE I AND INTERIM PHASE II
REPORT FOR FORMER HOUSATONIC RIVER
OXBOW AREAS A, B, C, J, AND K

VOLUME II OF II

General Electric Company
Pittsfield, Massachusetts

February 1996

BBL
BLASLAND, BOUCK & LEE, INC.
engineers & scientists

012098
GE

REPORT

OUTSIDE THE RIVER	
Site:	GE-OUNO
Break:	2.2
Other:	5811

MCP PHASE I AND INTERIM PHASE II REPORT FOR FORMER HOUSATONIC RIVER OXBOW AREAS A, B, C, J, AND K

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GE-P 012098

MCP PHASE I AND INTERIM PHASE II REPORT FOR FORMER HOUSATONIC
RIVER OXBOW AREAS A, B, C, J, AND K

VOLUME II OF II

GENERAL ELECTRIC COMPANY
PITTSFIELD, MASSACHUSETTS

FEBRUARY 1996

BLASLAND, BOUCK & LEE, INC.
6723 TOWPATH ROAD, BOX 66
SYRACUSE, NEW YORK 13214

MCP PHASE I AND INTERIM PHASE II REPORT FOR
FORMER Housatonic RIVER OXBOW AREAS
A, B, C, J, AND K

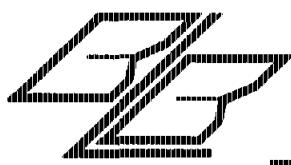
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Section 1

VOLATILE ORGANICS ANALYSIS (SOIL)

FP-1, 8-12'	Soil sample from Oxbow Area J, Soil Boring FP-1 at 8-12 feet
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FP-3, 4-8'	Soil sample from Oxbow Area J, Soil Boring FP-3 at 4-8 feet
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ROA011416	Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 14-16 feet
ROA012022	Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 20-22 feet
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ROB2B0002	Soil sample from Oxbow Area B, Groundwater Monitoring Well B-2 at 0-2 feet
ROB-DPA1	Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)
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ROC021214	Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet
ROC3B0204	Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet
ROC-DPA1	Soil sample from Oxbow Area C, Groundwater Monitoring Well 12-14 (Duplicate)

VOLATILE ORGANICS ANALYSIS (SOIL) (CONT'D)

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1
ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2
ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3
ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4
ROO1B1416 - Soil sample from Oxbow Area K, Soil Boring K1 at 14-16 feet
ROO2B0810 - Soil sample from Oxbow Area K, Soil Boring K2 at 8-10 feet

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: FP-1,8-12'
Lab Sample ID: JJ7107

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	.. 10 U	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
bromoform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	6 U
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	5 U
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	4 J -
chloroform	5 U	1,1,1-trichloroethane	5 U
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	1 J -
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: FP-2,4-8'
Lab Sample ID: JJ7108

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	10 U	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
bromoform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	6 J
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	5 U
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	3 J
chloroform	5 U	1,1,1-trichloroethane	5 U
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	5 U
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in ug/kg (ppb)

Sample Matrix: Soil

Client Sample ID: FP-3,4-8'
Lab Sample ID: JJ7109

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	10 U	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
bromoform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	5 J
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	5 U
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	3 J
chloroform	5 U	1,1,1-trichloroethane	5 U
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	5 U
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in ug/kg (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank
Lab Sample ID: VB1010

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	3 J	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
bromoform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	2 J -
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	1 J -
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	5 U
chloroform	5 U	1,1,1-trichloroethane	5 U
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	5 U
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

SOIL SURROGATE PERCENT RECOVERY SUMMARY

Sample No.	VOLATILE		
	Toluene-08 (81-117%)*	BFB (74-121%)*	1,2 Dichloroethane-04 (70-121%)*
Method Blank	95	94	94
FP-1,8-12'	105	80	96
FP-2,4-8'	103	87	94
FP-3,4-8'	106	93	84

*Values in parenthesis represent USEPA contract required QC limits.

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

- Client Project ID: GE Pittsfield

Job Number: GMIN 44313

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: YB-2, 4-8'
Lab Sample ID: JJ7170

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	10 U	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
bromoform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	3 J
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	5 U
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	1 J
chloroform	5 U	1,1,1-trichloroethane	4 J
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	5 U
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44313

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: YB-4, 0-4'
Lab Sample ID: JJ7171

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	10 U	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
chloroform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	2 J
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	5 U
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	5 U
chloroform	5 U	1,1,1-trichloroethane	5
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	5 U
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44313

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank
Lab Sample ID: VBI010

<u>Compound</u>		<u>Compound</u>	
acrolein	10 U	1,1-dichloroethene	5 U
acrylonitrile	3 J	trans-1,2-dichloroethene	5 U
benzene	5 U	1,2-dichloropropane	5 U
bromodichloromethane	5 U	cis-1,3-dichloropropene	5 U
bromoform	5 U	trans-1,3-dichloropropene	5 U
bromomethane	10 U	ethyl benzene	5 U
carbon tetrachloride	5 U	methylene chloride	2 J
chlorobenzene	5 U	1,1,2,2-tetrachloroethane	1 J
chloroethane	10 U	tetrachloroethene	5 U
2-chloroethylvinyl ether	10 U	toluene	5 U
chloroform	5 U	1,1,1-trichloroethane	5 U
chloromethane	10 U	1,1,2-trichloroethane	5 U
dibromochloromethane	5 U	trichloroethene	5 U
1,1-dichloroethane	5 U	vinyl chloride	10 U
1,2-dichloroethane	5 U		

U - Compound analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

-Client Project ID: GE Pittsfield

Job Number: GMIN 44313

SOIL SURROGATE PERCENT RECOVERY SUMMARY

Sample No.	VOLATILE		
	Toluene-08 (81-117%)*	BFB (74-121%)*	1,2 Dichloroethane-04 (70-121%)*
Method Blank	95	94	94
YB-2, 4-8'	96	90	83
YB-4, 0-4'	104	86	91

*Values in parenthesis represent USEPA contract required QC limits.

LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM.RTP	Contract: 500077	ROA010406
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 26
Matrix: (soil/water) SOIL	Lab Sample ID: 462129	
Sample wt/vol: 5.0 (g/mL) G	Lab File ID: GB062129C13	
Level: (low/med) LOW	Date Received: 11/08/91	
* Moisture: not dec. 8	Date Analyzed: 11/13/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chlороethane	11	U
75-09-2	Methylene Chloride	31	B
67-64-1	Acetone	23	B
75-15-0	Carbon Disulfide	5	U
75-35-4	1, 1-Dichloroethene	5	U
75-34-3	1, 1-Dichloroethane	5	U
540-59-0	1, 2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1, 2-Dichloroethane	5	U
78-93-3	2-Butanone	11	U
71-55-6	1, 1, 1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1, 2-Dichloropropane	5	U
10061-01-5	cis-1, 3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1, 1, 2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1, 3-Dichloropropene	5	U
110-75-8	2-Chloroethylvinylether	11	U
75-25-2	Bromoform	11	U
108-10-1	4-Methyl-2-Pantanone	16	U
591-78-6	2-Hexanone	16	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	11	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	U
74-88-4	Iodomethane	11	U

FORM I VOL

1/87 Rev.

107-02-8-----	Acrolein	98	U
107-13-1-----	Acrylonitrile	130	U
75-69-4-----	Trichlorofluoromethane	5	U
107-05-1-----	3-Chloropropene	16	U
76-13-1-----	1,1,2-Trichloro-1,2,2-trifluo	11	U
354-58-5-----	1,1,1-Trichloro-2,2,2-trifluo	11	U
74-95-3-----	Dibromomethane	11	U
4170-30-3-----	Crotonaldehyde	110	U
106-93-4-----	1,2-Dibromoethane	5	U
630-20-6-----	1,1,1,2-Tetrachloroethane	5	U
764-41-0-----	cis-1,4-Dichloro-2-butene	16	U
96-18-4-----	1,2,3-Trichloropropane	16	U
764-71-0-----	trans-1,4-Dichloro-2-butene	16	U
96-18-4-----	Ethylmethacrylate	11	U
96-12-8-----	1,2-Dibromo-3-chloropropane	11	U
97-63-2-----	Ethylmethacrylate	11	U

R0A010406

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA011214
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>26</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>462131</u>	
Sample wt/vol: <u>5.0</u> (g/mL) G	Lab File ID: <u>GH062131C13</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/08/91</u>	
# Moisture: not dec. <u>9</u>	Date Analyzed: <u>11/13/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG		Q
74-87-3	Chloromethane	11	U	
74-83-9	Bromomethane	5	U	
75-01-4	Vinyl Chloride	11	U	
75-00-3	Chloroethane	11	U	
75-09-2	Methylene Chloride	30	B	
67-64-1	Acetone	12	B	
75-15-0	Carbon Disulfide	5	U	
75-35-4	1,1-Dichloroethene	5	U	
75-34-3	1,1-Dichloroethane	5	U	
540-59-0	1,2-Dichloroethene (total)	5	U	
67-66-3	Chloroform	5	U	
107-06-2	1,2-Dichloroethane	5	U	
78-93-3	2-Butanone	11	U	
71-55-6	1,1,1-Trichloroethane	5	U	
56-23-5	Carbon Tetrachloride	5	U	
108-05-4	Vinyl Acetate	11	U	
75-27-4	Bromodichloromethane	5	U	
78-87-5	1,2-Dichloropropane	5	U	
10061-01-5	cis-1,3-Dichloropropene	5	U	
79-01-6	Trichloroethene	5	U	
124-48-1	Dibromochloromethane	5	U	
79-00-5	1,1,2-Trichloroethane	5	U	
71-43-2	Benzene	5	U	
10061-02-6	Trans-1,3-Dichloropropene	5	U	
110-75-8	2-Chloroethylvinylether	11	U	
75-25-2	Bromoform	11	U	
108-10-1	4-Methyl-2-Pentanone	16	U	
591-78-6	2-Hexanone	16	U	
127-18-4	Tetrachloroethene	5	U	
79-34-5	1,1,2,2-Tetrachloroethane	11	U	
108-88-3	Toluene	5	U	
108-90-7	Chlorobenzene	5	U	
100-41-4	Ethylbenzene	5	U	
100-42-5	Styrene	5	U	
1330-20-7	Total Xylenes	5	U	
74-88-4	Iodomethane	11	U	

FORM I VOA

1/87 Rev.

107-02-8	Acrolein	99	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	5	U
107-05-1	3-Chloropropene	16	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoropropane	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluoroethane	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	5	U
630-20-6	1,1,1,2-Tetrachloroethane	5	U
764-41-0	cis-1,4-Dichloro-2-butene	16	U
96-18-4	1,2,3-Trichloropropane	16	U
764-71-0	trans-1,4-Dichloro-2-butene	16	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Ethylmethacrylate	11	U

R04011214

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP

Contract: 500077

ROA011416

Lab Code: COMPU Case No.: 24105 SAS No.: SDG No.: 26

Matrix: (soil/water) SOIL Lab Sample ID: 462132

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH062132C13

Level: (low/med) LOW Date Received: 11/08/91

* Moisture: not dec. 7 Date Analyzed: 11/13/91

Column: (pack/cap) CAP Dilution Factor: 1.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/kg

CAS NO.	COMPOUND	Q
74-87-3	Chloromethane	11 U
74-83-9	Bromomethane	5 U
75-01-4	Vinyl Chloride	11 U
75-00-3	Chloroethane	11 U
75-09-2	Methylene Chloride	27 B
67-64-1	Acetone	12 B
75-15-0	Carbon Disulfide	5 U
75-35-4	1,1-Dichloroethene	5 U
75-34-3	1,1-Dichloroethane	5 U
540-59-0	1,2-Dichloroethene (total)	5 U
67-66-3	Chloroform	5 U
107-06-2	1,2-Dichloroethane	5 U
78-93-3	2-Butanone	11 U
71-55-6	1,1,1-Trichloroethane	5 U
56-23-5	Carbon Tetrachloride	5 U
108-05-4	Vinyl Acetate	11 U
75-27-4	Bromodichloromethane	5 U
78-87-5	1,2-Dichloropropane	5 U
10061-01-5	cis-1,3-Dichloropropene	5 U
79-01-6	Trichloroethene	5 U
124-48-1	Dibromoethylchloromethane	5 U
79-00-5	1,1,2-Trichloroethane	5 U
71-43-2	Benzene	5 U
10061-02-6	Trans-1,3-Dichloropropene	5 U
110-75-8	2-Chloroethylvinylether	11 U
75-25-2	Bromoform	11 U
108-10-1	4-Methyl-2-Pentanone	16 U
591-78-6	2-Hexanone	16 U
127-18-4	Tetrachloroethene	5 U
79-34-5	1,1,2,2-Tetrachloroethane	11 U
108-88-3	Toluene	5 U
108-90-7	Chlorobenzene	5 U
100-41-4	Ethylbenzene	5 U
100-42-5	Styrene	5 U
1330-20-7	Total Xylenes	5 U
74-88-4	Iodomethane	11 U

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107-02-8	Acrolein	97	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	5	U
107-05-1	3-Chloropropene	16	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	5	U
630-20-6	1,1,1,2-Tetrachloroethane	5	U
764-41-0	cis-1,4-Dichloro-2-butene	16	U
96-18-4	1,2,3-Trichloropropane	16	U
764-71-0	trans-1,4-Dichloro-2-butene	16	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Ethylmethacrylate	11	U

R0A011416

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROAO12022

Lab Name: COMPUCHEM RTP Contract: 500077
 Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 26
 Matrix: (soil/water) SOIL Lab Sample ID: 462133
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH062133A13
 Level: (low/med) LOW Date Received: 11/08/91
 * Moisture: not dec. .24 Date Analyzed: 11/13/91
 Column: (pack/cap) GAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
74-87-3	Chloromethane	13	U	
74-83-9	Bromomethane	7	U	
75-01-4	Vinyl Chloride	13	U	
75-00-3	Chloroethane	13	U	
75-09-2	Methylene Chloride	64	B	
67-64-1	Acetone	63	B	
75-15-0	Carbon Disulfide	7	U	
75-35-4	1,1-Dichloroethene	7	U	
75-34-3	1,1-Dichloroethane	7	U	
540-59-0	1,2-Dichloroethene (total)	7	U	
67-66-3	Chloroform	7	U	
107-06-2	1,2-Dichloroethane	7	U	
78-93-3	2-Butanone	7	J	
71-55-6	1,1,1-Trichloroethane	7	U	
56-23-5	Carbon Tetrachloride	7	U	
108-05-4	Vinyl Acetate	13	U	
75-27-4	Bromodichloromethane	7	U	
78-87-5	1,2-Dichloropropane	7	U	
10061-01-5	cis-1,3-Dichloropropene	7	U	
79-01-6	Trichloroethene	7	U	
124-48-1	Dibromochloromethane	7	U	
79-00-5	1,1,2-Trichloroethane	7	U	
71-43-2	Benzene	7	U	
10061-02-6	Trans-1,3-Dichloropropene	7	U	
110-75-8	2-Chloroethylvinylether	13	U	
75-25-2	Bromoform	13	U	
108-10-1	4-Methyl-2-Pentanone	20	U	
591-78-6	2-Hexanone	20	U	
127-18-4	Tetrachloroethene	7	U	
79-34-5	1,1,2,2-Tetrachloroethane	13	U	
108-88-3	Toluene	7	U	
108-90-7	Chlorobenzene	7	U	
100-41-4	Ethylbenzene	19		
100-42-5	Styrene	7	U	
1330-20-7	Total Xylenes	13		
74-88-4	Iodomethane	13	U	

107-02-8	Acrolein	120	U
107-13-1	Acrylonitrile	160	U
75-69-4	Trichlorofluoromethane	7	U
107-05-1	3-Chloropropene	20	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	13	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	13	U
74-95-3	Dibromomethane	13	U
4170-30-3	Crotonaldehyde	130	U
106-93-4	1,2-Dibromoethane	7	U
630-20-6	1,1,1,2-Tetrachloroethane	7	U
764-41-0	cis-1,4-Dichloro-2-butene	20	U
96-18-4	1,2,3-Trichloropropane	20	U
764-71-0	trans-1,4-Dichloro-2-butene	20	U
96-18-4	Ethylmethacrylate	13	U
96-12-8	1,2-Dibromo-3-chloropropane	13	U
97-63-2	Ethylmethacrylate	13	U

462133

FORM I VOA

1/87 Rev.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500077	ROA012224
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 26
Matrix: (soil/water) SOIL		Lab Sample ID: 462135
Sample wt/vol: 5.0 (g/mL) G		Lab File ID: GH062135C11
Level: (low/med) LOW		Date Received: 11/08/91
* Moisture: not dec. 10		Date Analyzed: 11/13/91
Column: (pack/cap) CAP		Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	Q
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	6	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chloroethane	11	U
75-09-2	Methylene Chloride	4.4	B
67-64-1	Acetone	1.6	B
75-15-0	Carbon Disulfide	6	U
75-35-4	1,1-Dichloroethene	6	U
75-34-3	1,1-Dichloroethane	6	U
540-59-0	1,2-Dichloroethene (total)	6	U
67-66-3	Chloroform	6	U
107-06-2	1,2-Dichloroethane	6	U
78-93-3	2-Butanone	11	U
71-55-6	1,1,1-Trichloroethane	6	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	6	U
78-87-5	1,2-Dichloropropane	6	U
10061-01-5	cis-1,3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1,1,2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1,3-Dichloropropene	6	U
110-75-8	2-Chloroethylvinylether	11	U
75-25-2	Bromoform	11	U
108-10-1	4-Methyl-2-Pentanone	17	U
591-78-6	2-Hexanone	17	U
127-18-4	Tetrachloroethene	6	U
79-34-5	1,1,2,2-Tetrachloroethane	11	U
108-88-3	Toluene	6	U
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	6	U
74-88-4	Iodomethane	11	U

FORM I VOA

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107-02-8	Acrolein	100	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	6	U
107-05-1	3-Chloropropene	17	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	6	U
630-20-6	1,1,1,2-Tetrachloroethane	6	U
764-41-0	cis-1,4-Dichloro-2-butene	17	U
96-18-4	1,2,3-Trichloropropane	17	U
764-71-0	trans-1,4-Dichloro-2-butene	17	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Ethylmethacrylate	11	U

R0A012324

LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500977	ROA-DFV1
Lab Code: COMPU	Case No.: 24105	SAS No.: 26
Matrix: (soil/water) SOIL		Lab Sample ID: 462157
Sample wt/vol: 5.0 (g/mL) G		Lab File ID: GH062157C13
Level: (low/med) LOW		Date Received: 11/08/91
* Moisture: not dec. 9		Date Analyzed: 11/13/91
Column: (pack/cap) CAP		Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chloroethane	11	U
75-09-2	Methylene Chloride	37	B
67-64-1	Acetone	17	B
75-15-0	Carbon Disulfide	5	U
75-35-4	1,1-Dichloroethene	5	U
75-34-3	1,1-Dichloroethane	5	U
540-59-0	1,2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1,2-Dichloroethane	5	U
78-93-3	2-Butanone	11	U
71-55-6	1,1,1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1,2-Dichloropropane	5	U
10061-01-5	cis-1,3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromoethylmethane	5	U
79-00-5	1,1,2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1,3-Dichloropropene	5	U
110-75-8	2-Chloroethylvinylether	11	U
75-25-2	Bromoform	11	U
108-10-1	4-Methyl-2-Pentanone	16	U
591-78-6	2-Hexanone	16	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1,1,2,2-Tetrachloroethane	11	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	U
74-88-4	Iodomethane	11	U

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	Acrolein	99	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	5	U
107-05-1	3-Chloropropene	16	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoropropane	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluoropropane	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	5	U
630-20-6	1,1,1,2-Tetrachloroethane	5	U
764-41-0	cis-1,4-Dichloro-2-butene	16	U
96-18-4	1,2,3-Trichloropropane	16	U
764-71-0	trans-1,4-Dichloro-2-butene	16	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Ethylmethacrylate	11	U

R0A-DPVI

FORM I VOA

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	<u>ROA2B0608</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>87</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465047</u>	
Sample wt/vol: <u>5.0</u> (g/mL) G	Lab File ID: <u>GH065047B13</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/21/91</u>	
% Moisture: not dec. <u>9</u>	Date Analyzed: <u>11/22/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chloroethane	11	U
75-09-2	Methylene Chloride	34	B
67-64-1	Acetone	17	B
75-15-0	Carbon Disulfide	5	U
75-35-4	1, 1-Dichloroethene	5	U
75-34-3	1, 1-Dichloroethane	5	U
540-59-0	1, 2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1, 2-Dichloroethane	5	U
78-93-3	2-Butanone	11	U
71-55-6	1, 1, 1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1, 2-Dichloropropane	5	U
10061-01-5	cis-1, 3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1, 1, 2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1, 3-Dichloropropene	5	U
110-75-8	2-Chloroethylvinylether	11	U
75-25-2	Bromoform	11	U
108-10-1	4-Methyl-2-Pentanone	16	U
591-78-6	2-Hexanone	16	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	11	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	U
74-88-4	Iodomethane	11	U

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107-02-8	Acrolein	99	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	5	U
107-05-1	3-Chloropropene	16	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	5	U
630-20-6	1,1,1,2-Tetrachloroethane	5	U
764-41-0	cis-1,4-Dichloro-2-butene	16	U
96-18-4	1,2,3-Trichloropropane	16	U
764-71-0	trans-1,4-Dichloro-2-butene	16	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Ethylmethacrylate	11	U

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA3B1214
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>577</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>473667</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>G</u>	Lab File ID: <u>GH073667B13</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>01/07/92</u>	
% Moisture: not dec. <u>9</u>	Date Analyzed: <u>01/10/92</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chloroethane	11	U
75-09-2	Methylene Chloride	26	B
67-64-1	Acetone	26	B
75-15-0	Carbon Disulfide	5	U
75-35-4	1,1-Dichloroethene	5	U
75-34-3	1,1-Dichloroethane	5	U
540-59-0	1,2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1,2-Dichloroethane	5	U
78-93-3	2-Butanone	11	U
71-55-6	1,1,1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1,2-Dichloropropane	5	U
10061-01-5	cis-1,3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1,1,2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1,3-Dichloropropene	5	U
110-75-8	2-Chloroethylvinylether	11	U
75-25-2	Bromoform	11	U
108-10-1	4-Methyl-2-Pentanone	16	U
591-78-6	2-Hexanone	16	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1,1,2,2-Tetrachloroethane	11	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	J
74-88-4	Iodomethane	11	U

FORM I VOA

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107-02-8	Acrolein	99	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	5	U
107-05-1	3-Chloropropene	16	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoropropane	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluoropropane	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	5	U
630-20-6	1,1,1,2-Tetrachloroethane	5	U
764-41-0	cis-1,4-Dichloro-2-butene	16	U
96-18-4	1,2,3-Trichloropropane	16	U
764-71-0	trans-1,4-Dichloro-2-butene	16	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Eethylmethacrylate	11	U

R0A3B1214

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUTECH, RTP	Contract: 500077	ROB1B0406
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 87
Matrix: (soil/water) SOIL	Lab Sample ID: 465524	
Sample wt/vol: 5.0 (g/mL) S	Lab File ID: GH065524B13	
Level: (low/med) LOW	Date Received: 11/22/91	
% Moisture: not dec. 17	Date Analyzed: 11/25/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
74-87-3	Chloromethane	12	U
74-83-9	Bromomethane	12	U
75-01-4	Vinyl Chloride	12	U
75-00-3	Chloroethane	12	U
75-09-2	Methylene Chloride	32	B
67-64-1	Acetone	23	B
75-15-0	Carbon Disulfide	12	U
75-35-4	1,1-Dichloroethene	12	U
75-34-3	1,1-Dichloroethane	12	U
540-59-0	1,2-Dichloroethene (total)	12	U
67-66-3	Chloroform	12	U
107-06-2	1,2-Dichloroethane	12	U
78-93-3	2-Butanone	12	U
71-55-6	1,1,1-Trichloroethane	12	U
56-23-5	Carbon Tetrachloride	12	U
75-27-4	Bromodichloromethane	12	U
78-87-5	1,2-Dichloropropane	12	U
10061-01-5	cis-1,3-Dichloropropene	12	U
79-01-6	Trichloroethene	12	U
124-48-1	Dibromochloromethane	12	U
79-00-5	1,1,2-Trichloroethane	12	U
71-43-2	Benzene	12	U
10061-02-6	Trans-1,3-Dichloropropene	12	U
75-25-2	Bromoform	12	U
108-10-1	4-Methyl-2-Pentanone	12	U
591-78-6	2-Hexanone	12	U
127-18-4	Tetrachloroethene	12	U
79-34-5	1,1,2,2-Tetrachloroethane	12	U
108-88-3	Toluene	4	J
108-90-7	Chlorobenzene	12	U
100-41-4	Ethylbenzene	12	U
100-42-5	Styrene	12	U
1330-20-7	Xylene (total)	12	U
74-88-4	Iodomethane	120	U
107-02-8	Acrolein	110	U
107-13-1	Acrylonitrile	140	U

FORM I VOA

1/87 Rev.

75-69-4-----	Trichlorofluoromethane	6	U
107-05-1-----	3-Chloropropene	6	U
76-13-1-----	1,1,2-Trichloro-1,2,2-trifluo	6	U
354-58-5-----	1,1,1-Trichloro-2,2,2-trifluo	6	U
74-95-3-----	Dibromomethane	6	U
4170-30-3-----	Crotonaldehyde	6	U
106-93-4-----	1,2-Dibromoethane	6	U
630-20-6-----	1,1,1,2-Tetrachloroethane	6	U
764-71-0-----	cis-1,4-Dichloro-2-butene	6	U
96-18-4-----	1,2,3-Trichloropropane	12	U
764-71-0-----	trans-1,4-Dichloro-2-butene	6	U
96-12-8-----	1,2-Dibromo-3-chloropropane	6	U

ROB1B0406

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	<u>ROB2B0002</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>87</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465877</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>G</u>	Lab File ID: <u>GH065877B12</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/23/91</u>	
# Moisture: not dec. <u>3.1</u>	Date Analyzed: <u>11/25/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	Chloromethane	15	U
74-83-9	Bromomethane	15	U
75-01-4	Vinyl Chloride	15	U
75-00-3	Chloroethane	15	U
75-09-2	Methylene Chloride	100	B
67-64-1	Acetone	46	B
75-15-0	Carbon Disulfide	15	U
75-35-4	1,1-Dichloroethene	15	U
75-34-3	1,1-Dichloroethane	15	U
540-59-0	1,2-Dichloroethene (total)	15	U
67-66-3	Chloroform	15	U
107-06-2	1,2-Dichloroethane	15	U
78-93-3	2-Butanone	15	U
71-55-6	1,1,1-Trichloroethane	15	U
56-23-5	Carbon Tetrachloride	15	U
75-27-4	Bromodichloromethane	15	U
78-87-5	1,2-Dichloropropane	15	U
10061-01-5	cis-1,3-Dichloropropene	15	U
79-01-6	Trichloroethene	15	U
124-48-1	Dibromochloromethane	15	U
79-00-5	1,1,2-Trichloroethane	15	U
71-43-2	Benzene	15	U
10061-02-6	Trans-1,3-Dichloropropene	15	U
75-25-2	Bromoform	15	U
108-10-1	4-Methyl-2-Pentanone	15	U
591-78-6	2-Hexanone	15	U
127-18-4	Tetrachloroethene	15	U
79-34-5	1,1,2,2-Tetrachloroethane	15	U
108-88-3	Toluene	7	J
108-90-7	Chlorobenzene	15	U
100-41-4	Ethylbenzene	5	J
100-42-5	Styrene	15	U
1330-20-7	Xylene (total)	12	J
74-88-4	Iodomethane	150	U
107-02-8	Acrolein	130	U
107-13-1	Acrylonitrile	180	U

FORM I VOA

1/87 Rev.

75-69-4	Trichlorofluoromethane	7	U
107-05-1	3-Chloropropene	7	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	7	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	7	U
74-95-3	Dibromomethane	7	U
4170-30-3	Crotonaldehyde	7	U
106-93-4	1,2-Dibromoethane	7	U
630-20-6	1,1,1,2-Tetrachloroethane	7	U
764-71-0	cis-1,4-Dichloro-2-butene	7	U
96-18-4	1,2,3-Trichloropropane	15	U
764-71-0	trans-1,4-Dichloro-2-butene	7	U
96-12-8	1,2-Dibromo-3-chloropropane	7	U

R0B2B0002

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500077	ROB-DPA1
Lab Code: COMEU	Case No.: 24105	SAS No.: 87
Matrix: (soil/water) SOIL	Lab Sample ID: 465549	
Sample wt/vol: 5.0 (g/mL) G	Lab File ID: GH065549B13	
Level: (low/med) LOW	Date Received: 11/22/91	
* Moisture: not dec. 17	Date Analyzed: 11/25/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	Chloromethane	12	U
74-83-9	Bromomethane	12	U
75-01-4	Vinyl Chloride	12	U
75-00-3	Chloroethane	12	U
75-09-2	Methylene Chloride	51	B
67-64-1	Acetone	24	B
75-15-0	Carbon Disulfide	12	U
75-35-4	1, 1-Dichloroethene	12	U
75-34-3	1, 1-Dichloroethane	12	U
540-59-0	1, 2-Dichloroethene (total)	12	U
67-66-3	Chloroform	12	U
107-06-2	1, 2-Dichloroethane	12	U
78-93-3	2-Butanone	12	U
71-55-6	1, 1, 1-Trichloroethane	12	U
56-23-5	Carbon Tetrachloride	12	U
75-27-4	Bromodichloromethane	12	U
78-87-5	1, 2-Dichloropropane	12	U
10061-01-5	cis-1, 3-Dichloropropene	12	U
79-01-6	Trichloroethene	12	U
124-48-1	Dibromochloromethane	12	U
79-00-5	1, 1, 2-Trichloroethane	12	U
71-43-2	Benzene	12	U
10061-02-6	Trans-1, 3-Dichloropropene	12	U
75-25-2	Bromoform	12	U
108-10-1	4-Methyl-2-Pentanone	12	U
591-78-6	2-Hexanone	12	U
127-18-4	Tetrachloroethene	12	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	12	U
108-88-3	Toluene	2	J
108-90-7	Chlorobenzene	12	U
100-41-4	Ethylbenzene	12	U
100-42-5	Styrene	12	U
1330-20-7	Xylene (total)	12	U
74-88-4	Iodomethane	120	U
107-02-8	Acrolein	110	U
107-13-1	Acrylonitrile	140	U

FORM I VOL

1/87 Rev.

75-69-4-----	Trichlorofluoromethane	6	U
107-05-1-----	3-Chloropropene	6	U
76-13-1-----	1,1,2-Trichloro-1,2,2-trifluor	6	U
354-58-5-----	1,1,1-Trichloro-2,2,2-trifluor	6	U
74-95-3-----	Dibromomethane	6	U
4170-30-3-----	Crotonaldehyde	6	U
106-93-4-----	1,2-Dibromoethane	6	U
630-20-6-----	1,1,1,2-Tetrachloroethane	6	U
764-71-0-----	cis-1,4-Dichloro-2-butene	6	U
96-18-4-----	1,2,3-Trichloropropane	12	U
764-71-0-----	trans-1,4-Dichloro-2-butene	6	U
96-12-8-----	1,2-Dibromo-3-chloropropane	6	U

ROB-DPAI

FORM I VOA

1/87 Rev.

LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500077	ROCO11012
Lab Code: CQMPUL	Case No.: 24105	SAS No.: _____ SDG No.: 26
Matrix: (soil/water) SOIL	Lab Sample ID: 461431	
Sample wt/vol: 5.0 (g/mL) G	Lab File ID: GH061431A13	
Level: (low/med) LOW	Date Received: 11/07/91	
* Moisture: not dec. 10	Date Analyzed: 11/14/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	6	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chloroethane	11	U
75-09-2	Methylene Chloride	28	B
67-64-1	Acetone	36	B
75-15-0	Carbon Disulfide	6	U
75-35-4	1, 1-Dichloroethene	6	U
75-34-3	1, 1-Dichloroethane	6	U
540-59-0	1, 2-Dichloroethene (total)	6	U
67-66-3	Chloroform	6	U
107-06-2	1, 2-Dichloroethane	6	U
78-93-3	2-Butanone	11	U
71-55-6	1, 1, 1-Trichloroethane	6	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	6	U
78-87-5	1, 2-Dichloropropane	6	U
10061-01-5	cis-1, 3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1, 1, 2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1, 3-Dichloropropene	6	U
110-75-8	2-Chloroethylvinylether	11	U
75-25-2	Bromoform	11	U
108-10-1	4-Methyl-2-Pentanone	17	U
591-78-6	2-Hexanone	17	U
127-18-4	Tetrachloroethene	6	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	11	U
108-88-3	Toluene	6	U
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	6	U
74-88-4	Iodomethane	11	U

FORM I VOA

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107-05-1	Acrolein	100	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	6	U
107-05-1	3-Chloropropene	17	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	11	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	6	U
630-20-6	1,1,1,2-Tetrachloroethane	6	U
764-41-0	cis-1,4-Dichloro-2-butene	17	U
96-18-4	1,2,3-Trichloropropane	17	U
764-71-0	trans-1,4-Dichloro-2-butene	17	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U
97-63-2	Ethylmethacrylate	11	U

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LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP Contract: 500077 ROC021214
 Lab Code: COMPU Case No.: 24105 SAS No.: SDG No.: 26
 Matrix: (soil/water) SOIL Lab Sample ID: 461148
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GR061148B13
 Level: (low/med) LOW Date Received: 11/06/91
 % Moisture: not dec. 18 Date Analyzed: 11/12/91
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
74-87-3	Chloromethane	12	U
74-83-9	Bromomethane	6	U
75-01-4	Vinyl Chloride	12	U
75-00-3	Chloroethane	12	U
75-09-2	Methylene Chloride	12	U
67-64-1	Acetone	58	B
75-15-0	Carbon Disulfide	48	B
75-35-4	1, 1-Dichloroethene	6	U
75-34-3	1, 1-Dichloroethane	6	U
540-59-0	1, 2-Dichloroethene (total)	6	U
67-66-3	Chloroform	6	U
107-06-2	1, 2-Dichloroethane	6	U
78-93-3	2-Butanone	6	U
71-55-6	1, 1, 1-Trichloroethane	12	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	6	U
75-27-4	Bromodichloromethane	12	U
78-87-5	1, 2-Dichloropropane	6	U
10061-01-5	cis-1, 3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1, 1, 2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1, 3-Dichloropropene	6	U
110-75-8	2-Chloroethylvinylether	6	U
75-25-2	Bromoform	12	U
108-10-1	4-Methyl-2-Pentanone	12	U
591-78-6	2-Hexanone	18	U
127-18-4	Tetrachloroethene	18	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	6	U
108-88-3	Toluene	12	U
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	6	U
74-88-4	Iodomethane	12	U

FORM I VOL

1/87 Rev.

	-Acrylene	110	U
107-13-1	-Acrylonitrile	150	U
75-69-4	-Trichlorofluoromethane	6	U
107-05-1	-3-Chloropropene	18	U
76-13-1	-1,1,2-Trichloro-1,2,2-trifluo	12	U
354-58-5	-1,1,1-Trichloro-2,2,2-trifluo	12	U
74-95-3	-Dibromomethane	12	U
4170-30-3	-Crotonaldehyde	120	U
106-93-4	-1,2-Dibromoethane	6	U
630-20-6	-1,1,1,2-Tetrachloroethane	6	U
764-41-0	-cis-1,4-Dichloro-2-butene	18	U
96-18-4	-1,2,3-Trichloropropane	18	U
764-71-0	-trans-1,4-Dichloro-2-butene	18	U
96-18-4	-Ethylmethacrylate	12	U
96-12-8	-1,2-Dibromo-3-chloropropane	12	U
97-63-2	-Ethylmethacrylate	12	U

R0C02/2/4

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROC3B0204
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>87</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465069</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>G</u>	Lab File ID: <u>GH065069B13</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/21/91</u>	
* Moisture: not dec. <u>8</u>	Date Analyzed: <u>11/22/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	-Chloromethane	11	U
74-83-9	-Bromomethane	5	U
75-01-4	-Vinyl Chloride	11	U
75-00-3	-Chloroethane	11	U
75-09-2	-Methylene Chloride	34	B
67-64-1	-Acetone	14	B
75-15-0	-Carbon Disulfide	5	U
75-35-4	-1,1-Dichloroethene	5	U
75-34-3	-1,1-Dichloroethane	5	U
540-59-0	-1,2-Dichloroethene (total)	5	U
67-66-3	-Chloroform	5	U
107-06-2	-1,2-Dichloroethane	5	U
78-93-3	-2-Butanone	11	U
71-55-6	-1,1,1-Trichloroethane	5	U
56-23-5	-Carbon Tetrachloride	5	U
108-05-4	-Vinyl Acetate	11	U
75-27-4	-Bromodichloromethane	5	U
78-87-5	-1,2-Dichloropropane	5	U
10061-01-5	-cis-1,3-Dichloropropene	5	U
79-01-6	-Trichloroethene	5	U
124-48-1	-Dibromochloromethane	5	U
79-00-5	-1,1,2-Trichloroethane	5	U
71-43-2	-Benzene	5	U
10061-02-6	-Trans-1,3-Dichloropropene	5	U
110-75-8	-2-Chloroethylvinylether	11	U
75-25-2	-Bromoform	11	U
108-10-1	-4-Methyl-2-Pentanone	16	U
591-78-6	-2-Hexanone	16	U
127-18-4	-Tetrachloroethene	5	U
79-34-5	-1,1,2,2-Tetrachloroethane	11	U
108-88-3	-Toluene	2	J
108-90-7	-Chlorobenzene	5	U
100-41-4	-Ethylbenzene	5	U
100-42-5	-Styrene	5	U
1330-20-7	-Total Xylenes	5	U
74-88-4	-Iodomethane	11	U

FORM I VOA

1/87 Rev.

107-02-8-----	Acrolein	98	U
107-13-1-----	Acrylonitrile	130	U
75-69-4-----	Trichlorofluoromethane	5	U
107-05-1-----	3-Chloropropene	16	U
76-13-1-----	1,1,2-Trichloro-1,2,2-trifluoropropane	11	U
354-58-5-----	1,1,1-Trichloro-2,2,2-trifluoropropane	11	U
74-95-3-----	Dibromomethane	11	U
4170-30-3-----	Crotonaldehyde	110	U
106-93-4-----	1,2-Dibromoethane	5	U
630-20-6-----	1,1,1,2-Tetrachloroethane	5	U
764-41-0-----	cis-1,4-Dichloro-2-butene	16	U
96-18-4-----	1,2,3-Trichloropropane	16	U
764-71-0-----	trans-1,4-Dichloro-2-butene	16	U
96-18-4-----	Ethylmethacrylate	11	U
96-12-8-----	1,2-Dibromo-3-chloropropane	11	U
97-63-2-----	Ethylmethacrylate	11	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPOCHEM RTP	Contract: 500027	ROC-DPAL
Lab Code: COMPEL	Case No.: 24105	SAS No.: SDG No.: 26
Matrix: (soil/water) SOIL	Lab Sample ID: 461185	
Sample wt/vol: 5.0 (g/mL) G	Lab File ID: GHQ61185B13	
Level: (low/med) LOW	Date Received: 11/06/91	
* Moisture: not dec. 46	Date Analyzed: 11/08/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
74-87-3	Chloromethane	19	U
74-83-9	Bromomethane	9	U
75-01-4	Vinyl Chloride	19	U
75-00-3	Chloroethane	19	U
75-09-2	Methylene Chloride	45	B
67-64-1	Acetone	44	
75-15-0	Carbon Disulfide	9	U
75-35-4	1, 1-Dichloroethene	9	U
75-34-3	1, 1-Dichloroethane	9	U
540-59-0	1, 2-Dichloroethene (total)	9	U
67-66-3	Chloroform	9	U
107-06-2	1, 2-Dichloroethane	9	U
78-93-3	2-Butanone	19	U
71-55-6	1, 1, 1-Trichloroethane	9	U
56-23-5	Carbon Tetrachloride	9	U
108-05-4	Vinyl Acetate	19	U
75-27-4	Bromodichloromethane	9	U
78-87-5	1, 2-Dichloropropane	9	U
10061-01-5	cis-1, 3-Dichloropropene	9	U
79-01-6	Trichloroethene	9	U
124-48-1	Dibromochloromethane	9	U
79-00-9	1, 1, 2-Trichloroethane	9	U
71-43-2	Benzene	9	U
10061-02-6	Trans-1, 3-Dichloropropene	9	U
110-75-8	2-Chloroethylvinylether	19	U
75-25-2	Bromoform	19	U
108-10-1	4-Methyl-2-Pentanone	28	U
591-78-6	2-Hexanone	28	U
127-18-4	Tetrachloroethene	9	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	19	U
108-68-3	Toluene	9	U
108-90-7	Chlorobenzene	9	U
100-41-4	Ethylbenzene	9	U
100-42-5	Styrene	9	U
1330-20-7	Total Xylenes	9	U
74-88-4	Iodomethane	19	U

FORM I VOL

1/87 Rev.

107-02-6	Acrolein	170	U
107-13-1	Acrylonitrile	220	U
75-69-4	Trichlorofluoromethane	9	U
107-05-1	3-Chloropropene	28	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoropropane	19	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluoropropane	19	U
74-95-3	Dibromomethane	19	U
4170-30-3	Crotonaldehyde	190	U
106-93-4	1,2-Dibromoethane	9	U
630-20-6	1,1,1,2-Tetrachloroethane	9	U
764-41-0	cis-1,4-Dichloro-2-butene	28	U
96-18-4	1,2,3-Trichloropropane	28	U
764-71-0	trans-1,4-Dichloro-2-butene	28	U
96-18-4	Ethylmethacrylate	19	U
96-12-8	1,2-Dibromo-3-chloropropane	19	U
97-63-2	Ethylmethacrylate	19	U

461185

FORM I VOA

1/87 Rev.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPLICHEM R&P	Contract: 500077	ROJ1S
Lab Code: QOMPU	Case No.: 24105	SAC No.: 424
Matrix: (soil/water) SOIL	Lab Sample ID: 469282	
Sample wt/vol: 5.0 (g/mL) G	Lab File ID: GR069282H13	
Level: (low/med) LOW	Date Received: 12/11/91	
# Moisture: not dec. 2.2	Date Analyzed: 12/16/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	Q
74-87-3	Chloromethane	13	U
74-83-9	Bromomethane	6	U
75-01-4	Vinyl Chloride	13	U
75-00-3	Chloroethane	13	U
75-09-2	Methylene Chloride	56	B
67-64-1	Acetone	23	
75-15-0	Carbon Disulfide	6	U
75-35-4	1,1-Dichloroethene	6	U
75-34-3	1,1-Dichloroethane	6	U
540-59-0	1,2-Dichloroethene (total)	6	U
67-66-3	Chloroform	6	U
107-06-2	1,2-Dichloroethane	6	U
78-93-3	2-Butanone	6	U
71-55-6	1,1,1-Trichloroethane	13	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	6	U
75-27-4	Bromodichloromethane	13	U
78-87-5	1,2-Dichloropropane	6	U
10061-01-5	cis-1,3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1,1,2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1,3-Dichloropropene	6	U
110-75-8	2-Chloroethylvinylether	13	U
75-25-2	Bromoform	13	U
108-10-1	4-Methyl-2-Pentanone	19	U
591-78-6	2-Hexanone	19	U
127-18-4	Tetrachloroethene	6	U
79-34-5	1,1,2,2-Tetrachloroethane	13	U
108-88-3	Toluene	6	U
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	6	U
74-88-4	Iodomethane	13	U

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107-02-8	Acrolein		120	U
107-13-1	Acrylonitrile		150	U
75-69-4	Trichlorofluoromethane		6	U
107-05-1	3-Chloropropene		19	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo		13	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo		13	U
74-95-3	Dibromomethane		13	U
4170-30-3	Crotonaldehyde		13	U
106-93-4	1,2-Dibromoethane		130	U
630-20-6	1,1,1,2-Tetrachloroethane		6	U
764-41-0	cis-1,4-Dichloro-2-butene		6	U
96-18-4	1,2,3-Trichloropropane		19	U
764-71-0	trans-1,4-Dichloro-2-butene		19	U
96-18-4	Ethylmethacrylate		19	U
96-12-8	1,2-Dibromo-3-chloropropene		13	U
			13	U

RDS 15

FORM I VOA

1/87 Rev.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM INC</u>	Contract: <u>500077</u>	<u>ROJ2S</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>424</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>469273</u>	
Sample wt/vol: <u>5.0</u> (g/mL) G	Lab File ID: <u>GH069273BL13</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/11/91</u>	
* Moisture: not dec. <u>15</u>	Date Analyzed: <u>12/13/91</u>	
Column: (pack/cap) <u>GAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
74-87-3	Chloromethane	12	U
74-83-9	Bromomethane	6	U
75-01-4	Vinyl Chloride	12	U
75-00-3	Chloroethane	12	U
75-09-2	Methylene Chloride	74	B
67-64-1	Acetone	39	
75-15-0	Carbon Disulfide	6	U
75-35-4	1, 1-Dichloroethene	6	U
75-34-3	1, 1-Dichloroethane	6	U
540-59-0	1, 2-Dichloroethene (total)	6	U
67-66-3	Chloform	6	U
107-06-2	1, 2-Dichloroethane	6	U
78-93-3	2-Butanone	12	U
71-55-6	1, 1, 1-Trichloroethane	6	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	12	U
75-27-4	Bromodichloromethane	6	U
78-87-5	1, 2-Dichloropropane	6	U
10061-01-5	cis-1, 3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1, 1, 2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1, 3-Dichloropropene	6	U
110-75-8	2-Chloroethylvinylether	12	U
75-25-2	Bromoform	12	U
108-10-1	4-Methyl-2-Pentanone	18	U
591-78-6	2-Hexanone	18	U
127-18-4	Tetrachloroethane	6	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	12	U
108-88-3	Toluene	6	U
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	6	U
74-88-4	Iodomethane	12	U

FORM I VOA

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107-02-8	Acrolein		110	U
107-13-1	Acrylonitrile		140	U
75-69-4	Trichlorofluoromethane		6	U
107-05-1	3-Chloropropene		18	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo		3	J
354-58-5	1,1,1-Trichloro-2,2,2-trifluo		12	U
74-95-3	Dibromomethane		12	U
4170-30-3	Crotonaldehyde		120	U
106-93-4	1,2-Dibromoethane		6	U
630-20-6	1,1,1,2-Tetrachloroethane		6	U
764-41-0	cis-1,4-Dichloro-2-butene		18	U
96-18-4	1,2,3-Trichloropropane		18	U
764-71-0	trans-1,4-Dichloro-2-butene		18	U
96-18-4	Ethylmethacrylate		12	U
96-12-8	1,2-Dibromo-3-chloropropane		12	U

035

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500027	ROJ3S
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 427
Matrix: (soil/water) SOIL	Lab Sample ID: 469291	
Sample wt/vol: 5.0 (g/mL) G	Lab File ID: GH069291B13	
Level: (low/med) LOW	Date Received: 12/11/91	
# Moisture: not dec. 9	Date Analyzed: 12/13/91	
Column: (pack/cap) CAP	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/Kg		
		Q		
74-87-3	Chloromethane	11	U	
74-83-9	Bromomethane	5	U	
75-01-4	Vinyl Chloride	11	U	
75-00-3	Chloroethane	11	U	
75-09-2	Methylene Chloride	55	B	
67-64-1	Acetone	28	U	
75-15-0	Carbon Disulfide	5	U	
75-35-4	1,1-Dichloroethene	5	U	
75-34-3	1,1-Dichloroethane	5	U	
540-59-0	1,2-Dichloroethene (total)	5	U	
67-66-3	Chloroform	5	U	
107-06-2	1,2-Dichloroethane	5	U	
78-93-3	2-Butanone	11	U	
71-55-6	1,1,1-Trichloroethane	5	U	
56-23-5	Carbon Tetrachloride	5	U	
108-05-4	Vinyl Acetate	11	U	
75-27-4	Bromodichloromethane	5	U	
78-87-5	1,2-Dichloropropane	5	U	
10061-01-5	cis-1,3-Dichloropropene	5	U	
79-01-6	Trichloroethene	5	U	
124-48-1	Dibromochloromethane	5	U	
79-00-5	1,1,2-Trichloroethane	5	U	
71-43-2	Benzene	5	U	
10061-02-6	Trans-1,3-Dichloropropene	5	U	
110-75-8	2-Chloroethylvinylether	11	U	
75-25-2	Bromoform	11	U	
108-10-1	4-Methyl-2-Pentanone	16	U	
591-78-6	2-Hexanone	16	U	
127-18-4	Tetrachloroethene	5	U	
79-34-5	1,1,2,2-Tetrachloroethane	11	U	
108-88-3	Toluene	5	U	
108-90-7	Chlorobenzene	5	U	
100-41-4	Ethylbenzene	5	U	
100-42-5	Styrene	5	U	
1330-20-7	Total Xylenes	5	U	
74-88-4	Iodomethane	11	U	

FORM I VOA

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107-02-8	Acrolein	99	U
107-13-1	Acrylonitrile	130	U
75-69-4	Trichlorofluoromethane	5	U
107-05-1	3-Chloropropene	16	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	2	J
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	11	U
74-95-3	Dibromomethane	11	U
4170-30-3	Crotonaldehyde	110	U
106-93-4	1,2-Dibromoethane	5	U
630-20-6	1,1,1,2-Tetrachloroethane	5	U
764-41-0	cis-1,4-Dichloro-2-butene	16	U
96-18-4	1,2,3-Trichloropropane	16	U
764-71-0	trans-1,4-Dichloro-2-butene	16	U
96-18-4	Ethylmethacrylate	11	U
96-12-8	1,2-Dibromo-3-chloropropane	11	U

POTS

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500077	ROJ4S
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 424
Matrix: (soil/water) SOIL		Lab Sample ID: 469257
Sample wt/vol: 5.0 (g/mL) G		Lab File ID: GH069257BL3
Level: (low/med) LOW		Date Received: 12/11/91
* Moisture: not dec. 33		Date Analyzed: 12/13/91
Column: (pack/cap) CAP		Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
74-87-3	Chloromethane	15	U
74-83-9	Bromomethane	7	U
75-01-4	Vinyl Chloride	15	U
75-00-3	Chloroethane	15	U
75-09-2	Methylene Chloride	87	B
67-64-1	Acetone	59	
75-15-0	Carbon Disulfide	7	U
75-35-4	1, 1-Dichloroethene	7	U
75-34-3	1, 1-Dichloroethane	7	U
540-59-0	1, 2-Dichloroethene (total)	7	U
67-66-3	Chloroform	7	U
107-06-2	1, 2-Dichloroethane	7	U
78-93-3	2-Butanone	15	U
71-55-6	1, 1, 1-Trichloroethane	7	U
56-23-5	Carbon Tetrachloride	7	U
108-05-4	Vinyl Acetate	15	U
75-27-4	Bromodichloromethane	7	U
78-87-5	1, 2-Dichloropropane	7	U
10061-01-5	cis-1, 3-Dichloropropene	7	U
79-01-6	Trichloroethene	7	U
124-48-1	Dibromochloromethane	7	U
79-00-5	1, 1, 2-Trichloroethane	7	U
71-43-2	Benzene	7	U
10061-02-6	Trans-1, 3-Dichloropropene	7	U
110-75-8	2-Chloroethylvinylether	15	U
75-25-2	Bromoform	15	U
108-10-1	4-Methyl-2-Pentanone	15	U
591-78-6	2-Hexanone	22	U
127-18-4	Tetrachloroethene	22	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	7	U
108-88-3	Toluene	15	U
108-90-7	Chlorobenzene	7	U
100-41-4	Ethylbenzene	7	U
100-42-5	Styrene	7	U
1330-20-7	Total Xylenes	7	U
74-88-4	Iodomethane	15	U

FORM I VOA

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107-02-8	Acrolein		130	U
107-13-1	Acrylonitrile		180	U
75-69-4	Trichlorofluoromethane		7	U
107-05-1	3-Chloropropene		22	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane		3	J
354-58-5	1,1,1-Trichloro-2,2,2-trifluoroethane		15	U
74-95-3	Dibromomethane		15	U
4170-30-3	Crotonaldehyde		150	U
106-93-4	1,2-Dibromoethane		7	U
630-20-6	1,1,1,2-Tetrachloroethane		7	U
764-41-0	cis-1,4-Dichloro-2-butene		22	U
96-18-4	1,2,3-Trichloropropane		22	U
764-71-0	trans-1,4-Dichloro-2-butene		22	U
96-18-4	Eethylmethacrylate		15	U
96-12-8	1,2-Dibromo-3-chloropropane		15	U

QJ34S

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	R001B1416
Lab Code: <u>COMPU</u>	Case No.: <u>22469</u>	SAS No.: _____ SDG No.: <u>01</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>396806</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>G</u>	Lab File ID: <u>GH096806B12</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>02/01/91</u>	
% Moisture: not dec. <u>15</u>	Date Analyzed: <u>02/05/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	Chloromethane	12	U
74-83-9	Bromomethane	6	U
75-01-4	Vinyl Chloride	12	U
75-00-3	Chloroethane	12	U
75-09-2	Methylene Chloride	33	B
67-64-1	Acetone	22	B
75-15-0	Carbon Disulfide	6	U
75-35-4	1,1-Dichloroethene	6	U
75-34-3	1,1-Dichloroethane	6	U
540-59-0	1,2-Dichloroethene (total)	6	U
67-66-3	Chloroform	6	U
107-06-2	1,2-Dichloroethane	6	U
78-93-3	2-Butanone	12	U
71-55-6	1,1,1-Trichloroethane	6	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	12	U
75-27-4	Bromodichloromethane	6	U
78-87-5	1,2-Dichloropropane	6	U
10061-01-5	cis-1,3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1,1,2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1,3-Dichloropropene	6	U
110-75-8	2-Chloroethylvinylether	12	U
75-25-2	Bromoform	12	U
108-10-1	4-Methyl-2-Pentanone	18	U
591-78-6	2-Hexanone	18	U
127-18-4	Tetrachloroethene	6	U
79-34-5	1,1,2,2-Tetrachloroethane	12	U
108-88-3	Toluene	6	U
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	6	U
74-88-4	Iodomethane	12	U

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107-02-8	Acrolein	110	U
107-13-1	Acrylonitrile	140	U
75-69-4	Trichlorofluoromethane	6	U
107-05-1	3-Chloropropene	18	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoropropane	12	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluoroethane	12	U
74-95-3	Dibromomethane	12	U
4170-30-3	Crotonaldehyde	120	U
106-93-4	1,2-Dibromoethane	6	U
630-20-6	1,1,1,2-Tetrachloroethane	6	U
764-71-0	cis-1,4-Dichloro-2-butene	18	U
96-18-4	1,2,3-Trichloropropane	18	
764-71-0	trans-1,4-Dichloro-2-butene	18	U
96-18-4	Eethylmethacrylate	12	U
96-12-8	1,2-Dibromo-3-chloropropane	12	U

FORM I VOA

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CCN 396806

LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	R002B0810
Lab Code: <u>COMPU</u>	Case No.: <u>22469</u>	SAS No.: _____ SDG No.: <u>01</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>396817</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>G</u>	Lab File ID: <u>GH096817B12</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>02/01/91</u>	
% Moisture: not dec. <u>18</u>	Date Analyzed: <u>02/05/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
74-87-3	-Chloromethane	12	U
74-83-9	-Bromomethane	6	U
75-01-4	-Vinyl Chloride	12	U
75-00-3	-Chloroethane	12	U
75-09-2	-Methylene Chloride	38	B
67-64-1	-Acetone	32	B
75-15-0	-Carbon Disulfide	6	U
75-35-4	-1,1-Dichloroethene	6	U
75-34-3	-1,1-Dichloroethane	6	U
540-59-0	-1,2-Dichloroethene (total)	6	U
67-66-3	-Chloroform	6	U
107-06-2	-1,2-Dichloroethane	6	U
78-93-3	-2-Butanone	12	U
71-55-6	-1,1,1-Trichloroethane	6	U
56-23-5	-Carbon Tetrachloride	6	U
108-05-4	-Vinyl Acetate	12	U
75-27-4	-Bromodichloromethane	6	U
78-87-5	-1,2-Dichloropropane	6	U
10061-01-5	-cis-1,3-Dichloropropene	6	U
79-01-6	-Trichloroethene	6	U
124-48-1	-Dibromochloromethane	6	U
79-00-5	-1,1,2-Trichloroethane	6	U
71-43-2	-Benzene	6	U
10061-02-6	-Trans-1,3-Dichloropropene	6	U
110-75-8	-2-Chloroethylvinylether	12	U
75-25-2	-Bromoform	12	U
108-10-1	-4-Methyl-2-Pentanone	18	U
591-78-6	-2-Hexanone	18	U
127-18-4	-Tetrachloroethene	6	U
79-34-5	-1,1,2,2-Tetrachloroethane	12	U
108-88-3	-Toluene	6	U
108-90-7	-Chlorobenzene	6	U
100-41-4	-Ethylbenzene	6	U
100-42-5	-Styrene	6	U
1330-20-7	-Total Xylenes	6	U
74-88-4	-Iodomethane	12	U

FORM I VOA

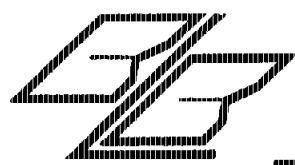
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107-02-8	Acrolein	110	U
107-13-1	Acrylonitrile	150	U
75-69-4	Trichlorofluoromethane	6	U
107-05-1	3-Chloropropene	18	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluo	12	U
354-58-5	1,1,1-Trichloro-2,2,2-trifluo	12	U
74-95-3	Dibromomethane	12	U
4170-30-3	Crotonaldehyde	120	U
106-93-4	1,2-Dibromoethane	6	U
630-20-6	1,1,1,2-Tetrachloroethane	6	U
764-71-0	cis-1,4-Dichloro-2-butene	18	U
96-18-4	1,2,3-Trichloropropane	18	U
764-71-0	trans-1,4-Dichloro-2-butene	18	U
96-18-4	Ethylmethacrylate	12	U
96-12-8	1,2-Dibromo-3-chloropropane	12	U

FORM I VOA

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CCN 396817



Section 2

SEMIVOLATILE ORGANICS ANALYSIS (SOIL)

FP-1, 8-12' -	Soil sample from Oxbow Area J, Soil Boring FP-1 at 8-12'
FP-2, 4-8' -	Soil sample from Oxbow Area J, Soil Boring FP-2 at 4-8'
FP-3, 4-8' -	Soil sample from Oxbow Area J, Soil Boring FP-3 at 4-8'
YB-2, 4-8' -	Soil sample from Oxbow Area J, Soil Boring YB-2 at 4-8'
YB-4, 0-4' -	Soil sample from Oxbow Area J, Soil Boring YB-4 at 0-4'
ROA012224 -	Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet
ROA2B0608 -	Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet
ROA3B1214 -	Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 12-14 feet
ROB1B0406 -	Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet
ROB-DPA1 -	Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)
ROB2B0002 -	Soil sample from Oxbow Area B, Groundwater Monitoring Well B-2 at 0-2 feet
ROC011012 -	Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet
ROC021214 -	Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet
ROC-DPA1 -	Soil sample from Oxbow Area C, Groundwater Monitoring Well 12-14 (Duplicate)
ROC3B0204 -	Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet
ROJ1S -	Soil sample from Oxbow Area J, Surficial Soil Sample J-1
ROJ2S -	Soil sample from Oxbow Area J, Surficial Soil Sample J-2
ROJ3S -	Soil sample from Oxbow Area J, Surficial Soil Sample J-3
ROJ4S -	Soil sample from Oxbow Area J, Surficial Soil Sample J-4
ROO1B1416 -	Soil sample from Oxbow Area K, Soil Boring K1 at 14-16 feet
ROO2B0810 -	Soil sample from Oxbow Area K, Soil Boring K2 at 8-10 feet

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: FP-1,8-12¹
Lab Sample ID: JJ7130

<u>Compound</u>		<u>Compound</u>	
acenaphthene	2,000 U	3,3'-dichlorobenzidine	4,000 U
acenaphthylene	2,000 U	diethyl phthalate	2,000 U
anthracene	2,000 U	dimethyl phthalate	2,000 U
benzidine	9,900 U	2,4-dinitrotoluene	2,000 U
benzo(a)anthracene	2,000 U	2,6-dinitrotoluene	2,000 U
benzo(b)fluoranthene	2,000 U	di-n-octylphthalate	2,000 U
benzo(k)fluoranthene	2,000 U	1,2-diphenylhydrazine ¹	2,000 U
benzo(a)pyrene	2,000 U	fluoranthene	350 J
benzo(g,h,i)perylene	2,000 U	fluorene	2,000 U
benzyl butyl phthalate	2,000 U	hexachlorobenzene	2,000 U
bis(2-chloroethoxy)methane	2,000 U	hexachlorobutadiene	2,000 U
bis(2-chloroethyl)ether	2,000 U	hexachlorocyclopentadiene	2,000 U
bis(2-chloroisopropyl)ether	2,000 U	hexachloroethane	2,000 U
bis(2-ethylhexyl)phthalate	2,000 U	indeno(1,2,3-cd)pyrene	2,000 U
4-bromophenyl phenyl ether	2,000 U	isophorone	2,000 U
2-chloronaphthalene	2,000 U	naphthalene	2,000 U
4-chlorophenyl phenyl ether	2,000 U	nitrobenzene	2,000 U
chrysene	2,000 U	N-nitrosodimethylamine	2,000 U
dibenzo(a,h)anthracene	2,000 U	N-nitrosodi-n-propylamine	2,000 U
di-n-butylphthalate	2,000 U	N-nitrosodiphenylamine ²	2,000 U
1,2-dichlorobenzene	2,000 U	phenanthrene	480 J
1,3-dichlorobenzene	2,000 U	pyrene	270 J
1,4-dichlorobenzene	2,000 U	1,2,4-trichlorobenzene	2,000 U

1 Screened for as Azobenzene

2 Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89

Date of Analysis: 10/12/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: FP-2,4-8'
Lab Sample ID: JJ7131

Compound

Compound

acenaphthene	1,900 U	3,3'-dichlorobenzidine	3,800 U
acenaphthylene	1,900 U	diethyl phthalate	1,900 U
anthracene	1,900 U	dimethyl phthalate	1,900 U
benzidine	9,500 U	2,4-dinitrotoluene	1,900 U
benzo(a)anthracene	260 J	2,6-dinitrotoluene	1,900 U
benzo(b)fluoranthene	1,900 U	di-n-octylphthalate	1,900 U
benzo(k)fluoranthene	1,900 U	1,2-diphenylhydrazine ¹	1,900 U
benzo(a)pyrene	200 J	fluoranthene	550 J
benzo(g,h,i)perylene	1,900 U	fluorene	1,900 U
benzyl butyl phthalate	1,900 U	hexachlorobenzene	1,900 U
bis(2-chloroethoxy)methane	1,900 U	hexachlorobutadiene	1,900 U
bis(2-chloroethyl)ether	1,900 U	hexachlorocyclopentadiene	1,900 U
bis(2-chloroisopropyl)ether	1,900 U	hexachloroethane	1,900 U
bis(2-ethylhexyl)phthalate	1,900 U	indeno(1,2,3-cd)pyrene	1,900 U
4-bromophenyl phenyl ether	1,900 U	isophorone	1,900 U
2-chloronaphthalene	1,900 U	naphthalene	1,900 U
4-chlorophenyl phenyl ether	1,900 U	nitrobenzene	1,900 U
chrysene	230 J	N-nitrosodimethylamine	1,900 U
dibenzo(a,h)anthracene	1,900 U	N-nitrosodi-n-propylamine	1,900 U
di-n-butylphthalate	1,900 U	N-nitrosodiphenylamine ²	1,900 U
1,2-dichlorobenzene	1,900 U	phenanthrene	480 J
1,3-dichlorobenzene	1,900 U	pyrene	420 J
1,4-dichlorobenzene	1,900 U	1,2,4-trichlorobenzene	1,900 U

1 Screened for as Azobenzene

2 Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89

Date of Analysis: 10/12/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: FP-3,4-8'
Lab Sample ID: JJ7132

<u>Compound</u>	<u>Compound</u>
acenaphthene	3,3'-dichlorobenzidine
acenaphthylene	diethyl phthalate
anthracene	dimethyl phthalate
benzidine	2,4-dinitrotoluene
benzo(a)anthracene	2,6-dinitrotoluene
benzo(b)fluoranthene	di-n-octylphthalate
benzo(k)fluoranthene	1,2-diphenylhydrazine ¹
benzo(a)pyrene	fluoranthene
benzo(g,h,i)perylene	fluorene
benzyl butyl phthalate	hexachlorobenzene
bis(2-chloroethoxy)methane	hexachlorobutadiene
bis(2-chloroethyl)ether	hexachlorocyclopentadiene
bis(2-chloroisopropyl)ether	hexachloroethane
bis(2-ethylhexyl)phthalate	indeno(1,2,3-cd)pyrene
4-bromophenyl phenyl ether	isophorone
2-chloronaphthalene	naphthalene
4-chlorophenyl phenyl ether	nitrobenzene
chrysene	N-nitrosodimethylamine
dibenzo(a,h)anthracene	N-nitrosodi-n-propylamine
di-n-butylphthalate	N-nitrosodiphenylamine ²
1,2-dichlorobenzene	phenanthrene
1,3-dichlorobenzene	pyrene
1,4-dichlorobenzene	1,2,4-trichlorobenzene

1 Screened for as Azobenzene

2 Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89
Date of Analysis: 10/12/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank
Lab Sample ID: BL4910

<u>Compound</u>		<u>Compound</u>	
acenaphthene	1,000 U	3,3'-dichlorobenzidine	2,000 U
acenaphthylene	1,000 U	diethyl phthalate	1,000 U
anthracene	1,000 U	dimethyl phthalate	1,000 U
benzidine	5,000 U	2,4-dinitrotoluene	1,000 U
benzo(a)anthracene	1,000 U	2,6-dinitrotoluene	1,000 U
benzo(b)fluoranthene	1,000 U	di-n-octylphthalate	1,000 U
benzo(k)fluoranthene	1,000 U	1,2-diphenylhydrazine ¹	1,000 U
benzo(a)pyrene	1,000 U	fluoranthene	1,000 U
benzo(g,h,i)perylene	1,000 U	fluorene	1,000 U
benzyl butyl phthalate	1,000 U	hexachlorobenzene	1,000 U
bis(2-chloroethoxy)methane	1,000 U	hexachlorobutadiene	1,000 U
bis(2-chloroethyl)ether	1,000 U	hexachlorocyclopentadiene	1,000 U
bis(2-chloroisopropyl)ether	1,000 U	hexachloroethane	1,000 U
bis(2-ethylhexyl)phthalate	310 J	indeno(1,2,3-cd)pyrene	1,000 U
4-bromophenyl phenyl ether	1,000 U	isophorone	1,000 U
2-chloronaphthalene	1,000 U	naphthalene	1,000 U
4-chlorophenyl phenyl ether	1,000 U	nitrobenzene	1,000 U
chrysene	1,000 U	N-nitrosodimethylamine	1,000 U
dibenzo(a,h)anthracene	1,000 U	N-nitrosodi-n-propylamine	1,000 U
di-n-butylphthalate	1,000 U	N-nitrosodiphenylamine ²	1,000 U
1,2-dichlorobenzene	1,000 U	phenanthrene	1,000 U
1,3-dichlorobenzene	1,000 U	pyrene	1,000 U
1,4-dichlorobenzene	1,000 U	1,2,4-trichlorobenzene	1,000 U

I Screened for as Azobenzene

2 Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89

Date of Analysis: 10/12/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

SOIL SURROGATE PERCENT RECOVERY SUMMARY

Sample No.	SEMI-VOLATILE					2,4,6 Tribromo- Phenol (18-122%)*
	Nitro-Benzene-D5 (23-120%)*	2-Fluoro-Biphenyl (30-116%)*	Terphenyl-D14 (18-137%)*	Phenol-D5 (24-113%)*	2-Fluoro- Phenol (26-121%)*	
FP-1,8-12'	71	69	66	55	61	74
FP-2,4-8'	60	61	64	46	52	69
FP-3,4-8'	54	60	57	40	42	43
Method Blank	78	75	74	55	64	66

*Values in parenthesis represent USEPA contract required QC limits.

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44313

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil -

Client Sample ID: YB-2, 4-8'
Lab Sample ID: JJ7182

<u>Compound</u>		<u>Compound</u>	
acenaphthene	2,000 U	3,3'-dichlorobenzidine	4,000 U
acenaphthylene	270 J	diethyl phthalate	2,000 U
anthracene	2,000 U	dimethyl phthalate	2,000 U
benzidine	10,000 U	2,4-dinitrotoluene	2,000 U
benzo(a)anthracene	300 J	2,6-dinitrotoluene	2,000 U
benzo(b)fluoranthene	380 J	di-n-octylphthalate	2,000 U
benzo(k)fluoranthene	460 J	1,2-diphenylhydrazine ¹	2,000 U
benzo(a)pyrene	370 J	fluoranthene	470 J
benzo(g,h,i)perylene	2,000 U	fluorene	2,000 U
benzyl butyl phthalate	2,000 U	hexachlorobenzene	2,000 U
bis(2-chloroethoxy)methane	2,000 U	hexachlorobutadiene	2,000 U
bis(2-chloroethyl)ether	2,000 U	hexachlorocyclopentadiene	2,000 U
bis(2-chloroisopropyl)ether	2,000 U	hexachloroethane	2,000 U
bis(2-ethylhexyl)phthalate	2,000 U	indeno(1,2,3-cd)pyrene	2,000 U
4-bromophenyl phenyl ether	2,000 U	isophorone	2,000 U
2-chloronaphthalene	2,000 U	naphthalene	2,000 U
4-chlorophenyl phenyl ether	2,000 U	nitrobenzene	2,000 U
chrysene	310 J	N-nitrosodimethylamine	2,000 U
dibenzo(a,h)anthracene	2,000 U	N-nitrosodi-n-propylamine	2,000 U
di-n-butylphthalate	2,000 U	N-nitrosodiphenylamine ²	2,000 U
1,2-dichlorobenzene	2,000 U	phenanthrene	290 J
1,3-dichlorobenzene	2,000 U	pyrene	700 J
1,4-dichlorobenzene	2,000 U	1,2,4-trichlorobenzene	2,000 U

1 Screened for as Azobenzene

2 Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89

Date of Analysis: 10/11/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44313

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: YB-4, 0-4'
Lab Sample ID: JJ7183

<u>Compound</u>		<u>Compound</u>	
acenaphthene	2,000 U	3,3'-dichlorobenzidine	4,000 U
acenaphthylene	420 J	diethyl phthalate	2,000 U
anthracene	260 J	dimethyl phthalate	2,000 U
benzidine	9,900 U	2,4-dinitrotoluene	2,000 U
benzo(a)anthracene	650 J	2,6-dinitrotoluene	2,000 U
benzo(b)fluoranthene	1,000 J	di-n-octylphthalate	2,000 U
benzo(k)fluoranthene	910 J	1,2-diphenylhydrazine ¹	2,000 U
benzo(a)pyrene	930 J	fluoranthene	890 J
benzo(g,h,i)perylene	770 J	fluorene	2,000 U
benzyl butyl phthalate	2,000 U	hexachlorobenzene	2,000 U
bis(2-chloroethoxy)methane	2,000 U	hexachlorobutadiene	2,000 U
bis(2-chloroethyl)ether	2,000 U	hexachlorocyclopentadiene	2,000 U
bis(2-chloroisopropyl)ether	2,000 U	hexachloroethane	2,000 U
bis(2-ethylhexyl)phthalate	2,000 U	Indeno(1,2,3-cd)pyrene	660 J
4-bromophenyl phenyl ether	2,000 U	isophorone	2,000 U
2-chloronaphthalene	2,000 U	naphthalene	2,000 U
4-chlorophenyl phenyl ether	2,000 U	nitrobenzene	2,000 U
chrysene	640 J	N-nitrosodimethylamine	2,000 U
dibenzo(a,h)anthracene	240 J	N-nitrosodi-n-propylamine	2,000 U
di-n-butylphthalate	2,000 U	N-nitrosodiphenylamine ²	630 J
1,2-dichlorobenzene	2,000 U	phenanthrene	430 J
1,3-dichlorobenzene	2,000 U	pyrene	940 J
1,4-dichlorobenzene	2,000 U	1,2,4-trichlorobenzene	2,000 U

¹ Screened for as Azobenzene

² Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89

Date of Analysis: 10/11/89

Geraghty & Miller, Inc.
October 16, 1989

Client Project ID: GE Pittsfield

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Job Number: GMIN 44313

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in $\mu\text{g}/\text{kg}$ (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank
Lab Sample ID: BL4910

<u>Compound</u>		<u>Compound</u>	
acenaphthene	1,000 U	3,3'-dichlorobenzidine	2,000 U
acenaphthylene	1,000 U	diethyl phthalate	1,000 U
anthracene	1,000 U	dimethyl phthalate	1,000 U
benzidine	5,000 U	2,4-dinitrotoluene	1,000 U
benzo(a)anthracene	1,000 U	2,6-dinitrotoluene	1,000 U
benzo(b)fluoranthene	1,000 U	di-n-octylphthalate	1,000 U
benzo(k)fluoranthene	1,000 U	1,2-diphenylhydrazine ¹	1,000 U
benzo(a)pyrene	1,000 U	fluoranthene	1,000 U
benzo(g,h,i)perylene	1,000 U	fluorene	1,000 U
benzyl butyl phthalate	1,000 U	hexachlorobenzene	1,000 U
bis(2-chloroethoxy)methane	1,000 U	hexachlorobutadiene	1,000 U
bis(2-chloroethyl)ether	1,000 U	hexachlorocyclopentadiene	1,000 U
bis(2-chloroisopropyl)ether	1,000 U	hexachloroethane	1,000 U
bis(2-ethylhexyl)phthalate	290 J	indeno(1,2,3-cd)pyrene	1,000 U
4-bromophenyl phenyl ether	1,000 U	isophorone	1,000 U
2-chloronaphthalene	1,000 U	naphthalene	1,000 U
4-chlorophenyl phenyl ether	1,000 U	nitrobenzene	1,000 U
chrysene	1,000 U	N-nitrosodimethylamine	1,000 U
dibenzo(a,h)anthracene	1,000 U	N-nitrosodi-n-propylamine	1,000 U
di-n-butylphthalate	1,000 U	N-nitrosodiphenylamine ²	200 J
1,2-dichlorobenzene	1,000 U	phenanthrene	1,000 U
1,3-dichlorobenzene	1,000 U	pyrene	1,000 U
1,4-dichlorobenzene	1,000 U	1,2,4-trichlorobenzene	1,000 U

¹ Screened for as Azobenzene

² Detected as Diphenylamine

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction: 10/09/89

Date of Analysis: 10/10/89

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44313

SOIL SURROGATE PERCENT RECOVERY SUMMARY

Sample No.	SEMI-VOLATILE					2,4,6 Tribromo-Phenol (18-122%)*
	Nitro-Benzene-05 (23-120%)*	2-Fluoro-Biphenyl (30-116%)*	Terphenyl-014 (18-137%)*	Phenol-05 (24-113%)*	2-Fluoro-Phenol (26-121%)*	
YB-2, 4-8"	80	78	96	57	65	77
YB-4, 0-4"	86	84	83	65	72	83
Method Blank	73	70	73	58	62	68

*Values in parenthesis represent USEPA contract required QC limits.

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE N

ROA012224

Lab Name: COMPUCHEM, RTP Contract: 500077
 Lab Code: COMPU Case No.: 24105 SAS No.: SDG No.: 27
 Matrix: (soil/water) SOIL Lab Sample ID: 462144
 Sample wt/vol: 30.5 (g/mL) G Lab File ID: GJ062144A20
 Level: (low/med) LOW Date Received: 11/08/91
 % Moisture: not dec. 10 dec. Date Extracted: 11/15/91
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 11/23/91
 GPC Cleanup: (Y/N) N pH: Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
62-75-9	N-Nitrosodimethylamine	360	U
110-86-1	Pyridine	360	U
97-63-2	Ethyl methacrylate	360	U
123-63-7	Paraldehyde	360	U
109-06-8	2-Picoline	720	U
10595-95-6	Nitrosomethylamine	360	U
66-27-3	Methyl methanesulfonate	360	U
55-18-5	N-Nitrosodiethylamine	360	U
62-50-0	Ethyl methanesulfonate	360	U
108-95-2	Phenol	360	U
62-53-3	Aniline	360	U
76-01-7	Pentachloroethane	360	U
111-44-4	bis(2-Chloroethyl)Ether	720	U
95-57-8	2-Chlorophenol	360	U
541-73-1	1,3-Dichlorobenzene	360	U
106-46-7	1,4-Dichlorobenzene	360	U
100-51-6	Benzyl Alcohol	360	U
95-50-1	1,2-Dichlorobenzene	360	U
95-48-7	2-Methylphenol	360	U
39638-32-9	bis(2-Chloroisopropyl)Ether	360	U
108-39-4	3-Methylphenol	360	U
106-44-5	4-Methylphenol	360	U
930-55-2	N-Nitrosopyrrolidine	360	U
59-89-2	N-Nitrosomorpholine	360	U
98-86-2	Acetophenone	360	U
621-64-7	N-Nitroso-Di-n-Propylamine	360	U
636-21-5	O-Toluidine hydrochloride	360	U
67-72-1	Hexachloroethane	360	U
98-95-3	Nitrobenzene	360	U
100-75-4	N-Nitrosopiperidine	360	U
78-59-1	Iscophorone	360	U
88-75-5	2-Nitrophenol	360	U
105-67-9	2,4-Dimethylphenol	360	U
108-70-3	1,3,5-Trichlorobenzene	360	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3	Benzal Chloride	360	U
65-85-0	Benzoic Acid	3600	U
111-91-1	bis(2-Chloroethoxy)Methane	360	U
120-83-2	2,4-Dichlorophenol	360	U
120-82-1	1,2,4-Trichlorobenzene	360	U
91-20-3	Naphthalene	360	U
106-47-8	4-Chloraniline	360	U
87-65-0	2,6-Dichlorophenol	720	U
95-54-5	o-Phenylenediamine	360	U
122-09-8	dimethylphenylethylamine	360	U
1888-71-7	Hexachloropropene	360	U
87-68-3	Hexachlorobutadiene	360	U
87-61-6	1,2,3-Trichlorobenzene	360	U
98-07-7	Benzotrichloride	720	U
924-16-3	N-Nitroso-di-n-butylamine	360	U
59-50-7	4-Chloro-3-Methylphenol	360	U
106-50-3	p-Phenylenediamine	360	U
94-59-7	Safrole	360	U
106-50-3	m-Phenylenediamine	360	U
91-57-6	2-Methylnaphthalene	360	U
90-12-0	1-Methylnaphthalene	360	U
95-94-3	1,2,4,5-Tetrachlorobenzene	360	U
634-90-2	1,2,3,5-Tetrachlorobenzene	360	U
77-47-4	Hexachlorocyclopentadiene	360	U
88-06-2	2,4,6-Trichlorophenol	720	U
95-95-4	2,4,5-Trichlorophenol	720	U
120-58-1	Icosafrole	720	U
91-58-7	2-Chloronaphthalene	360	U
90-13-1	1-Chloronaphthalene	360	U
634-66-2	1,2,3,4-Tetrachlorobenzene	360	U
88-74-4	2-Nitroaniline	360	U
130-15-4	1,4-Naphthoquinone	720	U
100-25-4	1,4-Dinitrobenzene	720	U
131-11-3	Dimethyl Phthalate	360	U
208-96-8	Acenaphthylene	360	U
606-20-2	2,6-Dinitrotoluene	360	U

462144

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE N

Lab Name: COMMERCIAL RTP	Contract: 590077	ROA012224
Lab Code: COMPU	Case No.: 24105	SAS No.: _____ SDG No.: 27
Matrix: (soil/water) SOIL	Lab Sample ID: 462144	
Sample wt/vol: 20.5 (g/mL) G	Lab File ID: GJ062144A20	
Level: (low/med) LOW	Date Received: 11/08/91	
# Moisture: not dec. 10 dec.	Date Extracted: 11/15/91	
Extraction: (SepF/Cont/Sonc) SQNC	Date Analyzed: 11/23/91	
GPC Cleanup: (Y/N) N	pH: _____	Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
99-09-2	3-Nitroaniline	720	U
83-32-9	Acenaphthene	360	U
51-28-5	2, 4-Dinitrophenol	1400	U
100-02-7	4-Nitrophenol	360	U
132-64-9	Dibenzofuran	360	U
121-14-2	2, 4-Dinitrotoluene	360	U
608-93-5	Pentachlorobenzene	360	U
91-59-8	2-Naphthylamine	720	U
134-32-7	1-Naphthylamine	720	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	720	U
84-66-2	Diethylphthalate	360	U
297-97-2	Zinophos	360	U
7005-72-3	4-Chlorophenyl-phenylether	360	U
86-73-7	Fluorene	360	U
100-01-6	4-Nitroaniline	360	U
99-55-8	5-Nitro-o-toluidine	720	U
122-66-7	1, 2-Diphenylhydrazine	720	U
534-52-1	4, 6-Dinitro-2-Methylphenol	360	U
86-30-6	N-Nitrosodiphenylamine (1)	1100	U
122-39-4	Diphenylamine	360	U
99-35-4	1, 3, 5-Trinitrobenzene	720	U
62-44-2	Phenacetin	360	U
101-55-3	4-Bromophenyl-phenylether	360	U
2303-16-4	Diallate	360	U
60-51-5	Dimethoate	360	U
118-74-1	Hexachlorobenzene	360	U
92-67-1	4-Aminobiphenyl	360	U
23950-58-5	Pronamide	360	U
87-86-5	Pentachlorophenol	360	U
82-68-8	Pentachloronitrobenzene	720	U
95-01-8	Phenanthrrene	360	U
120-12-7	Anthracene	360	U
84-74-2	Di-n-Butylphthalate	360	U
91-80-5	Methapyrilene	720	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	1700	U
206-44-0	fluoranthene	360	U
92-87-5	Benzidine	360	U
129-00-0	Pyrene	360	U
60-11-7	p-Dimethylaminoazobenzene	360	U
510-15-6	Chlorobenzilate	360	U
119-93-7	3,3'-Dimethylbenzidine	720	U
85-68-7	Butylbenzylphthalate	360	U
53-96-3	2-Acetylaminofluorene	360	U
101-14-4	Methylene-bis(2-Chloroaniline	360	U
91-94-1	3,3'-Dichlorobenzidine	360	U
119-90-4	3,3'-Dimethoxybenzidine	360	U
56-55-3	Benzo(a)Anthracene	360	U
218-01-9	Chrysene	360	U
117-81-7	bis(2-Ethylhexyl) Phthalate	360	U
117-84-0	Di-n-Octyl Phthalate	360	U
205-99-2	Benzo(b)Fluoranthene	360	U
57-97-6	7,12-Dimethylbenzanthracene	360	U
207-08-9	Benzo(k)Fluoranthene	360	U
50-32-8	Benzo(a)Pyrene	360	U
56-49-5	3-Methylchloranthrene	360	U
224-42-0	Dibenzo(a,j)acridine	360	U
193-39-5	Indeno(1,2,3-cd) Pyrene	360	U
53-70-3	Dibenz(a,h)Anthracene	360	U
191-24-2	Benzo(g,h,i)Perylene	360	U

(1) - Cannot be separated from Diphenylamine

5

4/6/2 1444

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA2B0608
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465049</u>	
Sample wt/vol: <u>30.2</u> (g/mL) <u>G</u>	Lab File ID: <u>GD065049B22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/21/91</u>	
% Moisture: not dec. <u>9</u> dec. _____	Date Extracted: <u>11/27/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/09/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>2.0</u>

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) UG/KG	Q
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62-75-9	N-Nitrosodimethylamine	720	U
110-86-1	Pyridine	720	U
97-63-2	Ethyl methacrylate	720	U
123-63-7	Paraldehyde	720	U
109-06-8	2-Picoline	1400	U
10595-95-6	Nitrosomethylethylamine	720	U
66-27-3	Methyl methanesulfonate	720	U
55-18-5	N-Nitrosodiethylamine	720	U
62-50-0	Ethyl methanesulfonate	720	U
108-95-2	Phenol	720	U
62-53-3	Aniline	720	U
76-01-7	Pentachloroethane	720	U
111-44-4	bis(2-Chloroethyl) Ether	1400	U
95-57-8	2-Chlorophenol	720	U
541-73-1	1,3-Dichlorobenzene	720	U
106-46-7	1,4-Dichlorobenzene	720	U
100-51-6	Benzyl Alcohol	720	U
95-50-1	1,2-Dichlorobenzene	720	U
95-48-7	2-Methylphenol	720	U
108-60-1	bis(2-Chloroisopropyl) Ether	720	U
108-39-4	3-Methylphenol	720	U
106-44-5	4-Methylphenol	720	U
930-55-2	N-Nitroso- <i>Di-n</i> -Propylamine	720	U
59-89-2	N-Nitrosomorpholine	720	U
98-86-2	Acetophenone	720	U
621-64-7	N-Nitroso- <i>Di-n</i> -Propylamine	720	U
636-21-5	o-Toluidine hydrochloride	720	U
67-72-1	Hexachloroethane	720	U
98-95-3	Nitrobenzene	720	U
100-75-4	N-Nitroscopiperidine	720	U
78-59-1	Isophorone	720	U
88-75-5	2-Nitrophenol	720	U
105-67-9	2,4-Dimethylphenol	720	U
108-70-3	1,3,5-Trichlorobenzene	720	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3-----	Benzal Chloride	720	U
65-85-0-----	Benzoic Acid	100	J
111-91-1-----	bis(2-Chloroethoxy)Methane	720	U
120-83-2-----	2, 4-Dichlorophenol	720	U
120-82-1-----	1, 2, 4-Trichlorobenzene	720	U
91-20-3-----	Naphthalene	2200	
106-47-8-----	4-Chloroaniline	720	U
87-65-0-----	2, 6-Dichlorophenol	1400	U
95-54-5-----	O-Phenylenediamine	720	U
122-09-8-----	dimethylphenylethylamine	720	U
1888-71-7-----	Hexachloropropene	720	U
87-68-3-----	Hexachlorobutadiene	720	U
87-61-6-----	1, 2, 3-Trichlorobenzene	720	U
98-07-7-----	Benzotrichloride	1400	U
924-16-3-----	N-Nitroso-di-n-butylamine	720	U
59-50-7-----	4-Chloro-3-Methylphenol	720	U
106-50-3-----	P-Phenylenediamine	720	U
94-59-7-----	Safrole	720	U
106-50-3-----	m-Phenylenediamine	720	U
91-57-6-----	2-Methylnaphthalene	930	
90-12-0-----	1-Methylnaphthalene	1900	
95-94-3-----	1, 2, 4, 5-Tetrachlorobenzene	720	U
634-90-2-----	1, 2, 3, 5-Tetrachlorobenzene	720	U
77-47-4-----	Hexachlorocyclopentadiene	720	U
88-06-2-----	2, 4, 6-Trichlorophenol	1400	U
95-95-4-----	2, 4, 5-Trichlorophenol	1400	U
120-58-1-----	Isosafrole	1400	U
91-58-7-----	2-Chloronaphthalene	720	U
90-13-1-----	1-Chloronaphthalene	720	U
634-66-2-----	1, 2, 3, 4-Tetrachlorobenzene	720	U
88-74-4-----	2-Nitroaniline	720	U
130-15-4-----	1, 4-Naphthoquinone	1400	U
100-25-4-----	1, 4-Dinitrobenzene	1400	U
131-11-3-----	Dimethyl Phthalate	720	U
208-96-8-----	Acenaphthylene	1000	
606-20-2-----	2, 6-Dinitrotoluene	720	U

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	<u>ROA2B0608</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465049</u>	
Sample wt/vol: <u>.30±.2</u> (g/mL) <u>G</u>	Lab File ID: <u>GD065049B22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/21/91</u>	
% Moisture: not dec. <u>2</u> dec. <u> </u>	Date Extracted: <u>11/27/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/09/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: <u> </u>	Dilution Factor: <u>2.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
99-09-2	3-Nitroaniline	1400	U
83-32-9	Acenaphthene	630	J
51-28-5	2, 4-Dinitrophenol	2800	U
100-02-7	4-Nitrophenol	720	U
132-64-9	Dibenzofuran	1100	
121-14-2	2, 4-Dinitrotoluene	720	U
608-93-5	Pentachlorobenzene	720	U
91-59-8	2-Naphthylamine	1400	U
134-32-7	1-Naphthylamine	1400	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	1400	U
84-66-2	Diethylphthalate	720	U
297-97-2	Zinophos	720	U
7005-72-3	4-Chlorophenyl-phenylether	720	U
86-73-7	Fluorene	2200	
100-01-6	4-Nitroaniline	1400	U
99-55-8	5-Nitro-o-toluidine	1400	U
122-66-7	1, 2-Diphenylhydrazine	720	U
534-52-1	4, 6-Dinitro-2-Methylphenol	2200	U
86-30-6	N-Nitrosodiphenylamine (1)	720	U
122-39-4	Diphenylamine	720	U
99-35-4	1, 3, 5-Trinitrobenzene	1400	U
62-44-2	Phenacetin	720	U
101-55-3	4-Bromophenyl-phenylether	720	U
2303-16-4	Diallate	720	U
60-51-5	Dimethoate	720	U
118-74-1	Hexachlorobenzene	720	U
92-67-1	4-Aminobiphenyl	720	U
23950-58-5	Pronamide	720	U
87-86-5	Pentachlorophenol	510	J
82-68-8	Pentachloronitrobenzene	720	U
85-01-8	Phenanthrene	5700	
120-12-7	Anthracene	1900	
84-74-2	Di-n-Butylphthalate	720	U
91-80-5	Methapyrilene	1400	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	3500	U
206-44-0	Fluoranthene	6700	
92-87-5	Benzidine	720	U
129-00-0	Pyrene	5300	
60-11-7	p-Dimethylaminoazobenzene	720	U
510-15-6	Chlorobenzilate	720	U
119-93-7	3,3'-Dimethylbenzidine	1400	U
85-68-7	Butylbenzylphthalate	720	U
53-96-3	2-Acetylaminofluorene	720	U
101-14-4	Methylene-bis(2-Chloroaniline	720	U
91-94-1	3,3'-Dichlorobenzidine	720	U
106-51-4	3,3'-Dimethoxybenzidine	720	U
56-55-3	Benzo(a)Anthracene	3000	
218-01-9	Chrysene	2700	
117-81-7	bis(2-Ethylhexyl)Phthalate	350	J
117-84-0	Di-n-Octyl Phthalate	720	U
205-99-2	Benzo(b)Fluoranthene	4000	
57-97-6	7,12-Dimethylbenzanthracene	720	U
207-08-9	Benzo(k)Fluoranthene	7000	
50-32-8	Benzo(a)Pyrene	2500	
56-49-5	3-Methylchloranthrene	720	U
224-42-0	Dibenzo(a,j)acridine	720	U
193-39-5	Indeno(1,2,3-cd)Pyrene	1100	
53-70-3	Dibenz(a,h)Anthracene	340	J
191-24-2	Benzo(g,h,i)Perylene	1100	

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name:	COMPUCHEM RTP	Contract:	500077	ROA3B1214		
Lab Code:	COMPEU	Case No.:	24105	SAS No.:	SDG No.:	578
Matrix:	(soil/water) SOIL	Lab Sample ID:	473668			
Sample wt/vol:	30.1 (g/mL) G	Lab File ID:	GD073668C22			
Level:	(low/med) LOW	Date Received:	01/07/92			
% Moisture:	not dec. 9 dec.	Date Extracted:	01/08/92			
Extraction:	(SepF/Cont/Sonc)	SONC	Date Analyzed:	01/11/92		
GPC Cleanup:	(Y/N) N	pH:	Dilution Factor:	10		

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
62-75-9	N-Nitrosodimethylamine	3600	U	
110-86-1	Pyridine	3600	U	
97-63-2	Ethyl methacrylate	3600	U	
123-63-7	Paraldehyde	3600	U	
109-06-8	2-Picoline	7200	U	
10595-95-6	Nitrosomethylethylamine	3600	U	
66-27-3	Methyl methanesulfonate	3600	U	
55-18-5	N-Nitrosodiethylamine	3600	U	
62-50-0	Ethyl methanesulfonate	3600	U	
108-95-2	Phenol	3600	U	
62-53-3	Aniline	3600	U	
76-01-7	Pentachloroethane	3600	U	
111-44-4	bis(2-Chloroethyl) Ether	7200	U	
95-57-8	2-Chlorophenol	3600	U	
541-73-1	1, 3-Dichlorobenzene	3600	U	
100-44-7	Benzyl Chloride	3600	U	
106-46-7	1, 4-Dichlorobenzene	3600	U	
100-51-6	Benzyl Alcohol	3600	U	
95-50-1	1, 2-Dichlorobenzene	3600	U	
95-48-7	2-Methylphenol	3600	U	
108-60-1	bis(2-Chloroisopropyl) Ether	3600	U	
108-39-4	3-Methylphenol	3600	U	
106-44-5	4-Methylphenol	3600	U	
930-55-2	N-Nitrosopyrrolidine	3600	U	
59-89-2	N-Nitrosomorpholine	3600	U	
98-86-2	Acetophenone	3600	U	
621-64-7	N-Nitroso-Di-n-Propylamine	3600	U	
636-21-5	O-Toluidine hydrochloride	3600	U	
67-72-1	Hexachloroethane	3600	U	
98-95-3	Nitrobenzene	3600	U	
100-75-4	N-Nitroscopiperidine	3600	U	
78-59-1	Isophorone	3600	U	
88-75-5	2-Nitrophenol	3600	U	
105-67-9	2, 4-Dimethylphenol	3600	U	

FORM I SV-1

1/87 Rev.

108-70-3	1, 3, 5-Trichlorobenzene	3600	U
98-87-3	Benzal Chloride	3600	U
65-85-0	Benzoic Acid	36000	U
111-91-1	bis(2-Chloroethoxy)Methane	3600	U
120-83-2	2, 4-Dichlorophenol	3600	U
120-82-1	1, 2, 4-Trichlorobenzene	3600	U
91-20-3	Naphthalene	23000	
106-47-8	4-Chloroaniline	3600	U
87-65-0	2, 6-Dichlorophenol	7200	U
95-54-5	c-Phenylenediamine	3600	U
122-09-8	dimethylphenylethylamine	3600	U
1888-71-7	Hexachloropropene	3600	U
87-68-3	Hexachlorobutadiene	3600	U
87-61-6	1, 2, 3-Trichlorobenzene	3600	U
98-07-7	Benzotrichloride	7200	U
924-16-3	N-Nitroso-di-n-butylamine	3600	U
59-50-7	4-Chloro-3-Methylphenol	3600	U
106-50-3	P-Phenylenediamine	3600	U
94-59-7	Safrole	3600	U
106-50-3	m-Phenylenediamine	3600	U
91-57-6	2-Methylnaphthalene	17000	
90-12-0	1-Methylnaphthalene	22000	
95-94-3	1, 2, 4, 5-Tetrachlorobenzene	3600	U
634-90-2	1, 2, 3, 5-Tetrachlorobenzene	3600	U
77-47-4	Hexachlorocyclopentadiene	3600	U
88-06-2	2, 4, 6-Trichlorophenol	7200	U
95-95-4	2, 4, 5-Trichlorophenol	7200	U
120-58-1	Isosafrole	7200	U
91-58-7	2-Chloronaphthalene	3600	U
90-13-1	1-Chloronaphthalene	3600	U
634-66-2	1, 2, 3, 4-Tetrachlorobenzene	3600	U
88-74-4	2-Nitroaniline	3600	U
130-15-4	1, 4-Naphthoquinone	7200	U
100-25-4	1, 4-Dinitrobenzene	7200	U
131-11-3	Dimethyl Phthalate	3600	U
208-96-8	Acenaphthylene	6100	
606-20-2	2, 6-Dinitrotoluene	3600	U

ROB Bla14

1C
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA3B1214
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SDG No.: <u>578</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>473668</u>	
Sample wt/vol: <u>30.1</u> (g/mL) G	Lab File ID: <u>GD073668C20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>01/07/92</u>	
% Moisture: not dec. <u>9</u> dec. <u> </u>	Date Extracted: <u>01/08/92</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>01/11/92</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: <u> </u>	Dilution Factor: <u>10</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	Q
99-09-2	3-Nitroaniline	7200	U
83-32-9	Acenaphthene	6100	
51-28-5	2, 4-Dinitrophenol	14000	U
100-02-7	4-Nitrophenol	3600	U
132-64-9	Dibenzofuran	7300	
121-14-2	2, 4-Dinitrotoluene	3600	U
608-93-5	Pentachlorobenzene	3600	U
91-59-8	2-Naphthylamine	7200	U
134-32-7	1-Naphthylamine	7200	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	7200	U
84-66-2	Diethylphthalate	3600	U
297-97-2	Zinophos	3600	U
7005-72-3	4-Chlorophenyl-phenylether	3600	U
86-73-7	Fluorene	17000	
100-01-6	4-Nitroaniline	7200	U
99-55-8	5-Nitro-o-toluidine	7200	U
122-66-7	1, 2-Diphenylhydrazine	3600	U
534-52-1	4, 6-Dinitro-2-Methylphenol	11000	U
86-30-6	N-Nitrosodiphenylamine (1)	3600	U
122-39-4	Diphenylamine	3600	U
99-35-4	1, 3, 5-Trinitrobenzene	7200	U
62-44-2	Phenacetin	3600	U
101-55-3	4-Bromophenyl-phenylether	3600	U
2303-16-4	Diallate	3600	U
60-51-5	Dimethoate	3600	U
118-74-1	Hexachlorobenzene	3600	U
92-67-1	4-Aminobiphenyl	3600	U
23950-58-5	Pronamide	3600	U
87-86-5	Pentachlorophenol	7200	U
82-68-8	Pentachloronitrobenzene	3600	U
85-01-8	Phenanthrene	59000	E
120-12-7	Anthracene	14000	
84-74-2	Di-n-Butylphthalate	3600	U
91-80-5	Methapyrilene	7200	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	18000	U
206-44-0	Fluoranthene	49000	
92-87-5	Benzidine	3600	U
129-00-0	Pyrene	42000	
60-11-7	p-Dimethylaminoazobenzene	3600	U
510-15-6	Chlorobenzilate	3600	U
119-93-7	3, 3'-Dimethylbenzidine	7200	U
85-68-7	Butylbenzylphthalate	3600	U
53-96-3	2-Acetylaminofluorene	3600	U
101-14-4	Methylene-bis(2-Chloroaniline	3600	U
91-94-1	3, 3'-Dichlorobenzidine	3600	U
106-51-4	3, 3'-Dimethoxybenzidine	3600	U
56-55-3	Benzo(a)Anthracene	17000	
218-01-9	Chrysene	18000	
117-81-7	bis(2-Ethylhexyl) Phthalate	880	J
117-84-0	Di-n-Octyl Phthalate	3600	U
205-99-2	Benzo(b) Fluoranthene	26000	X
57-97-6	7, 12-Dimethylbenzanthracene	3600	U
207-08-9	Benzo(k) Fluoranthene	26000	X
50-32-8	Benzo(a) Pyrene	15000	
56-49-5	3-Methylchloranthrene	3600	U
224-42-0	Dibenzo(a, j) acridine	3600	U
193-39-5	Indeno(1, 2, 3-cd) Pyrene	6600	
53-70-3	Dibenz(a, h) Anthracene	2100	J
191-24-2	Benzo(g, h, i) Perylene	7600	

(1) - Cannot be separated from Diphenylamine

ROAB1214

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB1B0406
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465527</u>	
Sample wt/vol: <u>10.3</u> (g/mL) <u>G</u>	Lab File ID: <u>GH065527A22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/22/91</u>	
% Moisture: not dec. <u>17</u> dec. _____	Date Extracted: <u>11/27/91</u>	
Extraction: (SepP/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/09/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg)	UG/KG	Q
62-75-9	-N-Nitrosodimethylamine	390	U	
110-86-1	-Pyridine	390	U	
97-63-2	-Ethyl methacrylate	390	U	
123-63-7	-Paraldehyde	390	U	
109-06-8	-2-Picoline	790	U	
10595-95-6	-Nitrosomethylethylamine	390	U	
66-27-3	-Methyl methanesulfonate	390	U	
55-18-5	-N-Nitrosodiethylamine	390	U	
62-50-0	-Ethyl methanesulfonate	390	U	
108-95-2	-Phenol	390	U	
62-53-3	-Aniline	390	U	
76-01-7	-Pentachloroethane	390	U	
111-44-4	-bis(2-Chloroethyl)Ether	790	U	
95-57-8	-2-Chlorophenol	390	U	
541-73-1	-1, 3-Dichlorobenzene	390	U	
106-46-7	-1, 4-Dichlorobenzene	390	U	
100-51-6	-Benzyl Alcohol	390	U	
95-50-1	-1, 2-Dichlorobenzene	390	U	
95-48-7	-2-Methylphenol	390	U	
108-60-1	-bis(2-Chloroisopropyl)Ether	390	U	
108-39-4	-3-Methylphenol	390	U	
106-44-5	-4-Methylphenol	390	U	
930-55-2	-N-Nitrosopyrrolidine	390	U	
59-89-2	-N-Nitrosomorpholine	390	U	
98-86-2	-Acetophenone	390	U	
621-64-7	-N-Nitroso-Di-n-Propylamine	390	U	
636-21-5	-o-Toluidine hydrochloride	390	U	
67-72-1	-Hexachloroethane	390	U	
98-95-3	-Nitrobenzene	390	U	
100-75-4	-N-Nitrosopiperidine	390	U	
78-59-1	-Isophorone	390	U	
88-75-5	-2-Nitrophenol	390	U	
105-67-9	-2, 4-Dimethylphenol	390	U	
108-70-3	-1, 3, 5-Trichlorobenzene	390	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3	Benzal Chloride	390	U
65-85-0	Benzoic Acid	3900	U
111-91-1	bis(2-Chloroethoxy)Methane	390	U
120-83-2	2,4-Dichlorophenol	390	U
120-82-1	1,2,4-Trichlorobenzene	390	U
91-20-3	Naphthalene	390	U
106-47-8	4-Chloroaniline	390	U
87-65-0	2,6-Dichlorophenol	790	U
95-54-5	o-Phenylenediamine	390	U
122-09-8	dimethylphenylethylamine	390	U
1888-71-7	Hexachloropropene	390	U
87-68-3	Hexachlorobutadiene	390	U
87-61-6	1,2,3-Trichlorobenzene	390	U
98-07-7	Benzotrichloride	790	U
924-16-3	N-Nitroso-di-n-butylamine	390	U
59-50-7	4-Chloro-3-Methylphenol	390	U
106-50-3	P-Phenylenediamine	390	U
94-59-7	Safrole	390	U
106-50-3	m-Phenylenediamine	390	U
91-57-6	2-Methylnaphthalene	390	U
90-12-0	1-Methylnaphthalene	50	J
95-94-3	1,2,4,5-Tetrachlorobenzene	390	U
634-90-2	1,2,3,5-Tetrachlorobenzene	390	U
77-47-4	Hexachlorocyclopentadiene	390	U
88-06-2	2,4,6-Trichlorophenol	790	U
95-95-4	2,4,5-Trichlorophenol	790	U
120-58-1	Isosafrole	790	U
91-58-7	2-Chloronaphthalene	390	U
90-13-1	1-Chloronaphthalene	390	U
634-66-2	1,2,3,4-Tetrachlorobenzene	390	U
88-74-4	2-Nitroaniline	390	U
130-15-4	1,4-Naphthoquinone	790	U
100-25-4	1,4-Dinitrobenzene	790	U
131-11-3	Dimethyl Phthalate	390	U
208-96-8	Acenaphthylene	160	J
606-20-2	2,6-Dinitrotoluene	390	U

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP	Contract: 500077	ROB1B0406
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 147
Matrix: (soil/water) SOIL	Lab Sample ID: 465527	
Sample wt/vol: 30.3 (g/mL) G	Lab File ID: GH065527A22	
Level: (low/med) LOW	Date Received: 11/22/91	
% Moisture: not dec. 12 dec.	Date Extracted: 11/27/91	
Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 12/09/91	
GPC Cleanup: (Y/N) N	pH:	Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
99-09-2	3-Nitroaniline	790	U	
83-32-9	Acenaphthene	50	J	
51-28-5	2,4-Dinitrophenol	1600	U	
100-02-7	4-Nitrophenol	390	U	
132-64-9	Dibenzofuran	390	U	
121-14-2	2,4-Dinitrotoluene	390	U	
608-93-5	Pentachlorobenzene	390	U	
91-59-8	2-Naphthylamine	790	U	
134-32-7	1-Naphthylamine	790	U	
58-90-2	2,3,4,6-Tetrachlorophenol	790	U	
84-66-2	Diethylphthalate	390	U	
297-97-2	Zinophos	390	U	
7005-72-3	4-Chlorophenyl-phenylether	390	U	
86-73-7	Fluorene	79	J	
100-01-6	4-Nitroaniline	790	U	
99-55-8	5-Nitro-o-toluidine	790	U	
122-66-7	1,2-Diphenylhydrazine	390	U	
534-52-1	4,6-Dinitro-2-Methylphenol	1200	U	
86-30-6	N-Nitrosodiphenylamine (1)	390	U	
122-39-4	Diphenylamine	390	U	
99-35-4	1,3,5-Trinitrobenzene	790	U	
62-44-2	Phenacetin	390	U	
101-55-3	4-Bromophenyl-phenylether	390	U	
2303-16-4	Diallate	390	U	
60-51-5	Dimethoate	390	U	
118-74-1	Hexachlorobenzene	390	U	
92-67-1	4-Aminobiphenyl	390	U	
23950-58-5	Pronamide	390	U	
87-86-5	Pentachlorophenol	720	J	
82-68-8	Pentachloronitrobenzene	390	U	
85-01-8	Phenanthrene	500		
120-12-7	Anthracene	190	J	
84-74-2	Di-n-Butylphthalate	390	U	
91-80-5	Methapyrilene	790	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	1900	U
206-44-0	Fluoranthene	760	
92-87-5	Benzidine	390	U
129-00-0	Pyrene	770	
60-11-7	p-Dimethylaminoazobenzene	390	U
510-15-6	Chlorobenzilate	390	U
119-93-7	3, 3'-Dimethylbenzidine	790	U
85-68-7	Butylbenzylphthalate	390	U
53-96-3	2-Acetylaminofluorene	390	U
101-14-4	Methylene-bis(2-Chloroaniline	390	U
91-94-1	3, 3'-Dichlorobenzidine	390	U
106-51-4	3, 3'-Dimethoxybenzidine	390	U
56-55-3	Benzo(a)Anthracene	510	
117-81-7	bis(2-Ethylhexyl)Phthalate	390	U
218-01-9	Chrysene	490	
117-84-0	Di-n-Octyl Phthalate	390	U
205-99-2	Benzo(b)Fluoranthene	1100	X
57-97-6	7, 12-Dimethylbenzanthracene	390	U
207-08-9	Benzo(k)Fluoranthene	1100	X
50-32-8	Benzo(a)Pyrene	660	
56-49-5	3-Methylchloranthrene	390	U
224-42-0	Dibenzo(a,j)acridine	390	U
193-39-5	Indeno(1,2,3-cd)Pyrene	330	J
53-70-3	Dibenz(a,h)Anthracene	120	J
191-24-2	Benzo(g,h,i)Perylene	350	J

(1) - Cannot be separated from Diphenylamine

465527

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB-DPA1DL
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465553</u>	
Sample wt/vol: <u>.30.2</u> (g/mL) G	Lab File ID: <u>GD065553C22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/22/91</u>	
* Moisture: not dec. <u>17</u> dec. _____	Date Extracted: <u>11/27/91</u>	
Extraction: (SepP/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/11/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>4.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
62-75-9	N-Nitrosodimethylamine	1600	U	
110-86-1	Pyridine	1600	U	
97-63-2	Ethyl methacrylate	1600	U	
123-63-7	Paraldehyde	1600	U	
109-06-8	2-Picoline	3200	U	
10595-95-6	Nitrosomethyl ethylamine	1600	U	
66-27-3	Methyl methanesulfonate	1600	U	
55-18-5	N-Nitrosodiethylamine	1600	U	
62-50-0	Ethyl methanesulfonate	1600	U	
108-95-2	Phenol	1600	U	
62-53-3	Aniline	1600	U	
76-01-7	Pentachloroethane	1600	U	
111-44-4	bis(2-Chloroethyl) Ether	3200	U	
95-57-8	2-Chlorophenol	1600	U	
541-73-1	1,3-Dichlorobenzene	1600	U	
106-46-7	1,4-Dichlorobenzene	1600	U	
100-51-6	Benzyl Alcohol	1600	U	
95-50-1	1,2-Dichlorobenzene	1600	U	
95-48-7	2-Methylphenol	1600	U	
108-60-1	bis(2-Chloroisopropyl) Ether	1600	U	
108-39-4	3-Methylphenol	1600	U	
106-44-5	4-Methylphenol	1600	U	
930-55-2	N-Nitroso pyrrolidine	1600	U	
59-89-2	N-Nitrosomorpholine	1600	U	
98-86-2	Acetophenone	1600	U	
621-64-7	N-Nitroso-Di-n-Propylamine	1600	U	
636-21-5	O-Toluidine hydrochloride	1600	U	
67-72-1	Hexachloroethane	1600	U	
98-95-3	Nitrobenzene	1600	U	
100-75-4	N-Nitrosopiperidine	1600	U	
78-59-1	Isophorone	1600	U	
88-75-5	2-Nitrophenol	1600	U	
105-67-9	2,4-Dimethylphenol	1600	U	
108-70-3	1,3,5-Trichlorobenzene	1600	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-67-3	Benzal Chloride		1600	U
65-85-0	Benzoic Acid		16000	U
111-91-1	bis(2-Chloroethoxy)Methane		1600	U
120-83-2	2,4-Dichlorophenol		1600	U
120-82-1	1,2,4-Trichlorobenzene		1600	U
91-20-3	Naphthalene		1700	D
106-47-8	4-Chloroaniline		1600	U
87-65-0	2,6-Dichlorophenol		3200	U
95-54-5	O-Phenylenediamine		1600	U
122-09-8	dimethylphenylethylamine		1600	U
1888-71-7	Hexachloropropene		1600	U
87-68-3	Hexachlorobutadiene		1600	U
87-61-6	1,2,3-Trichlorobenzene		1600	U
98-07-7	Benzotrichloride		1600	U
924-16-3	N-Nitroso-di-n-butylamine		3200	U
59-50-7	4-Chloro-3-Methylphenol		1600	U
106-50-3	P-Phenylenediamine		1600	U
94-59-7	Safrole		1600	U
106-50-3	m-Phenylenediamine		1600	U
91-57-6	2-Methylnaphthalene		1600	U
90-12-0	1-Methylnaphthalene		730	DJ
95-94-3	1,2,4,5-Tetrachlorobenzene		950	DJ
634-90-2	1,2,3,5-Tetrachlorobenzene		1600	U
77-47-4	Hexachlorocyclopentadiene		1600	U
88-06-2	2,4,6-Trichlorophenol		3200	U
95-95-4	2,4,5-Trichlorophenol		3200	U
120-58-1	Isosafrole		3200	U
91-58-7	2-Chloronaphthalene		1600	U
90-13-1	1-Chloronaphthalene		1600	U
634-66-2	1,2,3,4-Tetrachlorobenzene		1600	U
88-74-4	2-Nitroaniline		1600	U
130-15-4	1,4-Naphthoquinone		3200	U
100-25-4	1,4-Dinitrobenzene		3200	U
131-11-3	Dimethyl Phthalate		1600	U
208-96-8	Acenaphthylene		500	DJ
606-20-2	2,6-Dinitrotoluene		1600	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROB-DPA1DL

Lab Name: COMPUGHEM RTP	Contract: 500077	
Lab Code: COMPU	Case No.: 24105	SAS No.:
Matrix: (soil/water) SOIL	Lab Sample ID: 465553	
Sample wt/vol: 30.2 (g/mL) G	Lab File ID: GD065553C22	
Level: (low/med) LOW	Date Received: 11/22/91	
% Moisture: not dec. 12 dec.	Date Extracted: 11/27/91	
Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 12/11/91	
GPC Cleanup: (Y/N) N	pH:	Dilution Factor: 4.0

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	ug/kg

99-09-2	3-Nitroaniline	3200	U
83-32-9	Acenaphthene	2100	D
51-28-5	2, 4-Dinitrophenol	6200	U
100-02-7	4-Nitrophenol	1600	U
132-64-9	Dibenzofuran	1600	U
121-14-2	2, 4-Dinitrotoluene	1900	D
608-93-5	Pentachlorobenzene	1600	U
91-59-8	2-Naphthylamine	1600	U
134-32-7	1-Naphthylamine	3200	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	3200	U
84-66-2	Diethylphthalate	1600	U
297-97-2	Zinophos	1600	U
7005-72-3	4-Chlorophenyl-phenylether	1600	U
86-73-7	Fluorene	3400	D
100-01-6	4-Nitroaniline	3200	U
99-55-8	5-Nitro-o-toluidine	3200	U
122-66-7	1, 2-Diphenylhydrazine	1600	U
534-52-1	4, 6-Dinitro-2-Methylphenol	4700	U
86-30-6	N-Nitrosodiphenylamine (1)	1600	U
122-39-4	Diphenylamine	1600	U
99-35-4	1, 3, 5-Trinitrobenzene	3200	U
62-44-2	Phenacetin	1600	U
101-55-3	4-Bromophenyl-phenylether	1600	U
2303-16-4	Diallate	1600	U
60-51-5	Dimethoate	1600	U
118-74-1	Hexachlorobenzene	1600	U
92-67-1	4-Aminobiphenyl	1600	U
23950-58-5	Pronamide	1600	U
87-86-5	Pentachlorophenol	2300	DJ
82-68-8	Pentachloronitrobenzene	1600	U
85-01-8	Phenanthrrene	13000	D
120-12-7	Anthracene	10000	D
84-74-2	Di-n-Butylphthalate	1600	U
91-80-5	Methapyrilene	3200	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	7700	U
206-44-0	Fluoranthene	16000	D
92-87-5	Benzidine	1600	U
129-00-0	Pyrene	13000	D
60-11-7	p-Dimethylaminoazobenzene	1600	U
510-15-6	Chlorobenzilate	1600	U
119-93-7	3, 3'-Dimethylbenzidine	3200	U
85-68-7	Butylbenzylphthalate	1600	U
53-96-3	2-Acetylaminofluorene	1600	U
101-14-4	Methylene-bis(2-Chloroaniline)	1600	U
91-94-1	3, 3'-Dichlorobenzidine	1600	U
106-51-4	3, 3'-Dimethoxybenzidine	1600	U
56-55-3	Benzo(a)Anthracene	7300	D
218-01-9	Chrysene	6900	D
117-81-7	bis(2-Ethylhexyl) Phthalate	270	DJ
117-84-0	Di-n-Octyl Phthalate	1600	U
205-99-2	Benzo(b) Fluoranthene	13000	D
57-97-6	7, 12-Dimethylbenzanthracene	1600	U
207-08-9	Benzo(k) Fluoranthene	13000	D
50-32-8	Benzo(a) Pyrene	5700	D
56-49-5	3-Methylchloranthrene	1600	U
224-42-0	Dibenzo(a, j) acridine	1600	U
193-39-5	Indeno(1, 2, 3-cd) Pyrene	3100	D
53-70-3	Dibenz(a, h) Anthracene	880	DJ
191-24-2	Benzo(g, h, i) Perylene	3400	D

(1) - Cannot be separated from Diphenylamine

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB2B0002
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465879</u>	
Sample wt/vol: <u>.30.2</u> (g/mL) <u>G</u>	Lab File ID: <u>GJ065879B22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/23/91</u>	
% Moisture: not dec. <u>33</u> dec. _____	Date Extracted: <u>11/27/91</u>	
Extraction: (SepF/Cont/Sonic) <u>SONIC</u>	Date Analyzed: <u>12/09/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg)	UG/KG	Q
62-75-9	N-Nitrosodimethylamine	490	U	
110-86-1	Pyridine	490	U	
97-63-2	Ethyl methacrylate	490	U	
123-63-7	Paraldehyde	490	U	
109-06-8	2-Picoline	980	U	
10595-95-6	Nitrosomethylamine	490	U	
66-27-3	Methyl methanesulfonate	490	U	
55-18-5	N-Nitrosodiethylamine	490	U	
62-50-0	Ethyl methanesulfonate	490	U	
108-95-2	Phenol	490	U	
62-53-3	Aniline	490	U	
76-01-7	Pentachloroethane	490	U	
111-44-4	bis(2-Chloroethyl)Ether	69	J	
95-57-8	2-Chlorophenol	490	U	
541-73-1	1,3-Dichlorobenzene	490	U	
106-46-7	1,4-Dichlorobenzene	490	U	
100-51-6	Benzyl Alcohol	490	U	
95-50-1	1,2-Dichlorobenzene	490	U	
95-48-7	2-Methylphenol	490	U	
108-60-1	bis(2-Chloroisopropyl)Ether	490	U	
108-39-4	3-Methylphenol	50	J	
106-44-5	4-Methylphenol	50	J	
930-55-2	N-Nitroso-pyrrolidine	490	U	
59-89-2	N-Nitrosomorpholine	490	U	
98-86-2	Acetophenone	490	U	
621-64-7	N-Nitroso-Di-n-Propylamine	490	U	
636-21-5	o-Tolidine hydrochloride	490	U	
67-72-1	Hexachloroethane	490	U	
98-95-3	Nitrobenzene	490	U	
100-75-4	N-Nitrosopiperidine	490	U	
78-59-1	Iscophorone	490	U	
88-75-5	2-Nitrophenol	490	U	
105-67-9	2,4-Dimethylphenol	490	U	
108-70-3	1,3,5-Trichlorobenzene	490	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3	Benzal Chloride	490	U
65-85-0	Benzoic Acid	85	J
111-91-1	bis(2-Chloroethoxy)Methane	490	U
120-83-2	2,4-Dichlorophenol	490	U
120-82-1	1,2,4-Trichlorobenzene	490	U
91-20-3	Naphthalene	220	J
106-47-8	4-Chloroaniline	490	U
87-65-0	2,6-Dichlorophenol	980	U
95-54-5	o-Phenylenediamine	490	U
122-09-8	dimethylphenylethylamine	490	U
1888-71-7	Hexachloropropene	490	U
87-68-3	Hexachlorobutadiene	490	U
87-61-6	1,2,3-Trichlorobenzene	490	U
98-07-7	Benzotrichloride	980	U
924-16-3	N-Nitroso-di-n-butylamine	490	U
59-50-7	4-Chloro-3-Methylphenol	490	U
106-50-3	p-Phenylenediamine	490	U
94-59-7	Safrole	490	U
106-50-3	m-Phenylenediamine	490	U
91-57-6	2-Methylnaphthalene	110	J
90-12-0	1-Methylnaphthalene	210	J
95-94-3	1,2,4,5-Tetrachlorobenzene	490	U
634-90-2	1,2,3,5-Tetrachlorobenzene	490	U
77-47-4	Hexachlorocyclopentadiene	490	U
88-06-2	2,4,6-Trichlorophenol	980	U
95-95-4	2,4,5-Trichlorophenol	980	U
120-58-1	Icosafrole	980	U
91-58-7	2-Chloronaphthalene	490	U
90-13-1	1-Chloronaphthalene	490	U
634-66-2	1,2,3,4-Tetrachlorobenzene	490	U
88-74-4	2-Nitroaniline	490	U
130-15-4	1,4-Naphthoquinone	980	U
100-25-4	1,4-Dinitrobenzene	980	U
131-11-3	Dimethyl Phthalate	490	U
208-96-8	Acenaphthylene	750	
606-20-2	2,6-Dinitrotoluene	490	U

1C
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB2B0002
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465879</u>	
Sample wt/vol: <u>30.2</u> (g/mL) G	Lab File ID: <u>GJ065879B22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>11/23/91</u>	
% Moisture: not dec. <u>13</u> dec. _____	Date Extracted: <u>11/27/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/09/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
99-09-2	3-Nitroaniline	980	U
83-32-9	Acenaphthene	270	J
51-28-5	2, 4-Dinitrophenol	1900	U
100-02-7	4-Nitrophenol	490	U
132-64-9	Dibenzofuran	140	J
121-14-2	2, 4-Dinitrotoluene	490	U
608-93-5	Pentachlorobenzene	490	U
91-59-8	2-Naphthylamine	980	U
134-32-7	1-Naphthylamine	980	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	980	U
84-66-2	Diethylphthalate	490	U
297-97-2	Zinophos	490	U
7005-72-3	4-Chlorophenyl-phenylether	490	U
86-73-7	Fluorene	370	J
100-01-6	4-Nitroaniline	980	U
99-55-8	5-Nitro-o-toluidine	980	U
122-66-7	1, 2-Diphenylhydrazine	490	U
534-52-1	4, 6-Dinitro-2-Methylphenol	1500	U
86-30-6	N-Nitrosodiphenylamine (1)	490	U
122-39-4	Diphenylamine	490	U
99-35-4	1, 3, 5-Trinitrobenzene	980	U
62-44-2	Phenacetin	490	U
101-55-3	4-Bromophenyl-phenylether	490	U
2303-16-4	Diallate	490	U
60-51-5	Dimethoate	490	U
118-74-1	Hexachlorobenzene	490	U
92-67-1	4-Aminobiphenyl	490	U
23950-58-5	Pronamide	490	U
87-86-5	Pentachlorophenol	620	J
82-68-8	Pentachloronitrobenzene	490	U
85-01-8	Phenanthrone	2200	
120-12-7	Anthracene	710	
84-74-2	Di-n-Butylphthalate	85	J
91-80-5	Methapyrilene	980	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	2400	U
206-44-0	Fluoranthene	3600	
92-87-5	Benzidine	490	U
129-00-0	Pyrene	2500	
60-11-7	p-Dimethylaminoazobenzene	490	U
510-15-6	Chlorobenzilate	490	U
119-93-7	3,3'-Dimethylbenzidine	980	U
85-68-7	Butylbenzylphthalate	300	J
53-96-3	2-Acetylaminofluorene	490	U
101-14-4	Methylene-bis(2-Chloroaniline	490	U
91-94-1	3,3'-Dichlorobenzidine	490	U
106-51-4	3,3'-Dimethoxybenzidine	490	U
56-55-3	Benzo(a)Anthracene	1700	
218-01-9	Chrysene	1500	
117-81-7	bis(2-Ethylhexyl)Phthalate	330	J
117-84-0	Di-n-Octyl Phthalate	490	U
205-99-2	Benzo(b)Fluoranthene	4200	
57-97-6	7,12-Dimethylbenzanthracene	490	U
207-08-9	Benzo(k)Fluoranthene	4200	
50-32-8	Benzo(a)Pyrene	2100	
56-49-5	3-Methylchloranthrene	490	U
224-42-0	Dibenzo(a,j)acridine	490	U
193-39-5	Indeno(1,2,3-cd)Pyrene	970	
53-70-3	Dibenz(a,h)Anthracene	230	J
191-24-2	Benzo(g,h,i)Perylene	1200	

(1) -- Cannot be separated from Diphenylamine

1C
SEMITVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO

ROCO11012

Lab Name: COMPUCHEM RTP	Contract: 500077		
Lab Code: COMPU	Case No.: 24105	SAS No.: _____	SDG No.: 27
Matrix: (soil/water) SOIL	Lab Sample ID: 461432		
Sample wt/vol: 30.0 (g/mL) S	Lab File ID: G2D61432B20		
Level: (low/med) LOW	Date Received: 11/15/91		
† Moisture: not dec. 10 dec.	Date Extracted: 11/15/91		
Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 11/23/91		
GPC Cleanup: (Y/N) N	pH:	Dilution Factor: 5.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg		Q
		3700	U	
99-09-2	3-Nitroaniline	240	J ✓	
83-32-9	Acenaphthene	7200	U	
51-28-5	2, 4-Dinitrophenol	1800	U	
100-02-7	4-Nitrophenol	270	J ✓	
132-64-9	Dibenzofuran	1800	U	
121-14-2	2, 4-Dinitrotoluene	1800	U	
608-93-5	Pentachlorobenzene	1800	U	
91-59-8	2-Naphthylamine	3700	U	
134-32-7	1-Naphthylamine	3700	U	
58-90-2	2, 3, 4, 6-Tetrachlorophenol	3700	U	
84-66-2	Diethylphthalate	1800	U	
297-97-2	Zinophos	1800	U	
7005-72-3	4-Chlorophenyl-phenylether	1800	U	
86-73-7	Fluorene	1200	J ✓	
100-01-6	4-Nitroaniline	3700	U	
99-55-8	5-Nitro-o-toluidine	3700	U	
122-66-7	1, 2-Diphenylhydrazine	1800	U	
534-52-1	4, 6-Dinitro-2-Methylphenol	5500	U	
86-30-6	N-Nitrosodiphenylamine (1)	1800	U	
122-39-4	Diphenylamine	1800	U	
99-35-4	1, 3, 5-Trinitrobenzene	3700	U	
62-44-2	Phenacetin	1800	U	
101-55-3	4-Bromophenyl-phenylether	1800	U	
2303-16-4	Diallate	1800	U	
60-51-5	Dimethoate	1800	U	
118-74-1	Hexachlorobenzene	1800	U	
92-67-1	4-Aminobiphenyl	1800	U	
23950-58-5	Pronamide	1800	U	
87-86-5	Pentachlorophenol	3700	U	
82-68-8	Pentachloronitrobenzene	1800	U	
85-01-8	Phenanthrene	13000	J ✓	
120-12-7	Anthracene	1600	J ✓	
84-74-2	Di-n-Butylphthalate	1800	U	
91-80-5	Methapyrilene	3700	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

98-87-3	Benzal Chloride	1800	U
65-85-0	Benzoic Acid	18000	U
111-91-1	bis (2-Chloroethoxy)Methane	1800	U
120-83-2	2, 4-Dichlorophenol	1800	U
120-82-1	1, 2, 4-Trichlorobenzene	1800	U
91-20-3	Naphthalene	230	J
106-47-8	4-Chloroaniline	1800	U
87-65-0	2, 6-Dichlorophenol	3700	U
95-54-5	c-Phenylenediamine	1800	U
122-09-8	dimethylphenylethylamine	1800	U
1888-71-7	Hexachloropropene	1800	U
87-68-3	Hexachlorobutadiene	1800	U
87-61-6	1, 2, 3-Trichlorobenzene	1800	U
98-07-7	Benzotrichloride	3700	U
924-16-3	N-Nitroso-di-n-butylamine	1800	U
59-50-7	4-Chloro-3-Methylphenol	1800	U
106-50-3	p-Phenylenediamine	1800	U
94-59-7	Safrole	1800	U
106-50-3	m-Phenylenediamine	1800	U
91-57-6	2-Methylnaphthalene	1800	U
90-12-0	1-Methylnaphthalene	330	J
95-94-3	1, 2, 4, 5-Tetrachlorobenzene	1800	U
634-90-2	1, 2, 3, 5-Tetrachlorobenzene	1800	U
77-47-4	Hexachlorocyclopentadiene	1800	U
88-06-2	2, 4, 6-Trichlorophenol	3700	U
95-95-4	2, 4, 5-Trichlorophenol	3700	U
120-58-1	Icosafrole	3700	U
91-58-7	2-Chloronaphthalene	1800	U
90-13-1	1-Chloronaphthalene	1800	U
634-66-2	1, 2, 3, 4-Tetrachlorobenzene	1800	U
88-74-4	2-Nitroaniline	1800	U
130-15-4	1, 4-Naphthoquinone	3700	U
100-25-4	1, 4-Dinitrobenzene	3700	U
131-11-3	Dimethyl Phthalate	1800	U
208-96-8	Acenaphthylene	2200	
606-20-2	2, 6-Dinitrotoluene	1800	U

461432

FORM I SV-1

1/87 Rev.

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO

ROCO11012

Lab Name:	COMPUCHEM RTP	Contract:	500077
Lab Code:	COMPU	Case No.:	24105
Matrix:	(soil/water) SOIL	Lab Sample ID:	461432
Sample wt/vol:	30.0 (g/mL) G	Lab File ID:	G2D61432B20
Level:	(low/med) LOW	Date Received:	11/15/91
Moisture:	not dec. 10 dec.	Date Extracted:	11/15/91
Extraction:	(SepP/Cont/Sonc)	Date Analyzed:	11/23/91
GPC Cleanup:	(Y/N) N	pH:	Dilution Factor: 5.0

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) ug/kg	Q
62-75-9	N-Nitrosodimethylamine	1800	U
110-86-1	Pyridine	1800	U
97-63-2	Ethyl methacrylate	1800	U
123-63-7	Paraldehyde	1800	U
109-06-8	2-Picoline	3700	U
10595-95-6	Nitrosomethylamine	1800	U
66-27-3	Methyl methanesulfonate	1800	U
55-18-5	N-Nitrosodiethylamine	1800	U
62-50-0	Ethyl methanesulfonate	1800	U
108-95-2	Phenol	1800	U
62-53-3	Aniline	1800	U
76-01-7	Pentachloroethane	1800	U
111-44-4	bis(2-Chloroethyl)Ether	3700	U
95-57-8	2-Chlorophenol	1800	U
541-73-1	1,3-Dichlorobenzene	1800	U
106-46-7	1,4-Dichlorobenzene	1800	U
100-51-6	Benzyl Alcohol	1800	U
95-50-1	1,2-Dichlorobenzene	1800	U
95-48-7	2-Methylphenol	1800	U
39638-32-9	bis(2-Chloroisopropyl)Ether	1800	U
108-39-4	3-Methylphenol	1800	U
106-44-5	4-Methylphenol	1800	U
930-55-2	N-Nitrospyrrolidine	1800	U
59-89-2	N-Nitrosomorpholine	1800	U
98-86-2	Acetophenone	1800	U
621-64-7	N-Nitroso-Di-n-Propylamine	1800	U
636-21-5	O-Toluidine hydrochloride	1800	U
67-72-1	Hexachloroethane	1800	U
98-95-3	Nitrobenzene	1800	U
100-75-4	N-Nitrospiperidine	1800	U
78-59-1	Isophorone	1800	U
88-75-5	2-Nitrophenol	1800	U
105-67-9	2,4-Dimethylphenol	1800	U
108-70-3	1,3,5-Trichlorobenzene	1800	U

50-18-0	Cyclophosphamide	8900	U
206-44-0	Fluoranthene	20000	✓
92-87-5	Benzidine	1800	U
129-00-0	Pyrene	19000	✓
60-11-7	p-Dimethylaminoazobenzene	1800	U
510-15-6	Chlorobenzilate	1800	U
119-93-7	3, 3'-Dimethylbenzidine	3700	U
85-68-7	Butylbenzylphthalate	1800	U
53-96-3	2-Acetylaminofluorene	1800	U
101-14-4	Methylene-bis(2-Chloroaniline)	1800	U
91-94-1	3, 3'-Dichlorobenzidine	1800	U
119-90-4	3, 3'-Dimethoxybenzidine	1800	U
56-55-3	Benzo(a)Anthracene	11000	✓
218-01-9	Chrysene	13000	✓
117-81-7	bis(2-Ethylhexyl) Phthalate	260	J ✓
117-84-0	Di-n-Octyl Phthalate	1800	U
205-99-2	Benzo(b) Fluoranthene	20000	X ✓
57-97-6	7, 12-Dimethylbenzanthracene	1800	U
207-08-9	Benzo(k) Fluoranthene	20000	X ✓
50-32-8	Benzo(a) Pyrene	10000	✓
56-49-5	3-Methylchloranthrene	1800	U
224-42-0	Dibenzo(a, j) acridine	1800	U
193-39-5	Indeno(1, 2, 3-cd) Pyrene	3600	✓
53-70-3	Dibenz(a, h) Anthracene	1100	J ✓
191-24-2	Benzo(g, h, i) Perylene	3300	✓

(1) - Cannot be separated from Diphenylamine

441432

FORM I SV-3

1/87 Rev.

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE N

Lab Name: COMPUCHEM, RTP	Contract: 500077	ROC021214
Lab Code: COMPU	Case No.: 24105	SAS No.: SDG No.: 27
Matrix: (soil/water) SOIL		Lab Sample ID: 461152
Sample wt/vol: 30.0 (g/mL) G		Lab File ID: GRJ61152A20
Level: (low/med) LOW		Date Received: 11/06/91
% Moisture: not dec. 18 dec.		Date Extracted: 12/06/91
Extraction: (SepF/Cont/Sonc)	SQNC	Date Analyzed: 12/24/91
GPC Cleanup: (Y/N) Y	pH:	Dilution Factor: 0.50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
62-75-9	N-Nitrosodimethylamine	400	U
110-86-1	Pyridine	400	U
97-63-2	Ethyl methacrylate	400	U
123-63-7	Paraldehyde	400	U
109-06-8	2-Picoline	810	U
10595-95-6	Nitrosomethylethylamine	400	U
66-27-3	Methyl methanesulfonate	400	U
55-18-5	N-Nitrosodiethylamine	400	U
62-50-0	Ethyl methanesulfonate	400	U
108-95-2	Phenol	400	U
62-53-3	Aniline	400	U
76-01-7	Pentachloroethane	400	U
111-44-4	bis(2-Chloroethyl) Ether	400	U
95-57-8	2-Chlorophenol	810	U
541-73-1	1, 3-Dichlorobenzene	400	U
106-46-7	1, 4-Dichlorobenzene	400	U
100-51-6	Benzyl Alcohol	400	U
95-50-1	1, 2-Dichlorobenzene	400	U
95-48-7	2-Methylphenol	400	U
39638-32-9	bis(2-Chloroisopropyl) Ether	400	U
108-39-4	3-Methylphenol	400	U
106-44-5	4-Methylphenol	400	U
930-55-2	N-Nitrosopyrrolidine	400	U
59-89-2	N-Nitrosomorpholine	400	U
98-86-2	Acetophenone	400	U
621-64-7	N-Nitroso-Di-n-Propylamine	400	U
636-21-5	O-Tolidine hydrochloride	400	U
67-72-1	Hexachloroethane	400	U
98-95-3	Nitrobenzene	400	U
100-75-4	N-Nitroscopiperidine	400	U
78-59-1	Isophorone	400	U
88-75-5	2-Nitrophenol	400	U
105-67-9	2, 4-Dimethylphenol	400	U
108-70-3	1, 3, 5-Trichlorobenzene	400	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3	Benzal Chloride		400	U
65-85-0	Benzoic Acid		4000	U
111-91-1	bis(2-Chloroethoxy)Methane		400	U
120-83-2	2, 4-Dichlorophenol		400	U
120-82-1	1, 2, 4-Trichlorobenzene		400	U
91-20-3	Naphthalene		400	U
106-47-8	4-Chloroaniline		400	U
87-65-0	2, 6-Dichlorophenol		400	U
95-54-5	o-Phenylenediamine		810	U
122-09-8	dimethylphenylethylamine		400	U
1888-71-7	Hexachloropropene		400	U
87-68-3	Hexachlorobutadiene		400	U
87-61-6	1, 2, 3-Trichlorobenzene		400	U
98-07-7	Benzotrichloride		400	U
924-16-3	N-Nitroso-di-n-butylamine		810	U
59-50-7	4-Chloro-3-Methylphenol		400	U
106-50-3	p-Phenylenediamine		400	U
94-59-7	Safrole		400	U
106-50-3	m-Phenylenediamine		400	U
91-57-6	2-Methylnaphthalene		400	U
90-12-0	1-Methylnaphthalene		400	U
95-94-3	1, 2, 4, 5-Tetrachlorobenzene		400	U
634-90-2	1, 2, 3, 5-Tetrachlorobenzene		400	U
77-47-4	Hexachlorocyclopentadiene		400	U
88-06-2	2, 4, 6-Trichlorophenol		400	U
95-95-4	2, 4, 5-Trichlorophenol		810	U
120-58-1	Icosafrole		810	U
91-58-7	2-Chloronaphthalene		810	U
90-13-1	1-Chloronaphthalene		400	U
634-66-2	1, 2, 3, 4-Tetrachlorobenzene		400	U
88-74-4	2-Nitroaniline		400	U
130-15-4	1, 4-Naphthoquinone		400	U
100-25-4	1, 4-Dinitrobenzene		810	U
131-11-3	Dimethyl Phthalate		810	U
208-96-8	Acenaphthylene		400	U
606-20-2	2, 6-Dinitrotoluene		400	U

461152

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE N

ROC021214

Lab Name: COMPUCHEM, RTP Contract: 500077
 Lab Code: COMPU Case No.: 24105 SAS No.: SDG No.: 27
 Matrix: (soil/water) SOIL Lab Sample ID: 461152
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GRJ61152A20
 Level: (low/med) LOW Date Received: 11/06/91
 % Moisture: not dec. 18 dec. Date Extracted: 12/06/91
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/24/91
 GPC Cleanup: (Y/N) Y pH: Dilution Factor: 0.50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
99-09-2	3-Nitroaniline	810	U
83-32-9	Acenaphthene	400	U
51-28-5	2,4-Dinitrophenol	1600	U
100-02-7	4-Nitrophenol	400	U
132-64-9	Dibenzofuran	400	U
121-14-2	2,4-Dinitrotoluene	400	U
608-93-5	Pentachlorobenzene	400	U
91-59-8	2-Naphthylamine	810	U
134-32-7	1-Naphthylamine	810	U
58-90-2	2,3,4,6-Tetrachlorophenol	810	U
84-66-2	Diethylphthalate	400	U
297-97-2	Zinophos	400	U
7005-72-3	4-Chlorophenyl-phenylether	400	U
86-73-7	Fluorene	400	U
100-01-6	4-Nitroaniline	810	U
99-55-8	5-Nitro-o-toluidine	810	U
122-66-7	1,2-Diphenylhydrazine	400	U
534-52-1	4,6-Dinitro-2-Methylphenol	1200	U
86-30-6	N-Nitrosodiphenylamine (1)	400	U
122-39-4	Diphenylamine	400	U
99-35-4	1,3,5-Trinitrobenzene	810	U
62-44-2	Phenacetin	400	U
101-55-3	4-Bromophenyl-phenylether	400	U
2303-16-4	Diallate	400	U
60-51-5	Dimethoate	400	U
118-74-1	Hexachlorobenzene	400	U
92-67-1	4-Aminobiphenyl	400	U
23950-58-5	Pronamide	400	U
87-86-5	Pentachlorophenol	810	U
82-68-8	Pentachloronitrobenzene	400	U
85-01-8	Phenanthrene	210	J
120-12-7	Anthracene	230	J
84-74-2	Di-n-Butylphthalate	400	U
91-80-5	Methapyrilene	810	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	2000	U
206-44-0	Fluoranthene	340	J
92-87-5	Benzidine	400	U
129-00-0	Pyrene	230	J
60-11-7	p-Dimethylaminocazobenzene	400	U
510-15-6	Chlorobenzilate	400	U
119-93-7	3,3'-Dimethylbenzidine	810	U
85-68-7	Butylbenzylphthalate	400	U
53-96-3	2-Acetylaminofluorene	400	U
101-14-4	Methylene-bis(2-Chloroaniline)	400	U
91-94-1	3,3'-Dichlorobenzidine	400	U
119-90-4	3,3'-Dimethoxybenzidine	400	U
56-55-3	Benzo(a)Anthracene	180	J
218-01-9	Chrysene	150	J
117-81-7	bis(2-Ethylhexyl)Phthalate	49	J
117-84-0	Di-n-Octyl Phthalate	400	U
205-99-2	Benzo(b)Fluoranthene	140	J
57-97-6	7,12-Dimethylbenzanthracene	400	U
207-08-9	Benzo(k)Fluoranthene	140	J
50-32-8	Benzo(a)Pyrene	150	J
56-49-5	3-Methylchloranthrene	400	U
224-42-0	Dibenzo(a,j)acridine	400	U
193-39-5	Indeno(1,2,3-cd)Pyrene	400	U
53-70-3	Dibenz(a,h)Anthracene	400	U
191-24-2	Benzo(g,h,i)Perylene	400	U

(1) - Cannot be separated from Diphenylamine

4611S1

FORM I SV-3

1/87 Rev.

1B
SEMITOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC-DPA1

Lab Name: COMPUCHEM RTP	Contract: 500077		
Lab Code: COMPU	Case No.: 24105	SAS No.:	SDG No.: 27
Matrix: (soil/water) SOIL	Lab Sample ID: 461186		
Sample wt/vol: 30.1 (g/mL) G	Lab File ID: GH061186B20		
Level: (low/med) LOW	Date Received: 11/06/91		
# Moisture: not dec. 46 dec.	Date Extracted: 11/08/91		
Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 11/27/91		
GPC Cleanup: (Y/N) N	pH:	Dilution Factor: 1.0	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
62-75-9	N-Nitrosodimethylamine	610	U	
110-86-1	Pyridine	610	U	
97-63-2	Ethyl methacrylate	610	U	
123-63-7	Paraldehyde	610	U	
109-06-8	2-Picoline	1200	U	
10595-95-6	Nitrosomethylethylamine	610	U	
66-27-3	Methyl methanesulfonate	610	U	
55-18-5	N-Nitrosodiethylamine	610	U	
62-50-0	Ethyl methanesulfonate	610	U	
108-95-2	Phenol	610	U	
62-53-3	Aniline	610	U	
76-01-7	Pentachloroethane	610	U	
111-44-4	bis(2-Chloroethyl) Ether	1200	U	
95-57-8	2-Chlorophenol	610	U	
541-73-1	1, 3-Dichlorobenzene	610	U	
106-46-7	1, 4-Dichlorobenzene	610	U	
100-51-6	Benzyl Alcohol	610	U	
95-50-1	1, 2-Dichlorobenzene	610	U	
95-48-7	2-Methylphenol	610	U	
39638-32-9	bis(2-Chloroisopropyl) Ether	610	U	
108-39-4	3-Methylphenol	610	U	
106-44-5	4-Methylphenol	610	U	
930-55-2	N-Nitroso- <i>D</i> - <i>n</i> -Propylamine	610	U	
59-89-2	N-Nitrosomorpholine	610	U	
98-86-2	Acetophenone	610	U	
621-64-7	N-Nitroso- <i>D</i> - <i>n</i> -Propylamine	610	U	
636-21-5	o-Toluidine hydrochloride	610	U	
67-72-1	Hexachloroethane	610	U	
98-95-3	Nitrobenzene	610	U	
100-75-4	N-Nitroscopiperidine	610	U	
78-59-1	Isophorone	610	U	
88-75-5	2-Nitrophenol	610	U	
105-67-9	2, 4-Dimethylphenol	610	U	
108-70-3	1, 3, 5-Trichlorobenzene	610	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3	Benzal Chloride	610	U
65-85-0	Benzoic Acid	6100	U
111-91-1	bis(2-Chloroethoxy)Methane	610	U
120-83-2	2,4-Dichlorophenol	610	U
120-82-1	1,2,4-Trichlorobenzene	610	U
91-20-3	Naphthalene	610	U
106-47-8	4-Chloroaniline	610	U
87-65-0	2,6-Dichlorophenol	1200	U
95-54-5	O-Phenylenediamine	610	U
122-09-8	dimethylphenylethylamine	610	U
1888-71-7	Hexachloropropene	610	U
87-68-3	Hexachlorobutadiene	610	U
87-61-6	1,2,3-Trichlorobenzene	610	U
98-07-7	Benzotrichloride	1200	U
924-16-3	N-Nitroso-di-n-butylamine	610	U
59-50-7	4-Chloro-3-Methylphenol	610	U
106-50-3	P-Phenylenediamine	610	U
94-59-7	Safrole	610	U
106-50-3	m-Phenylenediamine	610	U
91-57-6	2-Methylnaphthalene	610	U
90-12-0	1-Methylnaphthalene	610	U
95-94-3	1,2,4,5-Tetrachlorobenzene	610	U
634-90-2	1,2,3,5-Tetrachlorobenzene	610	U
77-47-4	Hexachlorocyclopentadiene	610	U
88-06-2	2,4,6-Trichlorophenol	1200	U
95-95-4	2,4,5-Trichlorophenol	1200	U
120-58-1	Iosafrole	1200	U
91-58-7	2-Chloronaphthalene	610	U
90-13-1	1-Chloronaphthalene	610	U
634-66-2	1,2,3,4-Tetrachlorobenzene	610	U
88-74-4	2-Nitroaniline	610	U
130-15-4	1,4-Naphthoquinone	1200	U
100-25-4	1,4-Dinitrobenzene	1200	U
131-11-3	Dimethyl Phthalate	610	U
208-96-8	Acenaphthylene	610	U
606-20-2	2,6-Dinitrotoluene	610	U

46302-1

FORM I SV-1

1/87 Rev.

1C
SEMITVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPCU/CHM/LTE	Contract: 599977	ROC-DPA1
Lab Code: COMPCU	Case No.: 24105	SAS No.: _____ SDG No.: 27
Matrix: (soil/water) SOIL	Lab Sample ID: 461186	
Sample wt/vol: 30.1 (g/mL) g	Lab File ID: GH061186B20	
Level: (low/med) LOW	Date Received: 11/06/91	
* Moisture: not dec. 46 dec.	Date Extracted: 11/08/91	
Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 11/22/91	
GPC Cleanup: (Y/N) N	pH: _____	Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
99-09-2	3-Nitroaniline	1200	U
83-32-9	Acenaphthene	95	J
51-28-5	2, 4-Dinitrophenol	2400	U
100-02-7	4-Nitrophenol	610	U
132-64-9	Dibenzofuran	64	J
121-14-2	2, 4-Dinitrotoluene	610	U
608-93-5	Pentachlorobenzene	610	U
91-59-8	2-Naphthylamine	1200	U
134-32-7	1-Naphthylamine	1200	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	1200	U
84-66-2	Diethylphthalate	610	U
297-97-2	Zinophos	610	U
7005-72-3	4-Chlorophenyl-phenylether	610	U
86-73-7	Fluorene	140	J
100-01-6	4-Nitroaniline	1200	U
99-55-8	5-Nitro-o-toluidine	1200	U
122-66-7	1, 2-Diphenylhydrazine	610	U
534-52-1	4, 6-Dinitro-2-Methylphenol	1800	U
86-30-6	N-Nitrosodiphenylamine (1)	610	U
122-39-4	Diphenylamine	610	U
99-35-4	1, 3, 5-Trinitrobenzene	1200	U
62-44-2	Phenacetin	610	U
101-55-3	4-Bromophenyl-phenylether	610	U
2303-16-4	Diallate	610	U
60-51-5	Dimethoate	610	U
118-74-1	Hexachlorobenzene	610	U
92-67-1	4-Aminobiphenyl	610	U
23950-58-5	Pronamide	610	U
87-86-5	Pentachlorophenol	1200	U
82-68-8	Pentachloronitrobenzene	610	U
85-01-8	Phenanthrrene	1200	U
120-12-7	Anthracene	290	J
84-74-2	Di-n-Butylphthalate	130	J
91-80-5	Methapyrilene	1200	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide		3000	U
206-44-0	Fluoranthene		1500	
92-87-5	Benzidine		610	U
129-00-0	Pyrene		1100	
60-11-7	p-Dimethylaminoazobenzene		610	U
510-15-6	Chlorobenzilate		610	U
119-93-7	3, 3'-Dimethylbenzidine		1200	U
85-68-7	Butylbenzylphthalate		610	U
53-96-3	2-Acetylaminofluorene		610	U
101-14-4	Methylene-bis(2-Chloroaniline)		610	U
91-94-1	3, 3'-Dichlorobenzidine		610	U
119-90-4	3, 3'-Dimethoxybenzidine		610	U
56-55-3	Benzo(a)Anthracene		740	
218-01-9	Chrysene		710	
117-81-7	bis(2-Ethylhexyl) Phthalate		200	J
117-84-0	Di-n-Octyl Phthalate		610	U
205-99-2	Benzo(b) Fluoranthene		450	J
57-97-6	7, 12-Dimethylbenzanthracene		610	U
207-08-9	Benzo(k) Fluoranthene		280	J
50-32-8	Benzo(a) Pyrene		620	
56-49-5	3-Methylchloranthrene		610	U
224-42-0	Dibenzo(a, j) acridine		610	U
193-39-5	Indeno(1, 2, 3-cd) Pyrene		320	J
53-70-3	Dibenz(a, h) Anthracene		100	J
191-24-2	Benzo(g, h, i) Perylene		270	J

(1) - Cannot be separated from Diphenylamine

461186

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC3B0204DL

Lab Name: COMPUCHEM ETP

Contract: 500077

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 147

Matrix: (soil/water) SOIL

Lab Sample ID: 465088

Sample wt/vol: 30.4 (g/mL) G

Lab File ID: GD065088B22

Level: (low/med) LOW

Date Received: 11/21/91

± Moisture: not dec. 8 dec. _____

Date Extracted: 11/27/91

Extraction: (SepF/Cont/Sonc) SONC

Date Analyzed: 12/13/91

GPC Cleanup: (Y/N) N pH: _____

Dilution Factor: 10

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
62-75-9	-N-Nitrosodimethylamine	3500	U	
110-86-1	-Pyridine	3500	U	
97-63-2	-Ethyl methacrylate	3500	U	
123-63-7	-Paraldehyde	3500	U	
109-06-8	-2-Picoline	7100	U	
10595-95-6	-Nitrosomethylamine	3500	U	
66-27-3	-Methyl methanesulfonate	3500	U	
55-18-5	-N-Nitrosodiethylamine	3500	U	
62-50-0	-Ethyl methanesulfonate	3500	U	
108-95-2	-Phenol	3500	U	
62-53-3	-Aniline	3500	U	
76-01-7	-Pentachloroethane	3500	U	
111-44-4	-bis(2-Chloroethyl)Ether	7100	U	
95-57-8	-2-Chlorophenol	3500	U	
541-73-1	-1, 3-Dichlorobenzene	3500	U	
106-46-7	-1, 4-Dichlorobenzene	3500	U	
100-51-6	-Benzyl Alcohol	3500	U	
95-50-1	-1, 2-Dichlorobenzene	3500	U	
95-48-7	-2-Methylphenol	3500	U	
108-60-1	-bis(2-Chloroisopropyl)Ether	3500	U	
108-39-4	-3-Methylphenol	3500	U	
106-44-5	-4-Methylphenol	3500	U	
930-55-2	-N-Nitroso-pyrrolidine	3500	U	
59-89-2	-N-Nitrosomorpholine	3500	U	
98-86-2	-Acetophenone	3500	U	
621-64-7	-N-Nitroso-Di-n-Propylamine	3500	U	
636-21-5	-o-Toluidine hydrochloride	3500	U	
67-72-1	-Hexachloroethane	3500	U	
98-95-3	-Nitrobenzene	3500	U	
100-75-4	-N-Nitroso-piperidine	3500	U	
78-59-1	-Isophorone	3500	U	
88-75-5	-2-Nitrophenol	3500	U	
105-67-9	-2, 4-Dimethylphenol	3500	U	
108-70-3	-1, 3, 5-Trichlorobenzene	3500	U	

(1) - Cannot be separated from Diphenylamine
FORM I SV-4

1/87 Rev.

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC3B0204DL

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>		
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____	SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>465088</u>		
Sample wt/vol: <u>.30.4</u> (g/mL) <u>G</u>	Lab File ID: <u>GD065088B22</u>		
Level: (low/med) <u>LOW</u>	Date Received: <u>11/21/91</u>		
Moisture: not dec. <u>8</u> dec. _____	Date Extracted: <u>11/27/91</u>		
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/13/91</u>		
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>10</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
62-75-9	N-Nitrosodimethylamine	3500	U	
110-86-1	Pyridine	3500	U	
97-63-2	Ethyl methacrylate	3500	U	
123-63-7	Paraldehyde	3500	U	
109-06-8	2-Picoline	7100	U	
10595-95-6	Nitrosomethylethylamine	3500	U	
66-27-3	Methyl methanesulfonate	3500	U	
55-18-5	N-Nitrosodiethylamine	3500	U	
62-50-0	Ethyl methanesulfonate	3500	U	
108-95-2	Phenol	3500	U	
62-53-3	Aniline	3500	U	
76-01-7	Pentachloroethane	3500	U	
111-44-4	bis(2-Chloroethyl)Ether	7100	U	
95-57-8	2-Chlorophenol	3500	U	
541-73-1	1,3-Dichlorobenzene	3500	U	
106-46-7	1,4-Dichlorobenzene	3500	U	
100-51-6	Benzyl Alcohol	3500	U	
95-50-1	1,2-Dichlorobenzene	3500	U	
95-48-7	2-Methylphenol	3500	U	
108-60-1	bis(2-Chloroisopropyl)Ether	3500	U	
108-39-4	3-Methylphenol	3500	U	
106-44-5	4-Methylphenol	3500	U	
930-55-2	N-Nitrospyrrolidine	3500	U	
59-89-2	N-Nitrosomorpholine	3500	U	
98-86-2	Acetophenone	3500	U	
621-64-7	N-Nitroso-Di-n-Propylamine	3500	U	
636-21-5	o-Toluidine hydrochloride	3500	U	
67-72-1	Hexachloroethane	3500	U	
98-95-3	Nitrobenzene	3500	U	
100-75-4	N-Nitrosopiperidine	3500	U	
78-59-1	Isophorone	3500	U	
88-75-5	2-Nitrophenol	3500	U	
105-67-9	2,4-Dimethylphenol	3500	U	
108-70-3	1,3,5-Trichlorobenzene	3500	U	

(1) - Cannot be separated from Diphenylamine
FORM I SV-4

1/87 Rev.

98-87-3	Benzal Chloride	3500	U
65-85-0	Benzoic Acid	35000	U
111-91-1	bis(2-Chloroethoxy)Methane	3500	U
120-83-2	2,4-Dichlorophenol	3500	U
120-82-1	1,2,4-Trichlorobenzene	3500	U
91-20-3	Naphthalene	1900	DJ
106-47-8	4-Chloroaniline	3500	U
87-65-0	2,6-Dichlorophenol	7100	U
95-54-5	o-Phenylenediamine	3500	U
122-09-8	dimethylphenylethylamine	3500	U
1888-71-7	Hexachloropropene	3500	U
87-68-3	Hexachlorobutadiene	3500	U
87-61-6	1,2,3-Trichlorobenzene	7100	U
98-07-7	Benzotrichloride	3500	U
924-16-3	N-Nitroso-di-n-butylamine	3500	U
59-50-7	4-Chloro-3-Methylphenol	3500	U
106-50-3	p-Phenylenediamine	3500	U
94-59-7	Safrole	3500	U
106-50-3	m-Phenylenediamine	3500	DJ
91-57-6	2-Methylnaphthalene	1600	DJ
90-12-0	1-Methylnaphthalene	2500	DJ
95-94-3	1,2,4,5-Tetrachlorobenzene	3500	U
634-90-2	1,2,3,5-Tetrachlorobenzene	3500	U
77-47-4	Hexachlorocyclopentadiene	3500	U
88-06-2	2,4,6-Trichlorophenol	7100	U
95-95-4	2,4,5-Trichlorophenol	7100	U
120-58-1	Isosafrole	7100	U
91-58-7	2-Chloronaphthalene	3500	U
90-13-1	1-Chloronaphthalene	3500	U
634-66-2	1,2,3,4-Tetrachlorobenzene	3500	U
88-74-4	2-Nitroaniline	3500	U
130-15-4	1,4-Naphthoquinone	7100	U
100-25-4	1,4-Dinitrobenzene	7100	U
131-11-3	Dimethyl Phthalate	3500	U
208-96-8	Acenaphthylene	2900	DJ
606-20-2	2,6-Dinitrotoluene	3500	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC3B0204DL

Lab Name: COMPUCHEM RTP Contract: 500077
 Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 147
 Matrix: (soil/water) SOIL Lab Sample ID: 465088
 Sample wt/vol: 30.4 (g/mL) G Lab File ID: GD065088B22
 Level: (low/med) LOW Date Received: 11/21/91
 % Moisture: not dec. 8 dec. _____ Date Extracted: 11/27/91
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/13/91
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 10

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
99-09-2	3-Nitroaniline	7100	U
83-32-9	Acenaphthene	3100	DJ
51-28-5	2,4-Dinitrophenol	14000	U
100-02-7	4-Nitrophenol	3500	U
132-64-9	Dibenzofuran	2700	DJ
121-14-2	2,4-Dinitrotoluene	3500	U
608-93-5	Pentachlorobenzene	3500	U
91-59-8	2-Naphthylamine	7100	U
134-32-7	1-Naphthylamine	7100	U
58-90-2	2,3,4,6-Tetrachlorophenol	7100	U
84-66-2	Diethylphthalate	3500	U
297-97-2	Zincphos	3500	U
7005-72-3	4-Chlorophenyl-phenylether	3500	U
86-73-7	Fluorene	5400	D
100-01-6	4-Nitroaniline	7100	U
99-55-8	5-Nitro-o-toluidine	7100	U
122-66-7	1,2-Diphenylhydrazine	3500	U
534-52-1	4,6-Dinitro-2-Methylphenol	11000	U
86-30-6	N-Nitrosodiphenylamine (1)	3500	U
122-39-4	Diphenylamine	3500	U
99-35-4	1,3,5-Trinitrobenzene	7100	U
62-44-2	Phenacetin	3500	U
101-55-3	4-Bromophenyl-phenylether	3500	U
2303-16-4	Diallate	3500	U
60-51-5	Dimethoate	3500	U
118-74-1	Hexachlorobenzene	3500	U
92-67-1	4-Aminobiphenyl	3500	U
23950-58-5	Pronamide	3500	U
87-86-5	Pentachlorophenol	7100	U
82-68-8	Pentachloronitrobenzene	3500	U
95-01-8	Phenanthrrene	27000	D
120-12-7	Anthracene	10000	D
84-74-2	Di-n-Butylphthalate	3500	U
91-80-5	Methapyrilene	7100	U

(1) - Cannot be separated from Diphenylamine
FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	17000	U
206-44-0	Fluoranthene	41000	D
92-87-5	Benzidine	3500	U
129-00-0	Pyrene	43000	D
60-11-7	p-Dimethylaminoazobenzene	3500	U
510-15-6	Chlorobenzilate	3500	U
119-93-7	3,3'-Dimethylbenzidine	7100	U
85-68-7	Butylbenzylphthalate	3500	U
53-96-3	2-Acetylaminofluorene	3500	U
101-14-4	Methylene-bis(2-Chloroaniline)	3500	U
91-94-1	3,3'-Dichlorobenzidine	3500	U
106-51-4	3,3'-Dimethoxybenzidine	24000	D
56-55-3	Benzo(a)Anthracene	22000	D
218-01-9	Chrysene	3500	U
117-81-7	bis(2-Ethylhexyl) Phthalate	3500	U
117-84-0	Di-n-Octyl Phthalate	49000	D
205-99-2	Benzo(b)Fluoranthene	3500	U
57-97-6	7,12-Dimethylbenzanthracene	49000	D
207-08-9	Benzo(k)Fluoranthene	22000	D
50-32-8	Benzo(a)Pyrene	3500	U
56-49-5	3-Methylchloranthrene	3500	U
224-42-0	Dibenzo(a,j)acridine	13000	D
193-39-5	Indeno(1,2,3-cd)Pyrene	3600	D
53-70-3	Dibenz(a,h)Anthracene	12000	D
191-24-2	Benzo(g,h,i)Perylene		

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROJIS

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>		Lab Sample ID: <u>469285</u>
Sample wt/vol: <u>.30.2</u> (g/mL) G		Lab File ID: <u>GH069285B22</u>
Level: (low/med) <u>LOW</u>		Date Received: <u>12/11/91</u>
* Moisture: not dec. <u>.22</u> dec. _____		Date Extracted: <u>12/20/91</u>
Extraction: (SepF/Cont/Sonc) <u>SONC</u>		Date Analyzed: <u>01/02/92</u>
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) UG/KG	Q
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62-75-9	N-Nitrosodimethylamine	420	U
110-86-1	Pyridine	420	U
97-63-2	Ethyl methacrylate	420	U
123-63-7	Paraldehyde	420	U
109-06-8	2-Picoline	840	U
10595-95-6	Nitrosomethylethylamine	420	U
66-27-3	Methyl methanesulfonate	420	U
55-18-5	N-Nitrosodiethylamine	420	U
62-50-0	Ethyl methanesulfonate	420	U
108-95-2	Phenol	420	U
62-53-3	Aniline	420	U
76-01-7	Pentachloroethane	420	U
111-44-4	bis(2-Chloroethyl) Ether	840	U
95-57-8	2-Chlorophenol	420	U
541-73-1	1,3-Dichlorobenzene	420	U
106-46-7	1,4-Dichlorobenzene	420	U
100-51-6	Benzyl Alcohol	420	U
95-50-1	1,2-Dichlorobenzene	420	U
95-48-7	2-Methylphenol	420	U
108-60-1	bis(2-Chloroisopropyl) Ether	420	U
108-39-4	3-Methylphenol	420	U
106-44-5	4-Methylphenol	420	U
930-55-2	N-Nitrosopyrrolidine	420	U
59-89-2	N-Nitrosomorpholine	420	U
98-86-2	Acetophenone	420	U
621-64-7	N-Nitroso-Di-n-Propylamine	420	U
636-21-5	o-Toluidine hydrochloride	420	U
67-72-1	Hexachloroethane	420	U
98-95-3	Nitrobenzene	420	U
100-75-4	N-Nitrosopiperidine	420	U
78-59-1	Isophorone	420	U
88-75-5	2-Nitrophenol	420	U
105-67-9	2,4-Dimethylphenol	420	U
108-70-3	1,3,5-Trichlorobenzene	420	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

98-87-3-----	Benzal Chloride	420	U
65-85-0-----	Benzoic Acid	4200	U
111-91-1-----	bis(2-Chloroethoxy)Methane	420	U
120-83-2-----	2,4-Dichlorophenol	420	U
120-82-1-----	1,2,4-Trichlorobenzene	420	U
91-20-3-----	Naphthalene	420	U
106-47-8-----	4-Chloroaniline	420	U
87-65-0-----	2,6-Dichlorophenol	840	U
95-54-5-----	o-Phenylenediamine	420	U
122-09-8-----	dimethylphenylethylamine	420	U
1888-71-7-----	Hexachloropropene	420	U
87-68-3-----	Hexachlorobutadiene	420	U
87-61-6-----	1,2,3-Trichlorobenzene	420	U
98-07-7-----	Benzotrichloride	840	U
924-16-3-----	N-Nitroso-di-n-butylamine	420	U
59-50-7-----	4-Chloro-3-Methylphenol	420	U
106-50-3-----	P-Phenylenediamine	420	U
94-59-7-----	Safrole	420	U
106-50-3-----	m-Phenylenediamine	420	U
91-57-6-----	2-Methylnaphthalene	420	U
90-12-0-----	1-Methylnaphthalene	420	U
95-94-3-----	1,2,4,5-Tetrachlorobenzene	420	U
634-90-2-----	1,2,3,5-Tetrachlorobenzene	420	U
77-47-4-----	Hexachlorocyclopentadiene	420	U
88-06-2-----	2,4,6-Trichlorophenol	840	U
95-95-4-----	2,4,5-Trichlorophenol	840	U
120-58-1-----	Isosafrole	840	U
91-58-7-----	2-Chloronaphthalene	420	U
90-13-1-----	1-Chloronaphthalene	420	U
634-66-2-----	1,2,3,4-Tetrachlorobenzene	420	U
88-74-4-----	2-Nitroaniline	420	U
130-15-4-----	1,4-Naphthoquinone	840	U
100-25-4-----	1,4-Dinitrobenzene	840	U
131-11-3-----	Dimethyl Phthalate	420	U
208-96-8-----	Acenaphthylene	92	J
606-20-2-----	2,6-Dinitrotoluene	420	U

1C
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM, RTP</u>	Contract: <u>500077</u>	ROJ1S
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>147</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>469285</u>	
Sample wt/vol: <u>30.2</u> (g/mL) <u>G</u>	Lab File ID: <u>GH069285B22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/11/91</u>	
Moisture: not dec. <u>22</u> dec. _____	Date Extracted: <u>12/20/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>01/02/92</u>	
GPC Cleanup: (Y/N) <u>N</u> pH: _____	Dilution Factor: <u>1.00</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
99-09-2	3-Nitroaniline	840	U	
83-32-9	Acenaphthene	48	J	
51-28-5	2, 4-Dinitrophenol	1700	U	
100-02-7	4-Nitrophenol	420	U	
132-64-9	Dibenzofuran	420	U	
121-14-2	2, 4-Dinitrotoluene	420	U	
608-93-5	Pentachlorobenzene	420	U	
91-59-8	2-Naphthylamine	840	U	
134-32-7	1-Naphthylamine	840	U	
58-90-2	2, 3, 4, 6-Tetrachlorophenol	840	U	
84-66-2	Diethylphthalate	420	U	
297-97-2	Zinophos	420	U	
7005-72-3	4-Chlorophenyl-phenylether	420	U	
86-73-7	Fluorene	54	J	
100-01-6	4-Nitroaniline	840	U	
99-55-8	5-Nitro-o-toluidine	840	U	
122-66-7	1, 2-Diphenylhydrazine	420	U	
534-52-1	4, 6-Dinitro-2-Methylphenol	1300	U	
86-30-6	N-Nitrosodiphenylamine (1)	420	U	
122-39-4	Diphenylamine	420	U	
99-35-4	1, 3, 5-Trinitrobenzene	840	U	
62-44-2	Phenacetin	420	U	
101-55-3	4-Bromophenyl-phenylether	420	U	
2303-16-4	Diallate	420	U	
60-51-5	Dimethoate	420	U	
118-74-1	Hexachlorobenzene	420	U	
92-67-1	4-Aminobiphenyl	420	U	
23950-58-5	Pronamide	420	U	
87-86-5	Pentachlorophenol	840	U	
82-68-8	Pentachloronitrobenzene	420	U	
85-01-8	Phenanthrene	590	/	
120-12-7	Anthracene	130	J	
84-74-2	Di-n-Butylphthalate	420	U	
91-80-5	Methapyrilene	840	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	2000	U
206-44-0	Fluoranthene	1400	
92-87-5	Benzidine	420	U
129-00-0	Pyrene	950	
60-11-7	p-Dimethylaminoazobenzene	420	U
510-15-6	Chlorobenzilate	420	U
119-93-7	3,3'-Dimethylbenzidine	840	U
85-68-7	Butylbenzylphthalate	420	U
53-96-3	2-Acetylaminofluorene	420	U
101-14-4	Methylene-bis(2-Chloroaniline)	420	U
91-94-1	3,3'-Dichlorobenzidine	420	U
106-51-4	3,3'-Dimethoxybenzidine	420	U
56-55-3	Benzo(a)Anthracene	610	
218-01-9	Chrysene	750	
117-81-7	bis(2-Ethylhexyl)Phthalate	56	J
117-84-0	Di-n-Octyl Phthalate	420	U
205-99-2	Benzo(b)Fluoranthene	1500	
57-97-6	7,12-Dimethylbenzanthracene	420	U
207-08-9	Benzo(k)Fluoranthene	1500	
50-32-8	Benzo(a)Pyrene	700	
56-49-5	3-Methylchloranthrene	420	U
224-42-0	Dibenzo(a,j)acridine	420	U
193-39-5	Indeno(1,2,3-cd)Pyrene	360	J
53-70-3	Dibenz(a,h)Anthracene	130	J
191-24-2	Benzo(g,h,i)Perylene	430	

(1) -- Cannot be separated from Diphenylamine

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTPContract: 500077

ROJ2S

Lab Code: COMPU Case No.: 24105 SAS No.: SDG No.: 429Matrix: (soil/water) SOIL Lab Sample ID: 469275Sample wt/vol: 30.4 (g/mL) G Lab File ID: GH069275A20Level: (low/med) LOW Date Received: 12/11/91% Moisture: not dec. 15 dec. Date Extracted: 12/20/91Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/23/91GPC Cleanup: (Y/N) N pH: Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

99-09-2	3-Nitroaniline	770	U
83-32-9	Acenaphthene	52	J
51-28-5	2, 4-Dinitrophenol	1500	U
100-02-7	4-Nitrophenol	380	U
132-64-9	Dibenzofuran	380	U
121-14-2	2, 4-Dinitrotoluene	380	U
608-93-5	Pentachlorobenzene	380	U
91-59-8	2-Naphthylamine	770	U
134-32-7	1-Naphthylamine	770	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	770	U
84-66-2	Diethylphthalate	380	U
297-97-2	Zinophos	380	U
7005-72-3	4-Chlorophenyl-phenylether	380	U
86-73-7	Fluorene	58	J
100-01-6	4-Nitroaniline	770	U
99-55-8	5-Nitro-o-toluidine	770	U
122-66-7	1, 2-Diphenylhydrazine	380	U
534-52-1	4, 6-Dinitro-2-Methylphenol	1100	U
86-30-6	N-Nitrosodiphenylamine (1)	380	U
122-39-4	Diphenylamine	380	U
99-35-4	1, 3, 5-Trinitrobenzene	770	U
62-44-2	Phenacetin	380	U
101-55-3	4-Bromophenyl-phenylether	380	U
2303-16-4	Diallate	380	U
60-51-5	Dimethoate	380	U
118-74-1	Hexachlorobenzene	380	U
92-67-1	4-Aminobiphenyl	380	U
23950-58-5	Pronamide	380	U
87-86-5	Pentachlorophenol	770	U
82-68-8	Pentachloronitrobenzene	380	U
85-01-8	Phenanthrane	770	
120-12-7	Anthracene	140	J
84-74-2	Di-n-Butylphthalate	380	U
91-80-5	Methaprylene	770	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	1900	U
206-44-0	Fluoranthene	1000	
92-87-5	Benzidine	380	U
129-00-0	Pyrene	810	
60-11-7	p-Dimethylaminoazobenzene	380	U
510-15-6	Chlorobenzilate	380	U
119-93-7	3,3'-Dimethylbenzidine	770	U
85-68-7	Butylbenzylphthalate	380	U
53-96-3	2-Acetylaminofluorene	380	U
101-14-4	Methylene-bis(2-Chloroaniline	380	U
91-94-1	3,3'-Dichlorobenzidine	380	U
106-51-4	3,3'-Dimethoxybenzidine	380	U
56-55-3	Benzo(a)Anthracene	570	
117-81-7	bis(2-Ethylhexyl)Phthalate	380	U
218-01-9	Chrysene	700	
117-84-0	Di-n-Octyl Phthalate	380	U
205-99-2	Benzo(b)Fluoranthene	580	X
57-97-6	7,12-Dimethylbenzanthracene	380	U
207-08-9	Benzo(k)Fluoranthene	580	X
50-32-8	Benzo(a)Pyrene	450	
56-49-5	3-Methylchloranthrene	380	U
224-42-0	Dibenzo(a,j)acridine	380	U
193-39-5	Indeno(1,2,3-cd)Pyrene	320	J
53-70-3	Dibenz(a,h)Anthracene	97	J
191-24-2	Benzo(g,h,i)Perylene	280	J

(1) - Cannot be separated from Diphenylamine

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTPContract: 500077

ROJ3S

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 429Matrix: (soil/water) SOIL Lab Sample ID: 469296Sample wt/vol: 30.3 (g/mL) G Lab File ID: GJ069296A20Level: (low/med) LOW Date Received: 12/11/91% Moisture: not dec. 9 dec. _____ Date Extracted: 12/20/91Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/24/91GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

62-75-9-----	N-Nitrosodimethylamine	360	U
110-86-1-----	Pyridine	360	U
97-63-2-----	Ethyl methacrylate	360	U
123-63-7-----	Paraldehyde	360	U
109-06-8-----	2-Picoline	720	U
10595-95-6-----	Nitrosomethylamine	360	U
66-27-3-----	Methyl methanesulfonate	360	U
55-18-5-----	N-Nitrosodiethylamine	360	U
62-50-0-----	Ethyl methanesulfonate	360	U
108-95-2-----	Phenol	360	U
62-53-3-----	Aniline	360	U
76-01-7-----	Pentachloroethane	360	U
111-44-4-----	bis(2-Chloroethyl) Ether	720	U
95-57-8-----	2-Chlorophenol	360	U
541-73-1-----	1,3-Dichlorobenzene	360	U
100-44-7-----	Benzyl Chloride	360	U
106-46-7-----	1,4-Dichlorobenzene	360	U
100-51-6-----	Benzyl Alcohol	360	U
95-50-1-----	1,2-Dichlorobenzene	360	U
95-48-7-----	2-Methylphenol	360	U
108-60-1-----	bis(2-Chloroisopropyl) Ether	360	U
108-39-4-----	3-Methylphenol	360	U
106-44-5-----	4-Methylphenol	360	U
930-55-2-----	N-Nitroso-pyrrolidine	360	U
59-89-2-----	N-Nitrosomorpholine	360	U
98-86-2-----	Acetophenone	360	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	360	U
636-21-5-----	o-Toluidine hydrochloride	360	U
67-72-1-----	Hexachloroethane	360	U
98-95-3-----	Nitrobenzene	360	U
100-75-4-----	N-Nitrosopiperidine	360	U
78-59-1-----	Isophorone	360	U
88-75-5-----	2-Nitrophenol	360	U
105-67-9-----	2,4-Dimethylphenol	360	U

FORM I SV-1

1/87 Rev.

108-70-3	1, 3, 5-Trichlorobenzene	360	U
98-87-3	Benzal Chloride	360	U
65-65-0	Benzoic Acid	3600	U
111-91-1	bis(2-Chloroethoxy)Methane	360	U
120-83-2	2, 4-Dichlorophenol	360	U
120-82-1	1, 2, 4-Trichlorobenzene	360	U
91-20-3	Naphthalene	360	U
106-47-8	4-Chloroaniline	360	U
87-65-0	2, 6-Dichlorophenol	720	U
95-54-5	o-Phenylenediamine	360	U
122-09-8	dimethylphenylethylamine	360	U
1888-71-7	Hexachloropropene	360	U
87-68-3	Hexachlorobutadiene	360	U
87-61-6	1, 2, 3-Trichlorobenzene	360	U
98-07-7	Benzotrichloride	720	U
924-16-3	N-Nitroso-di-n-butylamine	360	U
59-50-7	4-Chloro-3-Methylphenol	360	U
106-50-3	p-Phenylenediamine	360	U
94-59-7	Safrole	360	U
106-50-3	m-Phenylenediamine	360	U
91-57-6	2-Methylnaphthalene	360	U
90-12-0	1-Methylnaphthalene	360	U
95-94-3	1, 2, 4, 5-Tetrachlorobenzene	360	U
634-90-2	1, 2, 3, 5-Tetrachlorobenzene	360	U
77-47-4	Hexachlorocyclopentadiene	360	U
88-06-2	2, 4, 6-Trichlorophenol	720	U
95-95-4	2, 4, 5-Trichlorophenol	720	U
120-58-1	Icosafrole	720	U
91-58-7	2-Chloronaphthalene	360	U
90-13-1	1-Chloronaphthalene	360	U
634-66-2	1, 2, 3, 4-Tetrachlorobenzene	360	U
88-74-4	2-Nitroaniline	360	U
130-15-4	1, 4-Naphthoquinone	720	U
100-25-4	1, 4-Dinitrobenzene	720	U
131-11-3	Dimethyl Phthalate	360	U
208-96-8	Acenaphthylene	360	U
606-20-2	2, 6-Dinitrotoluene	360	U

1C
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP Contract: 500077

ROJ3S

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 429Matrix: (soil/water) SOIL Lab Sample ID: 469296Sample wt/vol: 30.3 (g/mL) G Lab File ID: GJ069296A20Level: (low/med) LOW Date Received: 12/11/91t Moisture: not dec. 9 dec. _____ Date Extracted: 12/20/91Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/24/91GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) ug/kg Q

99-09-2	3-Nitroaniline	720	U
83-32-9	Acenaphthene	63	J
51-28-5	2, 4-Dinitrophenol	1400	U
100-02-7	4-Nitrophenol	360	U
132-64-9	Dibenzofuran	360	U
121-14-2	2, 4-Dinitrotoluene	360	U
608-93-5	Pentachlorobenzene	360	U
91-59-8	2-Naphthylamine	720	U
134-32-7	1-Naphthylamine	720	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	720	U
84-66-2	Diethylphthalate	360	U
297-97-2	Zinophos	360	U
7005-72-3	4-Chlorophenyl-phenylether	360	U
86-73-7	Fluorene	49	J
100-01-6	4-Nitroaniline	720	U
99-55-8	5-Nitro-o-toluidine	720	U
122-66-7	1, 2-Diphenylhydrazine	360	U
534-52-1	4, 6-Dinitro-2-Methylphenol	1100	U
86-30-6	N-Nitrosodiphenylamine (1)	360	U
122-39-4	Diphenylamine	360	U
99-35-4	1, 3, 5-Trinitrobenzene	720	U
62-44-2	Phenacetin	360	U
101-55-3	4-Bromophenyl-phenylether	360	U
2303-16-4	Diallate	360	U
60-51-5	Dimethoate	360	U
118-74-1	Hexachlorobenzene	360	U
92-67-1	4-Aminobiphenyl	360	U
23950-58-5	Pronamide	360	U
87-86-5	Pentachlorophenol	720	U
82-68-8	Pentachloronitrobenzene	360	U
85-01-8	Phenanthrene	630	
120-12-7	Anthracene	100	J
84-74-2	Di-n-Butylphthalate	360	U
91-80-5	Methapyrilene	720	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	1700	U
206-44-0	Fluoranthene	1200	
92-87-5	Benzidine	360	U
129-00-0	Pyrene	1000	
60-11-7	p-Dimethylaminocazobenzene	360	U
510-15-6	Chlorobenzilate	360	U
119-93-7	3,3'-Dimethylbenzidine	720	U
85-68-7	Butylbenzylphthalate	360	U
53-96-3	2-Acetylaminofluorene	360	U
101-14-4	Methylene-bis(2-Chloroaniline)	360	U
91-94-1	3,3'-Dichlorobenzidine	360	U
106-51-4	3,3'-Dimethoxybenzidine	360	U
56-55-3	Benzo(a)Anthracene	630	
117-81-7	bis(2-Ethylhexyl) Phthalate	53	J
218-01-9	Chrysene	640	
117-84-0	Di-n-Octyl Phthalate	360	U
205-99-2	Benzo(b)Fluoranthene	650	X
57-97-6	7,12-Dimethylbenzanthracene	360	U
207-08-9	Benzo(k)Fluoranthene	650	X
50-32-8	Benzo(a)Pyrene	600	
56-49-5	3-Methylchloranthrene	360	U
224-42-0	Dibenzo(a,j)acridine	360	U
193-39-5	Indeno(1,2,3-cd)Pyrene	290	J
53-70-3	Dibenz(a,h)Anthracene	88	J
191-24-2	Benzo(g,h,i)Perylene	350	J

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROJ4S
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>428</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>469259</u>	
Sample wt/vol: <u>30.2</u> (g/mL) G	Lab File ID: <u>GH069259A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/11/91</u>	
Moisture: not dec. <u>73</u> dec. _____	Date Extracted: <u>12/20/91</u>	
Extraction: (SepP/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>12/23/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
62-75-9	N-Nitrosodimethylamine	1200	U
110-86-1	Pyridine	1200	U
97-63-2	Ethyl methacrylate	1200	U
123-63-7	Paraldehyde	1200	U
109-06-8	2-Picoline	1200	U
10595-95-6	Nitrosomethylethylamine	2400	U
66-27-3	Methyl methanesulfonate	1200	U
55-18-5	N-Nitrosodiethylamine	1200	U
62-50-0	Ethyl methanesulfonate	1200	U
108-95-2	Phenol	1200	U
62-53-3	Aniline	1200	U
76-01-7	Pentachloroethane	1200	U
111-44-4	bis(2-Chloroethyl)Ether	2400	U
95-57-8	2-Chlorophenol	1200	U
541-73-1	1,3-Dichlorobenzene	1200	U
100-44-7	Benzyl Chloride	1200	U
106-46-7	1,4-Dichlorobenzene	1200	U
100-51-6	Benzyl Alcohol	1200	U
95-50-1	1,2-Dichlorobenzene	1200	U
95-48-7	2-Methylphenol	1200	U
108-60-1	bis(2-Chloroisopropyl)Ether	1200	U
108-39-4	3-Methylphenol	1200	U
106-44-5	4-Methylphenol	1200	U
930-55-2	N-Nitrosopyrrolidine	1200	U
59-89-2	N-Nitrosomorpholine	1200	U
98-86-2	Acetophenone	1200	U
621-64-7	N-Nitroso-Di-n-Propylamine	1200	U
636-21-5	o-Toluidine hydrochloride	1200	U
67-72-1	Hexachloroethane	1200	U
98-95-3	Nitrobenzene	1200	U
100-75-4	N-Nitrosopiperidine	1200	U
78-59-1	Isophorone	1200	U
88-75-5	2-Nitrophenol	1200	U
105-67-9	2,4-Dimethylphenol	1200	U

108-70-3	1, 3, 5-Trichlorobenzene	1200	U
98-87-3	Benzal Chloride	1200	U
65-85-0	Benzoic Acid	12000	U
111-91-1	bis(2-Chloroethoxy) Methane	1200	U
120-83-2	2, 4-Dichlorophenol	1200	U
120-82-1	1, 2, 4-Trichlorobenzene	1200	U
91-20-3	Naphthalene	150	J ✓
106-47-8	4-Chloroaniline	1200	U
87-65-0	2, 6-Dichlorophenol	2400	U
95-54-5	o-Phenylenediamine	1200	U
122-09-8	dimethylphenylethylamine	1200	U
1888-71-7	Hexachloropropene	1200	U
87-68-3	Hexachlorobutadiene	1200	U
87-61-6	1, 2, 3-Trichlorobenzene	1200	U
98-07-7	Benzotrichloride	2400	U
924-16-3	N-Nitroso-di-n-butylamine	1200	U
59-50-7	4-Chloro-3-Methylphenol	1200	U
106-50-3	p-Phenylenediamine	1200	U
94-59-7	Safrole	1200	U
106-50-3	m-Phenylenediamine	1200	U
91-57-6	2-Methylnaphthalene	1200	U
90-12-0	1-Methylnaphthalene	1200	U
95-94-3	1, 2, 4, 5-Tetrachlorobenzene	1200	U
634-90-2	1, 2, 3, 5-Tetrachlorobenzene	1200	U
77-47-4	Hexachlorocyclopentadiene	1200	U
88-06-2	2, 4, 6-Trichlorophenol	2400	U
95-95-4	2, 4, 5-Trichlorophenol	2400	U
120-58-1	Isosafrole	2400	U
91-58-7	2-Chloronaphthalene	1200	U
90-13-1	1-Chloronaphthalene	1200	U
634-66-2	1, 2, 3, 4-Tetrachlorobenzene	1200	U
88-74-4	2-Nitroaniline	1200	U
130-15-4	1, 4-Naphthoquinone	2400	U
100-25-4	1, 4-Dinitrobenzene	2400	U
131-11-3	Dimethyl Phthalate	1200	U
208-96-8	Acenaphthylene	250	J ✓
606-20-2	2, 6-Dinitrotoluene	1200	U

ROJ4S

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROJ4S

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>428</u>
Matrix: (soil/water) <u>SOIL</u>		Lab Sample ID: <u>469259</u>
Sample wt/vol: <u>30.2</u> (g/mL) G		Lab File ID: <u>GH069259A20</u>
Level: (low/med) <u>LOW</u>		Date Received: <u>12/11/91</u>
* Moisture: not dec. <u>22</u> dec. _____		Date Extracted: <u>12/20/91</u>
Extraction: (SepF/Cont/Sonc) <u>SONC</u>		Date Analyzed: <u>12/23/91</u>
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG Q

<u>99-09-2</u>	<u>3-Nitroaniline</u>	<u>2400</u>	<u>U</u>
<u>83-32-9</u>	<u>Acenaphthene</u>	<u>1200</u>	<u>U</u>
<u>51-28-5</u>	<u>2, 4-Dinitrophenol</u>	<u>4800</u>	<u>U</u>
<u>100-02-7</u>	<u>4-Nitrophenol</u>	<u>1200</u>	<u>U</u>
<u>132-64-9</u>	<u>Dibenzofuran</u>	<u>1200</u>	<u>U</u>
<u>121-14-2</u>	<u>2, 4-Dinitrotoluene</u>	<u>1200</u>	<u>U</u>
<u>608-93-5</u>	<u>Pentachlorobenzene</u>	<u>1200</u>	<u>U</u>
<u>91-59-8</u>	<u>2-Naphthylamine</u>	<u>2400</u>	<u>U</u>
<u>134-32-7</u>	<u>1-Naphthylamine</u>	<u>2400</u>	<u>U</u>
<u>58-90-2</u>	<u>2, 3, 4, 6-Tetrachlorophenol</u>	<u>2400</u>	<u>U</u>
<u>84-66-2</u>	<u>Diethylphthalate</u>	<u>1200</u>	<u>U</u>
<u>297-97-2</u>	<u>Zinophos</u>	<u>1200</u>	<u>U</u>
<u>7005-72-3</u>	<u>4-Chlorophenyl-phenylether</u>	<u>1200</u>	<u>U</u>
<u>86-73-7</u>	<u>Fluorene</u>	<u>140</u>	<u>J</u>
<u>100-01-6</u>	<u>4-Nitroaniline</u>	<u>2400</u>	<u>U</u>
<u>99-55-8</u>	<u>5-Nitro-o-toluidine</u>	<u>2400</u>	<u>U</u>
<u>122-66-7</u>	<u>1, 2-Diphenylhydrazine</u>	<u>1200</u>	<u>U</u>
<u>534-52-1</u>	<u>4, 6-Dinitro-2-Methylphenol</u>	<u>3600</u>	<u>U</u>
<u>86-30-6</u>	<u>N-Nitrosodiphenylamine (1)</u>	<u>1200</u>	<u>U</u>
<u>122-39-4</u>	<u>Diphenylamine</u>	<u>1200</u>	<u>U</u>
<u>99-35-4</u>	<u>1, 3, 5-Trinitrobenzene</u>	<u>2400</u>	<u>U</u>
<u>62-44-2</u>	<u>Phenacetin</u>	<u>1200</u>	<u>U</u>
<u>101-55-3</u>	<u>4-Bromophenyl-phenylether</u>	<u>1200</u>	<u>U</u>
<u>2303-16-4</u>	<u>Diallate</u>	<u>1200</u>	<u>U</u>
<u>60-51-5</u>	<u>Dimethoate</u>	<u>1200</u>	<u>U</u>
<u>118-74-1</u>	<u>Hexachlorobenzene</u>	<u>1200</u>	<u>U</u>
<u>92-67-1</u>	<u>4-Aminobiphenyl</u>	<u>1200</u>	<u>U</u>
<u>23950-58-5</u>	<u>Pronamide</u>	<u>1200</u>	<u>U</u>
<u>87-86-5</u>	<u>Pentachlorophenol</u>	<u>2400</u>	<u>U</u>
<u>82-68-8</u>	<u>Pentachloronitrobenzene</u>	<u>1200</u>	<u>U</u>
<u>85-01-8</u>	<u>Phenanthrene</u>	<u>1700</u>	
<u>120-12-7</u>	<u>Anthracene</u>	<u>180</u>	<u>J</u>
<u>84-74-2</u>	<u>Di-n-Butylphthalate</u>	<u>150</u>	<u>J</u>
<u>91-80-5</u>	<u>Methapyrilene</u>	<u>2400</u>	<u>U</u>

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

50-18-0	Cyclophosphamide	5900	U
206-44-0	Fluoranthene	2800	
92-87-5	Benzidine	1200	U
129-00-0	Pyrene	2400	
60-11-7	p-Dimethylaminoazobenzene	1200	U
510-15-6	Chlorobenzilate	1200	U
119-93-7	3,3'-Dimethylbenzidine	2400	U
85-68-7	Butylbenzylphthalate	1200	U
53-96-3	2-Acetylaminofluorene	1200	U
101-14-4	Methylene-bis(2-Chloroaniline)	1200	U
91-94-1	3,3'-Dichlorobenzidine	1200	U
106-51-4	3,3'-Dimethoxybenzidine	1200	U
56-55-3	Benzo(a)Anthracene	1500	
218-01-9	Chrysene	2200	
117-81-7	bis(2-Ethylhexyl) Phthalate	420	J
117-84-0	Di-n-Octyl Phthalate	1200	U
205-99-2	Benzo(b)Fluoranthene	3200	X
57-97-6	7,12-Dimethylbenzanthracene	1200	U
207-08-9	Benzo(k)Fluoranthene	3200	X
50-32-8	Benzo(a)Pyrene	1500	
56-49-5	3-Methylchloranthrene	1200	U
224-42-0	Dibenzo(a,j)acridine	1200	U
193-39-5	Indeno(1,2,3-cd)Pyrene	1200	U
53-70-3	Dibenz(a,h)Anthracene	1200	U
191-24-2	Benzo(g,h,i)Perylene	1200	U

(1) - Cannot be separated from Diphenylamine

R0545

18
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP Contract: 500077 R001B1416

Lab Code: COMPU Case No.: 22469 SAS No.: SDG No.: 03

Matrix: (soil/water) SOIL Lab Sample ID: 196810

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH096810C04

Level: (low/med) LOW Date Received: 02/01/91

Moisture: not dec. 15 dec. Date Extracted: 02/04/91

Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 02/07/91

SPC Cleanup: (Y/N) N pH: Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG

62-75-9-----	N-Nitrosodimethylamine	390	U
110-86-1-----	Pyridine	390	U
97-63-2-----	Ethyl methacrylate	390	U
123-63-7-----	Paraldehyde	390	U
109-06-8-----	2-Picoline	780	U
10595-95-6-----	Nitrosomethylethylamine	390	U
66-27-3-----	Methyl methanesulfonate	390	U
108-95-2-----	Phenol	390	U
55-18-5-----	N-Nitrosodiethylamine	390	U
62-50-5-----	Ethyl methanesulfonate	390	U
62-53-3-----	Aniline	390	U
76-01-7-----	Pentachloroethane	390	U
111-44-4-----	bis(2-Chloroethyl)Ether	780	U
95-57-8-----	2-Chlorophenol	390	U
541-73-1-----	1,3-Dichlorobenzene	390	U
100-44-7-----	Benzyl chloride	390	U
106-46-7-----	1,4-Dichlorobenzene	390	U
100-51-6-----	Benzyl Alcohol	390	U
95-50-1-----	1,2-Dichlorobenzene	390	U
95-48-7-----	2-Methylphenol	390	U
39608-32-9-----	bis(2-Chloroisopropyl)Ether	390	U
108-39-4-----	3-Methylphenol	390	U
106-44-5-----	4-Methylphenol	390	U
930-55-2-----	N-Nitrosopyrrolidine	390	U
59-89-2-----	N-Nitrosomorpholine	390	U
98-86-2-----	Acetophenone	390	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	390	U
636-21-5-----	o-Toluidine hydrochloride	390	U
67-72-1-----	Hexachloroethane	390	U
98-95-3-----	Nitrobenzene	390	U
100-75-4-----	N-Nitrosopiperidine	390	U
78-59-1-----	Isophorone	390	U
88-75-5-----	2-Nitrophenol	390	U
105-67-9-----	2,4-Dimethylphenol	390	U

FORM I SV-1

1/87 Rev.

22469 03 SAMPLE DATA SUMMARY

108-70-3-----	1,3,5-Trichlorobenzene	390	C
98-87-3-----	Benzal chloride	390	C
65-85-0-----	Benzoic Acid	3900	U
111-91-1-----	bis(2-Chloroethoxy)Methane	390	U
120-83-2-----	2,4-Dichlorophenol	390	U
120-82-1-----	1,2,4-Trichlorobenzene	390	U
91-20-3-----	Naphthalene	390	U
106-47-8-----	4-Chloroaniline	390	U
87-65-0-----	2,6-Dichlorophenol	780	U
95-54-5-----	O-Phenylenediamine	390	U
122-09-8-----	dimethylphenylethylamine	390	U
1888-71-7-----	Hexachloropropene	390	U
67-68-3-----	Hexachlorobutadiene	390	U
87-61-6-----	1,2,3-Trichlorobenzene	390	U
98-07-7-----	Benzotrifluoride	780	U
924-16-3-----	N-Nitroso-di-n-butylamine	390	U
59-50-7-----	4-Chloro-3-Methylphenol	390	U
106-50-3-----	P-Phenylenediamine	390	U
94-59-7-----	Safrole	390	U
106-50-3-----	m-Phenylenediamine	390	U
91-57-6-----	2-Methylnaphthalene	390	U
90-12-0-----	1-Methylnaphthalene	390	U
95-94-3-----	1,2,4,5-Tetrachlorobenzene	390	U
634-90-2-----	1,2,3,5-Tetrachlorobenzene	390	U
77-47-4-----	Hexachlorocyclopentadiene	390	U
88-06-2-----	2,4,6-Trichlorophenol	780	U
95-95-4-----	2,4,5-Trichlorophenol	780	U
120-58-1-----	Isosafrole	780	U
91-58-7-----	2-Chloronaphthalene	390	U
90-13-1-----	1-Chloronaphthalene	390	U
634-66-2-----	1,2,3,4-Tetrachlorobenzene	390	U
88-74-4-----	2-Nitroaniline	390	U
130-15-4-----	1,4-Naphthoquinone	780	U
100-25-4-----	1,4-Dinitrobenzene	780	U
131-11-3-----	Dimethyl Phthalate	390	U
208-96-8-----	Acenaphthylene	390	U

FORM I SV-1

1/87 Rev.

CCN 39681D

22469 03 SAMPLE DATA SUMMARY

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	R001B1416
Lab Code: <u>COMPU</u>	Case No.: <u>22469</u>	SAS No.: _____ SDG No.: <u>03</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>396810</u>	
Sample wt/vol: <u>30.0</u> (g/mL) G	Lab File ID: <u>GH096810C04</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>02/01/91</u>	
% Moisture: not dec. <u>15</u> dec. _____	Date Extracted: <u>02/04/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>02/07/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
99-09-2-----	3-Nitroaniline	780	U
83-32-9-----	Acenaphthene	390	U
51-28-5-----	2, 4-Dinitrophenol	1500	U
100-02-7-----	4-Nitrophenol	390	U
132-64-9-----	Dibenzofuran	390	U
121-14-2-----	2, 4-Dinitrotoluene	390	U
608-93-5-----	Pentachlorobenzene	390	U
134-32-7-----	2-Naphthylamine	780	U
606-20-2-----	2, 6-Dinitrotoluene	390	U
134-32-7-----	1-Naphthylamine	780	U
58-90-2-----	2, 3, 4, 6-Tetrachlorophenol	780	U
84-66-2-----	Diethylphthalate	390	U
297-97-2-----	Zinophos	390	U
7005-72-3-----	4-Chlorophenyl-phenylether	390	U
86-73-7-----	Fluorene	390	U
100-01-6-----	4-Nitroaniline	780	U
99-55-8-----	5-Nitro-o-toluidine	780	U
534-52-1-----	4, 6-Dinitro-2-Methylphenol	1200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	390	U
122-39-4-----	Diphenylamine	390	U
99-35-4-----	1, 3, 5-Trinitrobenzene	780	U
122-66-7-----	1, 2-Diphenylhydrazine	390	U
62-44-2-----	Phenacetin	390	U
101-55-3-----	4-Bromophenyl-phenylether	390	U
2303-16-4-----	Diallate	390	U
60-51-5-----	Dimethoate	390	U
118-74-1-----	Hexachlorobenzene	390	U
92-67-1-----	4-Aminobiphenyl	390	U
23950-58-5-----	Pronamide	390	U
87-86-5-----	Pentachlorophenol	780	U
82-68-8-----	Pentachloronitrobenzene	390	U
85-01-8-----	Phenanthrone	390	U
120-12-7-----	Anthracene	390	U
84-74-2-----	Di-n-Butylphthalate	390	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

91-80-5-----	Methapyrilene	780	U
50-18-0-----	Cyclophosphamide	1900	U
206-44-0-----	Fluoranthene	390	U
92-87-5-----	Benzidine	390	U
129-00-0-----	Pyrene	390	U
60-11-7-----	p-Dimethylaminoazobenzene	390	U
510-15-6-----	Chlorobenzilate	390	U
119-93-7-----	3,3'-Dimethylbenzidine	780	U
85-68-7-----	Butylbenzylphthalate	390	U
53-96-3-----	2-Acetylaminofluorene	390	U
101-14-4-----	Methylene-bis(2-chloroaniline)	390	U
91-94-1-----	3,3'-Dichlorobenzidine	390	U
106-51-4-----	3,3'-Dimethoxybenzidine	390	U
56-55-3-----	Benzo(a)Anthracene	390	U
218-01-9-----	Chrysene	390	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	390	U
117-84-0-----	Oi-n-Octyl Phthalate	390	U
205-99-2-----	Benzo(b)Fluoranthene	390	U
57-97-6-----	7,12-Dimethylbenzanthracene	390	U
207-08-9-----	Benzo(k)Fluoranthene	390	U
50-32-8-----	Benzo(a)Pyrene	390	U
56-49-5-----	3-Methylcholanthrene	390	U
224-42-0-----	Dibenzo(a,j)acridine	390	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	390	U
53-70-3-----	Dibenz(a,h)Anthracene	390	U
191-24-2-----	Benzo(g,h,i)Perylene	390	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM, RTP</u>	Contract: <u>500077</u>	R002B0810
Lab Code: <u>COMPU</u>	Case No.: <u>22469</u>	SAS No.: _____ SDG No.: <u>03</u>
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>396818</u>	
Sample wt/vol: <u>30.0</u> (g/mL) G	Lab File ID: <u>GH096818B04</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>02/01/91</u>	
% Moisture: not dec. <u>18</u> dec. _____	Date Extracted: <u>02/04/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>02/07/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/kg	Q
62-75-9-----	N-Nitrosodimethylamine	400	U
110-86-1-----	Pyridine	400	U
97-63-2-----	Ethyl methacrylate	400	U
123-63-7-----	Paraldehyde	400	U
109-06-8-----	2-Picoline	810	U
10595-95-6-----	Nitrosomethylethylamine	400	U
66-27-3-----	Methyl methanesulfonate	400	U
108-95-2-----	Phenol	400	U
55-18-5-----	N-Nitrosodiethylamine	400	U
62-50-5-----	Ethyl methanesulfonate	400	U
62-53-3-----	Aniline	400	U
76-01-7-----	Pentachloroethane	400	U
111-44-4-----	bis(2-Chloroethyl) Ether	810	U
95-57-8-----	2-Chlorophenol	400	U
541-73-1-----	1,3-Dichlorobenzene	400	U
100-44-7-----	Benzyl chloride	400	U
106-46-7-----	1,4-Dichlorobenzene	400	U
100-51-6-----	Benzyl Alcohol	400	U
95-50-1-----	1,2-Dichlorobenzene	400	U
95-48-7-----	2-Methylphenol	400	U
39638-32-9-----	bis(2-Chloroisopropyl) Ether	400	U
108-39-4-----	3-Methylphenol	400	U
106-44-5-----	4-Methylphenol	400	U
930-55-2-----	N-Nitroscypyrrrolidine	400	U
59-89-2-----	N-Nitrosomorpholine	400	U
98-86-2-----	Acetophenone	400	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	400	U
636-21-5-----	o-Tolidine hydrochloride	400	U
67-72-1-----	Hexachloroethane	400	U
98-95-3-----	Nitrobenzene	400	U
100-75-4-----	N-Nitrosopiperidine	400	U
78-59-1-----	Isophorone	400	U
88-75-5-----	2-Nitrophenol	400	U
105-67-9-----	2,4-Dimethylphenol	400	U

(1) - Cannot be separated from Diphenylamine
FORM I SV-4

1/87 Rev.

108-70-3-----	1,3,5-Trichlorobenzene	400	U
98-87-3-----	Benzal chloride	400	U
65-85-0-----	Benzoic Acid	4000	U
111-91-1-----	bis(2-Chloroethoxy)Methane	400	U
120-83-2-----	2,4-Dichlorophenol	400	U
120-82-1-----	1,2,4-Trichlorobenzene	400	U
91-20-3-----	Naphthalene	400	U
106-47-8-----	4-Chloraniline	400	U
87-65-0-----	2,6-Dichlorophenol	810	U
95-54-5-----	o-Phenylenediamine	400	U
122-09-8-----	dimethylphenylethylamine	400	U
1888-71-7-----	Hexachloropropene	400	U
87-68-3-----	Hexachlorobutadiene	400	U
87-61-6-----	1,2,3-Trichlorobenzene	400	U
98-07-7-----	Benzotrichloride	810	U
924-16-3-----	N-Nitroso-di-n-butylamine	400	U
59-50-7-----	4-Chloro-3-Methylphenol	400	U
106-50-3-----	p-Phenylenediamine	400	U
94-59-7-----	Safrole	400	U
106-50-3-----	m-Phenylenediamine	400	U
91-57-6-----	2-Methylnaphthalene	400	U
90-12-0-----	1-Methylnaphthalene	400	U
95-94-3-----	1,2,4,5-Tetrachlorobenzene	400	U
634-90-2-----	1,2,3,5-Tetrachlorobenzene	400	U
77-47-4-----	Hexachlorocyclopentadiene	400	U
88-06-2-----	2,4,6-Trichlorophenol	810	U
95-95-4-----	2,4,5-Trichlorophenol	810	U
120-58-1-----	Isosafrole	810	U
91-58-7-----	2-Chloronaphthalene	400	U
90-13-1-----	1-Chloronaphthalene	400	U
634-66-2-----	1,2,3,4-Tetrachlorobenzene	400	U
88-74-4-----	2-Nitroaniline	400	U
130-15-4-----	1,4-Naphthoquinone	810	U
100-25-4-----	1,4-Dinitrobenzene	810	U
131-11-3-----	Dimethyl Phthalate	400	U
208-96-8-----	Acenaphthylene	400	U

FORM I SV-1

1/87 Rev.

CCN 394818

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM, RTP</u>	Contract: <u>500077</u>	R002B0810
Lab Code: <u>COMPU</u>	Case No.: <u>22469</u>	SAS No.: _____
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID: <u>396818</u>	
Sample wt/vol: <u>30.0</u> (g/mL) G	Lab File ID: <u>GH096818B04</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>02/01/91</u>	
Moisture: not dec. <u>18</u> dec. _____	Date Extracted: <u>02/04/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SONC</u>	Date Analyzed: <u>02/07/91</u>	
CPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.00</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
99-09-2-----	3-Nitroaniline	810	U	
83-32-9-----	Acenaphthene	400	U	
51-28-5-----	2,4-Dinitrophenol	1600	U	
100-02-7-----	4-Nitrophenol	400	U	
132-64-9-----	Dibenzofuran	400	U	
121-14-2-----	2,4-Dinitrotoluene	400	U	
608-93-5-----	Pentachlorobenzene	400	U	
134-32-7-----	2-Naphthylamine	810	U	
606-20-2-----	2,6-Dinitrotoluene	400	U	
134-32-7-----	1-Naphthylamine	810	U	
58-90-2-----	2,3,4,6-Tetrachlorophenol	810	U	
84-66-2-----	Diethylphthalate	400	U	
297-97-2-----	Zinophos	400	U	
7005-72-3-----	4-Chlorophenyl-phenylether	400	U	
86-73-7-----	Fluorene	400	U	
100-01-6-----	4-Nitroaniline	810	U	
99-55-8-----	5-Nitro-o-toluidine	810	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	1200	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	400	U	
122-39-4-----	Diphenylamine	400	U	
99-35-4-----	1,3,5-Trinitrobenzene	810	U	
122-66-7-----	1,2-Diphenylhydrazine	400	U	
62-44-2-----	Phenacetin	400	U	
101-55-3-----	4-Bromophenyl-phenylether	400	U	
2303-16-4-----	Diallate	400	U	
60-51-5-----	Dimethoate	400	U	
118-74-1-----	Hexachlorobenzene	400	U	
92-67-1-----	4-Aminobiphenyl	400	U	
23950-58-5-----	Pronamide	400	U	
87-86-5-----	Pentachlorophenol	810	U	
82-68-8-----	Pentachloronitrobenzene	400	U	
85-01-8-----	Phenanthrene	53	J	
120-12-7-----	Anthracene	400	U	
84-74-2-----	Di-n-Butylphthalate	53	J	

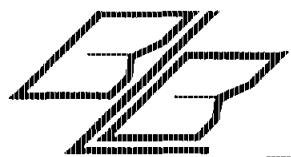
(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

91-80-5-----	Methapyrilene	810	U
50-18-0-----	Cyclophosphamide	2000	U
206-44-0-----	Fluoranthene	80	J
92-87-5-----	Benzidine	400	U
129-00-0-----	Pyrene	97	J
60-11-7-----	p-Dimethylaminoazobenzene	400	U
510-15-6-----	Chlorobenzilate	400	U
119-93-7-----	3,3'-Dimethylbenzidine	810	U
85-68-7-----	Butylbenzylphthalate	400	U
53-96-3-----	2-Acetylaminofluorene	400	U
101-14-4-----	Methylene-bis(2-chloroaniline	400	U
91-94-1-----	3,3'-Dichlorobenzidine	400	U
106-51-4-----	3,3'-Dimethoxybenzidine	400	U
56-55-3-----	Benzo(a)Anthracene	45	J
218-01-9-----	Chrysene	59	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	67	J
117-84-0-----	Di-n-Octyl Phthalate	400	U
205-99-2-----	Benzo(b)Fluoranthene	86	JX
57-97-6-----	7,12-Dimethylbenzanthracene	400	U
207-08-9-----	Benzo(k)Fluoranthene	86	JX
50-32-8-----	Benzo(a)Pyrone	42	J
56-49-5-----	3-Methylcholanthrene	400	U
224-42-0-----	Dibenzo(a,j)acridine	400	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	400	U
53-70-3-----	Dibenz(a,h)Anthracene	400	U
191-24-2-----	Benzo(g,h,i)Perylene	400	U

(1) - Cannot be separated from Diphenylamine



Section 3

PESTICIDES/PCB ANALYSIS (SOIL)

ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet

ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet

ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet

ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet

ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)

ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet

ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet

ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)

ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1

ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2

ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3

ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

ROO1B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet

ROO2B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

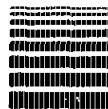
SAMPLE IDENTIFIER: R0A012224
COMPUCHEM SAMPLE NUMBER: 462141
DRY WEIGHT FACTOR: 1.11

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	3.9
2P. 4,4'-DDE	BDL	3.9
3P. 4,4'-DDT	BDL	3.9
4P. ALDRIN	BDL	1.1
5P. CHLORDANE	BDL	4.4
6P. DIELDRIN	BDL	1.7
7P. ENDOSULFAN I	BDL	1.7
8P. ENDOSULFAN II	BDL	3.9
9P. ENDOSULFAN SULFATE	BDL	2.2
10P. ENDRIN	BDL	2.8
11P. ENDRIN ALDEHYDE	BDL	1.1
12P. HEPTACHLOR	BDL	1.1
13P. HEPTACHLOR EPOXIDE	BDL	1.1
14P. KEPONE	BDL	1.1
15P. p,p'-METHOXYCHLOR	BDL	3.9
16P. PCB-1016	BDL	22
17P. PCB-1221	BDL	22
18P. PCB-1232	BDL	22
19P. PCB-1242	BDL	22
20P. PCB-1248	BDL	22
21P. PCB-1254	57	22
22P. PCB-1260	BDL	22
23P. TOXAPHENE	BDL	22
24P. ALPHA-BHC	BDL	1.1
25P. BETA-BHC	BDL	1.1
26P. DELTA-BHC	BDL	1.1
27P. GAMMA-BHC (Lindane)	BDL	1.1

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROA012224
COMPUCHEM SAMPLE NUMBER: 462141
DRY WEIGHT FACTOR: 1.11

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	† Recovery	Control Range †
Dibutylchloroendate	54	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: ROA2B0608
COMPUCHEM SAMPLE NUMBER: 465050
DRY WEIGHT FACTOR: 1.1

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	3.9
2P. 4,4'-DDE	BDL	3.9
3P. 4,4'-DDT	BDL	3.9
4P. ALDRIN	BDL	1.1
5P. CHLORDANE	BDL	4.4
6P. DIELDRIN	BDL	1.7
7P. ENDOSULFAN I	BDL	1.7
8P. ENDOSULFAN II	BDL	3.9
9P. ENDOSULFAN SULFATE	BDL	2.2
10P. ENDRIN	BDL	2.8
11P. ENDRIN ALDEHYDE	BDL	1.1
12P. HEPTACHLOR	BDL	1.1
13P. HEPTACHLOR EPOXIDE	BDL	1.1
14P. KEPONE	BDL	1.1
15P. p,p'-METHOXYCHLOR	BDL	3.9
16P. PCB-1016	BDL	22
17P. PCB-1221	BDL	22
18P. PCB-1232	BDL	22
19P. PCB-1242	BDL	22
20P. PCB-1243	BDL	22
21P. PCB-1254	BDL	22
22P. PCB-1260	BDL	22
23P. TOXAPHENE	BDL	22
24P. ALPHA-BHC	BDL	1.1
25P. BETA-BHC	BDL	1.1
26P. DELTA-BHC	BDL	1.1
27P. GAMMA-BHC (Lindane)	BDL	1.1

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: R0A2B0608
COMPUCHEM SAMPLE NUMBER: 465050
DRY WEIGHT FACTOR: 1.1

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroendate	150	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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APPENDIX VIII, IX - PESTICIDES, METHOD 8080

RESULTS REPORTED ON DRY WEIGHT BASIS

(Page 1)

SAMPLE IDENTIFIER: ROB1B0406
COMPUCHEM SAMPLE NUMBER: 465533
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	21
2P. 4,4'-DDE	BDL	21
3P. 4,4'-DDT	BDL	21
4P. ALDRIN	BDL	6
5P. CHLORDANE	BDL	24
6P. DIELDRIN	BDL	9
7P. ENDOSULFAN I	BDL	9
8P. ENDOSULFAN II	BDL	21
9P. ENDOSULFAN SULFATE	BDL	12
10P. ENDRIN	BDL	15
11P. ENDRIN ALDEHYDE	BDL	6
12P. HEPTACHLOR	BDL	6
13P. HEPTACHLOR EPOXIDE	BDL	6
14P. KEPONE	BDL	6
15P. p,p'-METHOXYCHLOR	BDL	21
16P. PCB-1016	BDL	120
17P. PCB-1221	BDL	120
18P. PCB-1232	BDL	120
19P. PCB-1242	BDL	120
20P. PCB-1248	BDL	120
21P. PCB-1254	5300	120
22P. PCB-1260	2800	120
23P. TOXAPHENE	BDL	120
24P. ALPHA-BHC	BDL	6
25P. BETA-BHC	BDL	6
26P. DELTA-BHC	BDL	6
27P. GAMMA-BHC (Lindane)	100	6

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

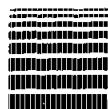
APPENDIX VIII, IX - PESTICIDES, METHOD 8060
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROB1B0406
COMPUCHEM SAMPLE NUMBER: 465533
DRY WEIGHT FACTOR: 1.21

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	† Recovery	Control Range †
Dibutylchlorendate	69	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: ROB2B0002
COMPUCHEM SAMPLE NUMBER: 465880
DRY WEIGHT FACTOR: 1.50

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	26
2P. 4,4'-DDE	BDL	26
3P. 4,4'-DDT	BDL	26
4P. ALDRIN	BDL	7.4
5P. CHLORDANE	BDL	30
6P. Dieldrin	BDL	11
7P. ENDOSULFAN I	BDL	11
8P. ENDOSULFAN II	BDL	26
9P. ENDOSULFAN SULFATE	BDL	15
10P. ENDRIN	BDL	19
11P. ENDRIN ALDEHYDE	BDL	7.4
12P. HEPTACHLOR	BDL	7.4
13P. HEPTACHLOR EPOXIDE	BDL	7.4
14P. KEPONE	BDL	7.4
15P. p,p'-METHOXYCHLOR	BDL	26
16P. PCB-1016	BDL	150
17P. PCB-1221	BDL	150
18P. PCB-1232	BDL	150
19P. PCB-1242	BDL	150
20P. PCB-1248	BDL	150
21P. PCB-1254	14000	150
22P. PCB-1260	BDL	150
23P. TOXAPENE	BDL	150
24P. ALPHA-BHC	BDL	7.4
25P. BETA-BHC	BDL	7.4
26P. DELTA-BHC	BDL	7.4
27P. GAMMA-BHC (Lindane)	BDL	7.4

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

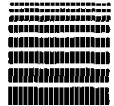
APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROBZB0002
COMPUCHEM SAMPLE NUMBER: 465880
DRY WEIGHT FACTOR: 1.5

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroendate	103	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: ROB-DPAI
COMPUCHEM SAMPLE NUMBER: 465556
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION ($\mu\text{g}/\text{kg}$)	DETECTION + LIMIT ($\mu\text{g}/\text{kg}$)
1P. 4,4'-DDD	BDL	42
2P. 4,4'-DDE	BDL	42
3P. 4,4'-DDT	BDL	42
4P. ALDRIN	BDL	12
5P. CHLORDANE	BDL	48
6P. DDELDRIN	BDL	18
7P. ENDOSULFAN I	BDL	18
8P. ENDOSULFAN II	BDL	42
9P. ENDOSULFAN SULFATE	BDL	24
10P. ENDRIN	BDL	30
11P. ENDRIN ALDEHYDE	BDL	12
12P. HEPTACHLOR	BDL	12
13P. HEPTACHLOR EPOXIDE	BDL	12
14P. KEPONE	BDL	12
15P. p,p'-METHOXYCHLOR	BDL	42
16P. PCB-1016	BDL	240
17P. PCB-1221	BDL	240
18P. PCB-1232	BDL	240
19P. PCB-1242	BDL	240
20P. PCB-1248	BDL	240
21P. PCB-1254	49000	240
22P. PCB-1260	BDL	240
23P. TOXAPENE	BDL	240
24P. ALPHA-BHC	BDL	12
25P. BETA-BHC	BDL	12
26P. DELTA-BHC	BDL	12
27P. GAMMA-BHC (Lindane)	BDL	12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

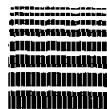
APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROB-DPA1
COMPUCHEM SAMPLE NUMBER: 465556
DRY WEIGHT FACTOR: 1.21

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroendate	103	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

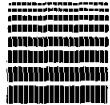
SAMPLE IDENTIFIER: ROC011012
COMPUCHEM SAMPLE NUMBER: 461433
DRY WEIGHT FACTOR: 1.11

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	97	3.9
2P. 4,4'-DDE	BDL	3.9
3P. 4,4'-DDT	BDL	3.9
4P. ALDRIN	BDL	1.1
5P. CHLORDANE	BDL	4.4
6P. DIELDRIN	BDL	1.7
7P. ENDOSULFAN I	BDL	1.7
8P. ENDOSULFAN II	BDL	3.9
9P. ENDOSULFAN SULFATE	BDL	2.2
10P. ENDRIN	BDL	2.8
11P. ENDRIN ALDEHYDE	BDL	1.1
12P. HEPTACHLOR	BDL	1.1
13P. HEPTACHLOR EPOXIDE	BDL	1.1
14P. KEPONE	BDL	1.1
15P. p,p'-METHOXYCHLOR	BDL	3.9
16P. PCB-1016	BDL	22
17P. PCB-1221	BDL	22
18P. PCB-1232	BDL	22
19P. PCB-1242	BDL	22
20P. PCB-1248	BDL	22
21P. PCB-1254	860	22
22P. PCB-1260	BDL	22
23P. TOXAPHENE	BDL	22
24P. ALPHA-BHC	BDL	1.1
25P. BETA-BHC	BDL	1.1
26P. DELTA-BHC	BDL	1.1
27P. GAMMA-BHC (Lindane)	BDL	1.1

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

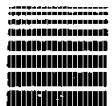
APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROC011012
COMPUCHEM SAMPLE NUMBER: 461433
DRY WEIGHT FACTOR: 1.11

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range †
Dibutylchloroendate	21	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: ROC021214
COMPUCHEM SAMPLE NUMBER: 461155
DRY WEIGHT FACTOR: 1.23

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	4.3
2P. 4,4'-DDE	BDL	4.3
3P. 4,4'-DDT	140	4.3
4P. ALDRIN	BDL	1.2
5P. CHLORDANE	BDL	4.9
6P. DIELDRIN	BDL	1.8
7P. ENDOSULFAN I	BDL	1.8
8P. ENDOSULFAN II	BDL	4.3
9P. ENDOSULFAN SULFATE	BDL	2.4
10P. ENDRIN	BDL	3.1
11P. ENDRIN ALDEHYDE	BDL	1.2
12P. HEPTACHLOR	BDL	1.2
13P. HEPTACHLOR EPOXIDE	BDL	1.2
14P. KEPONE	BDL	1.2
15P. p,p'-METHOXYCHLOR	BDL	4.3
16P. PCB-1016	BDL	24
17P. PCB-1221	BDL	24
18P. PCB-1232	BDL	24
19P. PCB-1242	BDL	24
20P. PCB-1248	BDL	24
21P. PCB-1254	BDL	24
22P. PCB-1260	260	24
23P. TOXAPHENE	BDL	24
24P. ALPHA-BHC	BDL	1.2
25P. BETA-BHC	BDL	1.2
26P. DELTA-BHC	23	1.2
27P. GAMMA-BHC (Lindane)	6.7	1.2

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROC021214
COMPUCHEM SAMPLE NUMBER: 461155
DRY WEIGHT FACTOR: 1.23

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroxendate	79	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

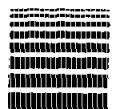
SAMPLE IDENTIFIER: ROC-DPA1
COMPUCHEM SAMPLE NUMBER: 461188
DRY WEIGHT FACTOR: 1.85

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	6.3
2P. 4,4'-DDE	BDL	6.3
3P. 4,4'-DDT	BDL	6.3
4P. ALDRIN	BDL	1.8
5P. CHLORDANE	BDL	7.2
6P. DIELDRIN	BDL	2.7
7P. ENDOSULFAN I	BDL	2.7
8P. ENDOSULFAN II	BDL	6.3
9P. ENDOSULFAN SULFATE	BDL	3.6
10P. ENDRIN	BDL	4.5
11P. ENDRIN ALDEHYDE	BDL	1.8
12P. HEPTACHLOR	BDL	1.8
13P. HEPTACHLOR EPOXIDE	BDL	1.8
14P. KEPONE	BDL	1.8
15P. p,p'-METHOXYCHLOR	BDL	6.3
16P. PCB-1016	BDL	36
17P. PCB-1221	BDL	36
18P. PCB-1232	BDL	36
19P. PCB-1242	BDL	36
20P. PCB-1248	BDL	36
21P. PCB-1254	BDL	36
22P. PCB-1260	1600	36
23P. TOXAPHENE	BDL	36
24P. ALPHA-BHC	BDL	1.8
25P. BETA-BHC	BDL	1.8
26P. DELTA-BHC	BDL	1.8
27P. GAMMA-BHC (Lindane)	BDL	1.8

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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(Page 2)

SAMPLE IDENTIFIER: ROC-DPAI
COMPUCHEM SAMPLE NUMBER: 461188
DRY WEIGHT FACTOR: 1.35

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroendate	73	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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RESULTS REPORTED ON DRY WEIGHT BASIS
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SAMPLE IDENTIFIER: ROC3E0204
COMPUCHEM SAMPLE NUMBER: 465093
DRY WEIGHT FACTOR: 1.08

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	3.8
2P. 4,4'-DDE	BDL	3.8
3P. 4,4'-DDT	BDL	3.8
4P. ALDRIN	BDL	1.1
5P. CHLORDANE	BDL	4.3
6P. Dieldrin	BDL	1.6
7P. ENDOSULFAN I	BDL	1.6
8P. ENDOSULFAN II	BDL	3.8
9P. ENDOSULFAN SULFATE	BDL	2.2
10P. ENDRIN	BDL	2.7
11P. ENDRIN ALDEHYDE	BDL	1.1
12P. HEPTACHLOR	BDL	1.1
13P. HEPTACHLOR EPOXIDE	BDL	1.1
14P. KEPONE	BDL	1.1
15P. p,p'-METHOXYCHLOR	BDL	3.8
16P. PCB-1016	BDL	22
17P. PCB-1221	BDL	22
18P. PCB-1232	BDL	22
19P. PCB-1242	BDL	22
20P. PCB-1248	BDL	22
21P. PCB-1254	BDL	22
22P. PCB-1260	BDL	22
23P. TOXAPHENE	BDL	22
24P. ALPHA-BHC	BDL	1.1
25P. BETA-BHC	BDL	1.1
26P. DELTA-BHC	BDL	1.1
27P. GAMMA-BHC (Lindane)	BDL	1.1

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

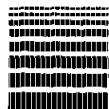
SAMPLE IDENTIFIER: ROC3B0204
COMPUCHEM SAMPLE NUMBER: 465093
DRY WEIGHT FACTOR: 1.08

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroendate	**	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.

**No surrogate recovery data available due to a matrix effect. See Quality Assurance Notice.



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APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: ROJIS
COMPUCHEM SAMPLE NUMBER: 469286
DRY WEIGHT FACTOR: 1.28

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	4.5
2P. 4,4'-DDE	BDL	4.5
3P. 4,4'-DDT	BDL	4.5
4P. ALDRIN	BDL	1.3
5P. CHLORDANE	BDL	5.1
6P. DIELDRIN	BDL	1.9
7P. ENDOSULFAN I	BDL	1.9
8P. ENDOSULFAN II	BDL	4.5
9P. ENDOSULFAN SULFATE	BDL	2.6
10P. ENDRIN	BDL	3.2
11P. ENDRIN ALDEHYDE	BDL	1.3
12P. HEPTACHLOR	BDL	1.3
13P. HEPTACHLOR EPOXIDE	BDL	1.3
14P. KEPONE	BDL	1.3
15P. p,p'-METHOXYCHLOR	BDL	4.5
16P. PCB-1016	BDL	26
17P. PCB-1221	BDL	26
18P. PCB-1232	BDL	26
19P. PCB-1242	BDL	26
20P. PCB-1248	BDL	26
21P. PCB-1254	BDL	26
22P. PCB-1260	970	26
23P. TOXAPHENE	BDL	26
24P. ALPHA-BHC	BDL	1.3
25P. BETA-BHC	BDL	1.3
26P. DELTA-BHC	BDL	1.3
27P. GAMMA-BHC (Lindane)	BDL	1.3

BDL = BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

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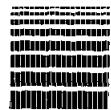
APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: R0J1S
COMPUCHEM SAMPLE NUMBER: 469286
DRY WEIGHT FACTOR: 1.28

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	† Recovery	Control Range †
Dibutylchloroendate	60	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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APPENDIX VIII, IX - PESTICIDES, METHOD 8080
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(Page 1)

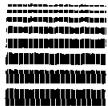
SAMPLE IDENTIFIER: ROJ2S
COMPUCHEM SAMPLE NUMBER: 469277
DRY WEIGHT FACTOR: 1.18

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	21
2P. 4,4'-DDE	BDL	21
3P. 4,4'-DDT	BDL	21
4P. ALDRIN	BDL	5.9
5P. CHLORDANE	BDL	24
6P. DIELDRIN	BDL	8.9
7P. ENDOSULFAN I	BDL	8.9
8P. ENDOSULFAN II	BDL	21
9P. ENDOSULFAN SULFATE	BDL	12
10P. ENDRIN	BDL	15
11P. ENDRIN ALDEHYDE	BDL	5.9
12P. HEPTACHLOR	BDL	5.9
13P. HEPTACHLOR EPOXIDE	BDL	5.9
14P. KEPONE	BDL	5.9
15P. p,p'-METHOXYCHLOR	BDL	21
16P. PCB-1016	BDL	120
17P. PCB-1221	BDL	120
18P. PCB-1232	BDL	120
19P. PCB-1242	BDL	120
20P. PCB-1248	BDL	120
21P. PCB-1254	BDL	120
22P. PCB-1260	440	120
23P. TOXAPHENE	BDL	120
24P. ALPHA-BHC	BDL	5.9
25P. BETA-BHC	BDL	5.9
26P. DELTA-BHC	BDL	5.9
27P. GAMMA-BHC (Lindane)	BDL	5.9

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROJ2S
COMPUCHEM SAMPLE NUMBER: 469277
DRY WEIGHT FACTOR: 1.18

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	† Recovery	Control Range ‡
Dibutylchlorendate	98	(20-150)†

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: ROJ3S
COMPUCHEM SAMPLE NUMBER: 469297
DRY WEIGHT FACTOR: 1.1

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	3.8
2P. 4,4'-DDE	BDL	3.8
3P. 4,4'-DDT	6.9	3.8
4P. ALDRIN	BDL	1.1
5P. CHLORDANE	BDL	4.4
6P. DIELDRIN	BDL	1.6
7P. ENDOSULFAN I	BDL	1.6
8P. ENDOSULFAN II	BDL	3.8
9P. ENDOSULFAN SULFATE	BDL	2.2
10P. ENDRIN	BDL	2.7
11P. ENDRIN ALDEHYDE	BDL	1.1
12P. HEPTACHLOR	BDL	1.1
13P. HEPTACHLOR EPOXIDE	BDL	1.1
14P. KEPONE	BDL	1.1
15P. p,p'-METHOXYCHLOR	BDL	3.8
16P. PCB-1016	BDL	22
17P. PCB-1221	BDL	22
18P. PCB-1232	BDL	22
19P. PCB-1242	BDL	22
20P. PCB-1248	BDL	22
21P. PCB-1254	BDL	22
22P. PCB-1260	BDL	22
23P. TOXAPHENE	BDL	22
24P. ALPHA-BHC	BDL	1.1
25P. BETA-BHC	BDL	1.1
26P. DELTA-BHC	BDL	1.1
27P. GAMMA-BHC (Lindane)	BDL	1.1

BDL= BELOW DETECTION LIMIT

* Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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(Page 2)

SAMPLE IDENTIFIER: ROJ3S
COMPUCHEM SAMPLE NUMBER: 469297
DRY WEIGHT FACTOR: 1.1

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	† Recovery	Control Range †
Dibutylchlorendate	110	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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(Page 1)

SAMPLE IDENTIFIER: ROJ4S
COMPUCHEM SAMPLE NUMBER: 469260
DRY WEIGHT FACTOR: 1.49

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	5.2
2P. 4,4'-DDE	BDL	5.2
3P. 4,4'-DDT	BDL	5.2
4P. ALDRIN	BDL	1.5
5P. CHLORDANE	BDL	5.9
6P. DIELDRIN	BDL	2.2
7P. ENDOSULFAN I	BDL	2.2
8P. ENDOSULFAN II	BDL	5.2
9P. ENDOSULFAN SULFATE	BDL	3
10P. ENDRIN	BDL	3.7
11P. ENDRIN ALDEHYDE	BDL	1.5
12P. HEPTACHLOR	BDL	1.5
13P. HEPTACHLOR EPOXIDE	BDL	1.5
14P. KEPONE	BDL	1.5
15P. p,p'-METHOXYCHLOR	BDL	5.2
16P. PCB-1016	BDL	30
17P. PCB-1221	BDL	30
18P. PCB-1232	BDL	30
19P. PCB-1242	BDL	30
20P. PCB-1248	BDL	30
21P. PCB-1254	BDL	30
22P. PCB-1260	710	30
23P. TOXAPHENE	BDL	30
24P. ALPHA-BHC	BDL	1.5
25P. BETA-BHC	BDL	1.5
26P. DELTA-BHC	BDL	1.5
27P. GAMMA-BHC (Lindane)	BDL	1.5

BDL= BELOW DETECTION LIMIT

* Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

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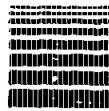
APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: ROJ4S
COMPUCHEM SAMPLE NUMBER: 469260
DRY WEIGHT FACTOR: 1.49

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchloroendate	68	(20-150)*

*Advisory surrogate: with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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(Page 1)

SAMPLE IDENTIFIER: R001B1416
COMPUCHEM SAMPLE NUMBER: 397161
DRY WEIGHT FACTOR: 1.18

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	4.1
2P. 4,4'DDE	BDL	4.1
3P. 4,4'-DDT	BDL	4.1
4P. ALDRIN	BDL	1.2
5P. CHLORDANE	BDL	4.7
6P. DIELDRIN	BDL	1.8
7P. ENDOSULFAN I	BDL	1.8
8P. ENDOSULFAN II	BDL	4.1
9P. ENDOSULFAN SULFATE	BDL	2.4
10P. ENDRIN	BDL	3
11P. ENDRIN ALDEHYDE	BDL	1.2
12P. HEPTACHLOR	BDL	1.2
13P. HEPTACHLOR EPOXIDE	BDL	1.2
14P. KEPONE	BDL	1.2
15P. p,p'-METHOXYCHLOR	BDL	4.1
16P. TOXAPHENE	BDL	24
17P. ALPHA-BEC	BDL	1.2
18P. BETA-BEC	BDL	1.2
19P. DELTA-BEC	BDL	1.2
20P. GAMMA-BEC (Lindane)	BDL	1.2

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

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APPENDIX VIII, IX - PESTICIDES, METHOD 8080
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(Page 2)

SAMPLE IDENTIFIER: R001B1416
COMPUCHEM SAMPLE NUMBER: 397161
DRY WEIGHT FACTOR: 1.18

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Dibutylchlorendate	94	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



COMPUCHEM
LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill/Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 1)

SAMPLE IDENTIFIER: R002B0810
COMPUCHEM SAMPLE NUMBER: 397165
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. 4,4'-DDD	BDL	4.2
2P. 4,4'-DDE	BDL	4.2
3P. 4,4'-DDT	BDL	4.2
4P. ALDRIN	BDL	1.2
5P. CHLORDANE	BDL	4.3
6P. DIELDRIN	BDL	1.3
7P. ENDOSULFAN I	BDL	1.8
8P. ENDOSULFAN II	BDL	4.2
9P. ENDOSULFAN SULFATE	BDL	2.4
10P. ENDRIN	BDL	3
11P. ENDRIN ALDEHYDE	BDL	1.2
12P. HEPTACHLOR	BDL	1.2
13P. HEPTACHLOR EPOXIDE	BDL	1.2
14P. KEPONE	BDL	1.2
15P. p,p'-METEOXYCHLOR	BDL	4.2
16P. TOXAPENE	BDL	24
17P. ALPHA-BHC	BDL	1.2
18P. BETA-BEC	BDL	1.2
19P. DELTA-BEC	BDL	1.2
20P. GAMMA-BHC (Lindane)	BDL	1.2

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variations from the nominal sample weight and dry weight.

(Continued)



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COMPOUND LIST

APPENDIX VIII, IX - PESTICIDES, METHOD 8080
RESULTS REPORTED ON DRY WEIGHT BASIS
(Page 2)

SAMPLE IDENTIFIER: R002B0810
COMPUCHEM SAMPLE NUMBER: 397165
DRY WEIGHT FACTOR: 1.21

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	† Recovery	Control Range †
Dibutylchlorendate	96	(20-150)*

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.

PCB ANALYSIS (SOIL) (CONT'D.)

YB-5, 4-8' - Soil sample from Oxbow Area J, Soil Boring YB-5 at 4-8 feet

ROA010002 to
ROA012224 - Soil samples from Oxbow Area A, Groundwater Monitoring Well A-1 at 0-24 feet (2-foot intervals)

ROA2B0002 to
ROA2B1416 - Soil samples from Oxbow Area A, Soil Boring A-2 at 0-16 feet (2-foot intervals)

ROA3B0002 to
ROA3B2022 - Soil samples from Oxbow Area A, Groundwater Monitoring Well A-3 at 0-22 feet (2-foot intervals)

ROA3-DP1 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 16-18 feet (Duplicate)

ROB1B0002 to
ROB1B1820 - Soil samples from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-20 feet (2-foot intervals)

ROB-DPP1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 8-10 feet (Duplicate)

ROB2B0002 to
ROB2B1618 - Soil samples from Oxbow Area B, Groundwater Monitoring Well B-2 at 0-18 feet (2-foot intervals)

ROB-DPP2 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 10-12 feet (Duplicate)

ROC010002 to
ROC012224 - Soil samples from Oxbow Area C, Groundwater Monitoring Well C-1 at 0-24 feet (2-foot intervals)

ROC-DPP1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 0-2 feet (Duplicate)

ROC020002 to
ROCO21820 - Soil samples from Oxbow Area C, Groundwater Monitoring Well C-2 at 0-20 feet (2-foot intervals)

ROC-DPP2 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 22-24 feet (Duplicate)

ROC3B0002 to
ROC3B1214 - Soil samples from Oxbow Area C, Soil Boring C-3 at 0-14 feet (2-foot intervals)

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1

ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2

ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

ROO1B0002 to
ROO1B1820 - Soil samples from Oxbow Area K, Soil Boring K1 at 0-20 feet (2-foot intervals)

ROO2B0002 to
ROO2B0810 - Soil samples from Oxbow Area K, Soil Boring K2 at 0-10 feet (2-foot intervals)

PCB ANALYSIS (SOIL) (CONT'D)

19-4-14A, (0-6") - Soil sample from Oxbow Area B, Surficial Soil Sample 19-4-14A, 0-6 inches.

19-4-14B, (0-6") - Soil sample from Oxbow Area B, Surficial Soil Sample 19-4-14B, 0-6 inches.

19-4-14C, (0-6") - Soil sample from Oxbow Area B, Surficial Soil Sample 19-4-14C, 0-6 inches.

19-4-14D, (0-6") - Soil sample from Oxbow Area B, Surficial Soil Sample 19-4-14D, 0-6 inches.

19-4-14E, (0-6") - Soil sample from Oxbow Area B, Surficial Soil Sample 19-4-14E, 0-6 inches.

OX-J-SS1 - Soil sample from Oxbow Area J, Surficial Sample OX-J-SS1 from 0-4 inches

OX-J-SS2 - Soil sample from Oxbow Area J, Surficial Sample OX-J-SS2 from 0-4 inches

OX-J-SS3 - Soil sample from Oxbow Area J, Surficial Sample OX-J-SS3 from 0-4 inches

OX-J-SS4 - Soil sample from Oxbow Area J, Surficial Sample OX-J-SS4 from 0-4 inches

OX-J-SS5 - Soil sample from Oxbow Area J, Surficial Sample OX-J-SS5 from 0-4 inches

OX-J-DUP - Duplicate Soil sample from Oxbow Area J, Surficial Sample OX-J-SS5 from 0-4 inches

OX-J-SS6 - Soil sample from Oxbow Area J, Surficial Sample OX-J-SS6 from 0-4 inches

A-3 - Soil sample from Oxbow Area A, Surficial Soil Sample A-3 from 0-4 inches

B-1 - Soil sample from Oxbow Area B, Surficial Soil Sample B-1 from 0-4 inches

B-2 - Soil sample from Oxbow Area B, Surficial Soil Sample B-2 from 0-4 inches

C-2 - Soil sample from Oxbow Area C, Surficial Soil Sample C-2 from 0-4 inches

C2-10N(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

C2-10S(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

PCB ANALYSIS (SOIL) (CONT'D)

— C2-10E(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

— C2-10W(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

— C2-20N(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

— C2-20S(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

— C2-20E(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

— C2-20W(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches

— C2-30E(0-6") - Soil sample from Oxbow Area C, Surficial Soil Sample at C-2 from 0-6 inches



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ANALYTICAL SERVICES

CERTIFICATE OF ANALYSIS

Geraghty & Miller, Inc.
125 East Bethpage Road
Plainview, NY 11803
ATTN: Dennis Colton

October 16, 1989

Job Number: GMIN 44299

P.O. Number: NY0360SS06

This is the Certificate of Analysis for the following samples:

Client Project ID: GE Pittsfield
Date Received by Lab: 10/06/89
Number of Samples: Twenty (20)
Sample Type: Soil

PCB's ANALYSIS

Results in mg/kg (ppm) dry weight

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
FP-1,0-4'	JJ7110	0.05 U	0.05 U	0.05 U	0.05 U
FP-1,4-8'	JJ7111	0.05 U	0.05 U	0.05 U	0.05 U
FP-1,8-12'	JJ7112	0.05 U	0.05 U	0.05 U	0.05 U
FP-1,12-16'	JJ7113	0.05 U	0.05 U	0.05 U	0.05 U

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

Date of Extraction: 10/09/89

Date of Analysis: 10/11/89

Reviewed and Approved:

Alyce Moore
Alyce Moore
Laboratory Manager

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44299

PCB's ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
FP-1,16-20'	JJ7114	0.05 U	0.05 U	0.05 U	0.05 U
FP-1,20-24'	JJ7115	0.05 U	0.05 U	0.05 U	0.05 U
FP-1,24-28'	JJ7116	0.05 U	0.05 U	0.05 U	0.05 U
FP-1,28-30'	JJ7117	0.05 U	0.05 U	0.05 U	0.05 U
FP-2,0-4'	JJ7118	0.64 *	12	0.8 U	13
FP-2,4-8'	JJ7119	0.05 U	0.05 U	0.05 U	0.05 U
FP-3,0-4'	JJ7120	0.05 U	2.4 *	0.41	2.8
FP-3,4-8'	JJ7121	0.05 U	0.33	0.05	0.38
FP-4,0-4'	JJ7122	0.05 U	0.05 U	0.19 *	0.19 *
FP-4,4-8'	JJ7123	0.05 U	0.05 U	0.05 U	0.05 U
SA-1,0-4'	JJ7124	0.05 U	0.16 *	0.09 *	0.25
SA-1,4-8'	JJ7125	0.05 U	0.05 U	0.05 U	0.05 U
SA-1,8-12'	JJ7126	0.05 U	0.05 U	0.05 *	0.05 *
SA-2,0-4'	JJ7127	0.05 U	0.10 *	0.10 *	0.20
SA-2,4-8'	JJ7128	0.05 U	0.08 *	0.05 *	0.13
SA-2,8-12'	JJ7129	0.05 U	0.05 U	0.05 U	0.05 U
Method Blank	BL4909	0.05 U	0.05 U	0.05 U	0.05 U

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

* - Sample exhibits alteration of standard Aroclor pattern.

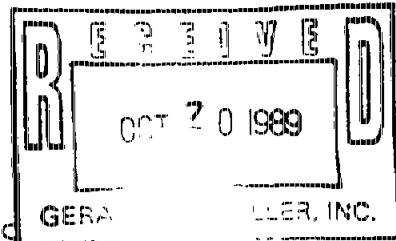
Date of Extraction: 10/09/89

Date of Analysis: 10/11/89



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CERTIFICATE OF ANALYSIS

Geraghty & Miller, Inc.
125 East Bethpage Road
Plainview, NY 11803
ATTN: Dennis Colton

October 16, 1989

Job Number: GMIN 44313

P.O. Number: NY0360SS06

This is the Certificate of Analysis for the following samples:

Client Project ID: GE Pittsfield
Date Received by Lab: 10/07/89
Number of Samples: Ten (10)
Sample Type: Soil

PCB's ANALYSIS

Results in mg/kg (ppm) dry weight

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
YB-1, 0-4'	JJ7172	0.05 U	0.60 *	0.35 *	0.95
YB-1, 4-8'	JJ7173	0.05 U	0.25 *	0.18 *	0.43
YB-2, 0-4'	JJ7174	0.05 U	1.2 *	1.1 *	2.3
YB-2, 4-8'	JJ7175	0.05 U	0.63 *	0.17 *	0.80

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

* - Sample exhibits alteration of standard Aroclor pattern.

Date of Extraction: 10/09/89

Date of Analysis: 10/12/89

Reviewed and Approved:

Alyce Moore
Alyce Moore
Laboratory Manager

Geraghty & Miller, Inc.
October 16, 1989

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE Pittsfield

Job Number: GMIN 44313

PCB's ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
YB-3, 0-4"	JJ7176	0.05 U	0.20 *	0.37 *	0.57
YB-3, 4-8"	JJ7177	0.05 U	0.05 U	0.05 U	0.05 U
YB-4, 0-4"	JJ7178	0.05 U	0.28 *	0.27 *	0.55
YB-4, 4-8"	JJ7179	0.05 U	0.05 U	0.05 U	0.05 U
YB-5, 0-4"	JJ7180	0.05 U	0.09 U	1.8	1.8
YB-5, 4-8"	JJ7181	0.05 U	0.05 U	0.08 *	0.08 *
Method Blank	BL4915	0.05 U	0.05 U	0.05 U	0.05 U

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

* - Sample exhibits alteration of standard Aroclor pattern.

Date of Extraction: 10/09/89

Date of Analysis: 10/12/89

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA010002	RR9997	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	Total <u>Aroclors</u>
ROA010204	RR9998	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor</u> <u>1016, 1232,</u> <u>1242† and/or 1248</u>	<u>Aroclor</u> <u>1254</u>	<u>Aroclor</u> <u>1260</u>	<u>Total</u> <u>Aroclors</u>
ROA010406	RR9999	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA010608	SS0001	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	<u>Aroclor</u> <u>1254</u>	<u>Aroclor</u> <u>1260</u>	Total Aroclors
ROA010810	SS0002	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA011012	SS0003	0.05 U	0.05 U	0.06 *	0.06

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA011214	SS0004	0.05 U	0.05 U	0.18 *	0.18

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROA011416	SS0005	0.09 U	0.13 U	0.89 *	0.89

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA011618	SS0006	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROA011820	SS0009	0.05 U	0.05 U	0.06 *	0.06

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA012022	SS0010	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 24, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05402/GE-Oxburn A

Job Number: GECP 49948

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROA012224	SS0011	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/15/91

Analysis Date: 11/22/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B0002	SS1163	0.05 U	0.21	0.17	0.38

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B0204	SS1164	0.05 U	1.8	1.3	3.1

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B0406	SS1165	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B0608	SSI166	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B0810	SS1167	0.05 U	0.05 U	0.19 *	0.19

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B1012	SS1168	0.05 U	0.11	0.24	0.35

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA2B1214	SS1169	0.05 U	0.16	0.19	0.35

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA2B1416	SS1170	0.07 U	0.74 U	3.4	3.4

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
February 3, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor <u>1016, 1232,</u> <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA3B0002	SS3401	0.27 U	7.3 *	18 *	25

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA3B0204	SS3402	0.05 U	0.09 U	3.0 *	3.0

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor</u> 1016, 1232, 1242† and/or 1248	<u>Aroclor</u> 1254	<u>Aroclor</u> 1260	<u>Total</u> <u>Aroclors</u>
ROA3B0406	SS3403	0.23 U	4.4 *	13 *	17

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA3B0608	SS3404	0.05 U	0.06 U	0.29 *	0.29

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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February 3, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA3B0810	SS3405	0.56 U	8.4 U	50 *	50

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROA3B1012	SS3406	0.05 U	0.13 U	0.87 *	0.87

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15/92 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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Client Project ID: AY05002/GE-Oxbow A

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA3B1214	SS3409	0.05 U	0.33 *	1.1 *	1.4

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern

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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor 1254	Aroclor 1260	Total Aroclors
ROA3B1416	SS3410	0.09 U	0.33 *	1.6 *	1.9

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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Client Project ID: AY05002/GE-Oxbow A

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total <u>Aroclors</u>
ROA3B1618	SS3411	0.09 U	0.17 U	1.9 *	1.9

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROA3B1820	SS3412	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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February 3, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor</u> <u>1016, 1232,</u> <u>1242† and/or 1248</u>	<u>Aroclor</u> <u>1254</u>	<u>Aroclor</u> <u>1260</u>	<u>Total</u> <u>Aroclors</u>
ROA3B2022	SS3413	0.05 U	0.18 *	0.44 *	0.62

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50353

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROA3-DP1	SS3414	0.15 U	0.20 U	3.0 *	3.0

Extraction Date: 01/09/92

Analysis Date: 01/14, 01/15 and 01/16/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B1B0002	SS1172	0.28 U	3.0 U	15	15

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROB1B0204	SS1173	0.47 U	19	4.1	23

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROB1B0406	SS1174	1.1 U	11 *	1.3 U	11

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B1B0608	SS1175	3.6 U**	1.3	4.9	6.2

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

** - Higher detection limit due to interference.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor		Aroclor	Aroclor	Total
		1016, 1232, 1242† and/or <u>1248</u>	<u>1254</u>			
ROB1B0810	SS1176	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B1B1012	SS1177	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B1B1214	SS1180	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROB1B1416	SS1181	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R0B1B1618	SS1182	0.17 U	1.9	0.75	2.6

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROB1B1820	SS1183	0.05 U	0.21 *	0.05 U	0.21

— Extraction Date: 12/02/91
Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total <u>Aroclors</u>
ROB-DPP1	SS1184	0.05 U	0.10	0.05 U	0.10

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROB2B0002	SS1200	4.3 U	140	42	180

Extraction Date: 11/27/91
Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B2B0204	SS1201	0.05 U	0.47	2.2	2.7

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
R0B2B0406	SS1202	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total <u>Aroclors</u>
R0B2B0608	SS1203	0.05 U	0.22	0.51	0.73

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROB2B0810	SS1206	0.05 U	0.05 U	0.09	0.09

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROB2B1012	SS1207	0.38 U	5.0	12	17

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B2B1214	SS1208	0.05 U	0.53	0.91	1.4

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
R0B2B1416	SS1209	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROB2B1618	SS1210	0.05 U	0.15	0.44	0.59

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE Oxbow B

Job Number: GECP 50077

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
R0B-DPP2	SS1211	0.05 U	0.61	1.2	1.8

Extraction Date: 11/27/91

Analysis Date: 12/13 and 12/16/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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December 31, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO10002	RR9792	0.05 U	0.79	0.05 U	0.79

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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December 31, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

lient Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total <u>Aroclors</u>
ROCO10204	RR9793	0.05 U	0.20 U	1.1	1.1

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO10406	RR9794	0.47 U	19	1.4 U	19

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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December 31, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor <u>1016, 1232,</u> <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROCO10608	RR9795	0.11 U	2.8	0.53	3.3

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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December 31, 1991

IT ANALYTICAL SERVICES
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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROCO10810	RR9796	0.22 U	8.7	0.57 U	8.7

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
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lient Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROCO11012	RR9799	0.39 U	11	1.4 U	11

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO11214	RR9800	1.2 U	57	2.7 U	57

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

f - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 31, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO11618	RR9802	1.1 U	49	2.8 U	49

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROCO11820	RR9803	0.24 U	13	0.62 U	13

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO12022	RR9804	0.44 U	13	1.5 U	13

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO12224	RR9805	0.11 U	6.6	0.30 U	6.6

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total <u>Aroclors</u>
ROC-DPPI	RR9806	0.05 U	0.21 *	0.17 *	0.38

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Samples exhibits alteration of standard Aroclor pattern.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor <u>1016, 1232,</u> <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROCO20002	RR9784	10 U	41 U	750	750

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO20406	RR9785	0.51 U	11	84	95

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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5815 MIDDLEBROOK PIKE
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lient Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO20608	RR9786	0.58 U	12	30	42

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROCO20810	RR9787	1.0 U	2.4 U	81	81

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO21012	RR9788	0.49 U	1.1 U	11	11

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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KNOXVILLE, TN

lient Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO21214	RR9789	0.26 U	3.8 U	22	26

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

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ient Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROCO21416	RR9790	2.9 U	11 U	150	150

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 31, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

lient Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor		Aroclor 1254	Aroclor 1260	Total Aroclors
		1016, 1232, 1242† and/or 1248	0.06 U			
ROCO21820	RR9791		0.06 U	0.22 U	5.3	5.3

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
December 31, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow C

Job Number: GECP 49925

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROC-DPP2	RR9807	0.22 U	6.7	1.2 U	6.7

Extraction Date: 11/15/91

Analysis Date: 11/18/91 through 11/21/91 and 11/25/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROC3B0002	SS1185	0.05 U	0.72	0.21	0.93

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROC3B0204	SS1186	0.05 U	0.34	0.11	0.45

Extraction Date: 12/02/91
Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total <u>Aroclors</u>
ROC3B0406	SS1187	0.05 U	0.06	0.07	0.13

-- Extraction Date: 12/02/91
Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROC3B0608	SS1188	0.28 U	19	4.9	24

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
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KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors 0.05 U
ROC3B0810	SS1189	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
ROC3B1012	SS1190	0.05 U	0.05 U	0.05 U	0.05 U

— Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 28, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow A

Job Number: GECP 50073

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, 1242† and/or <u>1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROC3B1214	SS1191	0.05 U	0.05 U	0.05 U	0.05 U

Extraction Date: 12/02/91

Analysis Date: 12/10 through 12/13/91 and 12/17/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 29, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow J

Job Number: GECP 50230

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, <u>1242† and/or 1248</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROJyS	SS2438	0.05 U	0.11 U	1.7 *	1.7

Extraction Date: 12/20/91

Analysis Date: 12/26 and 12/27/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
January 29, 1992

Client Project ID: AY05002/GE-Oxbow J

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Job Number: GECP 50230

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
ROI2/S	SS2441	0.05 U	0.05 U	0.53	0.53

Extraction Date: 12/20/91

Analysis Date: 12/26 and 12/27/91

—† — Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U — Compound was analyzed for but not detected. The number is the detection limit for the sample.

General Electric Company
January 29, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Oxbow J

Job Number: GECP 50230

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	Aroclor 1016, 1232, <u>1242† and/or 1243</u>	Aroclor <u>1254</u>	Aroclor <u>1260</u>	Total Aroclors
ROI49S	SS2443	0.05 U	0.11 U	1.9 *	1.9

Extraction Date: 12/20/91

Analysis Date: 12/26 and 12/27/91

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor</u> <u>1016, 1232,</u> <u>1242† and/or 1248</u>	<u>Aroclor</u> <u>1254</u>	<u>Aroclor</u> <u>1260</u>	<u>Total</u> <u>Aroclors</u>
R001B0002	PP7336	0.05 U	0.05 U	0.15 *	0.15

- † - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.
U - Compound was analyzed for but not detected. The number is the detection limit for the sample.
* - Sample exhibits alteration of standard Aroclor pattern.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B0204	PP7337	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B0406	PP7338	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B0608	PP7339	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B0810	PP7340	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R00181012	PP7343	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

-PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B1214	PP7344	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

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March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B1416	PP7345	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

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March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B1618	PP7346	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

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March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R001B1820	PP7347	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

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March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R002B0002	PP7348	0.05 U	0.05 U	0.07 *	0.07

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R002B0204	PP7349	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IIT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R002B0406	PP7350	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R002B0608	PP7351	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

General Electric Company
March 5, 1991

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: AY05002/GE-Housatonic River

Job Number: GECP 47699

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, 1242† and/or 1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>	<u>Total Aroclors</u>
R002B0810	PP7352	0.05 U	0.05 U	0.05 U	0.05 U

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 02/13/91

Date of Analysis: 02/14 to 02/20/91

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General Electric Company
September 9, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Project ID: GE-Housatonic River/101.97

Job Number: GECP 52153

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Aroclor 1016, 1232, †1242 and/or</u>			<u>Total Aroclors</u>	
		<u>1248</u>	<u>Aroclor 1254</u>	<u>Aroclor 1260</u>		
I7-2-32B, (0-6")	TT4372	0.45 U	2.6 U	24 *		24
I7-2-34B, (0-6")	TT4373	0.73 U	4.2 U	49 *		49
I8-24-5A, (0-6")	TT4374	0.48 U	2.8 U	38 *		38
I9-4-14A, (0-6")	TT4375	0.19 U	1.7 *	4.5 *		6.2
I9-4-14C, (0-6")	TT4376	0.51 U	35 *	12 *		47
I7-2-32A, (0-6")	TT4377	0.54 U	12 *	80 *		92
I7-2-34A, (0-6")	TT4378	0.13 U	0.58 U	12 *		12
I9-4-14B, (0-6")	TT4379	0.05 U	0.48 *	3.8 *		4.3
Method Blank	BLH1421	0.05 U	0.05 U	0.05 U		0.05 U

Extraction Date: 08/28/92

Analysis Date: 09/02/92

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
January 26, 1993

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: Housatonic River

Job Number: GECP 53182

PCBs ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID	Lab Sample ID	Aroclor 1016, 1232, †1242 and/or			Total Aroclors
		1248	Aroclor 1254	Aroclor 1260	
I9-4-14D (0"-6")	VV3777	0.08 U	0.19 U	4.3 *	4.3
I9-4-14D (6"-12")	VV3778	0.06 U	0.21 U	5.4 *	5.4
I9-4-14E (0"-6")	VV3779	0.05 U	0.17 U	3.5 *	3.5
I9-4-14E (6"-12")	VV3780	0.05 U	0.17 U	6.7 *	6.7
I7-2-32-C (0"-6")	VV3781	0.07 U	0.20 U	5.1 *	5.1
I7-2-32-C (6"-12")	VV3782	0.06 U	0.21 U	5.5 *	5.5
I7-2-32-D (0"-6")	VV3783	0.05 U	0.19 U	4.4 *	4.4
I7-2-32-D (6"-12")	VV3784	0.05 U	0.05 U	0.48 *	0.48
I7-2-34-C (0"-6")	VV3785	0.05 U	0.20 U	4.6 *	4.6
I7-2-34-C (6"-12")	VV3786	0.05 U	0.05 U	0.63 *	0.63
I7-2-34-D (0"-6")	VV3787	0.05 U	0.20 U	3.6 *	3.6
I7-2-34-D (6"-12")	VV3788	0.05 U	0.05 U	0.30 *	0.30
I7-3-7-C (0"-6")	VV3789	0.49 U	1.8 U	30 *	30
I7-3-7-C (6"-12")	VV3790	0.05 U	0.09 U	1.9 *	1.9
I7-3-7-D (0"-6")	VV3791	1.5 U	3.8 U	53 *	53

Extraction Date: 01/11/93

Analysis Date: 01/12, 01/13, 01/14, and 01/15/93

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

* - Sample exhibits alteration of standard Aroclor pattern.

General Electric Company
January 26, 1993

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Project ID: Housatonic River

Job Number: GECP 53182

SPIKED BLANK ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

Client Sample ID: Blank Spike
Lab Sample ID: H2887

<u>Compound</u>	<u>Conc.</u> <u>Spike Added</u>	<u>Conc.</u> <u>Blank Spike</u>	<u>%</u> <u>Rec.</u>
Aroclor 1242	1.0	0.77	77

Date of Extraction: 01/11/93
Date of Analysis: 01/14/93

000004

General Electric Company
September 9, 1992

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

Client Project ID: GE-Housatonic River/101.97

Job Number: GECP 52153

TOTAL ORGANIC CARBON ANALYSIS

Results in mg/kg (ppm)

Sample Matrix: Soil

<u>Client Sample ID</u>	<u>Lab Sample ID</u>	<u>Result</u>
Method Blank	P4201	1 U
I7-2-32B, (0-6")	TT4372	10,000
I7-2-34B, (0-6")	TT4373	19,000
I8-24-5A, (0-6")	TT4374	12,000
I9-4-14A, (0-6")	TT4375	14,000
I9-4-14C, (0-6")	TT4376	13,000

Date of Analysis: 09/01/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-J-SS1	Extraction Date:	09/16/94
Lab Sample ID:	AB9829	Analysis Date:	09/16, 19/94
Sample Matrix:	SOIL	Concentration Units:	µg/kg (ppb)

Aroclor 1016, 1232 †1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
120 U	240 U	630 +	630 +

+ - Positive result.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

- Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-J-SS2	Extraction Date:	09/16/94
Lab Sample ID:	AB9830	Analysis Date:	09/16, 19/94
Sample Matrix:	SOIL	Concentration Units:	µg/kg (ppb)

Aroclor 1016, 1232 †1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
240 U	470 U	1600 +	1600 +

+ - Positive result.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

* - Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-J-SSG	Extraction Date:	09/16/94
Lab Sample ID:	AB9831	Analysis Date:	09/16, 19/94
Sample Matrix:	SOIL	Concentration Units:	µg/kg (ppb)

Aroclor 1016, 1232 †1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
250 U	500 U	1500 +	1500 +

+ - Positive result.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

'' - Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-J-SS4	Extraction Date:	09/16/94
Lab Sample ID:	AB9832	Analysis Date:	09/16,19/94
Sample Matrix:	SOIL	Concentration Units:	µg/kg (ppb)

Aroclor 1016, 1232 +1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
260 U	510 U	1300 +	1300 +

+ - Positive result.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-I-SSS	Extraction Date:	09/16/94
Lab Sample ID:	AB9833	Analysis Date:	09/16, 19/94
Sample Matrix:	SOIL	Concentration Units:	$\mu\text{g}/\text{kg}$ (ppb)

Aroclor 1016, 1232 †1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
47 U	94 U	290 +	290 +

+ - Positive result.

* - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

† - Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-J-DUP	Extraction Date:	09/16/94
Lab Sample ID:	AB9835	Analysis Date:	09/16, 19/94
Sample Matrix:	SOIL	Concentration Units:	µg/kg (ppb)

Aroclor 1016, 1232 +1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
51 U	100 U	350 +	350 +

+ - Positive result.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

PCBs ANALYSIS

Laboratory Name:	QUANTERRA-KNOXVILLE	SDG Number:	N/A
Contract Name:	GENERAL ELECTRIC	Job Number:	1380
Client Sample ID:	OX-J-SS6	Extraction Date:	09/16/94
Lab Sample ID:	AB9834	Analysis Date:	09/16, 19/94
Sample Matrix:	SOIL	Concentration Units:	$\mu\text{g}/\text{kg}$ (ppb)

Aroclor 1016, 1232 †1242 &/or 1248	Aroclor 1254	Aroclor 1260	Total Aroclors
45 U	90 U	280 +	280 +

+ - Positive result.

† - Sample Aroclor pattern identified and/or calculated as Aroclor 1242.

- Compound was analyzed for but not detected. The number is the detection limit for the sample.



ANALYSIS INFORMATION

Table 1.0 Polychlorinated Biphenyls

Analysis EPA Method 8081		Extraction Method 3550		Analyst Amy Miine		Instrument GC-ECD		
Parameter	QC Lot:	Date Sampled	Date Extracted	Date Analyzed	Dilution Factor	Surr. %R	Total PCBs	MDL
Sample ID						(mg/Kg)	(mg/Kg)	
B-1	1010958081H2	10/10/95	10/11/95	10/11/95	10	108	4.18	0.100
B-2	1010958081H2	10/10/95	10/10/95	10/11/95	10	111	11.8	0.100
A-3	1010958081H2	10/10/95	10/11/95	10/11/95	1	104	0.397	0.100
C-2	1010958081H2	10/10/95	10/11/95	10/11/95	1000	DO	745	0.100
PS-DUP-1	1010958081H2	10/10/95	10/11/95	10/11/95	10	108	7.14	0.100

Note: Surrogate limit 41-131; % surrogate recovery limits are established guidelines suggested by the USEPA Contract Laboratory Statement of Work.

DO = Due to the high concentration of the sample, the surrogate standard was diluted beyond the normal quantitation range.



ANALYSIS INFORMATION

Table 1.0 Polychlorinated Biphenyls

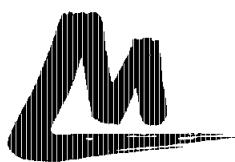
Analysis EPA Method 8081		Extraction Method 3550		Analyst ACM		Instrument GC-ECD		
Parameter	QC Lot:	Date Sampled	Date Extracted	Date Analyzed	Dilution Factor	Surr. %R	Total PCBs	MDL
						(mg/Kg)	(mg/Kg)	
Sample ID								
C2-10N (0-6")	1110958081-H	11/09/95	11/10/95	11/13/95	1000	DO	418	0.100
C2-10S (0-6")	1110958081-H	11/09/95	11/10/95	11/13/95	100	DO	36.8	0.100
C2-10E (0-6")	1110958081-H	11/09/95	11/10/95	11/13/95	100	DO	55.3	0.100
C2-10W(0-6")	1110958081-H	11/09/95	11/10/95	11/13/95	100	DO	22.3	0.100
C2-D-1 (0-6")	1110958081 H	11/09/95	11/10/95	11/13/95	100	DO	40.6	0.100
C2-20N (0-6")		11/09/95				HOLD		
C2-20S (0-6")		11/09/95				HOLD		
C2-20E (0-6")		11/09/95				HOLD		
C2-20W(0-6")		11/09/95				HOLD		

Note: Surrogate limit 41-131; % surrogate recovery limits are established guidelines suggested by the USEPA Contract Laboratory Statement of Work.

ND = Analyte of interest was non-detectable.

DO = Due to the high concentration of the sample, the surrogate standard was diluted beyond the normal quantitation range.

MDL = Method Detection Limit.



ANALYSIS INFORMATION

Table 1.0 Polychlorinated Biphenyls

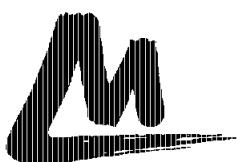
Analysis EPA Method 8081		Extraction Method 3550		Analyst ACM		Instrument GC-ECD		
Parameter	QC Lot:	Date Sampled	Date Extracted	Date Analyzed	Dilution Factor	Surr. %R	Total PCBs	MDL
							(mg/Kg)	(mg/Kg)
Sample ID								
C2-20N (0-6")	1110958081-H	11/09/95	11/15/95	11/15/95	10	107	5.58	0.100
C2-20S (0-6")	1110958081-H	11/09/95	11/15/95	11/15/95	10	106	7.55	0.100
C2-20E (0-6")	1110958081-H	11/09/95	11/15/95	11/15/95	100	DO	59.3	0.100

Note: Surrogate limit 41-131; % surrogate recovery limits are established guidelines suggested by the USEPA Contract Laboratory Statement of Work.

ND = Analyte of interest was not detected at or below the laboratory determined Analytical Method Detection Limit.

DO = The surrogate standard was diluted beyond the normal quantitation range of the analytical method.

MDL = Analytical Method Detection Limit.



ANALYSIS INFORMATION

Polychlorinated Biphenyls

Analysis Required EPA Method 8081	Extraction Method 3550A	Analyst JM	Instrument GC-ECD
Sample ID	C2-30E		MDL
Dilution Factor	10 (mg/Kg)		(mg/Kg)
Parameter PCBs	12.8		0.100

QC Lot:
1130958081-H1

ND = Analyte of interest was not detected at the laboratory determined Analytical Method Detection Limit.

MDL = Method Detection Limit

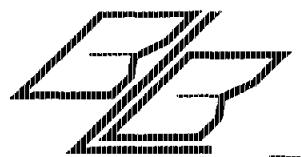
Polychlorinated Biphenyls

Analysis Required EPA Method 8081	Extraction Method 3510A	Analyst JM	Instrument GC-ECD
Sample ID	C2-RB-2		MDL
Dilution Factor	1 (ug/L)		(ug/L)
Parameter PCBs	ND		0.100

QC Lot:
1109958081-WV

ND = Analyte of interest was not detected at the laboratory determined Analytical Method Detection Limit.

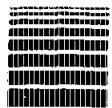
MDL = Method Detection Limit.



Section 4

ORGANOPHOSPHORUS PESTICIDES ANALYSIS (SOIL)

- ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-2 feet
- ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet
- ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)
- ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-2 at 0-2 feet
- ROC011012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet
- ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-1 at 2-4 feet
- ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1
- ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2
- ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3
- ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4
- R001B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet
- R002B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROA2B0608
COMPUCHEM SAMPLE NUMBER: 465051
DRY WEIGHT FACTOR: 1.10

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	11
2P. PHORATE	BDL	11
3P. DIMETHOATE	BDL	11
4P. DISULFOTON	BDL	11
5P. METHYL PARATHION	BDL	11
6P. PARATHION	BDL	11

BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	108	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROB1B0406
COMPUCHEM SAMPLE NUMBER: 465529
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEP)	BDL	12
2P. PHORATE	BDL	12
3P. DIMETHOATE	BDL	12
4P. DISULFOTON	BDL	12
5P. METHYL PARATHION	BDL	12
6P. PARATHION	BDL	12

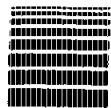
BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	116	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROB-DPA1
COMPUCHEM SAMPLE NUMBER: 465555
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	12
2P. PHORATE	BDL	12
3P. DIMETHOATE	BDL	12
4P. DISULFOTON	BDL	12
5P. METHYL PARATHION	BDL	12
6P. PARATHION	BDL	12

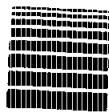
BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	106	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROB2B0002
COMPUCHEM SAMPLE NUMBER: 465881
DRY WEIGHT FACTOR: 1.50

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	15
2P. PROORATE	BDL	15
3P. DINETHOATE	BDL	15
4P. DISULFOTON	BDL	15
5P. METHYL PARATHION	BDL	15
6P. PARATHION	BDL	15

BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	104	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROC011012
COMPUCHEM SAMPLE NUMBER: 464370
DRY WEIGHT FACTOR: 1.11

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE (SULFOTEPP)	BDL	11
2P. PHORATE	BDL	11
3P. DIMETHOATE	BDL	11
4P. DISULFOTON	BDL	11
5P. METHYL PARATHION	BDL	11
6P. PARATHION	BDL	11

BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	66	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROC380204
COMPUCHEM SAMPLE NUMBER: 465099
DRY WEIGHT FACTOR: 1.08

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	11
2P. PHORATE	BDL	11
3P. DIMETHOATE	BDL	11
4P. DISULFOTON	BDL	11
5P. METHYL PARATHION	BDL	11
6P. PARATHION	BDL	11

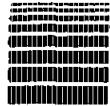
BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range †
Methidathion	64	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

†Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJLS
COMPUCHEM SAMPLE NUMBER: 469287
DRY WEIGHT FACTOR: 1.28

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	13
2P. PHORATE	BDL	13
3P. DIMETHOATE	BDL	13
4P. DISULFOTON	BDL	13
5P. METHYL PARATHION	BDL	13
6P. PARATHION	BDL	13

BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	106	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

*Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJ2S
COMPUCHEM SAMPLE NUMBER: 469278
DRY WEIGHT FACTOR: 1.18

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE (SULFOTEPP)	BDL	12
2P. PHORATE	BDL	12
3P. DIMETHOATE	BDL	12
4P. DISULFOTON	BDL	12
5P. METHYL PARATHION	BDL	12
6P. PARATHION	BDL	12

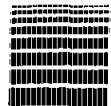
†
BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	104	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPOHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJ3S
COMPUCHEM SAMPLE NUMBER: 469299
DRY WEIGHT FACTOR: 1.10

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	11
2P. PEORATE	BDL	11
3P. DIMETHOATE	BDL	11
4P. DISULFOTON	BDL	11
5P. METHYL PARATHION	BDL	11
6P. PARATHION	BDL	11

BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	52	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJ4S
COMPUCHEM SAMPLE NUMBER: 469264
DRY WEIGHT FACTOR: 1.49

	CONCENTRATION ($\mu\text{g}/\text{kg}$)	DETECTION + LIMIT ($\mu\text{g}/\text{kg}$)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	15
2P. PBORATE	BDL	15
3P. DINETEAOATE	BDL	15
4P. DISULFOTON	BDL	15
5P. METHYL PARATHION	BDL	15
6P. PARATEION	BDL	15

BDL=BELOW DETECTION LIMIT

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	104	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

*Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: R001B1416
COMPUCHEN SAMPLE NUMBER: 398418
DRY WEIGHT FACTOR: 1.18

	CONCENTRATION ($\mu\text{g}/\text{kg}$)	DETECTION + LIMIT ($\mu\text{g}/\text{kg}$)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	11
2P. PHORATE	BDL	11
3P. DIMETHOATE	BDL	11
4P. DISULFOTON	BDL	11
5P. METHYL PARATHION	BDL	11
6P. PARATHION	BDL	11

BDL=BELOW DETECTION LIMIT

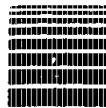
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range †
Methidathion	125 **	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

†Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.

**See Laboratory Notice # 1.



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COMPOUND LIST
APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: RCO2B0810
COMPUCHEM SAMPLE NUMBER: 398420
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION ($\mu\text{g}/\text{kg}$)	DETECTION + LIMIT ($\mu\text{g}/\text{kg}$)
1P. TETRAETHYLDITHIOPYROPHOSPHATE(SULFOTEPP)	BDL	12
2P. PHORATE	BDL	12
3P. DIMETHOATE	BDL	12
4P. DISULFOTON	BDL	12
5P. METHYL PARATHION	BDL	12
6P. PARATHION	BDL	12

BDL=BELOW DETECTION LIMIT

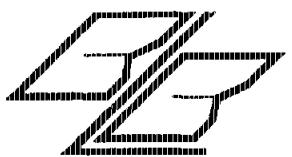
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
Methidathion	138 **	(60-120)*

*Advisory surrogate. See Quality Assurance Notice

+Detection limits have been adjusted to report variation from the nominal sample weight and dry weight.

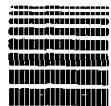
**See Laboratory Notice # 1.



Section 5

HERBICIDES ANALYSIS (SOIL)

- ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet
- ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet
- ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet
- ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet
- ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)
- ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet
- ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet
- ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)
- ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet
- ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1
- ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2
- ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3
- ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4
- R001B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet
- R002B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet



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RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROA012224
COMPUCHEM SAMPLE NUMBER: 462142
DRY WEIGHT FACTOR: 1.11

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	110
2. 2,4,5-TP (Silvex)	BDL	28
3. 2,4,5-T	BDL	28

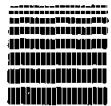
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	104	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
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RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROA2B0608
COMPUCHEM SAMPLE NUMBER: 465048
DRY WEIGHT FACTOR: 1.10

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	110
2. 2,4,5-TP (Silver)	BDL	27
3. 2,4,5-T	BDL	27

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range †
2,4-DB	118	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight, the dry weight, and dilution.

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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROB1B0406
COMPUCHEM SAMPLE NUMBER: 465525
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	120
2. 2,4,5-TP (Silvex)	BDL	30
3. 2,4,5-T	BDL	30

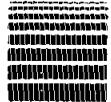
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties.[†] A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	78	(20-150)*

BDL=BELOW DETECTION LIMIT

[†]Detection limits have been adjusted to report variation from the nominal sample weight, the dry weight, and dilution.

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROB280002
COMPUCHEM SAMPLE NUMBER: 465878
DRY WEIGHT FACTOR: 1.50

	CONCENTRATION (ug/kg)	DETECTION LIMIT (ug/kg)
1. 2,4-D	BDL	150
2. 2,4,5-TP (Silvex)	BDL	37
3. 2,4,5-T	BDL	37

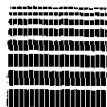
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

Recovery Control Range
2,4-DB 107 (20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight, the dry weight, and dilution.

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST

APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROB-DPA1
COMPUCHEM SAMPLE NUMBER: 465551
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	120
2. 2,4,5-TP (Silvex)	BDL	30
3. 2,4,5-T	BDL	30

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	95	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROC011012
COMPUCHEM SAMPLE NUMBER: 461434
DRY WEIGHT FACTOR: 1.11

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	110
2. 2,4,5-TP (Silvex)	BDL	28
3. 2,4,5-T	BDL	28

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	79	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROC021214
COMPUCHEM SAMPLE NUMBER: 461156
DRY WEIGHT FACTOR: 1.23

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	120
2. 2,4,5-TP (Silvex)	BDL	31
3. 2,4,5-T	BDL	31

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range †
2,4-DB	83	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

†Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 3150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROC-DPA1
COMPUCHEM SAMPLE NUMBER: 461189
DRY WEIGHT FACTOR: 1.85

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	180
2. 2,4,5-TP (Silvex)	BDL	46
3. 2,4,5-T	BDL	46

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	77	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST

APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROC3B0204

COMPUCHEM SAMPLE NUMBER: 465082

DRY WEIGHT FACTOR: 1.08

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	110
2. 2,4,5-TP (Silvex)	BDL	27
3. 2,4,5-T	BDL	27

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

Recovery Control Range
2,4-DB 85 (20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight, the dry weight, and dilution.

*Advisory surrogate: with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJIS
COMPUCHEM SAMPLE NUMBER: 469283
DRY WEIGHT FACTOR: 1.28

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	130
2. 2,4,5-TP (Silvex)	BDL	32
3. 2,4,5-T	BDL	32

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

† Recovery Control Range †		
2,4-DB	62	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJ2S
COMPUCHEM SAMPLE NUMBER: 469274
DRY WEIGHT FACTOR: 1.18

	CONCENTRATION (ug/kg)	DETECTION LIMIT (ug/kg)
1. 2,4-D	BDL	120
2. 2,4,5-TP (Silvex)	BDL	29
3. 2,4,5-T	BDL	29

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

Recovery	Control Range
78	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJ3S
COMPUCHEM SAMPLE NUMBER: 469295
DRY WEIGHT FACTOR: 1.10

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	110
2. 2,4,5-TP (Silvex)	BDL	27
3. 2,4,5-T	BDL	27

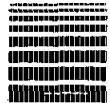
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	75	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



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COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: ROJ4S
COMPUCHEM SAMPLE NUMBER: 469258
DRY WEIGHT FACTOR: 1.49

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	150
2. 2,4,5-TP (Silvex)	BDL	37
3. 2,4,5-T	BDL	37

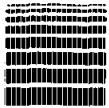
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	63	(20-150)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 20% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



**COMPUCHEM
LABORATORIES, INC.**

P.O. Box 12652 3308 Chapel Hill/Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

COMPOUND LIST

APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: R001B1416

COMPUCHEM SAMPLE NUMBER: 398419

DRY WEIGHT FACTOR: 1.18

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	220	120
2. 2,4,5-TP (Silvex)	51	29
3. 2,4,5-T	52	29

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	104	(16-124)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



**COMPUCHEM
LABORATORIES, INC.**

P.O. Box 12652 3308 Chapel Hill / Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

COMPOUND LIST
APPENDIX VIII, IX - HERBICIDES, METHOD 8150
RESULTS REPORTED ON DRY WEIGHT BASIS

SAMPLE IDENTIFIER: R002B0810
COMPUCHEM SAMPLE NUMBER: 398423
DRY WEIGHT FACTOR: 1.21

	CONCENTRATION (ug/kg)	DETECTION + LIMIT (ug/kg)
1. 2,4-D	BDL	120
2. 2,4,5-TP (Silver)	BDL	30
3. 2,4,5-T	BDL	30

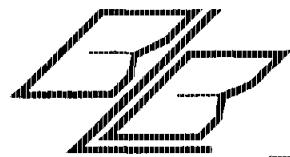
Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analyties. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	Recovery	Control Range
2,4-DB	94	(16-124)*

BDL=BELOW DETECTION LIMIT

*Detection limits have been adjusted to report variation from the nominal sample weight and the dry weight.

*Advisory surrogate; with the exception of dilutions recovery below 10% requires an action step (re-extraction and reanalysis). See Quality Assurance Notice.



Section 6

DIOXIN/FURAN ANALYSIS (SOIL)

ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet

ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet

ROA3B1214 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 12-14 feet

ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet

ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet

ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)

ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet

ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet

ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)

ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1

ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2

ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3

ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

ROO1B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet

ROO2B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet

OX-J-SS1 - Soil sample from Oxbow Area J, Soil Surficial Sample OX-J-SS1 at 0-4 inches

OX-J-SS2 - Soil sample from Oxbow Area J, Soil Surficial Sample OX-J-SS2 at 0-4 inches

OX-J-SS3 - Soil sample from Oxbow Area J, Soil Surficial Sample OX-J-SS3 at 0-4 inches

OX-J-SS4 - Soil sample from Oxbow Area J, Soil Surficial Sample OX-J-SS4 at 0-4 inches

OX-J-SS5 - Soil sample from Oxbow Area J, Soil Surficial Sample OX-J-SS5 at 0-4 inches

OX-J-DUP - Duplicate soil sample from Oxbow Area J, from location OX-J-SS5 at 0-4 inches

DIOXIN/FURAN ANALYSIS (SOIL) (CONT'D)

OX-J-SS6 -

**Soil sample from Oxbow Area J, Soil Surficial Sample OX-J-SS6 at 0-4
inches**

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 09/14/92

LABORATORY: ChemWest

Ticket# CW-8950

Project Name: General Electric

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	TOTAL ANALYTE QUANTITY FOUND (ppb or ng/g)											
					TCDD	TCDF	PoCDD	HxCDD	HpCDD	OCDD	TCDF	TCDF	PoCDF	HxCDF	HpCDF	OCDF
ROA012224 // 462174	8950	11/20/91	18:51	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit					0.012	0.012	0.017	0.031	0.020	0.060	0.0065	0.035	0.012	0.019	0.019	0.038

* = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTITATION REPORT

Ticket# CW-8950
 Project Name: General Electric

PAGE 2 of 2
 DATE: 09/14/92
 LABORATORY: ChemWest

CLIENT ID.	GC/MS CW#	GC/MS DATE	INST. TIME	ABSOLUTE % RECOVERY OF INTERNAL STANDARDS							SURROGATE % ACCURACY		
				ID.	*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-ODDD	*C-TCDF	*C-PeCDF		
ROA012224 // 462174	8950	11/20/91 18:51	CW-2	86.7	91.1	91.4	85.4	65.6	88.9	92.7	99.2	101	96.0
Detection Limit													

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HPCDF = 13C12-1234678-HpCDF

Approved by: _____ 

FORM I - QUANTITATION REPORT

PAGE 1 of 2

DATE: 09/15/92

LABORATORY: ChemWest

Ticket# CW-9008

Project Name: General Electric Company

CLIENT ID.	SAMPLE CW#	SIZE	GC/MS DATE	GC/MS TIME	INST. ID.	2378	TOTAL ANALYTE QUANTITY FOUND (ppb or ng/g)								
							TCDD	TCDF	PeCDD	HxCDD	HxCDF	HpCDD	HpCDF	OCDD	OCDF
ROA2B0608 // 465062	9008-1A	9.27 G	12/11/91	19:04	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.71	0.71	0.87	0.18	0.054	0.070	1.2	6.1	0.24	0.14
ROA2B0608 // 465062	9008-1B	1.05 G	12/11/91	19:46	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.11	0.11	0.28	0.30	0.29	0.34	0.13	0.44	0.095	0.15
ROC3B0204 // 465135	9008-2A	9.46 G	12/11/91	20:26	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.84	0.84	3.9	1.7	0.11	0.21	1.6	12.8	0.59	0.86
ROC3B0204 // 465135	9008-2B	0.97 G	12/11/91	21:08	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.38	0.38	0.28	0.17	0.28	0.72	0.49	2.8	0.16	0.14

* = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: J

FORM 1 - QUANTITATION REPORT

PAGE 2 of 2

DATE: 09/15/92

LABORATORY: ChemWest

Ticket# CW-9008

Project Name: General Electric Company

CLIENT ID.	GC/MS CW#	GC/MS DATE	INST. TIME	ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C1-TCDD	*C-HxCDD	*C-HpCDF
ROA2B0608 // 465062	9008-1A	12/11/91 19:04	CW-2	1,6	2,1	9.2	38.0	39.8	0.60	4.1		85.2	229	89.9
Detection Limit														
ROA2B0608 // 465062	9008-1B	12/11/91 19:46	CW-2	71.8	81.2	91.2	84.8	56.7	55.3	81.9		101	96.9	91.6
Detection Limit														
RO03B0204 // 465135	9008-2A	12/11/91 20:26	CW-2	1.0	0.40	1.2	14.1	20.1	0.23	1.0		72.1	518	89.1
Detection Limit														
RO03B0204 // 465135	9008-2B	12/11/91 21:08	CW-2	21.8	52.1	79.0	74.3	47.3	9.1	57.8		99.2	102	94.5
Detection Limit														

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____ 

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 09/28/92

LABORATORY: ChemWest

Ticket #: CW-9154

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppb or ng/g)

CLIENT ID.	CH#	GC/MS DATE	GC/MS TIME	INST. ID.	2378	2378											
					TCDD	TCDF	PeCDD	HxCDD	HxCDF	HpCDD	OCDD	OCDF	TCDF	PeCDF	HxCDF	HxCDD	HpCDF
ROA381214 // 473666	9154	01/22/92	12:10	CW-2	ND 0.11	ND 0.11	ND 0.057	ND 0.026	ND 0.062	ND 0.025*	ND 0.43	ND 1.6	ND 0.022	ND 0.082	ND 0.073	ND 0.065	
Detection Limit																	
ROA381214 // 473666 MS	9154-MS	01/22/92	13:21	CW-2	9.3	9.3	9.9	31.5	9.7	10.1	9.6	9.6	16.0	37.5	18.5	12.2	
Detection Limit																	
ROA381214 // 473666 MSD	9154-MSD	01/22/92	14:02	CW-2	10.1	10.1	9.8	32.2	10.4	10.4	10.0	10.0	16.9	38.1	19.9	12.7	
Detection Limit																	

* = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 2 of 2

DATE: 09/28/92

LABORATORY: ChemWest

Ticket# CH-9154
 Project Name: General Electric Company

CLIENT	CH#	GC/MS	GC/MS	INST.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY				
					DATE	TIME	ID.	*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C1-TCDD	*C-HxCDD
DET	9154	01/22/92	12:10	CH-2	8.6	30.1	70.8	71.4	46.9	2.5	32.8			103	108	96.9
DET	9154-MS	01/22/92	13:21	CH-2	7.5	27.6	69.4	82.0	50.6	2.4	28.9			92.6	114	91.1
DET	9154-MSD	01/22/92	14:02	CH-2	8.5	29.4	69.8	75.1	48.9	2.9	30.9			98.6	110	91.8

INTERNAL STANDARDS

C-TCDD = 13C12-2378-TCDD
 C-PeCDD = 13C12-12378-PeCDD
 C-HxCDD = 13C12-123678-HxCDD
 C-HpCDD = 13C12-1234678-HpCDD
 C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378 -TCDD
 *C-HxCDD = 13C12-1237 89-HxCDD
 *C-PeCDF = 13C12-1237 8-PeCDF
 *C-HpCDF = 13C12-1234 678-HpCDF

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

Ticket# CW-9018-RX

DATE: 09/16/92

Project Name: General Electric Company

LABORATORY: ChemWest

TOTAL ANALYTE QUANTITY FOUND

CLIENT ID.	SAMPLE CW#	SIZE	GC/MS DATE	GC/MS TIME	INST. ID.	2378	(ppb or ng/g)											
							TCDD	TCDF	PeCDD	HxCDD	HpCDD	OCDD	TCDF	TCDF	PeCDF	HxCDF	HpCDF	OCDF
ROB180406 // 465506	9018-1BRX	0.96 G	12/29/91	15:05	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.2
Detection Limit						0.096	0.096	0.21	0.28	0.17	0.41	0.10	0.44	1.1				
ROB-DPA1 // 465509	9018-2BRX	0.98 G	12/29/91	16:21	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.98	6.9	4.1	5.3
Detection Limit						0.21	0.21	0.16	0.12	0.16	0.57	0.31	1.8					

ND = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTITATIVE REPORT

PAGE 2 of 2

DATE: 09/16/92

LABORATORY: ChemWest

Ticket# CW-9018-RX

Project Name: General Electric Company

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*CI-TCDD	*C-HxCDD	*C-HpCDF
ROB1B0406 // 465506	9018-1BRX	12/29/91	15:05	CW-2	75.0	92.7	102	91.4	65.2	61.2	91.4	100	96.9	98.4
Detection Limit														
ROB-DPAI // 465509	9018-2BRX	12/29/91	16:21	CW-2	30.9	60.8	90.9	89.7	62.2	15.3	65.3	101	102	97.8
Detection Limit														

INTERNAL STANDARDS

C-TCDD = 13C12-2378-TCDD
 C-PeCDD = 13C12-12378-PeCDD
 C-HxCDD = 13C12-123678-HxCDD
 C-HpCDD = 13C12-1234678-HpCDD
 C-TCDF = 13C12-2378-TCDF

SURROGATES

*CI-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____

FORM 1 - DETECTION REPORT

PAGE 1 of 2

DATE: 09/16/92

LABORATORY: ChemWest

Ticket# CW-9020-RX

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppb or ng/g)

CLIENT ID.	SAMPLE CW#	SIZE	GC/MS DATE	INST. TIME	2378 ID.	2378												
						TCDD	TCDF	PeCDD	HxCDD	HxCDF	OCDD	OCDF	TCDF	PeCDF	HxCDF	HxCDF	OCDF	
ROB2B0002 // 465887	9020-ARX	6.84 G	12/29/91	17:15	CW-2	ND	ND	ND	ND	0.17	0.66	0.10	0.51	0.59	0.48	0.26	0.24	
Detection Limit						0.016	0.016	0.014	0.022									

S = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTIFICATION REPORT

PAGE 2 of 2

DATE: 09/16/92

LABORATORY: ChemWest

Ticket# CW-9020-RX

Project Name: General Electric Company

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C-IxCDD	*C-HxCDD	*C-HpCDF
ROB2B0002 // 465887	9020-ARX	12/29/91	17:15	CW-2	73.8	95.9	107	98.3	62.1	53.5	94.1	99.7	97.6	95.4
Detection Limit														

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C-IxCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____ 

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 09/14/92

LABORATORY: ChemWest

Ticket# CW-8942

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppb or ng/g)

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	2378												
					TCDD	TCDF	PeCDD	HxCDD	HxCDF	OCDD	OCDF	TCDF	TCDF	PeCDF	HxCDF	HxCDF	OCDF
ROCO11012 // 461439	8942	11/20/91	17:57	CW-2	ND	ND	ND	ND	ND	0.30	ND	ND	ND	0.38	0.41	ND	ND
Detection Limit					0.015	0.024	0.030	0.024	0.095	0.040	0.040			0.11	0.076		

a = MAXIMUM POSSIBLE CONCENTRATION

#C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

#C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

#C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 2 of 2

DATE: 09/14/92

LABORATORY: ChemWest

Ticket# CW-8942

Project Name: General Electric Company

CLIENT ID.	GC/MS CW#	GC/MS DATE	INST. TIME	ABSOLUTE % RECOVERY of INTERNAL STANDARDS								SURROGATE % ACCURACY
				ID.	*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	
ROC011012 // 461439	8942	11/20/91	17:57	CW-2	49.6	45.5	76.9	67.7	48.5	26.4	72.8	98.1
Detection Limit												98.5
												97.8

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____

FORM 1 - QUANTITATION REPORT

Ticket# CW-4725

Project Name: Compuchem

PAGE 2 of 2

DATE: 09/15/92

LABORATORY: ChemWest

CLIENT ID.	CW#	GC/HS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY OF INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*CI-TCDD	*C-HxCDD	*C-HpCDF
ROC021214 // 461182	8933-1	11/25/91	13:21	CW-2	87.9	89.7	92.3	82.4	66.0	73.3	89.2	97.8	98.7	102
Detection Limit														
ROC021214 // 461182 MS	8933-1MS	11/25/91	14:03	CW-2	91.7	89.4	94.5	90.0	68.4	89.4	93.7	98.3	98.1	92.8
Detection Limit														
ROC021214 // 461182 MSD	8933-1MSD	11/25/91	14:43	CW-2	89.1	86.4	91.2	88.5	65.4	80.9	88.8	95.3	101	93.1
Detection Limit														
ROC-DPA1 // 461194	8933-2	11/25/91	15:22	CW-2	96.2	93.8	97.9	93.2	74.4	96.5	97.0	98.5	98.2	95.0
Detection Limit														

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*CI-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____ J _____

FORM 1 - QUANTITATION REPORT

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DATE: 09/15/92

LABORATORY: ChemWest

Ticket# CW-8933

Project Name: Compuchem

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	TOTAL ANALYTE QUANTITY FOUND (ppb or ng/g)											
					2378			2378			2378			2378		
TCDD	TCDD	PeCDD	HxCDD	HxCDD	OCDD	TCDF	TCDF	PeCDF	HxCDF	HxCDF	OCDF					
ROC021214 // 461182	8933-1	11/25/91	13:21	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit					0.027	0.027	0.032	0.051	0.088	0.18*	0.026	0.041	0.032	0.061	0.046	0.13
ROC021214 // 461182 MS	8933-1MS	11/25/91	14:03	CW-2	10.7	10.7	11.0	33.4	10.7	11.0	10.8	10.8	21.8	41.2	20.8	13.5
Detection Limit					10.5	10.5	11.2	34.2	10.7	11.1	10.8	10.8	22.0	41.9	20.9	13.8
ROC-DPA1 // 461194	8933-2	11/25/91	15:22	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit					0.025	0.025	0.027	0.060	0.094	0.16	0.026	0.063	0.036	0.036	0.10	0.16

* = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 09/16/92

LABORATORY: ChemWest

Ticket# CW-9085

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppb or ng/g)

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	2378	2378											
						TCDD	TCDF	PeCDD	HxCDD	HxCDF	OCDD	TCDF	PeCDF	HxCDF	HxCDF	OCDF	
ROJ4S // 469246	9085-1	01/13/92	14:39	CW-2		ND	ND	1.6	8.5	6.7	2.0	0.23	16.8	57.2	22.5	2.6	0.31
Detection Limit						0.028	0.027										
ROJ2S // 469249	9085-2	01/13/92	15:19	CW-2		ND	ND	ND	ND	0.061	0.21	ND	ND	0.15	0.17	ND	0.045
Detection Limit						0.019	0.035	0.011	0.014			0.018	0.018				0.068
ROJ1S // 469251	9085-3	01/13/92	15:58	CW-2		ND	ND	ND	ND	0.17	0.94	0.047	0.22	ND	0.39	0.25	0.18
Detection Limit						0.012	0.012	0.015	0.017							0.055	
ROJ3S // 469254	9085-4	01/13/92	16:39	CW-2		ND	ND	ND	ND								
Detection Limit						0.023	0.023	0.019	0.012	0.019	0.043	0.016	0.072	0.0076	0.013	0.019	0.024

ND = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzofuran (12 carbons)

Approved by: J

FORM 1 - QUANTIFICATION REPORT

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DATE: 09/16/92

LABORATORY: ChemWest

Ticket# CW-9085

Project Name: General Electric Company

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C1-TCDD	*C-HxCDD	*C-HpCDF
ROJ4S // 469246 Detection Limit	9085-1	01/13/92	14:39	CW-2	75.6	79.0	78.8	63.0	39.4	59.8	80.7	99.3	93.9	94.2
ROJ2S // 469249 Detection Limit	9085-2	01/13/92	15:19	CW-2	81.3	86.7	88.2	74.2	47.7	77.0	89.3	100	97.0	95.3
ROJ1S // 469251 Detection Limit	9085-3	01/13/92	15:58	CW-2	73.3	81.0	74.7	61.0	44.2	72.0	81.0	100	97.2	97.2
ROJ3S // 469254 Detection Limit	9085-4	01/13/92	16:39	CW-2	52.7	79.0	84.4	73.1	50.6	35.8	80.5	99.7	97.8	96.7

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____

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DATE: 12/09/92

LABORATORY: ChemWest

Ticket# CW-7614

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppb or ng/g)

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	2378								2378							
					TCDD	TCDF	PeCDD	HxCDD	HxCDF	HpCDD	HpCDF	OCDD	OCDF	TCDF	PeCDF	HxCDF	HxCDF	HpCDF	OCDF	
R001B1416 // 398425	7614-1	02/28/91	16:26	CW-1	ND	ND														
Detection Limit					0.048	0.048	0.072	0.11	0.12	0.24	0.024	0.040	0.049	0.077	0.11	0.18				
R002B0810 // 398426	7614-2	02/28/91	17:04	CW-1	ND	ND														
Detection Limit					0.071	0.071	0.070	0.11	0.13	0.22	0.032	0.067	0.043	0.079	0.10	0.18				

ND = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by:



FORM 1 - QUANTITATION REPORT

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DATE: 12/09/92

LABORATORY: ChemWest

Ticket# CW-7614

Project Name: General Electric Company

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C1-TCDD	*C-HxCDD	*C-HpCDF
R001B1416 // 398425	7614-1	02/28/91	16:26	CW-1	65.1	61.0	69.7	71.9	49.7	67.5	70.9	101	100	88.0
Detection Limit														
R002B0810 // 398426	7614-2	02/28/91	17:04	CW-1	72.6	67.9	77.1	77.5	53.9	77.4	80.4	103	99.0	86.1
Detection Limit														

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HPCDF = 13C12-1234678-HpCDF

Approved by: _____ 

PCDD & PCDF

EPA METHOD 1613A

Sample ID: OX-J-SS1
 Lab ID: 13894-001-SA
 Matrix: Soil
 % Solid: 83

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.01 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N Ratio</u>	<u>Qualifier</u>
2,3,7,8-TCDD	ND	0.46			
Total TCDD	0.82		0.81	7:1	
1,2,3,7,8-PeCDD	ND	0.83			I
Total PeCDD	ND	1.2			I
1,2,3,4,7,8-HxCDD	1.1		1.29	>10:1	A
1,2,3,6,7,8-HxCDD	2.9		1.28	>10:1	
1,2,3,7,8,9-HxCDD	1.9		1.22	>10:1	A
Total HxCDD	24		1.22	>10:1	
1,2,3,4,6,7,8-HpCDD	50		1.04	>10:1	
Total HpCDD	100		1.03	>10:1	
OCDD	390		0.89	>10:1	
2,3,7,8-TCDF	6.8		0.77	>10:1	
1,2,7,8-TCDF	2.6		0.82	>10:1	
Total TCDF	59		0.74	>10:1	D
1,2,3,7,8-PeCDF	22.3		1.45	>10:1	A
2,3,4,7,8-PeCDF	9.8		1.56	>10:1	
Total PeCDF	130		1.74	>10:1	D
1,2,3,4,7,8-HxCDF	4.6		1.31	>10:1	
1,2,3,6,7,8-HxCDF	6.6		1.34	>10:1	
2,3,4,6,7,8-HxCDF	9.4		1.32	>10:1	
1,2,3,7,8,9-HxCDF	1.1		1.30	>10:1	A
Total HxCDF	130		1.30	>10:1	D
1,2,3,4,6,7,8-HpCDF	26		1.02	>10:1	D
1,2,3,4,7,8,9-HpCDF	2.1		1.08	>10:1	A
Total HpCDF	61		1.02	>10:1	D
OCDF	33		0.93	>10:1	

 Analyst: JM

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 Reviewer: DM

PCDD & PCDF
EPA METHOD 1613A

Sample ID: OX-J-SS2
 Lab ID: 13894-002-SA
 Matrix: Soil
 % Solid: 85

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.13 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N Ratio</u>	<u>Qualifier</u>
2,3,7,8-TCDD	ND	0.46			
Total TCDD	3.5		0.73	>10:1	
1,2,3,7,8-PeCDD	ND	2.1			I
Total PeCDD	14		1.52	>10:1	
1,2,3,4,7,8-HxCDD	1.9		1.32	>10:1	A
1,2,3,6,7,8-HxCDD	5.2		1.32	>10:1	
1,2,3,7,8,9-HxCDD	3.1		1.28	>10:1	
Total HxCDD	47		1.28	>10:1	
1,2,3,4,6,7,8-HpCDD	91		1.03	>10:1	
Total HpCDD	170		1.05	>10:1	
OCDD	860		0.89	>10:1	
2,3,7,8-TCDF	16		0.80	>10:1	
1,2,7,8-TCDF	7.8		0.81	>10:1	
Total TCDF	160		0.78	>10:1	D
1,2,3,7,8-PeCDF	5.9		1.55	>10:1	
2,3,4,7,8-PeCDF	16		1.58	>10:1	
Total PeCDF	260		1.60	>10:1	D
1,2,3,4,7,8-HxCDF	15		1.29	>10:1	
1,2,3,6,7,8-HxCDF	26		1.24	>10:1	D
2,3,4,6,7,8-HxCDF	16		1.35	>10:1	
1,2,3,7,8,9-HxCDF	2.9		1.33	>10:1	
Total HxCDF	350		1.33	>10:1	D
1,2,3,4,6,7,8-HpCDF	160		1.03	>10:1	D
1,2,3,4,7,8,9-HpCDF	6.0		1.08	>10:1	
Total HpCDF	300		1.03	>10:1	D
OCDF	100		0.94	>10:1	

 Analyst: Jewett

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 Reviewer: SM

PCDD & PCDF

EPA METHOD 1613A

Sample ID: OX-J-SS3
 Lab ID: 13894-003-SA
 Matrix: Soil
 % Solid: 79

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.13 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N Ratio</u>	<u>Qualifier</u>
2,3,7,8-TCDD	ND	0.32			
Total TCDD	4.6		0.83	>10:1	
1,2,3,7,8-PeCDD	ND	2.0			I
Total PeCDD	8.4		1.72	>10:1	
1,2,3,4,7,8-HxCDD	2.0		1.29	>10:1	A
1,2,3,6,7,8-HxCDD	7.4		1.30	>10:1	
1,2,3,7,8,9-HxCDD	3.8		1.41	>10:1	
Total HxCDD	57		1.25	>10:1	
1,2,3,4,6,7,8-HpCDD	110		1.04	>10:1	
Total HpCDD	250		1.03	>10:1	
OCDD	840		0.88	>10:1	
2,3,7,8-TCDF	37		0.80	>10:1	
1,2,7,8-TCDF	15		0.79	>10:1	
Total TCDF	320		0.76	>10:1	D
1,2,3,7,8-PeCDF	9.9		1.55	>10:1	
2,3,4,7,8-PeCDF	7.6		1.57	>10:1	
Total PeCDF	460		1.49	>10:1	D
1,2,3,4,7,8-HxCDF	18		1.27	>10:1	
1,2,3,6,7,8-HxCDF	30		1.28	>10:1	D
2,3,4,6,7,8-HxCDF	35		1.27	>10:1	
1,2,3,7,8,9-HxCDF	3.6		1.37	>10:1	
Total HxCDF	520		1.30	>10:1	D
1,2,3,4,6,7,8-HpCDF	110		1.04	>10:1	D
1,2,3,4,7,8,9-HpCDF	6.4		1.05	>10:1	
Total HpCDF	230		1.04	>10:1	D
OCDF	89		0.92	>10:1	

Analyst: QMM

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Reviewer: JM

PCDD & PCDF

EPA METHOD 1613A

Sample ID: OX-J-SS4
 Lab ID: 13894-004-SA
 Matrix: Soil
 % Solid: 86

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.30 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N</u>	<u>Qualifier</u>
2,3,7,8-TCDD	0.55		0.68	3:1	
Total TCDD	9.9		0.81	>10:1	
1,2,3,7,8-PeCDD	3.0		*	>10:1	MPC
Total PeCDD	11		1.67	>10:1	
1,2,3,4,7,8-HxCDD	4.3		1.18	>10:1	
1,2,3,6,7,8-HxCDD	23		1.31	>10:1	
1,2,3,7,8,9-HxCDD	6.8		1.21	>10:1	
Total HxCDD	130		1.21	>10:1	
1,2,3,4,6,7,8-HpCDD	680		1.04	>10:1	
Total HpCDD	2100		1.04	>10:1	
OCDD	6500		0.87	>10:1	
2,3,7,8-TCDF	35		0.75	>10:1	
1,2,7,8-TCDF	19		0.81	>10:1	
Total TCDF	320		0.82	>10:1	D
1,2,3,7,8-PeCDF	18		1.55	>10:1	
2,3,4,7,8-PeCDF	39		1.57	>10:1	
Total PeCDF	450		1.63	>10:1	D
1,2,3,4,7,8-HxCDF	36		1.27	>10:1	
1,2,3,6,7,8-HxCDF	32		1.26	>10:1	D
2,3,4,6,7,8-HxCDF	31		1.27	>10:1	
1,2,3,7,8,9-HxCDF	6.2		1.35	>10:1	
Total HxCDF	500		1.32	>10:1	D
1,2,3,4,6,7,8-HpCDF	150		1.03	>10:1	D
1,2,3,4,7,8,9-HpCDF	13		1.08	>10:1	
Total HpCDF	420		1.03	>10:1	D
OCDF	290		0.91	>10:1	

Analyst: JL

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Reviewer: BB

PCDD & PCDF

EPA METHOD 1613A

Sample ID: OX-J-SS5
 Lab ID: 13894-005-SA
 Matrix: Soil
 % Solid: 84

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.11 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N Ratio</u>	<u>Qualifier</u>
2,3,7,8-TCDD	ND	0.36			
Total TCDD	0.69		0.77	3:1	
1,2,3,7,8-PeCDD	ND	0.49			
Total PeCDD	ND	1.0			
1,2,3,4,7,8-HxCDD	0.68		1.25	>10:1	A
1,2,3,6,7,8-HxCDD	1.9		1.29	>10:1	A
1,2,3,7,8,9-HxCDD	1.1		1.23	>10:1	A
Total HxCDD	15		1.21	>10:1	
1,2,3,4,6,7,8-HpCDD	34		1.06	>10:1	
Total HpCDD	74		1.02	>10:1	
OCDD	260		0.88	>10:1	
2,3,7,8-TCDF	5.7		0.80	>10:1	
1,2,7,8-TCDF	3.2		0.82	>10:1	
Total TCDF	65		0.72	>10:1	D
1,2,3,7,8-PeCDF	2.0		1.54	>10:1	A
2,3,4,7,8-PeCDF	6.7		1.59	>10:1	
Total PeCDF	97		1.60	>10:1	D
1,2,3,4,7,8-HxCDF	4.3		1.31	>10:1	
1,2,3,6,7,8-HxCDF	6.0		1.25	>10:1	D
2,3,4,6,7,8-HxCDF	7.2		1.30	>10:1	D
1,2,3,7,8,9-HxCDF	0.99		1.24	>10:1	A
Total HxCDF	92		1.30	>10:1	D
1,2,3,4,6,7,8-HpCDF	24		1.03	>10:1	D
1,2,3,4,7,8,9-HpCDF	2.1		0.99	>10:1	A
Total HpCDF	52		1.03	>10:1	D
OCDF	26		0.91	>10:1	

 Analyst: J. M.

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 Reviewer: bly

PCDD & PCDF
EPA METHOD 1613A

Sample ID: OX-J-DUP
 Lab ID: 13894-008-SA
 Matrix: Soil
 % Solid: 88

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.18 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N Ratio</u>	<u>Qualifier</u>
2,3,7,8-TCDD	ND	0.28			
Total TCDD	1.1		0.77	8:1	
1,2,3,7,8-PeCDD	ND	0.45			
Total PeCDD	ND	0.91			
1,2,3,4,7,8-HxCDD	0.57		1.18	8:1	A
1,2,3,6,7,8-HxCDD	1.8		1.39	>10:1	A
1,2,3,7,8,9-HxCDD	0.90		1.09	>10:1	A
Total HxCDD	13		1.23	>10:1	
1,2,3,4,6,7,8-HpCDD	31		1.05	>10:1	
Total HpCDD	67		1.03	>10:1	
OCDD	240		0.89	>10:1	
2,3,7,8-TCDF	5.5		0.75	>10:1	
1,2,7,8-TCDF	3.0		0.78	>10:1	
Total TCDF	45		0.81	>10:1	D
1,2,3,7,8-PeCDF	1.8		1.54	>10:1	A
2,3,4,7,8-PeCDF	6.5		1.60	>10:1	
Total PeCDF	86		1.58	>10:1	D
1,2,3,4,7,8-HxCDF	3.2		1.20	>10:1	
1,2,3,6,7,8-HxCDF	5.1		1.26	>10:1	D
2,3,4,6,7,8-HxCDF	5.8		1.36	>10:1	D
1,2,3,7,8,9-HxCDF	0.81		1.29	>10:1	A
Total HxCDF	80		1.32	>10:1	D
1,2,3,4,6,7,8-HpCDF	19		1.04	>10:1	D
1,2,3,4,7,8,9-HpCDF	1.4		0.96	>10:1	A
Total HpCDF	43		1.04	>10:1	D
OCDF	22		0.93	>10:1	

Analyst: JL

PCDD & PCDF
EPA METHOD 1613A

Sample ID: OX-J-SS6
 Lab ID: 13894-006-SA
 Matrix: Soil
 % Solid: 88

Date Received: 9/15/94
 Date Extracted: 9/23/94
 Sample Amount: 10.06 g

ICAL ID: I1613A
 QC Lot: LC0923S
 Units: pg/g

<u>Compound</u>	<u>Conc.</u>	<u>D.L.</u>	<u>Ratio</u>	<u>S/N Ratio</u>	<u>Qualifier</u>
2,3,7,8-TCDD	ND	0.47			
Total TCDD	3.8		0.83	>10:1	
1,2,3,7,8-PeCDD	ND	0.71			
Total PeCDD	ND	1.9			
1,2,3,4,7,8-HxCDD	0.78		1.24	>10:1	A
1,2,3,6,7,8-HxCDD	2.2		1.23	>10:1	A
1,2,3,7,8,9-HxCDD	1.4		1.32	>10:1	A
Total HxCDD	20		1.35	>10:1	
1,2,3,4,6,7,8-HpCDD	37		1.02	>10:1	
Total HpCDD	100		0.98	>10:1	
OCDD	270		0.89	>10:1	
2,3,7,8-TCDF	13		0.80	>10:1	
1,2,7,8-TCDF	8.3		0.81	>10:1	
Total TCDF	110		0.73	>10:1	D
1,2,3,7,8-PeCDF	3.5		1.57	>10:1	
2,3,4,7,8-PeCDF	6.9		1.56	>10:1	
Total PeCDF	94		1.66	>10:1	D
1,2,3,4,7,8-HxCDF	4.8		1.24	>10:1	
1,2,3,6,7,8-HxCDF	6.5		1.31	>10:1	D
2,3,4,6,7,8-HxCDF	7.7		1.28	>10:1	D
1,2,3,7,8,9-HxCDF	1.1		1.26	>10:1	A
Total HxCDF	92		1.37	>10:1	D
1,2,3,4,6,7,8-HpCDF	22		1.03	>10:1	D
1,2,3,4,7,8,9-HpCDF	1.9		1.01	>10:1	A
Total HpCDF	46		1.03	>10:1	D
OCDF	22		0.92	>10:1	

 Analyst: JW

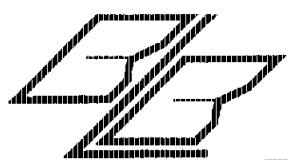
Page 1 of 2

 Reviewer: BB

DATA QUALIFIERS & ABBREVIATIONS

- A The amount detected is below the Method Quantitation Limit.
- B This compound was also detected in the blank.
- C The amount detected is less than five times the Method Quantitation Limit.
- D The amount reported is the maximum possible concentration.
- E The detection limit was raised above the Method Quantitation Limit due to chemical interferences.
- F This result has been confirmed on a DB-225 column.
- G This result has been confirmed on a SP-2331 column.
- H The signal-to-noise ratio is greater than 10:1.
- I Chemical Interference

Conc.	Concentration
D.L.	Detection Limit
NA	Not applicable
S/N	Signal-to-noise



Section 7

METALS ANALYSIS (SOIL)

ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet

ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet

ROA3B1214 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 12-14 feet

ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet

ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet

ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)

ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet

ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet

ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)

ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1

ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2

ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3

ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

ROO1B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet

ROO2B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIESContract: 7/88ROA012224Lab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937274Matrix (soil/water): SOILLab Sample ID: 462153Level (low/med): LOWDate Received: 11/08/91% Solids: 90.3Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	5590	E*	P	
7440-36-0	Antimony	4.3	B N	P	
7440-38-2	Arsenic	5.8	QN	F	
7440-39-3	Barium	24.1	E	P	
7440-41-7	Beryllium	.22	U	P	
7440-43-9	Cadmium	.55	U	P	
7440-70-2	Calcium	51600	E	P	
7440-47-3	Chromium	7.5	"	P	
7440-48-4	Cobalt	5.2	B	P	
7440-50-8	Copper	13.0		P	
7439-89-6	Iron	15100	E*	P	
7439-92-1	Lead	21.1	Q*	P	
7439-95-4	Magnesium	15100		P	
7439-96-5	Manganese	226	E	P	
7439-97-6	Mercury	.11	U	CV	
7440-02-0	Nickel	9.3		P	
7440-09-7	Potassium	207	B	P	
7782-49-2	Selenium	.46	B WN	F	
7440-22-4	Silver	.88	U N	P	
7440-23-5	Sodium	135	B	P	
7440-28-0	Thallium	.22	U WN	F	
7440-62-2	Vanadium	6.6		P	
7440-66-6	Zinc	43.5	EN*	P	
	Cyanide			NR	

Color Before: BROWN

Clarity Before: _____

Texture: MEDIUMColor After: YELLOW

Clarity After: _____

Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROA2B0608

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 937275Matrix (soil/water): SOIL Lab Sample ID: 465052Level (low/med): LOW Date Received: 11/21/91% Solids: 90.8Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	6120		P	
7440-36-0	Antimony	4.2	B	N	P
7440-38-2	Arsenic	6.5		QN	F
7440-39-3	Barium	27.6			P
7440-41-7	Beryllium	.29	B		P
7440-43-9	Cadmium	.55	U		P
7440-70-2	Calcium	57400			P
7440-47-3	Chromium	6.7			P
7440-48-4	Cobalt	7.0			P
7440-50-8	Copper	19.6			P
7439-89-6	Iron	17400	E		P
7439-92-1	Lead	16.3			F
7439-95-4	Magnesium	32900			P
7439-96-5	Manganese	446			P
7439-97-6	Mercury	.18		N*	CV
7440-02-0	Nickel	14.2			P
7440-09-7	Potassium	648			P
7782-49-2	Selenium	.36	B	WN	F
7440-22-4	Silver	.66	U	*	P
7440-23-5	Sodium	119	B		P
7440-28-0	Thallium	.22	U	W	F
7440-62-2	Vanadium	10.0			P
7440-66-6	Zinc	52.4	E		P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROA3B1214

Lab Name: COMPUCHEM LABORATORIESContract: SW-846Lab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937282Matrix (soil/water): SOILLab Sample ID: 473673Level (low/med): LOWDate Received: 01/07/92% Solids: 91.0Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	4980	*		P
7440-36-0	Antimony	6.6	U N		P
7440-38-2	Arsenic	5.7	Q*		F
7440-39-3	Barium	18.4	B *		P
7440-41-7	Beryllium	.15	B		P
7440-43-9	Cadmium	.55	U		P
7440-70-2	Calcium	15100	*		P
7440-47-3	Chromium	7.0	*		P
7440-48-4	Cobalt	6.1			P
7440-50-8	Copper	19.8			P
7439-89-6	Iron	12500			P
7439-92-1	Lead	28.8			P
7439-95-4	Magnesium	8650	*		P
7439-96-5	Manganese	376	*		P
7439-97-6	Mercury	.11	U *		CV
7440-02-0	Nickel	11.3			P
7440-09-7	Potassium	331	B		P
7782-49-2	Selenium	.44	U N		F
7440-22-4	Silver	1.1	U N		P
7440-23-5	Sodium	97.6	B		P
7440-28-0	Thallium	.33	U WN		F
7440-62-2	Vanadium	6.9	*		P
7440-66-6	Zinc	38.8	*		P
	Cyanide				NR

Color Before: GREY

Clarity Before: _____

Texture: MEDIUMColor After: YELLOW

Clarity After: _____

Artifacts: _____

Comments:

FORM 1.04 - PAGE 1PLEASE REFERENCE ENCLOSED NOTICE REGARDING "Q" FLAG IN COLUMN Q

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROB1B0406

Lab Name: <u>COMPUCHEM LABORATORIES</u>	Contract: <u>SW-846</u>	
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: _____ SDG No.: <u>937275</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>465531</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>11/22/91</u>	
% Solids: <u>85.0</u>		

Concentration Units (ug/L or mg/kg dry weight): MG/EG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	6950			P
7440-36-0	Antimony	4.4	B	N	P
7440-38-2	Arsenic	6.6		AN	F
7440-39-3	Barium	91.1			P
7440-41-7	Beryllium	.33	B		P
7440-43-9	Cadmium	.63			P
7440-70-2	Calcium	16100			P
7440-47-3	Chromium	15.1			P
7440-48-4	Cobalt	7.9			P
7440-50-8	Copper	333			P
7439-89-6	Iron	19800	E		P
7439-92-1	Lead	285		N	P
7439-95-4	Magnesium	4000			P
7439-96-5	Manganese	379			P
7439-97-6	Mercury	.37		N*	CV
7440-02-0	Nickel	23.1			P
7440-09-7	Potassium	599			P
7782-49-2	Selenium	.35	U	N	F
7440-22-4	Silver	1.1	B	*	P
7440-23-5	Sodium	159	B		P
7440-28-0	Thallium	.24	U	W	F
7440-62-2	Vanadium	13.9			P
7440-66-6	Zinc	342	E		P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROB2B0002

Lab Name: COMPUCHEM LABORATORIESContract: SW-846Lab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937275Matrix (soil/water): SOILLab Sample ID: 465882Level (low/med): LOWDate Received: 11/23/91% Solids: 91.8Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	5220			P
7440-36-0	Antimony	3.9	U	N	P
7440-38-2	Arsenic	5.1	AN		F
7440-39-3	Barium	37.7			P
7440-41-7	Beryllium	.21	B		P
7440-43-9	Cadmium	.80			P
7440-70-2	Calcium	8340			P
7440-47-3	Chromium	13.1			P
7440-48-4	Cobalt	5.1	B		P
7440-50-8	Copper	36.5			P
7439-89-6	Iron	11400	E		P
7439-92-1	Lead	94.2		N	P
7439-95-4	Magnesium	5950			P
7439-96-5	Manganese	190			P
7439-97-6	Mercury	.61		N*	CV
7440-02-0	Nickel	11.1			P
7440-09-7	Potassium	571			P
7782-49-2	Selenium	.38	B	N	F
7440-22-4	Silver	.77	B	*	P
7440-23-5	Sodium	90.5	B		P
7440-28-0	Thallium	.22	U	W	F
7440-62-2	Vanadium	10.5			P
7440-66-6	Zinc	135	E		P
	Cyanide				NR

Color Before: BROWN

Clarity Before: _____

Texture: MEDIUMColor After: YELLOW

Clarity After: _____

Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROB-DPA1

Lab Name: COMPUCHEM LABORATORIESContract: SW-846Lab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937275Matrix (soil/water): SOILLab Sample ID: 465558Level (low/med): LOWDate Received: 11/22/91t Solids: 85.0Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	7790			P
7440-36-0	Antimony	4.2	U	N	P
7440-38-2	Arsenic	4.8		AN	F
7440-39-3	Barium	68.1			P
7440-41-7	Beryllium	.41	B		P
7440-43-9	Cadmium	.59	U		P
7440-70-2	Calcium	3310			P
7440-47-3	Chromium	13.4			P
7440-48-4	Cobalt	8.8			P
7440-50-8	Copper	62.6			P
7439-89-6	Iron	15200	E		P
7439-92-1	Lead	97.5		N	P
7439-95-4	Magnesium	4280			P
7439-96-5	Manganese	273			P
7439-97-6	Mercury	.23		N*	CV
7440-02-0	Nickel	15.6			P
7440-09-7	Potassium	637			P
7782-49-2	Selenium	.41	B	N	F
7440-22-4	Silver	3.8	*		P
7440-23-5	Sodium	168	B		P
7440-28-0	Thallium	.24	U	W	F
7440-62-2	Vanadium	15.8			P
7440-66-6	Zinc	118	E		P
	Cyanide				NR

Color Before: BROWN

Clarity Before: _____

Texture: MEDIUMColor After: YELLOW

Clarity After: _____

Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIESContract: 7/88ROC011012Lab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937272Matrix (soil/water): SOILLab Sample ID: 461435Level (low/med): LOWDate Received: 11/07/91t Solids: 89.7Concentration Units (ug/L or mg/kg dry weight): MG/RG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	6550		P	
7440-36-0	Antimony	4.0	U	N	P
7440-38-2	Arsenic	4.3		F	
7440-39-3	Barium	36.5		P	
7440-41-7	Beryllium	.19	B		P
7440-43-9	Cadmium	.56	U		P
7440-70-2	Calcium	17200	*		P
7440-47-3	Chromium	9.1		P	
7440-48-4	Cobalt	6.6		P	
7440-50-8	Copper	287		N*	P
7439-89-6	Iron	16100	E		P
7439-92-1	Lead	104		N	P
7439-95-4	Magnesium	9560	*		P
7439-96-5	Manganese	351			P
7439-97-6	Mercury	.11	U		CV
7440-02-0	Nickel	12.6			P
7440-09-7	Potassium	435	B		P
7782-49-2	Selenium	.33	U	WN	F
7440-22-4	Silver	.67	U	N	P
7440-23-5	Sodium	111	B		P
7440-28-0	Thallium	.22	U	W	F
7440-62-2	Vanadium	11.5			P
7440-66-6	Zinc	187	E		P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88ROC021214Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 937272Matrix (soil/water): SOIL Lab Sample ID: 461159Level (low/med): LOW Date Received: 11/06/91% Solids: 81.5Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	6330			P
7440-36-0	Antimony	4.4	U	N	P
7440-38-2	Arsenic	3.6			F
7440-39-3	Barium	17.4	B		P
7440-41-7	Beryllium	.15	B		F
7440-43-9	Cadmium	.61	U		P
7440-70-2	Calcium	8050	*		P
7440-47-3	Chromium	8.3			P
7440-48-4	Cobalt	6.6			P
7440-50-8	Copper	15.3		N*	P
7439-89-6	Iron	15400	E		P
7439-92-1	Lead	28.9	A		F
7439-95-4	Magnesium	4820	*		P
7439-96-5	Manganese	223			P
7439-97-6	Mercury	.12	U		CV
7440-02-0	Nickel	13.1			P
7440-09-7	Potassium	404	B		P
7782-49-2	Selenium	.37	U	WN	F
7440-22-4	Silver	.74	U	N	P
7440-23-5	Sodium	102	B		P
7440-28-0	Thallium	.25	U	W	F
7440-62-2	Vanadium	7.7			P
7440-66-6	Zinc	51.4	E		P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIESContract: 7/88ROC-DPA1Lab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937272Matrix (soil/water): SOILLab Sample ID: 461190Level (low/med): LOWDate Received: 11/06/91% Solids: 54.0Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	9850			P
7440-36-0	Antimony	6.7	U	N	P
7440-38-2	Arsenic	4.8			F
7440-39-3	Barium	29.6	B		P
7440-41-7	Beryllium	.22	B		P
7440-43-9	Cadmium	.93	U		P
7440-70-2	Calcium	12400	*		P
7440-47-3	Chromium	12.0			P
7440-48-4	Cobalt	10.2			P
7440-50-8	Copper	18.0		N*	P
7439-89-6	Iron	20700	E		P
7439-92-1	Lead	33.3	A		F
7439-95-4	Magnesium	5740	*		P
7439-96-5	Manganese	298			P
7439-97-6	Mercury	.19	U		CV
7440-02-0	Nickel	17.7			P
7440-09-7	Potassium	534	B		P
7782-49-2	Selenium	.56	U	WN	F
7440-22-4	Silver	1.1	U	N	P
7440-23-5	Sodium	187	B		P
7440-28-0	Thallium	.37	U	W	F
7440-62-2	Vanadium	11.1			P
7440-66-6	Zinc	79.8	E		P
	Cyanide				NR

Color Before: BROWN

Clarity Before: _____

Texture: MEDIUMColor After: YELLOW

Clarity After: _____

Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROC3B0204

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 937275Matrix (soil/water): SOIL Lab Sample ID: 465105Level (low/med): LOW Date Received: 11/21/91% Solids: 91.5Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	8840			P
7440-36-0	Antimony	3.9	U	N	P
7440-38-2	Arsenic	4.9		N	F
7440-39-3	Barium	40.7			P
7440-41-7	Beryllium	.28	B		P
7440-43-9	Cadmium	.55	U		P
7440-70-2	Calcium	23100			P
7440-47-3	Chromium	8.6			P
7440-48-4	Cobalt	7.4			P
7440-50-8	Copper	123			P
7439-89-6	Iron	21200	E		P
7439-92-1	Lead	26.8			F
7439-95-4	Magnesium	14000			P
7439-96-5	Manganese	430			P
7439-97-6	Mercury	.11	U	N*	CV
7440-02-0	Nickel	16.4			P
7440-09-7	Potassium	772			P
7782-49-2	Selenium	.33	U	WN	F
7440-22-4	Silver	.66	U	*	P
7440-23-5	Sodium	101	B		P
7440-28-0	Thallium	.22	U	W	F
7440-62-2	Vanadium	14.0			P
7440-66-6	Zinc	67.3	E		P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIESContract: SW-846ROJ1SLab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 937278Matrix (soil/water): SOILLab Sample ID: 469288Level (low/med): LOWDate Received: 12/11/91% Solids: 78.2Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	9730	*	P	
7440-36-0	Antimony	8.9	B	N	P
7440-38-2	Arsenic	9.4	Q	F	
7440-39-3	Barium	57.3		P	
7440-41-7	Beryllium	.39	B		P
7440-43-9	Cadmium	1.2	U		P
7440-70-2	Calcium	6750	E		P
7440-47-3	Chromium	17.2		P	
7440-48-4	Cobalt	9.5	B		P
7440-50-8	Copper	30.8	N		P
7439-89-6	Iron	19600	*	P	
7439-92-1	Lead	97.8	*	P	
7439-95-4	Magnesium	5980		P	
7439-96-5	Manganese	517	N*		P
7439-97-6	Mercury	.20		CV	
7440-02-0	Nickel	17.7		P	
7440-09-7	Potassium	1070	B		P
7782-49-2	Selenium	1.0	U	F	
7440-22-4	Silver	1.5	U	N	P
7440-23-5	Sodium	145	B		P
7440-28-0	Thallium	.76	U	Q	F
7440-62-2	Vanadium	20.3			P
7440-66-6	Zinc	126			P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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PLEASE REFERENCE ENCLOSED NOTICE REGARDING "Q" FLAG IN COLUMN Q

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846ROJ2SLab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 937278Matrix (soil/water): SOIL Lab Sample ID: 469279Level (low/med): LOW Date Received: 12/11/91% Solids: 85.0Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	5670	*	P	
7440-36-0	Antimony	10.5	B	N	P
7440-38-2	Arsenic	21.9		A	F
7440-39-3	Barium	41.5	B		P
7440-41-7	Beryllium	.24	U		P
7440-43-9	Cadmium	1.2	U		P
7440-70-2	Calcium	9570		E	P
7440-47-3	Chromium	41.0			P
7440-48-4	Cobalt	9.4	B		P
7440-50-8	Copper	95.6		N	P
7439-89-6	Iron	68700	*		P
7439-92-1	Lead	121	*		P
7439-95-4	Magnesium	7150			P
7439-96-5	Manganese	854		N*	P
7439-97-6	Mercury	.60			CV
7440-02-0	Nickel	43.8			P
7440-09-7	Potassium	393	B		P
7782-49-2	Selenium	.92	U	W	F
7440-22-4	Silver	1.4	U	N	P
7440-23-5	Sodium	120	B		P
7440-28-0	Thallium	.69	U	W	F
7440-62-2	Vanadium	14.1			P
7440-66-6	Zinc	164			P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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FORM I - IN

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846ROJ3SLab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 937278Matrix (soil/water): SOIL Lab Sample ID: 469300Level (low/med): LOW Date Received: 12/11/91% Solids: 91.2Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	5500	*	P	
7440-36-0	Antimony	7.7	U	N	P
7440-38-2	Arsenic	5.5		A	P
7440-39-3	Barium	28.0	B		P
7440-41-7	Beryllium	.21	U		P
7440-43-9	Cadmium	1.1	U		P
7440-70-2	Calcium	8240	E		P
7440-47-3	Chromium	7.7			P
7440-48-4	Cobalt	5.6	B		P
7440-50-8	Copper	12.0		N	P
7439-89-6	Iron	14400	*		P
7439-92-1	Lead	13.5	*		F
7439-95-4	Magnesium	4590			P
7439-96-5	Manganese	214		N*	P
7439-97-6	Mercury	.11	U		CV
7440-02-0	Nickel	9.9			P
7440-09-7	Potassium	969	B		P
7782-49-2	Selenium	.87	U	W	F
7440-22-4	Silver	1.3	U	N	P
7440-23-5	Sodium	166	B		P
7440-28-0	Thallium	.65	U	W	F
7440-62-2	Vanadium	11.6			P
7440-66-6	Zinc	33.0			P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____Comments:
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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: <u>COMPUCHEM LABORATORIES</u>	Contract: <u>SW-846</u>	<u>ROJ4S</u>
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: _____ SDG No.: <u>937278</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>469265</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>12/11/91</u>	
% Solids: <u>66.9</u>		

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	10100	*	P	
7440-36-0	Antimony	11.1	B	N	P
7440-38-2	Arsenic	9.5		F	
7440-39-3	Barium	66.8		P	
7440-41-7	Beryllium	.30	B		P
7440-43-9	Cadmium	1.5	U		P
7440-70-2	Calcium	18100	E		P
7440-47-3	Chromium	17.8		P	
7440-48-4	Cobalt	14.8	B		P
7440-50-8	Copper	58.8		N	P
7439-89-6	Iron	44200	*		P
7439-92-1	Lead	195	*		P
7439-95-4	Magnesium	11500			P
7439-96-5	Manganese	987	N*		P
7439-97-6	Mercury	.21			CV
7440-02-0	Nickel	27.9			P
7440-09-7	Potassium	1120	B		P
7782-49-2	Selenium	1.2	U		F
7440-22-4	Silver	1.8	UN		P
7440-23-5	Sodium	174	B		P
7440-28-0	Thallium	.88	U	W	F
7440-62-2	Vanadium	27.3			P
7440-66-6	Zinc	266			P
	Cyanide				NR

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: YELLOW Clarity After: _____ Artifacts: _____

Comments:

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NORTHEASTERN ANALYTICAL CORPORATION

Report #1416

ANALYTICAL RESULTS

Metals

<u>Parameter</u>	910434-1 396814*	<u>Sample Designation</u>	
		Report Detection Limit*	
Aluminum, total	4,200	24	
Antimony, total	ND	1.2	
Arsenic, total	2.0	1.2	
Barium, total	ND	24	
Beryllium, total	ND	0.59	
Cadmium, total	ND	0.59	
Calcium, total	17,000	590	
Chromium, total	3.2	1.2	
Cobalt, total	ND	5.9	
Copper, total	11	3.0	
Iron, total	12,000	12	
Lead, total	ND	12	
Magnesium, total	9,800	590	
Manganese, total	300	1.8	
Mercury, total	ND	0.12	
Nickel, total	9.3	4.7	
Potassium, total	ND	590	
Selenium, total	ND	0.59	
Sodium, total	ND	590	
Silver, total	ND	1.2	
Thallium, total	ND	1.2	
Vanadium, total	5.9	5.9	
Zinc, total	38	2.4	
Units	(mg/kg)	(mg/kg)	

ND: Not Detected.

*: Calculated on a dry weight basis.



NORTHEASTERN ANALYTICAL CORPORATION

ANALYTICAL RESULTS

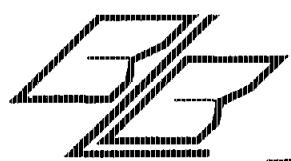
R092B0810

Metals

<u>Parameter</u>	<u>Sample Designation</u>	
	910434-2 396819*	Report Detection Limit*
Aluminum, total	2,900	24
Antimony, total	ND	1.2
Arsenic, total	ND	1.2
Barium, total	ND	24
Beryllium, total	ND	0.60
Cadmium, total	ND	0.60
Calcium, total	ND	600
Chromium, total	4.2	1.2
Cobalt, total	ND	6.0
Copper, total	ND	3.0
Iron, total	7,400	12
Lead, total	ND	12
Magnesium, total	1,300	600
Manganese, total	56	1.8
Mercury, total	ND	0.12
Nickel, total	ND	4.8
Potassium, total	ND	600
Selenium, total	ND	0.60
Sodium, total	ND	600
Silver, total	ND	1.2
Thallium, total	ND	1.2
Vanadium, total	ND	6.0
Zinc, total	19	2.4
Units	(mg/kg)	(mg/kg)

ND: Not Detected.

*: Calculated on a dry weight basis.



Section 8

CYANIDES ANALYSIS (SOIL)

ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet

ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet

ROA3B1214 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 12-14 feet

ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet

ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet

ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)

ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet

ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet

ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)

ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1

ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2

ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3

ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

ROO1B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet

ROO2B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet

OX-J-SS1 - Soil sample from Oxbow Area J, soil surficial sample OX-J-SS1 at 0-4 inches

OX-J-SS2 - Soil sample from Oxbow Area J, soil surficial sample OX-J-SS2 at 0-4 inches

OX-J-SS3 - Soil sample from Oxbow Area J, soil surficial sample OX-J-SS3 at 0-4 inches

OX-J-SS4 - Soil sample from Oxbow Area J, soil surficial sample OX-J-SS4 at 0-4 inches

OX-J-SS5 - Soil sample from Oxbow Area J, soil surficial sample OX-J-SS5 at 0-4 inches

OX-J-DUP - Duplicate soil sample from Oxbow Area J, at location OX-J-SS5 at 0-4 inches

CYANIDES ANALYSIS (SOIL) (CONT'D)

**OX-J-SS6 - Soil sample from Oxbow Area J, soil surficial sample OX-J-SS6 at 0-4
 inches**

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88ROA012224Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 410579Matrix (soil/water): SOIL Lab Sample ID: 462148Level (low/med): LOW Date Received: 11/08/91% Solids: 90.3Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury			NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc	.55	U		AS
	Cyanide				

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

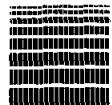
Lab Name: COMPUCHEM LABORATORIES Contract: 7/88ROA2B0608Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 410579Matrix (soil/water): SOIL Lab Sample ID: 465053Level (low/med): LOW Date Received: 11/21/91% Solids: 90.8Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury			NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc			NR	
	Cyanide	.55	U	AS	

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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COMPUCHEM
LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill/Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

COMPOUND LIST

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROA3B1214
COMPUCHEM SAMPLE NUMBER: 473677
DRY WEIGHT FACTOR: 1.10
PERCENT SOLID: 90.9

	CONCENTRATION (ng/kg)	DETECTION + LIMIT (ng/kg)
1. CYANIDE	BDL	0.55

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROB1B0406

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 10543AMatrix (soil/water): SOIL Lab Sample ID: 465536Level (low/med): LOW Date Received: 11/22/91% Solids: 82.9Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.60	U*		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: <u>COMPUCHEM LABORATORIES</u>	Contract: <u>7/88</u>	ROB2B0002
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: <u>SDG No.: 10543A</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>465683</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>11/23/91</u>	
% Solids: <u>66.7</u>		

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury			NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc			NR	
	Cyanide	.75	U *	AS	

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

- Lab Name: <u>COMPUCHEM LABORATORIES</u>	Contract: <u>7/88</u>	ROB-DPA1
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: <u>SDG No.: 10543A</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>465560</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>11/22/91</u>	
% Solids: <u>82.9</u>		

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.60	U *		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88

ROCO11012

Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 410543Matrix (soil/water): SOIL Lab Sample ID: 461437Level (low/med): LOW Date Received: 11/07/91% Solids: 85.0Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.59	U		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: <u>COMPUCHEM LABORATORIES</u>	Contract: <u>7/88</u>	<u>ROC021214</u>
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: _____ SDG No.: <u>410543</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>461162</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>11/06/91</u>	
% Solids: <u>81.5</u>		

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury			NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc			NR	
	Cyanide	.61	U	AS	

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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INORGANIC ANALYSIS DATA SHEET

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88ROC-DPA1Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 410543Matrix (soil/water): SOIL Lab Sample ID: 461192Level (low/med): LOW Date Received: 11/06/91t Solids: 54.0Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.93	U		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: <u>COMPUCHEM LABORATORIES</u>	Contract: <u>7/88</u>	<u>ROC3B0204</u>
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: _____ SDG No.: <u>410579</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>465114</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>11/21/91</u>	
% Solids: <u>92.0</u>		

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury			NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc			NR	
?	Cyanide	.54	U	AS	

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill / Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263
COMPOUND LIST

RESULTS REPORTED ON DRY WEIGHT BASIS USING TEE PERCENT SOLID

SAMPLE IDENTIFIER: ROJ15
COMPUCHEM SAMPLE NUMBER: 469289
DRY WEIGHT FACTOR: 1.28
PERCENT SOLID: 78.1

	CONCENTRATION (ng/kg)	DETECTION + LIMIT (ng/kg)
1. CYANIDE	1.3	0.64

BDL= BELOW DETECTION LIMIT

[†] Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST

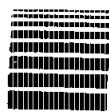
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJ2S
COMPUCHEM SAMPLE NUMBER: 469280
DRY WEIGHT FACTOR: 1.18
PERCENT SOLID: 84.7

	CONCENTRATION (mg/kg)	DETECTION + LIMIT (mg/kg)
1. CYANIDE	120	5.9

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST

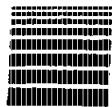
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJOS
COMPUCHEM SAMPLE NUMBER: 469301
DRY WEIGHT FACTOR: 1.10
PERCENT SOLID: 90.9

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. CYANIDE	BDL 0.55

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST

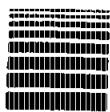
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJ4S
COMPUCHEM SAMPLE NUMBER: 469269
DRY WEIGHT FACTOR: 1.49
PERCENT SOLID: 67.1

	CONCENTRATION (mg/kg)	DETECTION + LIMIT (mg/kg)
1. CYANIDE	BDL	0.75

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST

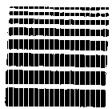
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: R001B1416
COMPUCHEM SAMPLE NUMBER: 396815
DRY WEIGHT FACTOR: 1.18
PERCENT SOLID: 84.7

	CONCENTRATION ($\mu\text{g}/\text{kg}$)	DETECTION + LIMIT ($\mu\text{g}/\text{kg}$)
1. CYANIDE	BDL	0.59

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: R002B0810
COMPUCHEM SAMPLE NUMBER: 396820
DRY WEIGHT FACTOR: 1.21
PERCENT SOLID: 82.6

	CONCENTRATION (ng/kg)	DETECTION + LIMIT (ng/kg)
I. CYANIDE	BDL	0.6

BDL= BELOW DETECTION LIMIT

[†] Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.

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INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM ENV. CORP. Contract: 3/90OX-J-SSILab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 292942Matrix (soil/water): SOIL Lab Sample ID: 639281Level (low/med): LOW Date Received: 09/15/94* Solids: 83.2Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.60	U		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM ENV. CORP. Contract: 3/90

OX-J-SS2

Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 292942Matrix (soil/water): SOIL Lab Sample ID: 639282Level (low/med): LOW Date Received: 09/15/94t Solids: 85.2Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury		.	NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc			NR	
	Cyanide	.59	U	AS	

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM ENV. CORP. Contract: 3/90OX-J-SS3Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 292942Matrix (soil/water): SOIL Lab Sample ID: 639277Level (low/med): LOW Date Received: 09/15/94% Solids: 81.3Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.62	U		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM ENV. CORP. Contract: 3/90OX-J-SS4Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 292942Matrix (soil/water): SOIL Lab Sample ID: 639283Level (low/med): LOW Date Received: 09/15/94* Solids: 78.8Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	63	U		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM ENV. CORP.Contract: 3/90OX-J-SSSLab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 292942Matrix (soil/water): SOILLab Sample ID: 639284Level (low/med): LOWDate Received: 09/15/94% Solids: 86.1Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc	.58	U		AS
	Cyanide				

Color Before: BROWN

Clarity Before: _____

Texture: MEDIUMColor After: COLORLESS

Clarity After: _____

Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM ENV. CORP.Contract: 3/90OX-J-DUPLab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 292942Matrix (soil/water): SOILLab Sample ID: 639286Level (low/med): LOWDate Received: 09/15/94% Solids: 86.2Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum			NR	
7440-36-0	Antimony			NR	
7440-38-2	Arsenic			NR	
7440-39-3	Barium			NR	
7440-41-7	Beryllium			NR	
7440-43-9	Cadmium			NR	
7440-70-2	Calcium			NR	
7440-47-3	Chromium			NR	
7440-48-4	Cobalt			NR	
7440-50-8	Copper			NR	
7439-89-6	Iron			NR	
7439-92-1	Lead			NR	
7439-95-4	Magnesium			NR	
7439-96-5	Manganese			NR	
7439-97-6	Mercury			NR	
7440-02-0	Nickel			NR	
7440-09-7	Potassium			NR	
7782-49-2	Selenium			NR	
7440-22-4	Silver			NR	
7440-23-5	Sodium			NR	
7440-28-0	Thallium			NR	
7440-62-2	Vanadium			NR	
7440-66-6	Zinc			NR	
	Cyanide	.58	U	AS	

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: <u>COMPUCHEM ENV. CORP.</u>	Contract: <u>3/90</u>	<u>OX-J-SS6</u>
Lab Code: <u>COMPU</u>	Case No.: <u>50007</u>	SAS No.: _____ SDG No.: <u>292942</u>
Matrix (soil/water): <u>SOIL</u>	Lab Sample ID: <u>639285</u>	
Level (low/med): <u>LOW</u>	Date Received: <u>09/15/94</u>	
# Solids: <u>88.7</u>		

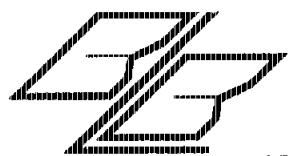
Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	.56	U		AS

Color Before: BROWN Clarity Before: _____ Texture: MEDIUMColor After: COLORLESS Clarity After: _____ Artifacts: _____

Comments:

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Section 9

PHENOL ANALYSIS (SOIL)

ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet

ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet

ROA3B1214 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 12-14 feet

ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet

ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet

ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)

ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet

ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet

ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)

ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet

ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1

ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2

ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3

ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

ROO1B1416 - Soil sample from Oxbow Area K, Soil Boring K-1 at 14-16 feet

ROO2B0810 - Soil sample from Oxbow Area K, Soil Boring K-2 at 8-10 feet



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COMPOUND LIST - CLASSICAL PARAMETERS

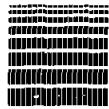
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROA012224
COMPUCHEM SAMPLE NUMBER: 462154
DRY WEIGHT FACTOR: 1.11
PERCENT SOLID: 90.1

CONCENTRATION (mg/kg)	DETECTION + LIMIT (mg/kg)
1. PHENOLS, TOTAL	BDL 0.11

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

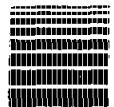
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROA012224DUP
COMPUCHEM SAMPLE NUMBER: 461165
DRY WEIGHT FACTOR: 1.11
PERCENT SOLID: 90.1

	DETECTION +	LIMIT
CONCENTRATION	(ng/kg)	(ng/kg)
I. PHENOLS, TOTAL	BDL	0.11

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROA2B0608
COMPUCHEM SAMPLE NUMBER: 465055
DRY WEIGHT FACTOR: 1.10
PERCENT SOLID: 90.9

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	3.6
	0.11

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUCHEM LIST - CLASSICAL PARAMETERS

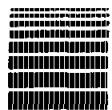
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROA3BL214
COMPUCHEM SAMPLE NUMBER: 473678
DRY WEIGHT FACTOR: 1.10
PERCENT SOLID: 90.9

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	0.93
	0.11

BDL= BELOW DETECTION LIMIT

* Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

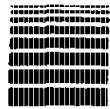
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: R0B1B0406
COMPUCHEM SAMPLE NUMBER: 465539
DRY WEIGHT FACTOR: 1.21
PERCENT SOLID: 82.6

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL 0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROB1B0406DUP

COMPUCHEM SAMPLE NUMBER: 465574

DRY WEIGHT FACTOR: 1.21

PERCENT SOLID: 82.6

	DETECTION +	LIMIT
CONCENTRATION	(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL	0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

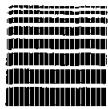
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROB2B0002
COMPUCHEM SAMPLE NUMBER: 465884
DRY WEIGHT FACTOR: 1.50
PERCENT SOLID: 66.7

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	0.31
	0.15

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

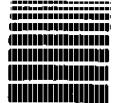
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROB-DPA1
COMPUCHEM SAMPLE NUMBER: 465563
DRY WEIGHT FACTOR: 1.21
PERCENT SOLID: 82.6

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL 0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROC011012
COMPUCHEM SAMPLE NUMBER: 461436
DRY WEIGHT FACTOR: 1.11
PERCENT SOLID: 90.1

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	0.22 0.11

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROC021214
COMPUCHEM SAMPLE NUMBER: 461160
DRY WEIGHT FACTOR: 1.23
PERCENT SOLID: 81.3

	DETECTION + CONCENTRATION (mg/kg)	LIMIT (mg/kg)
1. PHENOLS, TOTAL	BDL	0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROC-DPA1

COMPUCHEM SAMPLE NUMBER: 461191

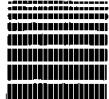
DRY WEIGHT FACTOR: 1.85

PERCENT SOLID: 54.1

CONCENTRATION (mg/kg)	DETECTION + LIMIT (mg/kg)
1. PHENOLS, TOTAL	BDL 0.19

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

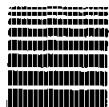
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROC3B0204
COMPUCHEM SAMPLE NUMBER: 465118
DRY WEIGHT FACTOR: 1.08
PERCENT SOLID: 92.6

	DETECTION +	LIMIT
CONCENTRATION	(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL	0.11

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJLS
COMPUCHEM SAMPLE NUMBER: 469290
DRY WEIGHT FACTOR: 1.28
PERCENT SOLID: 78.1

CONCENTRATION (mg/kg)	DETECTION + LIMIT (mg/kg)
1. PHENOLS, TOTAL	BDL 0.13

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

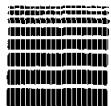
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJ2S
COMPUCHEM SAMPLE NUMBER: 469281
DRY WEIGHT FACTOR: 1.18
PERCENT SOLID: 84.7

	DETECTION +	LIMIT
CONCENTRATION	(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL	0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJ3S
COMPUCHEM SAMPLE NUMBER: 469302
DRY WEIGHT FACTOR: 1.10
PERCENT SOLID: 90.9

	DETECTION +
CONCENTRATION	LIMIT
(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL 0.11

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

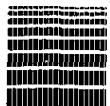
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: ROJ4S
COMPUCHEM SAMPLE NUMBER: 469272
DRY WEIGHT FACTOR: 1.49
PERCENT SOLID: 67.1

	DETECTION + CONCENTRATION (mg/kg)	LIMIT (mg/kg)
1. PHENOLS, TOTAL	0.29	0.15

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: R001B1416

COMPUCHEM SAMPLE NUMBER: 396816

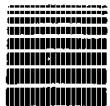
DRY WEIGHT FACTOR: 1.18

PERCENT SOLID: 84.7

	CONCENTRATION (ng/kg)	DETECTION + LIMIT (ng/kg)
1. PHENOLS, TOTAL	BDL	0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



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COMPOUND LIST - CLASSICAL PARAMETERS

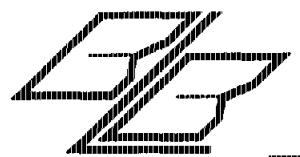
RESULTS REPORTED ON DRY WEIGHT BASIS USING THE PERCENT SOLID

SAMPLE IDENTIFIER: R002E0810
COMPUCHEM SAMPLE NUMBER: 396821
DRY WEIGHT FACTOR: 1.21
PERCENT SOLID: 82.6

	DETECTION +	LIMIT
CONCENTRATION	(mg/kg)	(mg/kg)
1. PHENOLS, TOTAL	BDL	0.12

BDL= BELOW DETECTION LIMIT

+ Detection limits have been adjusted to report variation from the nominal sample weight and the percent solid.



Section 10

SULFIDE ANALYSIS (SOIL)

- ROA012244 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-1 at 22-24 feet
- ROA3B1214 - Soil sample from Oxbow Area A, Groundwater Monitoring Well A-3 at 12-14 feet
- ROB1B0406 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet
- ROB-DPA1 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 4-6 feet (Duplicate)
- ROB2B0002 - Soil sample from Oxbow Area B, Groundwater Monitoring Well B-1 at 0-2 feet
- ROA2B0608 - Soil sample from Oxbow Area A, Soil Boring A-2 at 6-8 feet
- ROC3B0204 - Soil sample from Oxbow Area C, Soil Boring C-3 at 2-4 feet
- ROCO11012 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-1 at 10-12 feet
- ROCO21214 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet
- ROC-DPA1 - Soil sample from Oxbow Area C, Groundwater Monitoring Well C-2 at 12-14 feet (Duplicate)
- ROJ1S - Soil sample from Oxbow Area J, Surficial Soil Sample J-1
- ROJ2S - Soil sample from Oxbow Area J, Surficial Soil Sample J-2
- ROJ3S - Soil sample from Oxbow Area J, Surficial Soil Sample J-3
- ROJ4S - Soil sample from Oxbow Area J, Surficial Soil Sample J-4

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 11/15/91

Case: 8950
Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
462161/ROA012224	8950-1	90	BRL	11.1

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	52.0000	104.0
LQCSD	LQCSD	50.0	49.6000	99.2

Relative % Difference = 4.7

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by:

Date Reported:

09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 01/09/92

Case: 9154

Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
473665/R0A3B1214	9154-1	91	BRL	11.0

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	49.4400	98.9
LQCSD	LQCSD	50.0	49.0400	98.1

Relative % Difference = 0.8

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Matrix Spike	9154-1MS	54.9	62.2418	113.3
Matrix Spike DUP	9154-1MSD	54.9	62.2418	113.3

Relative % Difference = 0.0

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: *[Signature]*

Date Reported:

09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 11/30/91

Case: 9018
Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
465589/R0B1B0406	9018-1	83	BRL	12.0
465591/R0B-DPA1	9018-2	83	BRL	12.0

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	46.9312	93.9
LQCSD	LQCSD	50.0	47.7312	95.5

Relative % Difference = 1.7

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: fmj

Date Reported:
09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 11/30/91

Case: 9020

Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount	
			Detected (MG/KG)	RL (MG/KG)
465886/ROB2B0002	9020-1	67	BRL	14.9

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount	
			Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	46.9312	93.9
LQCSD	LQCSD	50.0	47.7312	95.5

Relative % Difference = 1.7

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: EJK

Date Reported:

09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 11/30/91

Case: 9008
Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
465058/R0A2B0608	9008-1	91	BRL	11.0
465127/ROC3B0204	9008-2	92	BRL	10.9

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	46.9312	93.9
LQCSD	LQCSD	50.0	47.7312	95.5

Relative % Difference = 1.7

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: Eny

Date Reported:
09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 11/13/91

Case: 8942

Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
461438/R0C011012	8942-1	90	92.4	11.1

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	53.6000	107.2
LQCSD	LQCSD	50.0	55.2000	110.4

Relative % Difference = 2.9

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: EMK

Date Reported:

09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 11/13/91

Case: 8933

Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
461180/ROC021214	8933-1	82	25.4	12.2
461193/ROC-DPA1	8933-2	54	34.1	18.5

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	53.6000	107.2
LQCSD	LQCSD	50.0	55.2000	110.4

Relative % Difference = 2.9

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: 

Date Reported:
09/16/92

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 9030

Date(s) Analyzed: 12/18/91

Case: 9085
Matrix: Soil

Client ID	CHEMWEST ID	% Solids	Amount Detected (MG/KG)	RL (MG/KG)
469246/R0J4S	9085-1	67	BRL	14.9
469249/R0J2S	9085-2	85	65.0	11.8
469251/R0J1S	9085-3	78	BRL	12.8
469254/R0J3S	9085-4	91	BRL	11.0
469312/R0JEB1	9085-5	100	BRL	10.0

Client ID	CHEMWEST ID	Spike Conc. (MG/KG)	Amount Detected (MG/KG)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	50.0	53.6800	107.4
LQCSD	LQCSD	50.0	54.4800	109.0

Relative % Difference = 1.5

BRL: Below Reporting Limit.

RL: Reporting Limit.

The reporting limit for the Method Blank is 10.0 MG/KG.

Approved by: FDA

Date Reported:
09/16/92

REV5:12.91



Section 11

VOLATILE ORGANICS ANALYSIS (GROUNDWATER)

WP-1 - Groundwater sample from Oxbow Area K, Well Point 1
WP-2 - Groundwater sample from Oxbow Area K, Well Point 2
WP-3 - Groundwater sample from Oxbow Area J, Well Point 3
WP-7 - Groundwater sample from Oxbow Area C, Well Point 7
WP-8 - Groundwater sample from Oxbow Area C, Well Point 8
WP-9 - Groundwater sample from Oxbow Area A, Well Point 9
ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 1 OF 19

Sample Description: WP-1 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	ND
benzene	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	ND
chlorobenzene	ND	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	vinyl chloride ¹	ND
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Approved by Alice R. Meier
Laboratory Manager

Title


CERTIFICATE OF ANALYSIS

TO: **Geraghty & Miller, Inc.**
 ATTN: Bill Gray
 125 E. Bethpage Road
 Plainview, NY 11803

DATE REPORTED: November 15, 1988.
 PROJECT CODE: GMIN 42142
 ORDER NUMBER:
 PAGE 2 OF 19

Sample Description: WP-2 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	ND
benzene ¹	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	ND
chlorobenzene	ND	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	vinyl chloride ¹	ND
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed

Alyce S. Moore
 Approved by _____ Laboratory Manager _____


CERTIFICATE OF ANALYSIS

TO: **Geraghty & Miller, Inc.**
 ATTN: Bill Gray
 125 E. Bethpage Road
 Plainview, NY 11803

DATE REPORTED: **November 15, 1988**
 PROJECT CODE: **GMIN 42142**
 ORDER NUMBER:
 PAGE **3** OF **19**

Sample Description: WP-3 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	18
benzene	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	ND
chlorobenzene	ND	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	v vinyl chloride ¹	<10
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Alice S. Mason
 Approved by **Laboratory Manager**



ANALYTICAL SERVICES



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
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Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 2 OF 25

Sample Description: WP-7 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)	<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	ND
benzene ¹	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	ND
chlorobenzene	ND	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	vinyl chloride ¹	ND
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Alice R. Moore
Approved by _____
Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 3 OF 25

Sample Description: WP-8 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	ND
benzene	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	ND
chlorobenzene	11	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	vinyl chloride ¹	ND
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Approved by

Laboratory Manager

Title

Alice R. Maece



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CERTIFICATE OF ANALYSIS

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Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 4 OF 25

Sample Description: WP-9 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)	<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	<5
benzene	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	<5
chlorobenzene	<5	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	vinyl chloride ¹	<10
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Ryan R. Moore
Approved by _____
Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

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Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 6 OF 25

Sample Description: Trfp Blank (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	ND
benzene	ND	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	<5
chlorobenzene	<5	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	ND	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	ND
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	ND	vinyl chloride ¹	ND
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Approved by Alan R. Nease Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 5 OF 25

Sample Description: 62 (Water)

VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)	<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)
acrolein ¹	ND	1,1-dichloroethene	ND
acrylonitrile ¹	ND	trans-1,2-dichloroethene	140
benzene	20	1,2-dichloropropane	ND
bromodichloromethane	ND	cis-1,3-dichloropropene	ND
bromoform	ND	trans-1,3-dichloropropene	ND
bromomethane ¹	ND	ethyl benzene	ND
carbon tetrachloride	ND	methylene chloride	<5
chlorobenzene	530	1,1,2,2-tetrachloroethane	ND
chloroethane ¹	48	tetrachloroethene	ND
2-chloroethylvinyl ether ¹	ND	toluene	12
chloroform	ND	1,1,1-trichloroethane	ND
chloromethane ¹	ND	1,1,2-trichloroethane	ND
dibromochloromethane	ND	trichloroethene	ND
1,1-dichloroethane	20	vinyl chloride ¹	300
1,2-dichloroethane	ND		

Remarks: 5 = Quantitation Limit

ND = Not detected

< = Detected but at a level less than the quantitation limit.

¹ = This component has a quantitation limit two (2) times that listed.

Approved by

Laboratory Manager:

Title

Alys R. Moore

LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM.RTP

Contract: 500077

ROA1G

Lab Code: COMPU Case No.: 24105

SAS No.: _____

SDG No.: 0336

Matrix: (soil/water) WATER

Lab Sample ID: 468162

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN068162A03

Level: (low/med) LOW

Date Received: 12/06/91

Moisture: not dec.

Date Analyzed: 12/11/91

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

74-87-3-----Chloromethane	10	U
74-83-9-----Bromomethane	5	U
75-01-4-----Vinyl Chloride	10	U
75-00-3-----Chloroethane	10	U
75-09-2-----Methylene Chloride	4	BJ
67-64-1-----Acetone	10	U
75-15-0-----Carbon Disulfide	5	U
75-35-4-----1,1-Dichloroethene	5	U
75-34-3-----1,1-Dichloroethane	5	U
540-59-0-----1,2-Dichloroethene (total)	5	U
67-66-3-----Chloroform	5	U
107-06-2-----1,2-Dichloroethane	5	U
78-93-3-----2-Butanone	10	U
71-55-6-----1,1,1-Trichloroethane	5	U
56-23-5-----Carbon Tetrachloride	5	U
108-05-4-----Vinyl Acetate	10	U
75-27-4-----Bromodichloromethane	5	U
78-87-5-----1,2-Dichloropropane	5	U
10061-01-5-----cis-1,3-Dichloropropene	5	U
79-01-6-----Trichloroethene	5	U
124-48-1-----Dibromochloromethane	5	U
79-00-5-----1,1,2-Trichloroethane	5	U
71-43-2-----Benzene	4	J
10061-02-6-----Trans-1,3-Dichloropropene	5	U
75-25-2-----Bromoform	10	U
108-10-1-----4-Methyl-2-Pentanone	15	U
591-78-6-----2-Hexanone	15	U
127-18-4-----Tetrachloroethene	5	U
79-34-5-----1,1,2,2-Tetrachloroethane	10	U
108-88-3-----Toluene	5	U
108-90-7-----Chlorobenzene	5	U
100-41-4-----Ethylbenzene	11	U
100-42-5-----Styrene	5	U
1330-20-7-----Total Xylenes	3	J

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA1G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>0336</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468162</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CN068162A03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/06/91</u>	
# Moisture: not dec.	Date Analyzed: <u>12/11/91</u>	
Column (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	<u>ROA3G</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>585</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>474974</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CN074974A03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>01/15/92</u>	
Moisture: not dec.	Date Analyzed: <u>01/17/92</u>	
Column: (pack/cap) <u>GAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3	Chloromethane	10	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	10	U
75-00-3	Chloroethane	10	U
75-09-2	Methylene Chloride	11	B
67-64-1	Acetone	10	U
75-15-0	Carbon Disulfide	5	U
75-35-4	1,1-Dichloroethene	5	U
75-34-3	1,1-Dichloroethane	5	U
540-59-0	1,2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1,2-Dichloroethane	5	U
78-93-3	2-Butanone	10	U
71-55-6	1,1,1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
108-05-4	Vinyl Acetate	10	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1,2-Dichloropropane	5	U
10061-01-5	cis-1,3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1,1,2-Trichloroethane	5	U
71-43-2	Benzene	3	J
10061-02-6	Trans-1,3-Dichloropropene	5	U
75-25-2	Bromoform	10	U
108-10-1	4-Methyl-2-Pentanone	15	U
591-78-6	2-Hexanone	15	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1,1,2,2-Tetrachloroethane	10	U
108-88-3	Toluene	3	J
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	3	J
74-88-4	Iodomethane	10	U
107-02-8	Acrolein	90	U

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM,RTP</u>	Contract: <u>500077</u>	<u>ROA3G</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>585</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>474974</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CN074974A03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>01/15/92</u>	
% Moisture: not dec.	Date Analyzed: <u>01/17/92</u>	
Column (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB1G
Lab Code: <u>CONPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>0336</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468152</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CN068152C03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/06/91</u>	
# Moisture: not dec.	Date Analyzed: <u>12/11/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L Q	
		10	U
74-87-3	Chloromethane	10	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	10	U
75-00-3	Chloroethane	10	U
75-09-2	Methylene Chloride	5	BJ
67-64-1	Acetone	10	U
75-15-0	Carbon Disulfide	5	U
75-35-4	1,1-Dichloroethene	5	U
75-34-3	1,1-Dichloroethane	5	U
540-59-0	1,2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1,2-Dichloroethane	5	U
78-93-3	2-Butanone	10	U
71-55-6	1,1,1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1,2-Dichloropropane	5	U
10061-01-5	cis-1,3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1,1,2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1,3-Dichloropropene	5	U
75-25-2	Bromoform	10	U
108-10-1	4-Methyl-2-Pentanone	15	U
591-78-6	2-Hexanone	15	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1,1,2,2-Tetrachloroethane	10	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	<u>ROB1G</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>0336</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468152</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CN068152C03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/06/91</u>	
# Moisture: not dec.	Date Analyzed: <u>12/11/91</u>	
Column (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CONCENTRATION UNITS:
Number TICs found: 0 (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM.RTP Contract: 500077 ROB2G
 Lab Code: COMPP Case No.: 24105 SAS No.: SDG No.: 0336
 Matrix: (soil/water) WATER Lab Sample ID: 468160
 Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN068160C03
 Level: (low/med) LOW Date Received: 12/06/91
 * Moisture: not dec. Date Analyzed: 12/11/91
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) ug/L	Q
74-87-3	Chloromethane	10	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	10	U
75-00-3	Chloroethane	10	U
75-09-2	Methylene Chloride	2	BJ
67-64-1	Acetone	10	U
75-15-0	Carbon Disulfide	1	J
75-35-4	1, 1-Dichloroethene	5	U
75-34-3	1, 1-Dichloroethane	5	U
540-59-0	1, 2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1, 2-Dichloroethane	5	U
78-93-3	2-Butanone	5	U
71-55-6	1, 1, 1-Trichloroethane	10	U
56-23-5	Carbon Tetrachloride	5	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1, 2-Dichloropropene	5	U
10061-01-5	cis-1, 3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1, 1, 2-Trichloroethane	5	U
71-43-2	Benzene	10	U
10061-02-6	Trans-1, 3-Dichloropropene	5	U
75-25-2	Bromoform	10	U
108-10-1	4-Methyl-2-Pantanone	15	U
591-78-6	2-Hexanone	15	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1, 1, 2, 2-Tetrachloroethane	10	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	1	J

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ROB2G

Lab Name: COMPUCHEM,REP..... Contract: 500077.....

Lab Code: COMPU..... Case No.: 24105..... SAS No.: SDG No.: Q336.....

Matrix: (soil/water) WATER..... Lab Sample ID: 468160.....

Sample wt/vol: 5.0 (g/mL) ML..... Lab File ID: CN068160C03.....

Level: (low/med) LOW..... Date Received: 12/06/91.....

% Moisture: not dec. Date Analyzed: 12/11/91.....

Column (pack/cap) CAR..... Dilution Factor: 1.0.....

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
.....

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM, RTP</u>	Contract: <u>500077</u>	ROC1G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>0336</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468504</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CR068504B03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/07/91</u>	
Moisture: not dec.	Date Analyzed: <u>12/12/91</u>	
Column: (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/L	Q
74-87-3	-Chloromethane		10	U
74-83-9	-Bromomethane		5	U
75-01-4	-Vinyl Chloride		10	U
75-00-3	-Chloroethane		10	U
75-09-2	-Methylene Chloride		4	BJ
67-64-1	-Acetone		10	U
75-15-0	-Carbon Disulfide		4	J
75-35-4	-1,1-Dichloroethene		5	U
75-34-3	-1,1-Dichloroethane		5	U
540-59-0	-1,2-Dichloroethene (total)		5	U
67-66-3	-Chloroform		5	U
107-06-2	-1,2-Dichloroethane		5	U
78-93-3	-2-Butanone		10	U
71-55-6	-1,1,1-Trichloroethane		5	U
56-23-5	-Carbon Tetrachloride		5	U
75-27-4	-Bromodichloromethane		5	U
78-87-5	-1,2-Dichloropropane		5	U
10061-01-5	-cis-1,3-Dichloropropene		5	U
79-01-6	-Trichloroethene		5	U
124-48-1	-Dibromochloromethane		5	U
79-00-5	-1,1,2-Trichloroethane		5	U
71-43-2	-Benzene		5	U
10061-02-6	-Trans-1,3-Dichloropropene		5	U
75-25-2	-Bromoform		10	U
108-10-1	-4-Methyl-2-Pentanone		15	U
591-78-6	-2-Hexanone		15	U
127-18-4	-Tetrachloroethene		5	U
79-34-5	-1,1,2,2-Tetrachloroethane		10	U
108-88-3	-Toluene		5	U
108-90-7	-Chlorobenzene		5	U
100-41-4	-Ethylbenzene		5	U
100-42-5	-Styrene		5	U
1330-20-7	-Total Xylenes		5	U

1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM.RTP</u>	Contract: <u>500077</u>	<u>ROC1G</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>0336</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468504</u>	
Sample wt/vol: <u>5.0</u> (g/mL) <u>ML</u>	Lab File ID: <u>CR068504B03</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/07/91</u>	
Moisture: not dec.	Date Analyzed: <u>12/12/91</u>	
Column (pack/cap) <u>CAP</u>	Dilution Factor: <u>1.0</u>	

CONCENTRATION UNITS:
 Number TICs found: 0 (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP Contract: 500077

ROC2G

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 0336

Matrix: (soil/water) WATER Lab Sample ID: 468586

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CR068586B03

Level: (low/med) LOW Date Received: 12/07/91

† Moisture: not dec. Date Analyzed: 12/12/91

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	10	U
74-83-9	Bromomethane	5	U
75-01-4	Vinyl Chloride	10	U
75-00-3	Chloroethane	10	U
75-09-2	Methylene Chloride	3	BJ
67-64-1	Acetone	10	U
75-15-0	Carbon Disulfide	5	U
75-35-4	1,1-Dichloroethene	5	U
75-34-3	1,1-Dichloroethane	5	U
540-59-0	1,2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1,2-Dichloroethane	5	U
78-93-3	2-Butanone	10	U
71-55-6	1,1,1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1,2-Dichloropropane	5	U
10061-01-5	cis-1,3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1,1,2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1,3-Dichloropropene	5	U
75-25-2	Bromoform	10	U
108-10-1	4-Methyl-2-Pentanone	15	U
591-78-6	2-Hexanone	15	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1,1,2,2-Tetrachloroethane	10	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: COMPEL CHEM ETC	Contract: 5000272	ROC2G
Lab Code: COMPU	Case No.: 24105	SAS No.: _____ SDG No.: 0336
Matrix: (soil/water) WATER	Lab Sample ID: 468586	
Sample wt/vol: 5.0 (g/mL) ML	Lab File ID: CR068586B03	
Level: (low/med) LOW	Date Received: 12/07/91	
% Moisture: not dec.	Date Analyzed: 12/12/91	
Column (pack/cap) CAP	Dilution Factor: 1.0	

CONCENTRATION UNITS:
Number TICs found: 0 (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROJ1G

Lab Name: COMPUCHEM, RTP Contract: 500077

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 489

Matrix: (soil/water) WATER Lab Sample ID: 472794

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN072794A03

Level: (low/med) LOW Date Received: 12/20/91

% Moisture: not dec. Date Analyzed: 01/02/92

Column: (pack/cap) CAP Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	5	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	10	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	5	U
75-35-4-----	1,1-Dichloroethene	5	U
75-34-3-----	1,1-Dichloroethane	5	U
540-59-0-----	1,2-Dichloroethene (total)	5	U
67-66-3-----	Chloroform	5	U
107-06-2-----	1,2-Dichloroethane	5	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	5	U
56-23-5-----	Carbon Tetrachloride	5	U
108-05-4-----	Vinyl Acetate	10	U
75-27-4-----	Bromodichloromethane	5	U
78-87-5-----	1,2-Dichloropropane	5	U
10061-01-5-----	cis-1,3-Dichloropropene	5	U
79-01-6-----	Trichloroethene	5	U
124-48-1-----	Dibromochloromethane	5	U
79-00-5-----	1,1,2-Trichloroethane	5	U
71-43-2-----	Benzene	5	U
10061-02-6-----	Trans-1,3-Dichloropropene	5	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	15	U
591-78-6-----	2-Hexanone	15	U
127-18-4-----	Tetrachloroethene	5	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	5	U
108-90-7-----	Chlorobenzene	5	U
100-41-4-----	Ethylbenzene	5	U
100-42-5-----	Styrene	5	U
1330-20-7-----	Total Xylenes	5	U
74-88-4-----	Iodomethane	10	U
107-02-8-----	Acrolein	90	U

107-13-1-----	Acrylonitrile	120	U
75-69-4-----	Trichlorofluoromethane	5	U
107-05-1-----	3-Chloropropene	15	U
74-95-3-----	Dibromomethane	10	U
106-93-4-----	1, 2-Dibromoethane	5	U
630-20-6-----	1, 1, 1, 2-Tetrachloroethane	5	U
96-18-4-----	1, 2, 3-Trichloropropane	15	U
764-71-0-----	trans-1, 4-Dichloro-2-butene	15	U
97-63-2-----	Ethylmethacrylate	10	U
96-12-8-----	1, 2-Dibromo-3-chloropropane	10	U



Section 12

SEMICVOLATILE ORGANICS ANALYSIS (GROUNDWATER)

WP-1 - Groundwater sample from Oxbow Area K, Well Point 1
WP-2 - Groundwater sample from Oxbow Area K, Well Point 2
WP-3 - Groundwater sample from Oxbow Area J, Well Point 3
WP-7 - Groundwater sample from Oxbow Area C, Well Point 7
WP-8 - Groundwater sample from Oxbow Area C, Well Point 8
WP-9 - Groundwater sample from Oxbow Area A, Well Point 9
ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1


CERTIFICATE OF ANALYSIS

TO **Geraghty & Miller, Inc.**
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 6 OF 19

Sample Description: WP-1 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acenaphthene	ND	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine ¹	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,i)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	<10	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
dibenzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
d1-n-butylphthalate	ND	N-nitrosodiphenylamine ²	ND
1,2-dichlorobenzene	ND	phenanthrene	ND
1,3-dichlorobenzene	ND	pyrene	ND
1,4-dichlorobenzene*	ND	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

¹ Screened for as Azobenzene ² Detected as Diphenylamine

Alyce R. Massie
Approved by _____ Laboratory Manager



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO:
Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 8 OF 19

Sample Description: WP-2 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acenaphthene	ND	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine ¹	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,i)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	ND	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
dibenzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
di-n-butylphthalate	ND	N-nitrosodiphenylamine ²	ND
1,2-dichlorobenzene	ND	phenanthrene	ND
1,3-dichlorobenzene	ND	pyrene	ND
1,4-dichlorobenzene	ND	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

¹ Screened for as Azobenzene ² Detected as Diphenylamine

Alice F. Moore
Approved by _____
Laboratory Manager

Title


CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
 ATTN: Bill Gray
 125 E. Bethpage Road
 Plainview, NY 11803

DATE REPORTED: November 15, 1988
 PROJECT CODE: GMIN 42142
 ORDER NUMBER:
 PAGE 10 OF 19

Sample Description: WP-3 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acenaphthene	ND	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine ¹	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,i)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	ND	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
dibenzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
di-n-butylphthalate	ND	N-nitrosodiphenylamine ²	ND
1,2-dichlorobenzene	ND	phenanthrene	ND
1,3-dichlorobenzene	ND	pyrene	ND
1,4-dichlorobenzene	ND	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

¹ Screened for as Azobenzene ² Detected as Diphenylamine

Alyce S. Massie
 Approved by _____ Laboratory Manager


CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 9 OF 25

Sample Description: WP-7 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)	<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)
acenaphthene	ND	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine†	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,i)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	ND	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
dibenzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
di-n-butylphthalate	ND	N-nitrosodiphenylamine ²	<10
1,2-dichlorobenzene	ND	phenanthrene	ND
1,3-dichlorobenzene	ND	pyrene	ND
1,4-dichlorobenzene	ND	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

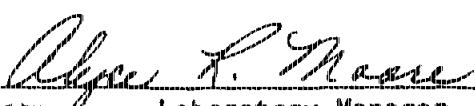
ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

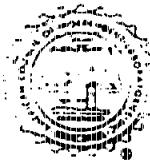
¹ Screened for as Azobenzene ² Detected as Diphenylamine

Approved by 
Glynn F. Moore
Laboratory Manager



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 11 OF 25

Sample Description: WP-8 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)	<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)
acenaphthene	<10	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine ¹	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,f)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	ND	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
dibenzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
di-n-butylphthalate	ND	N-nitrosodiphenylamine ²	ND
1,2-dichlorobenzene	ND	phenanthrene	ND
1,3-dichlorobenzene	ND	pyrene	ND
1,4-dichlorobenzene	ND	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

¹ Screened for as Azobenzene ² Detected as Diphenylamine

Approved by *Alcey L. Moore*
Laboratory Manager

Title


CERTIFICATE OF ANALYSIS

TO **Geraghty & Miller, Inc.**
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 13 OF 25

Sample Description: WP-9 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acenaphthene	ND	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine ¹	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,i)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	ND	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
dibenzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
di-n-butylphthalate	ND	N-nitrosodiphenylamine ²	<10
1,2-dichlorobenzene	ND	phenanthrene	ND
1,3-dichlorobenzene	ND	pyrene	ND
1,4-dichlorobenzene	ND	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

¹ Screened for as Azobenzene ² Detected as Diphenylamine

Alyce F. Moore
Approved by _____
Laboratory Manager


CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
 ATTN: Bill Gray
 125 E. Bethpage Road
 Plainview, NY 11803

DATE REPORTED: November 18, 1988
 PROJECT CODE: GMIN 42163
 ORDER NUMBER:
 PAGE 15 OF 25

Sample Description: 62 (Water)

BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
acenaphthene	13	3,3'-dichlorobenzidine **	ND
acenaphthylene	ND	diethyl phthalate	ND
anthracene	ND	dimethyl phthalate	ND
benzidine *	ND	2,4-dinitrotoluene	ND
benzo(a)anthracene	ND	2,6-dinitrotoluene	ND
benzo(b)fluoranthene	ND	di-n-octylphthalate	ND
benzo(k)fluoranthene	ND	1,2-diphenylhydrazine†	ND
benzo(a)pyrene	ND	fluoranthene	ND
benzo(g,h,i)perylene	ND	fluorene	ND
benzyl butyl phthalate	ND	hexachlorobenzene	ND
bis(2-chloroethoxy)methane	ND	hexachlorobutadiene	ND
bis(2-chloroethyl)ether	ND	hexachlorocyclopentadiene	ND
bis(2-chloroisopropyl)ether	ND	hexachloroethane	ND
bis(2-ethylhexyl)phthalate	<10	indeno(1,2,3-cd)pyrene	ND
4-bromophenyl phenyl ether	ND	isophorone	ND
2-chloronaphthalene	ND	naphthalene	ND
4-chlorophenyl phenyl ether	ND	nitrobenzene	ND
chrysene	ND	N-nitrosodimethylamine	ND
di benzo(a,h)anthracene	ND	N-nitrosodi-n-propylamine	ND
di-n-butylphthalate	ND	N-nitrosodiphenylamine‡	ND
1,2-dichlorobenzene	<10	phenanthrene	ND
1,3-dichlorobenzene	43	pyrene	ND
1,4-dichlorobenzene	92	1,2,4-trichlorobenzene	ND

Remarks: 10 = Quantitation limit.

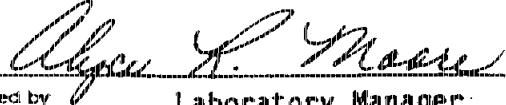
ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit five (5) times that listed.

** = This compound has a quantitation limit two (2) times that listed.

† Screened for as Azobenzene ‡ Detected as Diphenylamine


 Approved by Alyce J. Moore
 Title Laboratory Manager

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA1G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>226</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468169</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH068169A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/06/91</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>12/11/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>12/20/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
62-75-9	N-Nitrosodimethylamine	10	U
110-86-1	Pyridine	10	U
97-63-2	Ethyl methacrylate	10	U
109-06-8	2-Picoline	20	U
10595-95-6	Nitrosomethylethylamine	10	U
66-27-3	Methyl methanesulfonate	10	U
55-18-5	N-Nitrosodiethylamine	10	U
62-50-0	Ethyl methanesulfonate	10	U
108-95-2	Phenol	10	U
62-53-3	Aniline	10	U
76-01-7	Pentachloroethane	10	U
111-44-4	bis(2-Chloroethyl)Ether	20	U
95-57-8	2-Chlorophenol	10	U
541-73-1	1, 3-Dichlorobenzene	10	U
106-46-7	1, 4-Dichlorobenzene	10	U
100-51-6	Benzyl Alcohol	10	U
95-50-1	1, 2-Dichlorobenzene	10	U
95-48-7	2-Methylphenol	10	U
39638-32-9	bis(2-Chloroisopropyl)Ether	10	U
108-39-4	3-Methylphenol	10	U
106-44-5	4-Methylphenol	10	U
930-55-2	N-Nitrosopyrrolidine	10	U
59-89-2	N-Nitrosomorpholine	10	U
98-86-2	Acetophenone	10	U
621-64-7	N-Nitroso-Di-n-Propylamine	10	U
636-21-5	c-Toluidine hydrochloride	10	U
67-72-1	Hexachloroethane	10	U
98-95-3	Nitrobenzene	10	U
100-75-4	N-Nitrosopiperidine	10	U
78-59-1	Isophorone	10	U
88-75-5	2-Nitrophenol	10	U
105-67-9	2, 4-Dimethylphenol	10	U
65-85-0	Benzoic Acid	100	U
111-91-1	bis(2-Chloroethoxy)Methane	10	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

120-83-2-----	2, 4-Dichlorophenol	1.0	U
120-82-1-----	1, 2, 4-Trichlorobenzene	1.0	U
91-20-3-----	Naphthalene	34	
106-47-8-----	4-Chloroaniline	1.0	U
87-65-0-----	2, 6-Dichlorophenol	2.0	U
122-09-8-----	dimethylphenylethylamine	1.0	U
1888-71-7-----	Hexachloropropene	1.0	U
87-68-3-----	Hexachlorobutadiene	1.0	U
924-16-3-----	N-Nitroso-di-n-butylamine	1.0	U
59-50-7-----	4-Chloro-3-Methylphenol	1.0	U
106-50-3-----	P-Phenylenediamine	1.0	U
94-59-7-----	Safrole	1.0	U
91-57-6-----	2-Methylnaphthalene	1.3	
95-94-3-----	1, 2, 4, 5-Tetrachlorobenzene	1.0	U
77-47-4-----	Hexachlorocyclopentadiene	1.0	U
88-06-2-----	2, 4, 6-Trichlorophenol	2.0	U
95-95-4-----	2, 4, 5-Trichlorophenol	2.0	U
120-58-1-----	Icosafrole	2.0	U
91-58-7-----	2-Chloronaphthalene	1.0	U
68-74-4-----	2-Nitroaniline	1.0	U
130-15-4-----	1, 4-Naphthoquinone	2.0	U
131-11-3-----	Dimethyl Phthalate	1.0	U
208-96-8-----	Acenaphthylene	1	J
606-20-2-----	2, 6-Dinitrotoluene	1.0	U

463169

FORM I SV-1

1/87 Rev.

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM, RTP</u>	Contract: <u>500077</u>	ROA1G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>226</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468169</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH068169A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/06/91</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>12/11/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>12/20/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L Q	
99-09-2	3-Nitroaniline	20	U
83-32-9	Acenaphthene	15	
51-28-5	2,4-Dinitrophenol	40	U
100-02-7	4-Nitrophenol	10	U
132-64-9	Dibenzofuran	1	J
121-14-2	2,4-Dinitrotoluene	10	U
608-93-5	Pentachlorobenzene	10	U
91-59-8	2-Naphthylamine	20	U
134-32-7	1-Naphthylamine	20	U
58-90-2	2,3,4,6-Tetrachlorophenol	20	U
84-66-2	Diethylphthalate	10	U
297-97-2	Zinophos	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U
86-73-7	Fluorene	11	
100-01-6	4-Nitroaniline	20	U
99-55-8	5-Nitro-o-toluidine	20	U
122-66-7	1,2-Diphenylhydrazine	10	U
534-52-1	4,6-Dinitro-2-Methylphenol	30	U
86-30-6	N-Nitrosodiphenylamine (1)	10	U
122-39-4	Diphenylamine	10	U
99-35-4	1,3,5-Trinitrobenzene	20	U
62-44-2	Phenacetin	10	U
101-55-3	4-Bromophenyl-phenylether	10	U
2303-16-4	Diallate	10	U
60-51-5	Dimethoate	10	U
118-74-1	Hexachlorobenzene	10	U
92-67-1	4-Aminobiphenyl	10	U
23950-58-5	Pronamide	10	U
87-86-5	Pentachlorophenol	20	U
82-68-8	Pentachloronitrobenzene	10	U
85-01-8	Phenanthrrene	12	
120-12-7	Anthracene	4	J
84-74-2	Di-n-Butylphthalate	10	U
91-80-5	Methapyrilene	20	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0	Fluoranthene	2	J
92-87-5	Benzidine	10	J
129-00-0	Pyrene	4	J
60-11-7	p-Dimethylaminoazobenzene	10	U
510-15-6	Chlorobenzilate	10	U
119-93-7	3,3'-Dimethylbenzidine	20	U
85-68-7	Butylbenzylphthalate	10	U
53-96-3	2-Acetylaminofluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	10	U
218-01-9	Chrysene	10	U
117-81-7	bis(2-Ethylhexyl) Phthalate	10	U
117-84-0	Di-n-Octyl Phthalate	10	U
205-99-2	Benzo(b)Fluoranthene	10	U
57-97-6	7,12-Dimethylbenzanthracene	10	U
207-08-9	Benzo(k)Fluoranthene	10	U
50-32-8	Benzo(a)Pyrene	10	U
56-49-5	3-Methylcholanthrene	10	U
193-39-5	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3	Dibenz(a,h)Anthracene	10	U
191-24-2	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

10-31-9

1B
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	<u>ROA3G</u>
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>588</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>474981</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH074981A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>01/15/92</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>01/16/92</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>01/16/92</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>		Q
62-75-9	-N-Nitrosodimethylamine	10	U	
110-86-1	-Pyridine	10	U	
97-63-2	-Ethyl methacrylate	10	U	
109-06-8	-2-Picoline	10	U	
10595-95-6	-Nitrosomethylamine	20	U	
66-27-3	-Methyl methanesulfonate	10	U	
55-18-5	-N-Nitrosodiethylamine	10	U	
62-50-5	-Ethyl methanesulfonate	10	U	
108-95-2	-Phenol	10	U	
62-53-3	-Aniline	10	U	
76-01-7	-Pentachloroethane	10	U	
111-44-4	-bis(2-Chloroethyl) Ether	20	U	
95-57-8	-2-Chlorophenol	10	U	
541-73-1	-1, 3-Dichlorobenzene	10	U	
106-46-7	-1, 4-Dichlorobenzene	10	U	
100-51-6	-Benzyl Alcohol	10	U	
95-50-1	-1, 2-Dichlorobenzene	10	U	
95-48-7	-2-Methylphenol	8	J	
39638-32-9	-bis(2-Chloroisopropyl) Ether	10	U	
108-39-4	-3-Methylphenol	10	U	
106-44-5	-4-Methylphenol	10	U	
930-55-2	-N-Nitroso-N-propylamine	10	U	
59-89-2	-N-Nitrosomorpholine	10	U	
98-86-2	-Acetophenone	10	U	
621-64-7	-N-Nitroso-Di-n-Propylamine	10	U	
636-21-5	-o-Tolidine hydrochloride	10	U	
67-72-1	-Hexachloroethane	10	U	
98-95-3	-Nitrobenzene	10	U	
100-75-4	-N-Nitrosopiperidine	10	U	
78-59-1	-Isophorone	10	U	
88-75-5	-2-Nitrophenol	10	U	
105-67-9	-2, 4-Dimethylphenol	4	J	
65-85-0	-Benzoic Acid	3	J	
111-91-1	-bis(2-Chloroethoxy) Methane	10	U	

FORM I SV-1

1/87 Rev.

120-83-2	2, 4-Dichlorophenol	10	U
120-82-1	1, 2, 4-Trichlorobenzene	10	U
91-20-3	Naphthalene	3	J
106-47-8	4-Chloroaniline	10	U
87-65-0	2, 6-Dichlorophenol	20	U
122-09-8	dimethylphenylethylamine	10	U
1888-71-7	Hexachloropropene	10	U
87-68-3	Hexachlorobutadiene	10	U
924-16-3	N-Nitroso-di-n-butylamine	10	U
59-50-7	4-Chloro-3-Methylphenol	10	U
106-50-3	P-Phenylenediamine	10	U
94-59-7	Safrole	10	U
91-57-6	2-Methylnaphthalene	10	U
95-94-3	1, 2, 4, 5-Tetrachlorobenzene	10	U
77-47-4	Hexachlorocyclopentadiene	10	U
88-06-2	2, 4, 6-Trichlorophenol	10	U
95-95-4	2, 4, 5-Trichlorophenol	20	U
120-58-1	Isosafrole	20	U
91-58-7	2-Chloronaphthalene	10	U
88-74-4	2-Nitroaniline	10	U
130-15-4	1, 4-Naphthoquinone	20	U
131-11-3	Dimethyl Phthalate	10	U
208-96-8	Acenaphthylene	1	J
606-20-2	2, 6-Dinitrotoluene	10	U

FORM I SV-1

1/87 Rev.

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROA3G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>588</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>474981</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH074981A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>01/15/92</u>	
Moisture: not dec. _____ dec. _____	Date Extracted: <u>01/16/92</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>01/16/92</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>		Q
99-09-2	3-Nitroaniline	20	U	
83-32-9	Acenaphthene	36	U	
51-28-5	2, 4-Dinitrophenol	40	U	
100-02-7	4-Nitrophenol	10	U	
132-64-9	Dibenzofuran	10	U	
121-14-2	2, 4-Dinitrotoluene	10	U	
608-93-5	Pentachlorobenzene	10	U	
134-32-7	2-Naphthylamine	20	U	
134-32-7	1-Naphthylamine	20	U	
58-90-2	2, 3, 4, 6-Tetrachlorophenol	20	U	
84-66-2	Diethylphthalate	10	U	
297-97-2	Zinophos	10	U	
7005-72-3	4-Chlorophenyl-phenylether	10	U	
86-73-7	Fluorene	15	U	
100-01-6	4-Nitroaniline	20	U	
99-55-8	5-Nitro-o-toluidine	20	U	
122-66-7	1, 2-Diphenylhydrazine	10	U	
534-52-1	4, 6-Dinitro-2-Methylphenol	30	U	
86-30-6	N-Nitrosodiphenylamine (1)	10	U	
122-39-4	Diphenylamine	10	U	
99-35-4	1, 3, 5-Trinitrobenzene	20	U	
62-44-2	Phenacetin	10	U	
101-55-3	4-Bromophenyl-phenylether	10	U	
2303-16-4	Diallate	10	U	
60-51-5	Dimethoate	10	U	
118-74-1	Hexachlorobenzene	10	U	
92-67-1	4-Aminobiphenyl	10	U	
23950-58-5	Pronamide	10	U	
87-86-5	Pentachlorophenol	20	U	
82-68-8	Pentachloronitrobenzene	10	U	
85-01-8	Phenanthrene	7	J	
120-12-7	Anthracene	11	U	
84-74-2	Di-n-Butylphthalate	10	U	
91-80-5	Methapyrilene	20	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0	Fluoranthene	16	
92-87-5	Benzidine	10	U
129-00-0	Pyrene	12	
140-57-8	Aramite	20	U
60-11-7	p-Dimethylaminoazobenzene	10	U
510-15-6	Chlorobenzilate	10	U
119-93-7	3, 3'-Dimethylbenzidine	10	U
85-68-7	Butylbenzylphthalate	20	U
53-96-3	2-Acetylaminofluorene	10	U
91-94-1	3, 3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	10	J
117-81-7	bis(2-Ethylhexyl)Phthalate	5	BJ
218-01-9	Chrysene	3	J
117-84-0	Di-n-Octyl Phthalate	6	BJ
205-99-2	Benzo(b)Fluoranthene	10	U
57-97-6	7, 12-Dimethylbenzanthracene	7	JX
207-08-9	Benzo(k)Fluoranthene	10	U
50-32-8	Benzo(a)Pyrene	7	JX
56-49-5	3-Methylcholanthrene	4	J
193-39-5	Indeno(1, 2, 3-cd) Pyrene	10	U
53-70-3	Dibenz(a, h)Anthracene	1	J
191-24-2	Benzo(g, h, i)Perylene	10	U
		2	J

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROB1G

Lab Name: COMPUCHEM RTP Contract: 500077

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 226

Matrix: (soil/water) WATER Lab Sample ID: 468167

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH068167A20

Level: (low/med) LOW Date Received: 12/06/91

% Moisture: not dec. _____ dec. _____ Date Extracted: 12/11/91

Extraction: (SepF/Cont/Sonc) SEPF Date Analyzed: 12/20/91

GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
62-75-9	-N-Nitrosodimethylamine	10	U
110-86-1	-Pyridine	10	U
97-63-2	-Ethyl methacrylate	10	U
109-06-8	-2-Picoline	20	U
10595-95-6	-Nitrosomethyl ethylamine	10	U
66-27-3	-Methyl methanesulfonate	10	U
55-18-5	-N-Nitrosodiethylamine	10	U
62-50-0	-Ethyl methanesulfonate	10	U
108-95-2	-Phenol	10	U
62-53-3	-Aniline	10	U
76-01-7	-Pentachloroethane	10	U
111-44-4	-bis(2-Chloroethyl) Ether	20	U
95-57-8	-2-Chlorophenol	10	U
541-73-1	-1, 3-Dichlorobenzene	10	U
106-46-7	-1, 4-Dichlorobenzene	10	U
100-51-6	-Benzyl Alcohol	10	U
95-50-1	-1, 2-Dichlorobenzene	10	U
95-48-7	-2-Methylphenol	10	U
39638-32-9	-bis(2-Chloroisopropyl) Ether	10	U
108-39-4	-3-Methylphenol	10	U
106-44-5	-4-Methylphenol	10	U
930-55-2	-N-Nitrosopyrrolidine	10	U
59-89-2	-N-Nitrosomorpholine	10	U
98-86-2	-Acetophenone	10	U
621-64-7	-N-Nitroso-Di-n-Propylamine	10	U
636-21-5	-o-Tolidine hydrochloride	10	U
67-72-1	-Hexachloroethane	10	U
98-95-3	-Nitrobenzene	10	U
100-75-4	-N-Nitrosopiperidine	10	U
78-59-1	-Isophorone	10	U
88-75-5	-2-Nitrophenol	10	U
105-67-9	-2, 4-Dimethylphenol	10	U
65-85-0	-Benzoic Acid	100	U
111-91-1	-bis(2-Chloroethoxy) Methane	10	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-65-0-----	2,6-Dichlorophenol	20	U
122-09-8-----	dimethylphenylethylamine	10	U
1888-71-7-----	Hexachloropropene	10	U
87-68-3-----	Hexachlorobutadiene	10	U
924-16-3-----	N-Nitroso-di-n-butylamine	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
106-50-3-----	P-Phenylenediamine	10	U
94-59-7-----	Safrole	10	U
91-57-6-----	2-Methylnaphthalene	10	U
95-94-3-----	1,2,4,5-Tetrachlorobenzene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	20	U
95-95-4-----	2,4,5-Trichlorophenol	20	U
120-58-1-----	Isosafrole	20	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	10	U
130-15-4-----	1,4-Naphthoquinone	20	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U

4/23/67

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB1G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>226</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468167</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH068167A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/06/91</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>12/11/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>12/20/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L Q	
99-09-2	3-Nitroaniline	20	U
83-32-9	Acenaphthene	10	U
51-28-5	2,4-Dinitrophenol	40	U
100-02-7	4-Nitrophenol	10	U
132-64-9	Dibenzofuran	10	U
121-14-2	2,4-Dinitrotoluene	10	U
608-93-5	Pentachlorobenzene	10	U
91-59-8	2-Naphthylamine	20	U
134-32-7	1-Naphthylamine	20	U
58-90-2	2,3,4,6-Tetrachlorophenol	20	U
84-66-2	Diethylphthalate	10	U
297-97-2	Zinophos	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U
86-73-7	Fluorene	10	U
100-01-6	4-Nitroaniline	20	U
99-55-8	5-Nitro-o-toluidine	20	U
122-66-7	1,2-Diphenylhydrazine	10	U
534-52-1	4,6-Dinitro-2-Methylphenol	30	U
86-30-6	N-Nitrosodiphenylamine (1)	10	U
122-39-4	Diphenylamine	10	U
99-35-4	1,3,5-Trinitrobenzene	20	U
62-44-2	Phenacetin	10	U
101-55-3	4-Bromophenyl-phenylether	10	U
2303-16-4	Diallate	10	U
60-51-5	Dimethoate	10	U
118-74-1	Hexachlorobenzene	10	U
92-67-1	4-Aminobiphenyl	10	U
23950-58-5	Pronamide	10	U
87-86-5	Pentachlorophenol	20	U
82-68-8	Pentachloronitrobenzene	10	U
85-01-8	Phenanthrone	10	U
120-12-7	Anthracene	10	U
84-74-2	Di-n-Butylphthalate	10	U
91-80-5	Methapyrilene	20	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0	Fluoranthene	10	U
92-87-5	Benzidine	10	U
129-00-0	Pyrene	10	U
60-11-7	p-Dimethylaminoazobenzene	10	U
510-15-6	Chlorobenzilate	10	U
119-93-7	3,3'-Dimethylbenzidine	20	U
85-68-7	Butylbenzylphthalate	10	U
53-96-3	2-Acetylaminofluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	10	U
218-01-9	Chrysene	10	U
117-81-7	bis(2-Ethylhexyl)Phthalate	10	U
117-84-0	Di-n-Octyl Phthalate	10	U
205-99-2	Benzo(b)Fluoranthene	10	U
57-97-6	7,12-Dimethylbenzanthracene	10	U
207-08-9	Benzo(k)Fluoranthene	10	U
50-32-8	Benzo(a)Pyrene	10	U
56-49-5	3-Methylcholanthrene	10	U
193-39-5	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3	Dibenz(a,h)Anthracene	10	U
191-24-2	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

44-2167

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROB2G

Lab Name: COMPUCHEM RTP Contract: 500077

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 226

Matrix: (soil/water) WATER Lab Sample ID: 468168

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GR068168B22

Level: (low/med) LOW Date Received: 12/07/91

% Moisture: not dec. _____ dec. _____ Date Extracted: 12/12/91

Extraction: (SepF/Cont/Sonc) SEPF Date Analyzed: 12/26/91

GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/L	Q
62-75-9	-N-Nitrosodimethylamine	10	U	
110-86-1	-Pyridine	10	U	
97-63-2	-Ethyl methacrylate	10	U	
109-06-8	-2-Picoline	20	U	
10595-95-6	-Nitrosomethylethylamine	10	U	
66-27-3	-Methyl methanesulfonate	10	U	
55-18-5	-N-Nitrosodiethylamine	10	U	
62-50-0	-Ethyl methanesulfonate	10	U	
108-95-2	-Phenol	10	U	
62-53-3	-Aniline	10	U	
76-01-7	-Pentachloroethane	10	U	
111-44-4	-bis(2-Chloroethyl) Ether	20	U	
95-57-8	-2-Chlorophenol	10	U	
541-73-1	-1,3-Dichlorobenzene	10	U	
106-46-7	-1,4-Dichlorobenzene	10	U	
100-51-6	-Benzyl Alcohol	10	U	
95-50-1	-1,2-Dichlorobenzene	10	U	
95-48-7	-2-Methylphenol	10	U	
39638-32-9	-bis(2-Chloroisopropyl) Ether	10	U	
108-39-4	-3-Methylphenol	10	U	
106-44-5	-4-Methylphenol	10	U	
930-55-2	-N-Nitrosopyrrolidine	10	U	
59-89-2	-N-Nitrosomorpholine	10	U	
98-86-2	-Acetophenone	10	U	
621-64-7	-N-Nitroso-Di-n-Propylamine	10	U	
636-21-5	-o-Tolidine hydrochloride	10	U	
67-72-1	-Hexachloroethane	10	U	
98-95-3	-Nitrobenzene	10	U	
100-75-4	-N-Nitrosopiperidine	10	U	
78-59-1	-Isophorone	10	U	
88-75-5	-2-Nitrophenol	10	U	
105-67-9	-2,4-Dimethylphenol	10	U	
65-85-0	-Benzoic Acid	100	U	
111-91-1	-bis(2-Chloroethoxy) Methane	10	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

120-83-2-----	2, 4-Dichlorophenol	10	U
120-82-1-----	1, 2, 4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-65-0-----	2, 6-Dichlorophenol	20	U
122-09-8-----	dimethylphenylethylamine	10	U
1888-71-7-----	Hexachloropropene	10	U
87-68-3-----	Hexachlorobutadiene	10	U
924-16-3-----	N-Nitroso-di-n-butylamine	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
106-50-3-----	p-Phenylenediamine	10	U
94-59-7-----	Safrole	10	U
91-57-6-----	2-Methylnaphthalene	10	U
95-94-3-----	1, 2, 4, 5-Tetrachlorobenzene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2, 4, 6-Trichlorophenol	20	U
95-95-4-----	2, 4, 5-Trichlorophenol	20	U
120-58-1-----	Isosafrole	20	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	10	U
130-15-4-----	1, 4-Naphthoquinone	20	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2, 6-Dinitrotoluene	10	U

463168

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	ROB2G
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468168</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GR068168B22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/07/91</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>12/12/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>12/26/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/L	Q
99-09-2-----	3-Nitroaniline	20	U	
83-32-9-----	Acenaphthene	10	U	
51-28-5-----	2,4-Dinitrophenol	40	U	
100-02-7-----	4-Nitrophenol	10	U	
132-64-9-----	Dibenzofuran	10	U	
121-14-2-----	2,4-Dinitrotoluene	10	U	
608-93-5-----	Pentachlorobenzene	10	U	
91-59-8-----	2-Naphthylamine	20	U	
134-32-7-----	1-Naphthylamine	20	U	
58-90-2-----	2,3,4,6-Tetrachlorophenol	20	U	
84-66-2-----	Diethylphthalate	10	U	
297-97-2-----	Zinophos	10	U	
7005-72-3-----	4-Chlorophenyl-phenylether	10	U	
86-73-7-----	Fluorene	10	U	
100-01-6-----	4-Nitroaniline	20	U	
99-55-8-----	5-Nitro-o-toluidine	20	U	
122-66-7-----	1,2-Diphenylhydrazine	10	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	30	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U	
122-39-4-----	Diphenylamine	10	U	
99-35-4-----	1,3,5-Trinitrobenzene	20	U	
62-44-2-----	Phenacetin	10	U	
101-55-3-----	4-Bromophenyl-phenylether	10	U	
2303-16-4-----	Diallate	10	U	
60-51-5-----	Dimethoate	10	U	
118-74-1-----	Hexachlorobenzene	10	U	
92-67-1-----	4-Aminobiphenyl	10	U	
23950-58-5-----	Pronamide	10	U	
87-86-5-----	Pentachlorophenol	20	U	
82-68-8-----	Pentachloronitrobenzene	10	U	
85-01-8-----	Phenanthrene	10	U	
120-12-7-----	Anthracene	10	U	
84-74-2-----	Di-n-Butylphthalate	10	U	
91-80-5-----	Methapyrilene	20	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0	Fluoranthene	10	U
92-87-5	Benzidine	10	U
129-00-0	Pyrene	10	U
60-11-7	p-Dimethylaminoazobenzene	10	U
510-15-6	Chlorobenzilate	10	U
119-93-7	3,3'-Dimethylbenzidine	20	U
85-68-7	Butylbenzylphthalate	10	U
53-96-3	2-Acetylaminofluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	10	U
218-01-9	Chrysene	10	U
117-81-7	bis(2-Ethylhexyl)Phthalate	2	J
117-84-0	Di-n-Octyl Phthalate	10	U
205-99-2	Benzo(b)Fluoranthene	10	U
57-97-6	7,12-Dimethylbenzanthracene	10	U
207-08-9	Benzo(k)Fluoranthene	10	U
50-32-8	Benzo(a)Pyrene	10	U
56-49-5	3-Methylcholanthrene	10	U
193-39-5	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3	Dibenz(a,h)Anthracene	10	U
191-24-2	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

448'68

FORM I SV-3

1/87 Rev.

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC1G

Lab Name: COMPUCHEM RTP

Contract: 500077

Lab Code: COMPU Case No.: 24105

SAS No.: _____

SDG No.: 226

Matrix: (soil/water) WATER

Lab Sample ID: 468513

Sample wt/vol: 800 (g/mL) ML

Lab File ID: GH068513A20

Level: (low/med) LOW

Date Received: 12/07/91

% Moisture: not dec. _____ dec. _____

Date Extracted: 12/11/91

Extraction: (SepF/Cont/Sonc) SEPF

Date Analyzed: 12/21/91

GPC Cleanup: (Y/N) N pH: _____

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
62-75-9	N-Nitrosodimethylamine	12	U
110-86-1	Pyridine	12	U
97-63-2	Ethyl methacrylate	12	U
109-06-8	2-Picoline	25	U
10595-95-6	Nitrosomethylamine	12	U
66-27-3	Methyl methanesulfonate	12	U
55-18-5	N-Nitrosodiethylamine	12	U
62-50-0	Ethyl methanesulfonate	12	U
108-95-2	Phenol	12	U
62-53-3	Aniline	12	U
76-01-7	Pentachloroethane	12	U
111-44-4	bis(2-Chloroethyl) Ether	25	U
95-57-8	2-Chlorophenol	12	U
541-73-1	1,3-Dichlorobenzene	12	U
106-46-7	1,4-Dichlorobenzene	12	U
100-51-6	Benzyl Alcohol	12	U
95-50-1	1,2-Dichlorobenzene	12	U
95-48-7	2-Methylphenol	12	U
39638-32-9	bis(2-Chloroisopropyl) Ether	12	U
108-39-4	3-Methylphenol	12	U
106-44-5	4-Methylphenol	12	U
930-55-2	N-Nitrosopyrrolidine	12	U
59-89-2	N-Nitrosomorpholine	12	U
98-86-2	Acetophenone	12	U
621-64-7	N-Nitroso-Di-n-Propylamine	12	U
636-21-5	o-Tolidine hydrochloride	12	U
67-72-1	Hexachloroethane	12	U
98-95-3	Nitrobenzene	12	U
100-75-4	N-Nitrosopiperidine	12	U
78-59-1	Isophorone	12	U
88-75-5	2-Nitrophenol	12	U
105-67-9	2,4-Dimethylphenol	12	U
65-85-0	Benzoic Acid	120	U
111-91-1	bis(2-Chloroethoxy) Methane	12	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

120-83-2-----	2, 4-Dichlorophenol	12	U
120-82-1-----	1, 2, 4-Trichlorobenzene	12	U
91-20-3-----	Naphthalene	12	U
106-47-8-----	4-Chloroaniline	12	U
87-65-0-----	2, 6-Dichlorophenol	25	U
122-09-8-----	dimethylphenylethylamine	12	U
1888-71-7-----	Hexachloropropene	12	U
87-68-3-----	Hexachlorobutadiene	12	U
924-16-3-----	N-Nitroso-di-n-butylamine	12	U
59-50-7-----	4-Chloro-3-Methylphenol	12	U
106-50-3-----	P-Phenylenediamine	12	U
94-59-7-----	Safrole	12	U
91-57-6-----	2-Methylnaphthalene	12	U
95-94-3-----	1, 2, 4, 5-Tetrachlorobenzene	12	U
77-47-4-----	Hexachlorocyclopentadiene	12	U
88-06-2-----	2, 4, 6-Trichlorophenol	25	U
95-95-4-----	2, 4, 5-Trichlorophenol	25	U
120-58-1-----	Isosafrole	25	U
91-58-7-----	2-Chloronaphthalene	12	U
88-74-4-----	2-Nitroaniline	12	U
130-15-4-----	1, 4-Naphthoquinone	25	U
131-11-3-----	Dimethyl Phthalate	12	U
208-96-8-----	Acenaphthylene	12	U
606-20-2-----	2, 6-Dinitrotoluene	12	U

468513

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC1G

Lab Name: COMPUCHEM, RTP

Contract: 500077

Lab Code: COMPU Case No.: 24105 SAS No.: _____ SDG No.: 226

Matrix: (soil/water) WATER

Lab Sample ID: 468513

Sample wt/vol: 800 (g/mL) ML

Lab File ID: GH068513A20

Level: (low/med) LOW

Date Received: 12/07/91

* Moisture: not dec. _____ dec. _____

Date Extracted: 12/11/91

Extraction: (SepF/Cont/Sonc) SEPF

Date Analyzed: 12/21/91

GPC Cleanup: (Y/N) N pH: _____

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
---------	----------	---	---

99-09-2	3-Nitroaniline	25	U
83-32-9	Acenaphthene	12	U
51-28-5	2, 4-Dinitrophenol	50	U
100-02-7	4-Nitrophenol	12	U
132-64-9	Dibenzofuran	12	U
121-14-2	2, 4-Dinitrotoluene	12	U
608-93-5	Pentachlorobenzene	12	U
91-59-8	2-Naphthylamine	25	U
134-32-7	1-Naphthylamine	25	U
58-90-2	2, 3, 4, 6-Tetrachlorophenol	25	U
84-66-2	Diethylphthalate	12	U
297-97-2	Zinophos	12	U
7005-72-3	4-Chlorophenyl-phenylether	12	U
86-73-7	Fluorene	12	U
100-01-6	4-Nitroaniline	25	U
99-55-8	5-Nitro-o-toluidine	25	U
122-66-7	1, 2-Diphenylhydrazine	12	U
534-52-1	4, 6-Dinitro-2-Methylphenol	38	U
86-30-6	N-Nitrosodiphenylamine (1)	12	U
122-39-4	Diphenylamine	12	U
99-35-4	1, 3, 5-Trinitrobenzene	25	U
62-44-2	Phenacetin	12	U
101-55-3	4-Bromophenyl-phenylether	12	U
2303-16-4	Diallate	12	U
60-51-5	Dimethoate	12	U
118-74-1	Hexachlorobenzene	12	U
92-67-1	4-Aminobiphenyl	12	U
23950-58-5	Pronamide	12	U
87-86-5	Pentachlorophenol	25	U
82-68-8	Pentachloronitrobenzene	12	U
85-01-8	Phenanthrene	12	U
120-12-7	Anthracene	12	U
84-74-2	Di-n-Butylphthalate	12	U
91-80-5	Methapryrilene	25	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0-----	Fluoranthene	12	U
92-87-5-----	Benzidine	12	U
129-00-0-----	Pyrene	12	U
60-11-7-----	p-Dimethylaminoazobenzene	12	U
510-15-6-----	Chlorobenzilate	12	U
119-93-7-----	3,3'-Dimethylbenzidine	25	U
85-68-7-----	Butylbenzylphthalate	12	U
53-96-3-----	2-Acetylaminofluorene	12	U
91-94-1-----	3,3'-Dichlorobenzidine	12	U
56-55-3-----	Benzo(a)Anthracene	12	U
218-01-9-----	Chrysene	12	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	12	U
117-84-0-----	Di-n-Octyl Phthalate	12	U
205-99-2-----	Benzo(b)Fluoranthene	12	U
57-97-6-----	7,12-Dimethylbenzanthracene	12	U
207-08-9-----	Benzo(k)Fluoranthene	12	U
50-32-8-----	Benzo(a)Pyrene	12	U
56-49-5-----	3-Methylcholanthrene	12	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	12	U
53-70-3-----	Dibenz(a,h)Anthracene	12	U
191-24-2-----	Benzo(g,h,i)Perylene	12	U

(1) - Cannot be separated from Diphenylamine

443513

FORM I SV-3

1/87 Rev.

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC2G

Lab Name: COMPUCHEM, RTP

Contract: 500077

Lab Code: COMPU

Case No.: 24105

SAS No.: _____

SDG No.: 226

Matrix: (soil/water) WATER

Lab Sample ID: 468590

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH068590A20

Level: (low/med) LOW

Date Received: 12/07/91

% Moisture: not dec. dec.

Date Extracted: 12/11/91

Extraction: (SepF/Cont/Sonc) SEPF

Date Analyzed: 12/21/91

GPC Cleanup: (Y/N) N pH:

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/L

62-75-9	-N-Nitrosodimethylamine	10	U
110-86-1	-Pyridine	10	U
97-63-2	-Ethyl methacrylate	10	U
109-06-8	-2-Picoline	20	U
10595-95-6	-Nitrosomethylethylamine	10	U
66-27-3	-Methyl methanesulfonate	10	U
55-18-5	-N-Nitrosodiethylamine	10	U
62-50-0	-Ethyl methanesulfonate	10	U
108-95-2	-Phenol	10	U
62-53-3	-Aniline	10	U
76-01-7	-Pentachloroethane	10	U
111-44-4	-bis(2-Chloroethyl)Ether	20	U
95-57-8	-2-Chlorophenol	10	U
541-73-1	-1,3-Dichlorobenzene	10	U
106-46-7	-1,4-Dichlorobenzene	10	U
100-51-6	-Benzyl Alcohol	10	U
95-50-1	-1,2-Dichlorobenzene	10	U
95-48-7	-2-Methylphenol	10	U
39638-32-9	-bis(2-Chloroisopropyl)Ether	10	U
108-39-4	-3-Methylphenol	10	U
106-44-5	-4-Methylphenol	10	U
930-55-2	-N-Nitroso- <i>p</i> -Rolidine	10	U
59-89-2	-N-Nitrosomorpholine	10	U
98-86-2	-Acetophenone	10	U
621-64-7	-N-Nitroso-Di-n-Propylamine	10	U
636-21-5	-o-Toluidine hydrochloride	10	U
67-72-1	-Hexachloroethane	10	U
98-95-3	-Nitrobenzene	10	U
100-75-4	-N-Nitroso-piperidine	10	U
78-59-1	-Isophorone	10	U
88-75-5	-2-Nitrophenol	10	U
105-67-9	-2,4-Dimethylphenol	10	U
65-85-0	-Benzoic Acid	100	U
111-91-1	-bis(2-Chloroethoxy)Methane	10	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-4

1/87 Rev.

120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-65-0-----	2,6-Dichlorophenol	20	U
122-09-8-----	dimethylphenylethylamine	10	U
1888-71-7-----	Hexachloropropene	10	U
87-68-3-----	Hexachlorobutadiene	10	U
924-16-3-----	N-Nitroso-di-n-butylamine	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
106-50-3-----	P-Phenylenediamine	10	U
94-59-7-----	Safrole	10	U
91-57-6-----	2-Methylnaphthalene	10	U
95-94-3-----	1,2,4,5-Tetrachlorobenzene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	20	U
95-95-4-----	2,4,5-Trichlorophenol	20	U
120-58-1-----	Isosafrole	20	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	10	U
130-15-4-----	1,4-Naphthoquinone	20	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U

4/85/70

FORM I SV-1

1/87 Rev.

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROC2G

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>226</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>468590</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH068590A20</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/07/91</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>12/11/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>12/21/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>		Q
99-09-2	3-Nitroaniline	20	U	
83-32-9	Acenaphthene	10	U	
51-28-5	2, 4-Dinitrophenol	40	U	
100-02-7	4-Nitrophenol	10	U	
132-64-9	Dibenzofuran	10	U	
121-14-2	2, 4-Dinitrotoluene	10	U	
608-93-5	Pentachlorobenzene	10	U	
91-59-8	2-Naphthylamine	20	U	
134-32-7	1-Naphthylamine	20	U	
58-90-2	2, 3, 4, 6-Tetrachlorophenol	20	U	
84-66-2	Diethylphthalate	10	U	
297-97-2	Zinophos	10	U	
7005-72-3	4-Chlorophenyl-phenylether	10	U	
86-73-7	Fluorene	2	J	
100-01-6	4-Nitroaniline	20	U	
99-55-8	5-Nitro-o-toluidine	20	U	
122-66-7	1, 2-Diphenylhydrazine	10	U	
534-52-1	4, 6-Dinitro-2-Methylphenol	30	U	
86-30-6	N-Nitrosodiphenylamine (1)	10	U	
122-39-4	Diphenylamine	10	U	
99-35-4	1, 3, 5-Trinitrobenzene	20	U	
62-44-2	Phenacetin	10	U	
101-55-3	4-Bromophenyl-phenylether	10	U	
2303-16-4	Diallate	10	U	
60-51-5	Dimethoate	10	U	
118-74-1	Hexachlorobenzene	10	U	
92-67-1	4-Aminobiphenyl	10	U	
23950-58-5	Pronamide	10	U	
87-86-5	Pentachlorophenol	20	U	
82-68-8	Pentachloronitrobenzene	10	U	
85-01-8	Phenanthrone	4	J	
120-12-7	Anthracene	1	J	
84-74-2	Di-n-Butylphthalate	10	U	
91-80-5	Methapyrilene	20	U	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0	Fluoranthene	4	J
92-87-5	Benzidine	10	U
129-00-0	Pyrene	4	J
60-11-7	p-Dimethylaminoazobenzene	10	U
510-15-6	Chlorobenzilate	10	U
119-93-7	3,3'-Dimethylbenzidine	20	U
85-68-7	Butylbenzylphthalate	10	U
53-96-3	2-Acetylaminofluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	2	J
218-01-9	Chrysene	2	J
117-81-7	bis(2-Ethylhexyl)Phthalate	10	U
117-84-0	Di-n-Octyl Phthalate	10	U
205-99-2	Benzo(b)Fluoranthene	3	JX
57-97-6	7,12-Dimethylbenzanthracene	10	U
207-08-9	Benzo(k)Fluoranthene	3	JX
50-32-8	Benzo(a)Pyrrene	2	J
56-49-5	3-Methylcholanthrene	10	U
193-39-5	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3	Dibenz(a,h)Anthracene	10	U
191-24-2	Benzo(g,h,i)Perylene	1	J

(1) - Cannot be separated from Diphenylamine

481570

1B
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROJ1G

Lab Name: <u>COMPUCHEM RTP</u>	Contract: <u>500077</u>	
Lab Code: <u>COMPU</u>	Case No.: <u>24105</u>	SAS No.: _____ SDG No.: <u>540</u>
Matrix: (soil/water) <u>WATER</u>	Lab Sample ID: <u>472796</u>	
Sample wt/vol: <u>1000</u> (g/mL) <u>ML</u>	Lab File ID: <u>GH072796A22</u>	
Level: (low/med) <u>LOW</u>	Date Received: <u>12/20/91</u>	
% Moisture: not dec. _____ dec. _____	Date Extracted: <u>12/26/91</u>	
Extraction: (SepF/Cont/Sonc) <u>SEPF</u>	Date Analyzed: <u>12/28/91</u>	
GPC Cleanup: (Y/N) <u>N</u>	pH: _____	Dilution Factor: <u>1.0</u>

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/L Q

62-75-9-----	N-Nitrosodimethylamine	10	U
110-86-1-----	Pyridine	10	U
97-63-2-----	Ethyl methacrylate	10	U
109-06-8-----	2-Picoline	20	U
10595-95-6-----	Nitrosomethylethylamine	10	U
66-27-3-----	Methyl methanesulfonate	10	U
55-18-5-----	N-Nitrosodiethylamine	10	U
62-50-0-----	Ethyl methanesulfonate	10	U
108-95-2-----	Phenol	10	U
62-53-3-----	Aniline	10	U
76-01-7-----	Pentachloroethane	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	20	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1, 3-Dichlorobenzene	10	U
106-46-7-----	1, 4-Dichlorobenzene	10	U
100-51-6-----	Benzyl Alcohol	10	U
95-50-1-----	1, 2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
39638-32-9-----	bis(2-Chloroisopropyl) Ether	10	U
108-39-4-----	3-Methylphenol	10	U
106-44-5-----	4-Methylphenol	10	U
930-55-2-----	N-Nitrosopyrrolidine	10	U
59-89-2-----	N-Nitrosomorpholine	10	U
98-86-2-----	Acetophenone	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
636-21-5-----	o-Tolidine hydrochloride	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
100-75-4-----	N-Nitrosopiperidine	10	U
78-59-1-----	Iscophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2, 4-Dimethylphenol	10	U
65-85-0-----	Benzoic Acid	100	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U

120-83-2	2, 4-Dichlorophenol	10	U
120-82-1	1, 2, 4-Trichlorobenzene	10	U
91-20-3	Naphthalene	10	U
106-47-8	4-Chloroaniline	10	U
87-65-0	2, 6-Dichlorophenol	20	U
122-09-8	dimethylphenylethylamine	10	U
1888-71-7	Hexachloropropene	10	U
87-68-3	Hexachlorobutadiene	10	U
924-16-3	N-Nitroso-di-n-butylamine	10	U
59-50-7	4-Chloro-3-Methylphenol	10	U
106-50-3	P-Phenylenediamine	10	U
94-59-7	Safrole	10	U
91-57-6	2-Methylnaphthalene	10	U
95-94-3	1, 2, 4, 5-Tetrachlorobenzene	10	U
77-47-4	Hexachlorocyclopentadiene	10	U
88-06-2	2, 4, 6-Trichlorophenol	20	U
95-95-4	2, 4, 5-Trichlorophenol	20	U
120-58-1	Isosafrole	20	U
91-58-7	2-Chloronaphthalene	10	U
88-74-4	2-Nitroaniline	10	U
130-15-4	1, 4-Naphthoquinone	20	U
131-11-3	Dimethyl Phthalate	10	U
208-96-8	Acenaphthyliene	10	U
606-20-2	2, 6-Dinitrotoluene	10	U

1C
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ROJ1G

Lab Name: COMPUCHEM, RTP

Contract: 500077

Lab Code: COMPU Case No.: 24105 SAS No.: _____

SDG No.: 540

Matrix: (soil/water) WATER

Lab Sample ID: 472796

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH072796A22

Level: (low/med) LOW

Date Received: 12/20/91

% Moisture: not dec. _____ dec. _____

Date Extracted: 12/26/91

Extraction: (SepF/Cont/Sonc) SEPE

Date Analyzed: 12/28/91

GPC Cleanup: (Y/N) N pH: _____

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg)	UG/L	Q
99-09-2	3-Nitroaniline	20	U	
83-32-9	Acenaphthene	10	U	
51-28-5	2, 4-Dinitrophenol	40	U	
100-02-7	4-Nitrophenol	10	U	
132-64-9	Dibenzofuran	10	U	
121-14-2	2, 4-Dinitrotoluene	10	U	
608-93-5	Pentachlorobenzene	10	U	
91-59-8	2-Naphthylamine	20	U	
134-32-7	1-Naphthylamine	20	U	
58-90-2	2, 3, 4, 6-Tetrachlorophenol	20	U	
84-66-2	Diethylphthalate	10	U	
297-97-2	Zinophos	10	U	
7005-72-3	4-Chlorophenyl-phenylether	10	U	
86-73-7	Fluorene	10	U	
100-01-6	4-Nitroaniline	20	U	
99-55-8	5-Nitro-o-toluidine	20	U	
122-66-7	1, 2-Diphenylhydrazine	10	U	
534-52-1	4, 6-Dinitro-2-Methylphenol	30	U	
86-30-6	N-Nitrosodiphenylamine (1)	10	U	
122-39-4	Diphenylamine	10	U	
99-35-4	1, 3, 5-Trinitrobenzene	20	U	
62-44-2	Phenacetin	10	U	
101-55-3	4-Bromophenyl-phenylether	10	U	
2303-16-4	Diallate	10	U	
60-51-5	Dimethoate	10	U	
118-74-1	Hexachlorobenzene	10	U	
92-67-1	4-Aminobiphenyl	10	U	
23950-58-5	Pronamide	10	U	
87-86-5	Pentachlorophenol	20	U	
82-68-8	Pentachloronitrobenzene	10	U	
85-01-8	Phenanthrene	10	U	
120-12-7	Anthracene	10	U	
84-74-2	Di-n-Butylphthalate	10	U	
91-80-5	Methapyrilene	20	U	

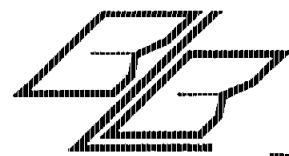
(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

206-44-0	Fluoranthene	10	U
92-87-5	Benzidine	10	U
129-00-0	Pyrene	10	U
60-11-7	p-Dimethylaminoazobenzene	10	U
510-15-6	Chlorobenzilate	10	U
119-93-7	3,3'-Dimethylbenzidine	20	U
85-68-7	Butylbenzylphthalate	10	U
53-96-3	2-Acetylaminofluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	10	U
218-01-9	Chrysene	10	U
117-81-7	bis(2-Ethylhexyl) Phthalate	1	BJ
117-84-0	Di-n-Octyl Phthalate	10	U
205-99-2	Benzo(b)Fluoranthene	10	U
57-97-6	7,12-Dimethylbenzanthracene	10	U
207-08-9	Benzo(k)Fluoranthene	10	U
50-32-8	Benzo(a)Pyrene	10	U
56-49-5	3-Methylcholanthrene	10	U
193-39-5	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3	Dibenz(a,h)Anthracene	10	U
191-24-2	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine



Section 13

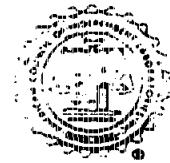
PESTICIDES / PCB ANALYSIS (GROUNDWATER)

WP-1 - Groundwater sample from Oxbow Area K, Well Point 1
WP-2 - Groundwater sample from Oxbow Area K, Well Point 2
WP-3 - Groundwater sample from Oxbow Area J, Well Point 3
WP-7 - Groundwater sample from Oxbow Area C, Well Point 7
WP-8 - Groundwater sample from Oxbow Area C, Well Point 8
WP-9 - Groundwater sample from Oxbow Area A, Well Point 9
ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 14 OF 19

Sample Description: WP-1 (Water)

PESTICIDES AND PCB's - PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)	<u>Compound</u>	<u>Concentration</u> ($\mu\text{g/liter}$)
aldrin	ND	endrin aldehyde ¹	ND
α -BHC	ND	heptachlor	ND (5.9)**
β -BHC	ND (14)**	heptachlor epoxide	ND
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND (1.1)**
δ -BHC	ND (0.7)**	PCB-(Aroclor)-1254 ¹	ND (1.9)*
chlordane ¹	ND	PCB-(Aroclor)-1221 ¹	ND (1.1)**
4,4'-DDT	ND	PCB-(Aroclor)-1232 ¹	ND (1.7)**
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND (2.0)**
4,4'-DDD	ND	PCB-(Aroclor)-1260 ¹	ND
dieldrin	ND	PCB-(Aroclor)-1016 ¹	ND
α -endosulfan	ND	toxaphene ¹	ND (3.3)*
β -endosulfan	ND		
endosulfan sulfate	ND		
endrin	ND		

Remarks: 0.5 = Quantitation limit

ND = Not detected

¹ = This component has a quantitation limit two (2) times that listed.

[†]Sample exhibits alteration of standard Aroclor pattern.

*Elevated detection limit due to presence of Aroclor 1260.

**Elevated quantitation limit listed in parenthesis due to sample matrix interferences.

Alyce S. Greene
Approved by _____
Laboratory Manager

Title _____



ANALYTICAL SERVICES



5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401

CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 15 OF 19

Sample Description: WP-2 (Water)

PESTICIDES AND PCB's - PRIORITY POLLUTANT ANALYSIS

Compound	Concentration (μ g/liter)	Compound	Concentration (μ g/liter)
aldrin	ND	endrin aldehyde ¹	ND
α -BHC	ND	heptachlor	ND (2:3)***
β -BHC	ND (5.4)**	heptachlor epoxide	ND
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND
δ -BHC	ND	PCB-(Aroclor)-1254 ¹	ND (1.1)**
chlordane ¹	ND	PCB-(Aroclor)-1221 ¹	ND
4,4'-DDT	ND	PCB-(Aroclor)-1232 ¹	ND
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND (1.5)**
4,4'-DDD	ND	PCB-(Aroclor)-1260 ¹	ND (1.3)**
dieldrin	ND	PCB-(Aroclor)-1016 ¹	ND
α -endosulfan	ND	toxaphene ¹	ND (1.5)**
β -endosulfan	ND		
endosulfan sulfate	ND		
endrin	ND		

Remarks: 0.5 = Quantitation limit

ND = Not detected

¹ = This component has a quantitation limit two (2) times that listed.

**Elevated quantitation limit listed in parenthesis due to sample matrix interferences.

Approved by Alice J. Nease Laboratory Manager

Title



ANALYTICAL SERVICES



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CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 16 OF 19

Sample Description: WP-3 (Water)

PESTICIDES AND PCB's - PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
aldrin	ND	endrin aldehyde ¹	ND
α -BHC	ND	heptachlor	ND (0.7)**
β -BHC	ND (1.6)**	heptachlor epoxide	ND
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND
δ -BHC	ND	PCB-(Aroclor)-1254 ¹	ND
chlordane ¹	ND	PCB-(Aroclor)-1221 ¹	ND
4,4'-DDT	ND	PCB-(Aroclor)-1232 ¹	ND
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND
4,4'-DDD	ND	PCB-(Aroclor)-1260 ¹	ND
dieldrin	ND	PCB-(Aroclor)-1016 ¹	ND
α -endosulfan	ND	toxaphene ¹	ND
β -endosulfan	ND		
endosulfan sulfate	ND		
endrin	ND		

Remarks: 0.5 = Quantitation limit

ND = Not detected

¹ = This component has a quantitation limit two (2) times that listed.

**Elevated quantitation limit listed in parenthesis due to sample matrix interferences.

Approved by *Alice R. Moore* Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 18 OF 25

Sample Description: WP-7 (Water)

PESTICIDES AND PCB's - PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
aldrin	ND (1.2)*	endrin aldehyde ¹	ND (2.4)***
α -BHC	ND	heptachlor ¹	ND (0.8)**
β -BHC	ND (1.9)*	heptachlor epoxide ¹	ND (0.5)***
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND (0.5)***
δ -BHC	ND	PCB-(Aroclor)-1254 ¹	ND (32)***
chlordane ¹	ND (3.7)***	PCB-(Aroclor)-1221 ¹	ND
4,4'-DDT	ND (2.1)***	PCB-(Aroclor)-1232 ¹	ND
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND (6.2)*
4,4'-DDD	ND	PCB-(Aroclor)-1260 ¹	ND (0.5)***
dieldrin	ND (0.8)***	PCB-(Aroclor)-1016 ¹	ND
α -endosulfan	ND	toxaphene ¹	ND (3.8)***
β -endosulfan	ND (0.8)***		
endosulfan sulfate	ND (1.4)***		
endrin	ND (2.1)***		

Remarks: 0.5 = Quantitation limit

ND = Not detected

¹ = This component has a quantitation limit two (2) times that listed.

[†]Sample exhibits alteration of standard Aroclor pattern.

*Elevated quantitation limit listed in parenthesis due to sample matrix interferences.

**Elevated detection limit listed in parenthesis due to presence of Aroclor.

Approved by *Alice S. Meier* Laboratory Manager

Title


CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
 ATTN: Bill Gray
 125 E. Bethpage Road
 Plainview, NY 11803

DATE REPORTED: November 18, 1988
 PROJECT CODE: GMIN 42163
 ORDER NUMBER:
 PAGE 19 OF 25

Sample Description: WP-8 (Water)

PESTICIDES AND PCB's - PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
aldrin	ND	endrin aldehyde ¹	ND
α -BHC	ND	heptachlor	ND
β -BHC	ND	heptachlor epoxide	ND
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND
δ -BHC	ND	PCB-(Aroclor)-1254 ¹	ND
chlordane ¹	ND	PCB-(Aroclor)-1221 ¹	ND
4,4'-DDT	ND	PCB-(Aroclor)-1232 ¹	ND
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND (1.8)**
4,4'-DDD	ND	PCB-(Aroclor)-1260 ¹	ND
dieldrin	ND	PCB-(Aroclor)-1016 ¹	ND
α -endosulfan	ND	toxaphene ¹	ND
γ -endosulfan	ND		
endosulfan sulfate	ND		
endrin	ND		

Remarks: 0.5 = Quantitation limit

ND = Not detected

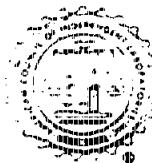
¹ = This component has a quantitation limit two (2) times that listed.

^tSample exhibits alteration of standard Aroclor pattern.

**Elevated detection limit listed in parenthesis due to presence of Aroclor.

Approved by Alice R. New Laboratory Manager

Title


CERTIFICATE OF ANALYSIS

TO **Geraghty & Miller, Inc.**
 ATTN: Bill Gray
 125 E. Bethpage Road
 Plainview, NY 11803

DATE REPORTED: November 18, 1988
 PROJECT CODE: GMIN 42163
 ORDER NUMBER:
 PAGE 20 OF 25

Sample Description: WP-9 (Water)

PESTICIDES AND PCB's - PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
aldrin	ND (1.1)*	endrin aldehyde ¹	ND (2.4)**
α -BHC	ND	heptachlor	ND (0.8)*
β -BHC	ND (1.8)*	heptachlor epoxide	ND
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND
δ -BHC	ND	PCB-(Aroclor)-1254 ¹	ND (39)**
chlordane ¹	ND (3.2)**	PCB-(Aroclor)-1221 ¹	ND
4,4'-DDT	ND (2.6)**	PCB-(Aroclor)-1232 ¹	ND (3.8)*
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND (4.3)*
4,4'-DDD	ND (1.3)**	PCB-(Aroclor)-1260 ¹	ND
dieldrin	ND (1.4)**	PCB-(Aroclor)-1016 ¹	ND (1.6)*
α -endosulfan	ND	toxaphene ¹	ND (3.8)**
β -endosulfan	ND (1.4)**		
endosulfan sulfate	ND (1.6)**		
endrin	ND (3.2)**		

Remarks: 0.5 = Quantitation limit

ND = Not detected

¹ = This component has a quantitation limit two (2) times that listed.

^tSample exhibits alteration of standard Aroclor pattern.

*Elevated quantitation limit listed in parenthesis due to sample matrix interferences.

**Elevated detection limit listed in parenthesis due to presence of Aroclor.

Ryan R. Meier
 Approved by Ryan R. Meier
 Laboratory Manager



ANALYTICAL SERVICES

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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 21 OF 25

Sample Description: 62 (Water)

PESTICIDES AND PCB'S - PRIORITY POLLUTANT ANALYSIS

Compound	Concentration (μ g/liter)	Compound	Concentration (μ g/liter)
aldrin	ND	endrin aldehyde ¹	ND
α -BHC	ND	heptachlor	ND
β -BHC	ND	heptachlor epoxide	ND
γ -BHC (lindane)	ND	PCB-(Aroclor)-1242 ¹	ND
δ -BHC	ND	PCB-(Aroclor)-1254 ¹	ND (2.0)**
chlor dane ¹	ND	PCB-(Aroclor)-1221 ¹	ND
4,4'-DDT	ND	PCB-(Aroclor)-1232 ¹	ND
4,4'-DDE	ND	PCB-(Aroclor)-1248 ¹	ND
4,4'-DDD	ND	PCB-(Aroclor)-1260 ¹	ND
dieldrin	ND	PCB-(Aroclor)-1016 ¹	ND
α -endosulfan	ND	toxaphene ¹	ND (1.4)**
β -endosulfan	ND		
endosulfan sulfate	ND		
endrin	ND		

Remarks: 0.5 = Quantitation limit

ND = Not detected

¹ = This component has a quantitation limit two (2) times that listed.

**Elevated quantitation limit listed in parenthesis due to sample matrix interferences.

Approved by Alice S. Rose
Laboratory Manager

Title

COMPOUND LIST

PESTICIDES, METHOD 8080
(Page 1)SAMPLE IDENTIFIER: ROA1G
COMPUCHEM® SAMPLE NUMBER: 468220

		CONCENTRATION (ug/L)	DETECTION LIMIT (ug/L)
1P.	4, 4'-DDD	BDL	0.10
2P.	4, 4'-DDE	BDL	0.10
3P.	4, 4'-DDT	BDL	0.10
4P.	ALDRIN	BDL	0.03
5P.	CHLORDANE	BDL	0.12
6P.	DIELDRIN	BDL	0.03
7P.	ENDOSULFAN I	BDL	0.05
8P.	ENDOSULFAN II	BDL	0.10
9P.	ENDOSULFAN SULFATE	BDL	0.05
10P.	ENDRIN	BDL	0.05
11P.	ENDRIN ALDEHYDE	BDL	0.03
12P.	HEPTACHLOR	BDL	0.03
13P.	HEPTACHLOR EPOXIDE	BDL	0.03
14P.	KEPONE	BDL	0.30
15P.	METHOXYCHLOR	BDL	0.30
16P.	PCB-1016	BDL	0.50
17P.	PCB-1221	BDL	0.50
18P.	PCB-1232	BDL	0.50
19P.	PCB-1242	BDL	0.50
20P.	PCB-1248	BDL	0.50
21P.	PCB-1254	BDL	0.50
22P.	PCB-1260	BDL	0.50
23P.	TOXAPHENE	BDL	1.00
24P.	ALPHA-BHC	BDL	0.03
25P.	BETA-BHC	BDL	0.03
26P.	DELTA-BHC	BDL	0.03
27P.	GAMMA-BHC	BDL	0.03

(Continued)

COMPOUND LIST

PESTICIDES, METHOD 8080
(Page 2)

SAMPLE IDENTIFIER: ROA1G
COMPUCHEM® SAMPLE NUMBER: 468220

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Dibutylchlorendate	93	(48-136)*

*Advisory surrogate; with the exception of dilutions recovery below 20% requires action step (re-extraction and re-analysis). See Quality Assurance Notice.

BDL=BELOW DETECTION LIMIT

COMPOUND LIST

PESTICIDES, METHOD 8080

(Page 1)

SAMPLE IDENTIFIER: ROC1G
 COMPUCHEM® SAMPLE NUMBER: 468518

	CONCENTRATION (ug/L)	DETECTION LIMIT (ug/L)
1P. 4,4'-DDD	BDL	0.10
2P. 4,4'-DDE	BDL	0.10
3P. 4,4'-DDT	BDL	0.10
4P. ALDRIN	BDL	0.03
5P. CHLORDANE	BDL	0.12
6P. DIELDRIN	BDL	0.03
7P. ENDOSULFAN I	BDL	0.05
8P. ENDOSULFAN II	BDL	0.10
9P. ENDOSULFAN SULFATE	BDL	0.05
10P. ENDRIN	BDL	0.05
11P. ENDRIN ALDEHYDE	BDL	0.03
12P. HEPTACHLOR	BDL	0.03
13P. HEPTACHLOR EPOXIDE	BDL	0.03
14P. KEPONE	BDL	0.30
15P. METHOXYCHLOR	BDL	0.30
16P. PCB-1016	BDL	0.50
17P. PCB-1221	BDL	0.50
18P. PCB-1232	BDL	0.50
19P. PCB-1242	BDL	0.50
20P. PCB-1248	BDL	0.50
21P. PCB-1254	BDL	0.50
22P. PCB-1260	BDL	0.50
23P. TOXAPHENE	BDL	1.00
24P. ALPHA-BHC	BDL	0.03
25P. BETA-BHC	BDL	0.03
26P. DELTA-BHC	BDL	0.03
27P. GAMMA-BHC	BDL	0.03

(Continued)

COMPOUND LIST

PESTICIDES, METHOD 8080
(Page 2)

SAMPLE IDENTIFIER: ROC1G
COMPUCHEM® SAMPLE NUMBER: 468518

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Dibutylchlorendate	93	(48-136)*

*Advisory surrogate; with the exception of dilutions recovery below 20% requires action step (re-extraction and re-analysis). See Quality Assurance Notice.

BDL=BELOW DETECTION LIMIT

PCB ANALYSIS (GROUNDWATER)

ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3

ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2

ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2

COMPOUND LIST
APPENDIX VIII, IX - PCBs, METHOD 8080

SAMPLE IDENTIFIER: ROA3G
COMPUCHEM® SAMPLE NUMBER: 474982

	CONCENTRATION ($\mu\text{g/L}$)	DETECTION LIMIT ($\mu\text{g/L}$)
1P. PCB-1016	BDL	0.53
2P. PCB-1221	BDL	0.53
3P. PCB-1232	BDL	0.53
4P. PCB-1242	BDL	0.53
5P. PCB-1248	BDL	0.53
6P. PCB-1254	BDL	0.53
7P. PCB-1260	BDL	0.53

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	% Recovery	Control Range %
Dibutylchlorendate	106	(48-136)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires action step (re-extraction and re-analysis).

BDL=BELOW DETECTION LIMIT

†Detection limits have been adjusted to report variation from the nominal sample weight.

COMPOUND LIST
APPENDIX VIII, IX - PCBs, METHOD 8080

SAMPLE IDENTIFIER: ROB2G
COMPUCHEM® SAMPLE NUMBER: 468172

	CONCENTRATION (μ g/L)	DETECTION LIMIT (μ g/L)
1P. PCB-1016	BDL	0.50
2P. PCB-1221	BDL	0.50
3P. PCB-1232	BDL	0.50
4P. PCB-1242	BDL	0.50
5P. PCB-1248	BDL	0.50
6P. PCB-1254	BDL	0.50
7P. PCB-1260	BDL	0.50

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	% Recovery	Control Range %
Dibutylchlorendate	95	(48-136)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires action step (re-extraction and re-analysis).

BDL=BELOW DETECTION LIMIT

COMPOUND LIST
APPENDIX VIII, IX - PCBs, METHOD 8080

SAMPLE IDENTIFIER: ROC2G
COMPUCHEM® SAMPLE NUMBER: 468591

	CONCENTRATION (μ g/L)	DETECTION LIMIT (μ g/L)
1P. PCB-1016	BDL	0.50
2P. PCB-1221	BDL	0.50
3P. PCB-1232	BDL	0.50
4P. PCB-1242	BDL	0.50
5P. PCB-1248	BDL	0.50
6P. PCB-1254	BDL	0.50
7P. PCB-1260	BDL	0.50

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	% Recovery	Control Range %
Dibutylchlorendate	93	(48-136)*

*Advisory surrogate; with the exception of dilutions recovery below 10% requires action step (re-extraction and re-analysis).

BDL=BELOW DETECTION LIMIT

COMPOUND LIST

PESTICIDES, METHOD 8080
(Page 2)

SAMPLE IDENTIFIER: ROJ1G
COMPUCHEM® SAMPLE NUMBER: 472797

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Dibutylchlorendate	122	(48-136)*

*Advisory surrogate; with the exception of dilutions recovery below 20% requires action step (re-extraction and re-analysis).

BDL=BELOW DETECTION LIMIT



Section 14

ORGANOPHOSPHORUS PESTICIDES ANALYSIS (GROUNDWATER)

ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1

COMPOUND LIST

APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140

SAMPLE IDENTIFIER: ROA1G
 COMPUCHEM® SAMPLE NUMBER: 468222

	<u>CONCENTRATION</u> <u>(ug/L)</u>	<u>DETECTION</u> <u>LIMIT</u> <u>(ug/L)</u>
1P. DIMETHOATE	BDL	0.50
2P. DISULFOTON	BDL	0.50
3P. METHYL PARATHION	BDL	0.50
4P. PARATHION	BDL	0.50
5P. PHORATE	BDL	0.50
6P. TETRAETHYLDITHIOPYROPHOSPHATE (SULFOTEPP)	BDL	0.50

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Methidathion	88	(60-120)*

*Advisory surrogate only.

BDL=BELOW DETECTION LIMIT

COMPOUND LIST

APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140

SAMPLE IDENTIFIER: ROB1G
 COMPUCHEM® SAMPLE NUMBER: 468221

	<u>CONCENTRATION</u> <u>(ug/L)</u>	<u>DETECTION</u> <u>LIMIT</u> <u>(ug/L)</u>
1P. DIMETHOATE	BDL	0.50
2P. DISULFOTON	BDL	0.50
3P. METHYL PARATHION	BDL	0.50
4P. PARATHION	BDL	0.50
5P. PHORATE	BDL	0.50
6P. TETRAETHYLDITHIOPYROPHOSPHATE (SULFOTEPP)	BDL	0.50

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Methidathion	99	(60-120)*

*Advisory surrogate only.

BDL=BELOW DETECTION LIMIT

COMPOUND LIST

APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140

SAMPLE IDENTIFIER: ROC1G
 COMPUCHEM® SAMPLE NUMBER: 468517

	<u>CONCENTRATION</u> (<u>ug/L</u>)	<u>DETECTION</u> <u>LIMIT</u> (<u>ug/L</u>)
1P. DIMETHOATE	BDL	0.50
2P. DISULFOTON	BDL	0.50
3P. METHYL PARATHION	BDL	0.50
4P. PARATHION	BDL	0.50
5P. PHORATE	BDL	0.50
6P. TETRAETHYLDITHIOPYROPHOSPHATE (SULFOTEPP)	BDL	0.50

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Methidathion	74	(60-120)*

*Advisory surrogate only.

BDL=BELOW DETECTION LIMIT

COMPOUND LIST

APPENDIX VIII, IX - ORGANOPHOSPHORUS PESTICIDES, METHOD 8140

SAMPLE IDENTIFIER: ROJ1G
 COMPUCHEM® SAMPLE NUMBER: 472798

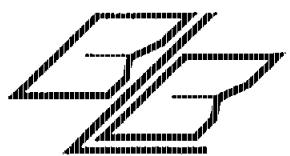
	<u>CONCENTRATION</u> <u>(ug/L)</u>	<u>DETECTION</u> <u>LIMIT</u> <u>(ug/L)</u>
1P. DIMETHOATE	BDL	0.50
2P. DISULFOTON	BDL	0.50
3P. METHYL PARATHION	BDL	0.50
4P. PARATHION	BDL	0.50
5P. PHORATE	BDL	0.50
6P. TETRAETHYLDITHIOPYROPHOSPHATE (SULFOTEPP)	BDL	0.50

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
Methidathion	111	(60-120)*

*Advisory surrogate only.

BDL=BELOW DETECTION LIMIT



Section 15

HERBICIDES ANALYSIS (GROUNDWATER)

ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1

COMPOUND LIST

APPENDIX VIII, IX - HERBICIDES, METHOD 8150

SAMPLE IDENTIFIER: ROA1G
COMPUCHEM® SAMPLE NUMBER: 468224

	CONCENTRATION (ug/L)	DETECTION LIMIT (ug/L)
1. 2,4-D	BDL	4.0
2. 2,4,5-TP(Silvex)	BDL	1.0
3. 2,4,5-T	BDL	1.0

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

<u>% Recovery</u>	<u>Control Range %</u>
73	(24-154)*

BDL=BELOW DETECTION LIMIT

*Advisory surrogate only; if recovery < 20%, re-methylation extract; minimum 10% recovery required after re-methylation. See Quality Assurance Notice.

COMPOUND LIST

APPENDIX VIII, IX - HERBICIDES, METHOD 8150

SAMPLE IDENTIFIER: ROBIG
COMPUCHEM® SAMPLE NUMBER: 468223

	CONCENTRATION (<u>ug/L</u>)	DETECTION LIMIT (<u>ug/L</u>)
1. 2,4-D	BDL	4.0
2. 2,4,5-TP(Silvex)	BDL	1.0
3. 2,4,S-T	BDL	1.0

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	<u>% Recovery</u>	<u>Control Range %</u>
2,4-DB	58	(24-154)*

BDL=BELOW DETECTION LIMIT

*Advisory surrogate only; if recovery < 20%, re-methylation extract; minimum 10% recovery required after re-methylation. See Quality Assurance Notice.

COMPOUND LIST

APPENDIX VIII, IX - HERBICIDES, METHOD 8150

SAMPLE IDENTIFIER: ROC1G
COMPUCHEM® SAMPLE NUMBER: 468525

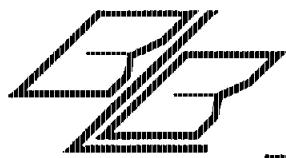
	CONCENTRATION (ug/L)	DETECTION LIMIT (ug/L)
1. 2,4-D	BDL	4.0
2. 2,4,5-TP(Silvex)	BDL	1.0
3. 2,4,5-T	BDL	1.0

Surrogate Recovery - Introduced at the beginning of the extraction, the surrogate standard is a select compound that analytically mimics the response of certain analytes. A known concentration of this surrogate is added to the sample and a percent recovery is calculated. This recovery acts as a barometer of extraction efficiency and analytical response for the individual sample.

	% Recovery	Control Range %
2,4-DB	53	(24-154)*

BDL=BELOW DETECTION LIMIT

*Advisory surrogate only; if recovery < 20%, re-methylation extract; minimum 10% recovery required after re-methylation. See Quality Assurance Notice.



Section 16

DIOXIN / FURAN ANALYSIS (GROUNDWATER)

ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1

FORM 1 - QUANTITATION REPORT

PAGE 2 of 2

DATE: 01/06/92

LABORATORY: ChemWest

Ticket# CW-9068

Project Name: General Electric Company

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS						SURROGATE % ACCURACY			
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C1-TCDD	*C-HxCDD	*C-HpCDF
Method Blank Detection Limit	9068-1MB	12/31/91	12:10	CW-1	79.9	83.8	79.7	63.6	37.7	82.2	84.2	98.6	98.5	102
ROFIG // 468214 Detection Limit	9068-1	12/31/91	12:46	CW-1	82.1	82.7	80.0	68.1	39.7	80.9	83.2	99.7	99.5	95.6
ROB1G // 468215 Detection Limit	9068-2	12/31/91	13:23	CW-1	82.4	79.1	72.2	58.4	35.0	82.3	83.8	99.6	98.9	100
ROB2G // 468216 Detection Limit	9068-3	12/31/91	14:01	CW-1	74.1	76.0	70.7	56.6	32.1	76.5	79.4	101	97.5	101
ROA1G // 468217 Detection Limit	9068-4	12/31/91	14:40	CW-1	83.4	84.6	77.7	63.4	40.0	80.4	87.3	99.9	100	97.7
RODP1G // 468218	9068-5	12/31/91	15:20	CW-1	73.6	76.8	69.7	50.4	27.3	74.9	78.5	99.5	95.9	97.1

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HpCDF = 13C12-1234678-HpCDF

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

Ticket# CW-9187

Project Name: General Electric Company

DATE: 02/06/92

LABORATORY: ChemWest

TOTAL ANALYTE QUANTITY FOUND

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	2378											
					TCDD	TCDF	PeCDD	HxCDD	HxCDF	OCDD	OCDF	TCDF	PeCDF	HxCDF	HxCDF	OCDF
Method Blank	9187-MBRX	02/04/92	17:24	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit					1.6	1.6	2.4	3.9	2.4	5.9	0.64	4.1	1.1	2.0	3.2	5.1
ROA3G // 474990	9187-RX	02/04/92	18:08	CW-2	ND	ND	ND	ND	ND	ND	ND	ND	ND	9.2	ND	ND
Detection Limit					1.5	1.5	2.5	4.3	3.3	5.0	3.7	12.9	1.8	5.3	6.8	

@ = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTIFICATION REPORT

PAGE 2 of 2

DATE: 02/06/92

LABORATORY: ChemWest

Ticket# CW-9187

Project Name: General Electric Company

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	ABSOLUTE % RECOVERY of INTERNAL STANDARDS							SURROGATE % ACCURACY		
					*C-TCDD	*C-PeCDD	*C-HxCDD	*C-HpCDD	*C-OCDD	*C-TCDF	*C-PeCDF	*C1-TCDD	*C-HxCDD	*C-HpCDF
Method Blank	9187-MBRX	02/04/92	17:24	CW-2	78.6	81.0	81.5	67.0	46.2	81.8	80.4	98.3	98.6	96.5
Detection Limit												99.6	97.0	93.1
ROA3G // 474990	9187-RX	02/04/92	18:08	CW-2	43.8	62.9	67.0	59.6	43.4	28.3	60.3			
Detection Limit														

INTERNAL STANDARDS

*C-TCDD = 13C12-2378-TCDD
 *C-PeCDD = 13C12-12378-PeCDD
 *C-HxCDD = 13C12-123678-HxCDD
 *C-HpCDD = 13C12-1234678-HpCDD
 *C-TCDF = 13C12-2378-TCDF

SURROGATES

*C1-TCDD = 37CL4-2378-TCDD
 *C-HxCDD = 13C12-123789-HxCDD
 *C-PeCDF = 13C12-12378-PeCDF
 *C-HPCDF = 13C12-1234678-HpCDF

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 01/16/92

LABORATORY: ChemWest

Ticket# CW-9092

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppt or ng/L)

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	2378		2378							
					TCDD	TCDF	PeCDD	HxCDD	HxCDF	OCDD	TCDF	PeCDF	HxCDF	HxCDF
Method Blank	9077-1MB	12/31/91	15:59	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit					0.18	0.18	0.26	0.27	0.45	0.51	0.10	0.36	0.25	0.26
ROC1G // 469599	9092	01/02/92	11:46	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit					0.17	0.17	0.34	0.58	0.70	1.1	0.23	0.34	0.21	0.40

ND = MAXIMUM POSSIBLE CONCENTRATION

#C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

#C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

#C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 01/09/92

LABORATORY: ChemWest

Ticket# CW-9079

Project Name: General Electric Company

TOTAL ANALYTE QUANTITY FOUND

(ppt or ng/L)

CLIENT ID.	CW#	GC/MS DATE	GC/MS TIME	INST. ID.	2378	2378											
						TCDD	TCDD	PeCDD	HxCDD	HxCDD	OCDD	TCDF	TCDF	PeCDF	HxCDF	HxCDF	OCDF
Method Blank	9077-1MB	12/31/91	15:59	CW-1		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.18	0.18	0.26	0.27	0.45	0.51	0.10	0.36	0.25	0.26	0.43	0.48
ROEB1 // 468581	9077-2	01/02/92	09:56	CW-1		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.15	0.15	0.23	0.50	0.45	0.68	0.084	0.36	0.21	0.31	0.36	0.67
ROC2G // 468599	9079-3	01/02/92	10:32	CW-1		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.22	0.22	0.40	0.39	0.74	0.78	0.20	0.31	0.24	0.30	1.3	0.54
ROE1G // 468619	9079-4	01/02/92	11:09	CW-1		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Detection Limit						0.26	0.26	0.25	0.58	0.51	0.74	0.10	0.38	0.30	0.34	0.35	0.48

* = MAXIMUM POSSIBLE CONCENTRATION

*C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

*C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

*C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: 

FORM 1 - QUANTITATION REPORT

PAGE 1 of 2

DATE: 01/16/92

LABORATORY: ChemWest

Ticket# CW-9136

Project Name: General Electric Company

CLIENT ID.	CW#	TOTAL ANALYTE QUANTITY FOUND (ppt or ng/L)													
		GC/MS DATE	GC/MS TIME	INST. ID.	2378		2378		OCDD	TCDF	TCDF	PeCDF	HxCDF	HpCDF	OCDF
					TCDD	TCDD	PeCDD	HxCDD							
Method Blank Detection Limit	9136-1MB	01/02/92	20:44	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
					0.25	0.25	0.48	0.40	0.74	0.69	0.10	0.41	0.29	0.22	0.45
RN10G // 472740 Detection Limit	9136-1	01/03/92	09:56	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
					0.16	0.16	0.35	0.52	0.63	0.63	0.15	0.32	0.27	0.81	1.5
RNDPG // 472742 Detection Limit	9136-2	01/03/92	10:34	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
					0.36	0.36	0.53	0.52	0.63	0.89	0.15	0.49	0.32	0.29	0.55
RN09G // 472744 Detection Limit	9136-3	01/03/92	11:11	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
					0.51	0.51	0.78	0.60	1.1	1.0	0.31	0.63	0.74	0.44	0.70
RN11G // 472746 Detection Limit	9136-4	01/03/92	11:50	CW-1	ND	ND	ND	ND	ND	4.1	ND	ND	ND	ND	ND
					0.28	0.28	0.27	0.56	1.3		0.35	0.58	0.70	1.4	1.1
ROJ1G // 472748 Detection Limit	9136-5	01/03/92	12:35	CW-1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
					0.33	0.33	0.62	0.53	0.63	0.75	0.14	0.43	0.29	0.32	0.52
															0.79

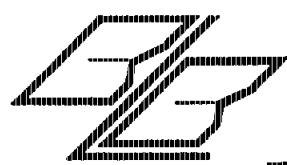
a = MAXIMUM POSSIBLE CONCENTRATION

#C-TCDD: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzodioxin (12 carbons)

#C-TCDF: Carbon 13 labeled 2,3,7,8-tetrachlorodibenzofuran (12 carbons)

#C-OCDD: Carbon 13 labeled octachlorodibenzodioxin (12 carbons)

Approved by: _____



Section 17

METALS ANALYSIS (GROUNDWATER)

WP-1 - Groundwater Sample from Oxbow Area K, Well Point 1
WP-2 - Groundwater Sample from Oxbow Area K, Well Point 2
WP-3 - Groundwater Sample from Oxbow Area J, Well Point 3
WP-7 - Groundwater Sample from Oxbow Area C, Well Point 7
WP-8 - Groundwater Sample from Oxbow Area C, Well Point 8
WP-9 - Groundwater Sample from Oxbow Area A, Well Point 9
ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 18 OF 19

Sample Description: Five (5) water samples received October 14, 1988

Concentration units are mg/liter (ppm)

PRIORITY POLLUTANT METALS

	WP-1	WP-2	WP-3	WP-5	WP-6
Antimony	<0.03	<0.03	<0.03	<0.03	<0.03
Arsenic	<0.03	<0.03	<0.03	<0.03	<0.03
Beryllium	<0.001	<0.001	<0.001	<0.001	<0.001
Cadmium	<0.005	<0.005	<0.005	<0.005	<0.005
Chromium	<0.01	<0.01	<0.01	<0.01	<0.01
Copper	<0.01	0.01	0.01	0.03	0.01
Lead	0.04	<0.03	<0.03	<0.03	<0.03
Mercury	<0.001	<0.001	<0.001	<0.001	<0.001
Nickel	<0.02	<0.02	<0.02	<0.02	<0.02
Selenium	<0.06	<0.06	<0.06	<0.06	<0.06
Silver	<0.005	<0.005	<0.005	<0.005	<0.005
Thallium	<0.03	<0.03	<0.05*	<0.03	<0.05*
Zinc	2.5	1.3	5.2	3.0	5.3

*Detection limits higher than normal due to sample matrix interferences.

Approved by *Allyn S. Russell*
Laboratory Manager

Title



ANALYTICAL SERVICES



5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401

CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 22 OF 25

Sample Description: Two (2) water samples received October 20, 1988

Concentration units are mg/liter (ppm)

PRIORITY POLLUTANT METALS

	<u>WP-4</u>	<u>WP-7</u>
Antimony	<0.03	<0.03
Arsenic	<0.03	<0.03
Beryllium	<0.001	<0.001
Cadmium	<0.005	<0.005
Chromium	<0.01	<0.01
Copper	<0.01	0.01
Lead	<0.03	<0.03
Mercury	<0.001	<0.001
Nickel	<0.02	<0.02
Selenium	<0.06	<0.06
Silver	<0.005	<0.005
Thallium	<0.04	<0.04
Zinc	0.93	0.36

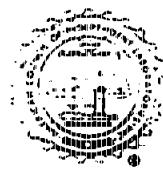
Allyn R. Moore
Approved by Laboratory Manager

Title



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER: 23 _____ OF 25
PAGE _____

Sample Description: Two (2) water samples received October 20, 1988

Concentration units are mg/liter (ppm)

PRIORITY POLLUTANT METALS

	<u>WP-8</u>	<u>WP-9</u>
Antimony	<0.03	<0.03
Arsenic	<0.03	<0.03
Beryllium	<0.001	<0.001
Cadmium	<0.005	<0.005
Chromium	<0.01	<0.01
Copper	<0.01	<0.01
Lead	<0.03	<0.03
Mercury	<0.001	<0.001
Nickel	<0.02	<0.02
Selenium	<0.06	<0.06
Silver	0.005 -	<0.005
Thallium	<0.04	<0.04
Zinc	0.42 -	0.36 -

Approved by

Ryan S. Meier Laboratory Manager

Title

**CERTIFICATE OF ANALYSIS**

TO **Geraghty & Miller, Inc.**
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 24 OF 25

Sample Description: One (1) water sample received October 20, 1988

Concentration units are mg/liter (ppm)

PRIORITY POLLUTANT METALS62

Antimony	<0.03
Arsenic	<0.03
Beryllium	<0.001
Cadmium	<0.005
Chromium	<0.01
Copper	<0.01
Lead	<0.03
Mercury	<0.001
Nickel	0.02
Selenium	<0.06
Silver	0.006
Thallium	<0.04
Zinc	0.044

Approved by Alice S. Moore Laboratory Manager

Title

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROA1G

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 93629AMatrix (soil/water): WATER Lab Sample ID: 468181Level (low/med): LOW Date Received: 12/06/91± Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	105	B		P
7440-36-0	Antimony	36.0	U		P
7440-38-2	Arsenic	5.0	U		F
7440-39-3	Barium	125	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	5.0	U		P
7440-79-2	Calcium	79400			P
7440-47-3	Chromium	4.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	6.0	U		P
7439-89-6	Iron	7290			P
7439-92-1	Lead	2.0	U		F
7439-95-4	Magnesium	25800			P
7439-96-5	Manganese	730			P
7439-97-6	Mercury	.20	U		CV
7440-92-0	Nickel	8.0	U		P
7440-99-7	Potassium	6710			P
7782-49-2	Selenium	3.0	U		F
7440-22-4	Silver	6.0	U		P
7440-23-5	Sodium	52300			P
7440-28-0	Thallium	2.0	U	WN	F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	35.8			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

FORM 1.04 - PAGE 1

U.S. EPA - SW-846

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

ROA3G

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846

Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 936308

Matrix (soil/water): WATER Lab Sample ID: 474994

Level (low/med): LOW Date Received: 01/15/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	24900		P	
7440-36-0	Antimony	60.0	U		P
7440-38-2	Arsenic	35.1	*		P
7440-39-3	Barium	154	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	5.0	U		P
7440-70-7	Calcium	108000			P
7440-47-3	Chromium	32.2			P
7440-48-4	Cobalt	10.3	B		P
7440-50-8	Copper	112			P
7439-89-6	Iron	56900			P
7439-91-1	Lead	112			P
7439-95-4	Magnesium	50500			P
7439-96-5	Manganese	3070			P
7439-97-6	Mercury	.20	U		CV
7440-02-0	Nickel	46.9			P
7440-09-7	Potassium	11800			P
7782-49-2	Selenium	4.0	U	QN	P
7440-22-4	Silver	10.0	U		P
7440-23-5	Sodium	222000			P
7440-78-0	Thallium	3.0	U	WN	P
7440-82-2	Vanadium	34.3	B		P
7440-66-6	Zinc	181			P
	Cyanide				NR

Color Before: BROWN Clarity Before: CLOUDY Texture: _____Color After: YELLOW Clarity After: CLEAR Artifacts: _____

Comments:

FORM 1.04 - PAGE 1PLEASE REFERENCE ENCLOSED NOTICE REGARDING "Q" FLAG IN COLUMN Q

FORM I - IN

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846ROB1GLab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 93629AMatrix (soil/water): WATER Lab Sample ID: 468178Level (low/med): LOW Date Received: 12/06/91t Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	103	B		P
7440-36-0	Antimony	36.0	U		P
7440-38-2	Arsenic	5.0	U W		F
7440-39-3	Barium	75.0	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	5.0	U		P
7440-70-2	Calcium	86100			P
7440-47-3	Chromium	4.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	6.0	U		P
7439-89-6	Iron	3700			P
7439-92-1	Lead	3.2			F
7439-95-4	Magnesium	30300			P
7439-96-5	Manganese	419			P
7439-97-6	Mercury	.20	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	2980	B		P
7782-49-2	Selenium	3.0	U W		F
7440-22-4	Silver	6.0	U		P
7440-23-5	Sodium	26700			P
7440-28-0	Thallium	2.0	U WN		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	28.3			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846

ROB2G

Lab Code: COMPU Case No.: 59007 SAS No.: _____ SDG No.: 93629AMatrix (soil/water): WATER Lab Sample ID: 468180Level (low/med): LOW Date Received: 12/06/91t Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	143	B		P
7440-36-0	Antimony	36.0	U		P
7440-38-2	Arsenic	5.0	U		F
7440-39-3	Barium	12.2	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-2	Cadmium	5.0	U		P
7440-70-2	Calcium	68700			P
7440-47-3	Chromium	4.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	6.0	U		P
7439-89-6	Iron	3010			P
7439-92-1	Lead	2.0	U		P
7439-95-4	Magnesium	24800			P
7439-96-5	Manganese	1670			P
7439-97-6	Mercury	20	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	1830	B		P
7782-49-2	Selenium	3.0	U		F
7440-22-4	Silver	6.0	U		P
7440-23-5	Sodium	35400			P
7440-28-0	Thallium	2.0	U	WN	F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	37.9			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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FORM I - IN

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846ROC1GLab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 93629AMatrix (soil/water): WATER Lab Sample ID: 468527Level (low/med): LOW Date Received: 12/07/91t Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	2520		P	
7440-16-0	Antimony	36.0	U	P	
7440-18-2	Arsenic	5.0	U	W	F
7440-39-3	Barium	79.0	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	5.0	U		P
7440-70-2	Calcium	135000		P	
7440-47-3	Chromium	4.8	B	P	
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	10.1	B		P
7439-89-6	Iron	14500		P	
7439-92-1	Lead	11.9			F
7432-95-4	Magnesium	28000		P	
7432-96-5	Manganese	2390		P	
7432-97-6	Mercury	.20	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5990		P	
7782-42-2	Selenium	3.0	U		F
7440-22-4	Silver	6.0	U		P
7440-23-5	Sodium	42700		P	
7440-28-0	Thallium	2.0	U	WN	F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	24.5			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: SW-846Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 93629AMatrix (soil/water): WATER Lab Sample ID: 468592Level (low/med): LOW Date Received: 12/07/91t Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	231			P
7440-36-0	Antimony	36.0	U		P
7440-38-2	Arsenic	5.0	U		F
7440-39-3	Barium	94.9	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	5.0	U		P
7440-70-2	Calcium	71800			P
7440-47-3	Chromium	4.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	6.0	U		P
7439-89-6	Iron	2210			P
7439-92-1	Lead	2.0	U		F
7439-95-4	Magnesium	13600			P
7439-96-5	Manganese	602			P
7439-97-6	Mercury	.20	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	7010			P
7782-49-2	Selenium	3.0	U		F
7440-22-4	Silver	6.0	U		P
7440-23-5	Sodium	41500			P
7440-28-0	Thallium	2.0	U	NN	F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	35.1			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIESContract: 7/88ROJIGLab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 936302Matrix (soil/water): WATERLab Sample ID: 472802Level (low/med): LOWDate Received: 12/20/91% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	17300	N*	P	
7440-36-0	Antimony	60.0	U	P	
7440-38-2	Arsenic	7.4	B W	F	
7440-39-3	Barium	120	B	P	
7440-41-7	Beryllium	1.0	U	P	
7440-43-9	Cadmium	5.0	U	P	
7440-70-2	Calcium	61400		P	
7440-47-3	Chromium	21.1	*	P	
7440-48-4	Cobalt	14.2	B	P	
7440-50-8	Copper	35.0		P	
7439-89-6	Iron	29100		P	
7439-92-1	Lead	12.6	N*	F	
7439-95-4	Magnesium	32700		P	
7439-96-5	Manganese	3320		F	
7439-97-6	Mercury	.20	U N	CV	
7440-02-0	Nickel	27.9	B	F	
7440-09-7	Potassium	5590		P	
7782-49-2	Selenium	4.0	U W	F	
7440-22-4	Silver	10.0	U	P	
7440-23-5	Sodium	16700	E	P	
7440-28-0	Thallium	3.0	U WN	F	
7440-62-2	Vanadium	19.7	B	P	
7440-65-6	Zinc	119		P	
	Cyanide				NR

Color Before: BROWNClarity Before: CLOUDY

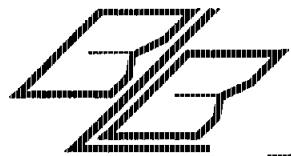
Texture: _____

Color After: YELLOWClarity After: CLEAR

Artifacts: _____

Comments:

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Section 18

CYANIDES ANALYSIS (GROUNDWATER)

WP-1 - Groundwater sample from Oxbow Area K, Well Point 1
WP-2 - Groundwater sample from Oxbow Area K, Well Point 2
WP-3 - Groundwater sample from Oxbow Area J, Well Point 3
WP-7 - Groundwater sample from Oxbow Area C, Well Point 7
WP-8 - Groundwater sample from Oxbow Area C, Well Point 8
WP-9 - Groundwater sample from Oxbow Area A, Well Point 9
ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 19 OF 19

Sample Description: Five (5) water samples received October 14, 1988

Concentration units are mg/liter (ppm)

	<u>Phenols</u>	<u>Cyanide</u>
WP-1	<0.01	<0.01
WP-2	<0.01	<0.01
WP-3	<0.01	<0.01
WP-5	<0.01	<0.01
WP-6	0.04	<0.01

Roger L. Moore
Approved by **Laboratory Manager**

Title



ANALYTICAL SERVICES



5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401

CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 25 OF 25

Sample Description: Five (5) water samples received October 20, 1988

Concentration units are mg/liter (ppm)

	<u>Cyanide</u>	<u>Phenols</u>
WP-4	<0.01	<0.01
WP-7	<0.01	<0.01
WP-8	<0.01	0.03
WP-9	<0.01	0.02
62	<0.01	0.05

Alice A. Koen
Approved by _____
Laboratory Manager

Title _____

1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIESContract: 7/88ROA1GLab Code: COMPUCase No.: 50007

SAS No.: _____

SDG No.: 75298AMatrix (soil/water): WATERLab Sample ID: 468188Level (low/med): LOWDate Received: 12/06/91% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
.	Cyanide	10.0	U		AS

Color Before: COLORLESSClarity Before: CLEAR

Texture: _____

Color After: COLORLESSClarity After: CLEAR

Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 3/90ROA3GLab Code: COMPU Case No.: 10510 SAS No.: _____ SDG No.: 475300Matrix (soil/water): WATER Lab Sample ID: 474986Level (low/med): LOW Date Received: 01/15/92% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	10.0	U		AS

Color Before: BROWN Clarity Before: CLOUDY Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88
 Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 75298A
 Matrix (soil/water): WATER Lab Sample ID: 468186
 Level (low/med): LOW Date Received: 12/06/91
 % Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	10.0	U		AS

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88 ROB2G

Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 75298A

Matrix (soil/water): WATER Lab Sample ID: 468187

Level (low/med): LOW Date Received: 12/06/91

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	10.0	U		AS

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88 ROC1G

Lab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 75298A

Matrix (soil/water): WATER Lab Sample ID: 468534

Level (low/med): LOW Date Received: 12/07/91

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	10.0	U		AS

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM LABORATORIES Contract: 7/88 ROC2GLab Code: COMPU Case No.: 50007 SAS No.: _____ SDG No.: 75298AMatrix (soil/water): WATER Lab Sample ID: 468593Level (low/med): LOW Date Received: 12/07/91% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	10.0	U		AS

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

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1
INORGANIC ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

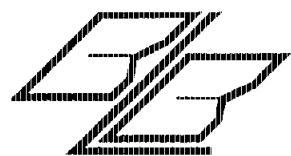
Lab Name: COMPUCHEM LABORATORIES Contract: 7/88 ROJ1GLab Code: COMPU Case No.: 33090 SAS No.: _____ SDG No.: 75299BMatrix (soil/water): WATER Lab Sample ID: 472804Level (low/med): LOW Date Received: 12/20/91% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum				NR
7440-36-0	Antimony				NR
7440-38-2	Arsenic				NR
7440-39-3	Barium				NR
7440-41-7	Beryllium				NR
7440-43-9	Cadmium				NR
7440-70-2	Calcium				NR
7440-47-3	Chromium				NR
7440-48-4	Cobalt				NR
7440-50-8	Copper				NR
7439-89-6	Iron				NR
7439-92-1	Lead				NR
7439-95-4	Magnesium				NR
7439-96-5	Manganese				NR
7439-97-6	Mercury				NR
7440-02-0	Nickel				NR
7440-09-7	Potassium				NR
7782-49-2	Selenium				NR
7440-22-4	Silver				NR
7440-23-5	Sodium				NR
7440-28-0	Thallium				NR
7440-62-2	Vanadium				NR
7440-66-6	Zinc				NR
	Cyanide	10.0	U		AS

Color Before: BROWN Clarity Before: CLOUDY Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

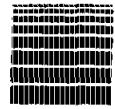
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Section 19

SULFIDES ANALYSIS (GROUNDWATER)

ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1



COMPUCHEM
LABORATORIES, INC.

SAMPLE CORRELATION SHEET

<u>CHEMWEST ID</u>	<u>COMPUCHEM ID</u>	<u>CLIENT ID</u>
9068-1	468208	ROFIG
9068-2	468210	ROB1G
9068-3	468211	ROB2G
9068-4	468212	ROA1G
9068-5	468213	RODPIG

SAMPLE CORRELATION SHEET

CHEMWEST ID

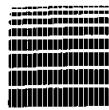
9187-1

COMPUCHEM ID

474993

CLIENT ID

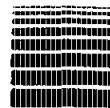
ROA3G



COMPUCHEM
LABORATORIES, INC.

SAMPLE CORRELATION SHEET

<u>CHEMWEST ID</u>	<u>COMPUCHEM ID</u>	<u>CLIENT ID</u>
9079-1	468548	ROC1G
9079-2	468584	ROEB1
9079-3	468603	ROC2G
9079-4	468620	ROE1G



COMPUCHEM
LABORATORIES, INC.

SAMPLE CORRELATION SHEET

CHEMWEST I.D.

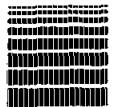
9136-1
9136-2
9136-3
9136-4
9136-5

COMPUCHEM I.D.

472739
472741
472743
472745
472747

CLIENT I.D.

RN10G
RNDPG
RN09G
RN11G
ROJ1G



COMPUCHEM
LABORATORIES, INC.

SULFIDE

Date(s) Analyzed: 12/11/91

Case: 9068

Matrix: Water

Client ID	CHEMWEST ID	Amount Detected (MG/L)
468208	9068-1	3.9
468210	9068-2	BRL
468211	9068-3	BRL
468212	9068-4	BRL
468213	9068-5	4.0

Client ID	CHEMWEST ID	Spike Conc. (MG/L)	Amount Detected (MG/L)	% Rec.
Method Blank	MB		BRL	
MBS	MBS	4.0	3.5680	89.2%
MBSD	MBSD	4.0	3.5360	88.4%

Relative % Difference = 0.9%

The reporting limit for Sulfide is 1.0 mg/L.

BRL: Below Reporting Limit.

Approved by: V.H.

Date Reported:
12/17/91

REV5:12.91

CHEMWEST ANALYTICAL LABORATORIES
SULFIDE
EPA METHOD 376.1

Date(s) Analyzed: 01/21/92

Case: 9187 PG1
Matrix: Water

Client ID	CHEMWEST ID	Amount Detected (MG/L)	RL (MG/L)
474993	9187-1	3.3	1.0

Client ID	CHEMWEST ID	Spike Conc. (MG/L)	Amount Detected (MG/L)	% Rec.
Method Blank	MB		BRL	
LQCS	LQCS	5.0	5.0960	101.9%
LQCSD	LQCSD	5.0	5.1360	102.7%

Relative % Difference = 0.8%

Client ID	CHEMWEST ID	Spike Conc. (MG/L)	Amount Detected (MG/L)	% Rec.
Matrix Spike	9187-1MS	5.0	10.0560	135.2%
Matrix Spike DUP	9187-1MSD	5.0	10.0160	134.4%

Relative % Difference = 0.6%

BRL: Below Reporting Limit.

Date Reported:

01/23/92

REV5:12.91

Approved by: V.H.



COMPUCHEM
LABORATORIES, INC.

SULFIDE

Date(s) Analyzed: 12/16/91

Case: 9079
Matrix: Water

Client ID	CHEMWEST ID	Amount Detected (MG/L)
468548	9079-1	5.2
468584	9079-2	BRL
468603	9079-3	2.2
468620	9079-4	BRL

Client ID	CHEMWEST ID	Spike Conc. (MG/L)	Amount Detected (MG/L)	% Rec.
Method Blank	MB		BRL	
MBS	MBS	5.0	5.0320	100.6%
MBSD	MBSD	5.0	5.1520	103.0%

Relative % Difference = 2.4%

The reporting limit for Sulfide is 1.0 mg/L.

BRL: Below Reporting Limit.

Approved by: V.H.

Date Reported:
12/17/91

REV5:12.91



COMPUCHEM
LABORATORIES, INC.

SULFIDE

Date(s) Analyzed: 12/27/91

Case : 9136

Matrix: Water

Client ID	CHEMWEST ID	Amount Detected (MG/L)
472739	9136-1	BRL
472741	9136-2	BRL
472743	9136-3	BRL
472745	9136-4	3.2
472745	9136-4DP	3.2
472747	9136-5	BRL

Client ID	CHEMWEST ID	Spike Conc. (MG/L)	Amount Detected (MG/L)	% Rec.
Method Blank	MB		BRL	
MBS	MBS	5.0	5.4400	108.8
MBSD	MBSD	5.0	5.5200	110.4

Relative % Difference = 1.5%

Client ID	CHEMWEST ID	Spike Conc. (MG/L)	Amount Detected (MG/L)	% Rec.
Matrix Spike	9136-4MS	5.0	9.2800	121.6
Matrix Spike Dup.	9136-4MSD	5.0	9.9600	135.2

Relative % Difference = 10.6%

The reporting limit for Sulfide is 1.0 mg/L.

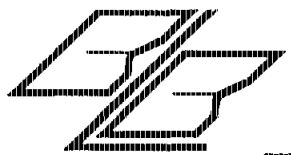
BRL: Below Reporting Limit.

Approved by: V.H.

Date Reported:

12/31/91

REV5:12.91



Section 20

PHENOLS ANALYSIS (GROUNDWATER)

WP-1 - Groundwater sample from Oxbow Area K, Well Point 1
WP-2 - Groundwater sample from Oxbow Area K, Well Point 2
WP-3 - Groundwater sample from Oxbow Area J, Well Point 3
WP-7 - Groundwater sample from Oxbow Area C, Well Point 7
WP-8 - Groundwater sample from Oxbow Area C, Well Point 8
WP-9 - Groundwater sample from Oxbow Area A, Well Point 9
ROA1G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-1
ROA3G - Groundwater sample from Oxbow Area A, Groundwater Monitoring Well A-3
ROB1G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-1
ROB2G - Groundwater sample from Oxbow Area B, Groundwater Monitoring Well B-2
ROC1G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-1
ROC2G - Groundwater sample from Oxbow Area C, Groundwater Monitoring Well C-2
ROJ1G - Groundwater sample from Oxbow Area J, Groundwater Monitoring Well J-1



ANALYTICAL SERVICES

5815 Middlebrook Pike • Knoxville, Tennessee 37921 • 615-588-6401



CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 19 OF 19

Sample Description: Five (5) water samples received October 14, 1988

Concentration units are mg/liter (ppm)

	<u>Phenols</u>	<u>Cyanide</u>
WP-1	<0.01	<0.01
WP-2	<0.01	<0.01
WP-3	<0.01	<0.01
WP-5	<0.01	<0.01
WP-6	0.04	<0.01

Roger L. Moore
Approved by Laboratory Manager

Title



ANALYTICAL SERVICES



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 25 OF 25

Sample Description: Five (5) water samples received October 20, 1988

Concentration units are mg/liter (ppm)

	<u>Cyanide</u>	<u>Phenols</u>
WP-4	<0.01	<0.01
WP-7	<0.01	<0.01
WP-8	<0.01	0.03
WP-9	<0.01	0.02
62	<0.01	0.05

Alice F. Moore
Approved by Laboratory Manager
Title _____



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 7 OF 19

Sample Description: WP-1 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
chloro-3-methylphenol	ND	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.

Allyn R. Moore
Approved by Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 9 OF 19

Sample Description: WP-2 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

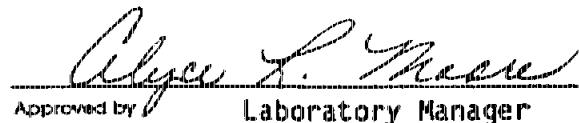
<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
4-chloro-3-methylphenol	ND	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.


Approved by Alyce J. Moore
Laboratory Manager



ANALYTICAL SERVICES

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CERTIFICATE OF ANALYSIS

TO: Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 15, 1988
PROJECT CODE: GMIN 42142
ORDER NUMBER:
PAGE 11 OF 19

Sample Description: WP-3 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
4-chloro-3-methylphenol	ND	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.

Alice S. Moore
Approved by Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 10 OF 25

Sample Description: WP-7 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Compound	Concentration (μ g/liter)	Compound	Concentration (μ g/liter)
4-chloro-3-methylphenol	<10	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.

Alce T. Moore
Approved by **Laboratory Manager**

Title



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 12 OF 25

Sample Description: WP-8 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>	<u>Compound</u>	<u>Concentration</u> <u>(μg/liter)</u>
4-chloro-3-methylphenol	ND	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.

Approved by Alyce F. Moore Laboratory Manager

Title



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CERTIFICATE OF ANALYSIS

TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 14 OF 25

Sample Description: WP-9 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
4-chloro-3-methylphenol	<10	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.

Approved by

Laboratory Manager

Title



ANALYTICAL SERVICES

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TO Geraghty & Miller, Inc.
ATTN: Bill Gray
125 E. Bethpage Road
Plainview, NY 11803

DATE REPORTED: November 18, 1988
PROJECT CODE: GMIN 42163
ORDER NUMBER:
PAGE 16 OF 25

Sample Description: 62 (Water)

ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Compound	Concentration ($\mu\text{g/liter}$)	Compound	Concentration ($\mu\text{g/liter}$)
4-chloro-3-methylphenol	ND	2-nitrophenol	ND
2-chlorophenol	ND	4-nitrophenol*	ND
2,4-dichlorophenol	ND	pentachlorophenol*	ND
2,4-dimethylphenol	ND	phenol	ND
2,4-dinitrophenol *	ND	2,4,6-trichlorophenol	ND
2-methyl-4,6-dinitrophenol *	ND		

Remarks: 10 = Quantitation limit.

ND = Not detected.

< = Detected but at a level less than the quantitation limit.

* = This compound has a quantitation limit of five (5) times that listed.

Alyce R. Mass
Approved by Laboratory Manager
Title _____

COMPOUND LIST - CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROA1G
COMPUCHEM SAMPLE NUMBER: 468198

	CONCENTRATION <u> </u> <u>(mg/L)</u>	DETECTION LIMIT <u> </u> <u>(mg/L)</u>
1. PHENOLS, TOTAL	0.018	0.010

COMPOUND LIST - CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROA3G
COMPUCHEM SAMPLE NUMBER: 474989

	<u>CONCENTRATION</u> <u>(mg/L)</u>	<u>DETECTION LIMIT</u> <u>(mg/L)</u>
1. PHENOLS, TOTAL	0.082	0.010

COMPOUND LIST - CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROB1G
COMPUCHEM SAMPLE NUMBER: 468192

	CONCENTRATION (mg/L)	DETECTION LIMIT (mg/L)
1. PHENOLS, TOTAL	BDL	0.010

BDL = BELOW DETECTION LIMITS

COMPOUND LIST -- CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROB2G
COMPUCHEM SAMPLE NUMBER: 468197

	CONCENTRATION (mg/L)	DETECTION LIMIT (mg/L)
1. PHENOLS, TOTAL	BDL	0.010

BDL = BELOW DETECTION LIMITS

COMPOUND LIST -- CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROC1G
COMPUCHEM SAMPLE NUMBER: 468538

	CONCENTRATION <u>(mg/L)</u>	DETECTION LIMIT <u>(mg/L)</u>
1. PHENOLS, TOTAL	0.019	0.010

COMPOUND LIST - CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROC2G
COMPUCHEM SAMPLE NUMBER: 468595

	CONCENTRATION <u>(mc/L)</u>	DETECTION LIMIT <u>(mc/L)</u>
1. PHENOLS, TOTAL	BDL	0.010

BDL = BELOW DETECTION LIMITS

COMPOUND LIST - CLASSICAL PARAMETERS

SAMPLE IDENTIFIER: ROJ1G
COMPUCHEM SAMPLE NUMBER: 472806

	CONCENTRATION <u>(mg/L)</u>	DETECTION LIMIT <u>(mg/L)</u>
1. PHENOLS, TOTAL	BDL	0.010

BDL = BELOW DETECTION LIMITS



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engineers & scientists