BACKGROUND DOCUMENT

REPORT ON REVISIONS TO 5TH EDITION AP-42 CHAPTER 15 - ORDNANCE DETONATION

EMISSION FACTORS DEVELOPED BASED ON EXPLODING ORDNANCE EMISSION STUDY PHASE II SERIES 2 TESTING CONDUCTED AT ABERDEEN PROVING GROUND, MARYLAND

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NOTICE

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TABLE OF CONTENTS

| | | | Page |
|-----|------|---|------|
| 1.0 | INT | RODUCTION | 1 |
| 2.0 | CON | MPOUNDS MEASURED AND EMISSION MEASUREMENT METHODS | 2 |
| | 2.1 | Carbon Monoxide, Carbon Dioxide, Nitrogen Oxides, and Sulfur Dioxide | |
| | 2.2 | Sulfur Hexafluoride, Ammonia, Methane, Carbon Monoxide, and Carbon Dioxide | |
| | 2.3 | Total Suspended Particulate | |
| | 2.4 | Particulate Matter with an Aerodynamic Diameter Less than or Equal to 10 or 2.5 M | |
| | 2.5 | Metals | |
| | 2.6 | Volatile Organic Compounds | |
| | 2.7 | Semivolatile Organic Compounds and Polycyclic Aromatic Hydrocarbons | |
| | 2.8 | Dioxin and Furan Compounds | |
| | 2.9 | Acid Gases | |
| | 2.10 | Hydrogen Cyanide | 7 |
| | 2.11 | Aldehydes and Carbonyl Compounds | |
| | 2.12 | Hydrocarbons, n-Hexane, and Sulfur Hexafluoride | |
| | 2.13 | · | |
| | 2.14 | e | |
| | 2.15 | Hexavalent Chromium | |
| 3.0 | TES | T DATA ANALYSIS AND RATING | 8 |
| | 3.1 | EPA Guidance Regarding Test Data Quality Ratings | |
| | 3.2 | Analysis of Test Data | |
| | 3.3 | Test Data Quality Ratings | 14 |
| 4.0 | EMI | SSION FACTOR CALCULATIONS | 16 |
| | 4.1 | Correction of Sample Volume to Standard Conditions | |
| | 4.2 | Incorporation of Analytical Detection-Limits to the Test Data | |
| | 4.3 | Determination of Background Concentration | |
| | 4.4 | Determination of Test Compound Concentration | |
| | 4.5 | Determination of Background-Corrected Concentration | 19 |
| | 4.6 | Determination of Dilution-Corrected Concentration | 21 |
| | 4.7 | Determination of Mass of Compound Released | 21 |
| | 4.8 | Determination of Emission Factors | 21 |
| | 4.9 | Determination of Average Emission Factors | 21 |
| 5.0 | EMI | SSION FACTOR RATINGS | 22 |
| | 5.1 | Emission Factor Ratings Assigned – Based on Ordnance-Specific Test Data | |
| | 5.2 | Emission Factor Ratings Assigned – Based on All Available Test Data | |
| 6.0 | REF | ERENCES | 32 |

TABLE OF CONTENTS (cont.)

| APPENDIX B: | NEW AP-42 SECTIONS FOR ORDNANCE INCLUDED IN EXPLODING ORDNANCE EMISSION STUDY PHASE II SERIES 2 TESTING AT ABERDEEN PROVING GROUND, MARYLAND |
|-------------|--|
| APPENDIX C: | PUBLIC COMMENTS ON AP-42 SECTIONS FOR ORDNANCE INCLUDED IN PHASE II SERIES 2 TESTING AT ABERDEEN PROVING GROUND, MARYLAND |
| | LIST OF TABLES |

| | LIST OF TABLES |
|---------------------------------|--|
| No. | <u>Title</u> Page |
| 1 2 3 4 5 6 7 | Ordnance for Which Emission Factors Were Developed |
| | LIST OF FIGURES |
| No. | <u>Title</u> <u>Page</u> |
| 1 | Calculation of background-corrected concentration |
| | |

1.0 INTRODUCTION

Due to the lack of credible data concerning emissions from training ordnance when used in their tactical configurations, the U.S. Army Environmental Center (USAEC) established a program to quantify emissions from the detonation of ordnance. This document presents background information concerning the development of air emission factors for four types of ordnance used during training exercises at U.S. Army installations. The air emission factors were developed from test data collected by USAEC. Ordnance for which emission factors have been developed and their corresponding AP-42 sections are identified in Table 1. To help readers easily find those emission factors of interest, the ordnance are organized according to their Department of Defense Identification Code (DODIC).

| DODIC | DODIC Ordnance Description | |
|-------------------------------|-----------------------------------|---------|
| G911 | G911 MK3A2 Offensive Hand Grenade | |
| K145 M18A1 Antipersonnel Mine | | 15.7.3 |
| L709 M25 Target-Hit Simulator | | 15.8.19 |
| M032 | 1-Pound Demolition Block Charge | 15.9.4 |

TABLE 1 ORDNANCE FOR WHICH EMISSION FACTORS WERE DEVELOPED

The emission factors described in this document are based on data obtained during testing conducted at Aberdeen Proving Ground, Maryland, as presented in the final test report titled *Report No.2* for the Exploding Ordnance Emission Study Phase II¹ and the document titled Detailed Test Plan No. 2 for the Exploding Ordnance Emission Study Phase II.² These documents were supplemented by additional data provided by the testing contractor.³ For each ordnance, two test runs were conducted. One ordnance was detonated per run. Source test protocols were developed by USAEC before any testing was conducted and were reviewed by the U.S. Environmental Protection Agency's (EPA's) Emission Measurement Center. The tests were conducted between April 2 and April 4, 2001.

The compounds that were measured include carbon monoxide (CO), carbon dioxide (CO₂), oxides of nitrogen (NO_X), sulfur dioxide (SO₂), ammonia (NH₃), methane (CH₄), total suspended particulate (TSP), particulate matter with an aerodynamic diameter less than or equal to 10 microns (PM-10), particulate matter with an aerodynamic diameter less than or equal to 2.5 microns (PM-2.5), metals, volatile organic compounds (VOC), semivolatile organic compounds (SVOC), dioxin and furan compounds, acid gases, hydrogen cyanide (HCN), aldehydes and carbonyl compounds, hydrocarbons, polycyclic aromatic hydrocarbons (PAH), energetic materials, total organic carbon (TOC), and hexavalent chromium. Within each of the AP-42 sections, only emission factors for criteria pollutants, carbon dioxide, hazardous air pollutants (as defined by §112(b)(1) of the *Clean Air Act* [CAA]), and toxic chemicals (as defined by §313 of the *Emergency Planning and Community Right-to-Know Act* [EPCRA]) are presented.

The emission factors were developed on a "per item" basis and on a "per net explosive weight (NEW)" basis. Users should choose the appropriate emission factor to estimate emissions based upon the data available; either factor is equally valid. The NEW of each ordnance tested is provided in the corresponding AP-42 section and in Table 2.

TABLE 2 ORDNANCE NET EXPLOSIVE WEIGHT

| DODIC | Ordnance Description | NEW (lb/item) ^a |
|--------------------------------------|-------------------------------|----------------------------|
| G911 | MK3A2 Offensive Hand Grenade | 4.94 E-01 |
| K145 | M18A1 Antipersonnel Mine | |
| L709 | L709 M25 Target-Hit Simulator | |
| M032 1-Pound Demolition Block Charge | | 1.00 |

^a NEW values were obtained from Reference 1.

This document includes five sections in addition to this Introduction. Section 2 of this document identifies the compounds measured during the test program and describes the emission measurement methods used. Section 3 includes a discussion of the emission factor final test report and ratings for the test data contained therein. Section 4 describes the calculations and methodologies used to develop emission factors for each type of compound measured. Section 5 describes the methodology used to rate the emission factors and provides emission factor ratings for each type of compound measured. Section 6 includes a complete list of the references cited in this document.

There are two appendices included with this document. Appendix A identifies, by ordnance type, all of the compounds for which analyses were performed and the emission factors that were developed. [Note: Compounds present in the method blank at greater than 50 percent of test levels are excluded from Appendix A as described in Section 3.2.4.] Appendix A also identifies the minimum detection levels associated with all compounds that were not detected. Emission factors and minimum detection levels presented in Appendix A were determined from the most accurate method if two sampling or analytical methods were used to measure one compound. Appendix B presents the new AP-42 sections for the four ordnance that were tested.

In addition to this document, there are electronic databases available on the web (http://www.epa.gov/ttn/chief/ap42/index.html) that contain the data used in the development of the emission factors. The general procedures that were followed to develop these emission factors can be found at the same web address under the title *Procedures for Preparing Emission Factor Documents*.⁴

2.0 COMPOUNDS MEASURED AND EMISSION MEASUREMENT METHODS

The USAEC Exploding Ordnance Phase II testing was conducted in the Large Octagon Test Chamber (LOTC) located at the Aberdeen Test Center (ATC), Maryland. The LOTC is an eight-sided steel structure with flat ends. Internally, the eight sides were increased to 16 sides by welding steel plates over each wall seam. This design strengthened the LOTC and removed sharp interior corners that would inhibit gaseous mixing during the testing. Additional plates were added to the upper and lower internal corners to create a more parabolic shape. The interior volume is approximately 2,400 cubic feet. Interior wall-to-wall distance is 14.5 feet, with a floor-to-ceiling height of 14 feet.

The measurement assembly and instrument enclosure are mounted on rollers in vertical tracks. After ordnance detonation, the assembly is lowered by an electric hoist from the top of the LOTC through the center of the ceiling and into the chamber. A metal cover on top of the assembly seals the interior when the assembly has been lowered into the LOTC. The LOTC has two flanged viewing ports for video recording and two ports with hinged internal doors through which the mixing fans are inserted.

A number of different test methods were employed to collect and analyze the emission data that were used to develop emission factors for detonation of ordnance. Table 3 identifies each emission test

method used; bracketed information identifies the purpose of using the method. The emissions data were collected using EPA test methods published in Title 40 of the Code of Federal Regulations, Part 50 (40 CFR 50); 40 CFR 60; 40 CFR 63; and in *Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air.*⁵ Some of the sample analytical procedures used were from EPA Office of Solid Waste (OSW) publication SW-846, *Test Methods for Evaluating Solid Waste*, *Physical/Chemical Methods*⁶ and the *National Institute for Occupational Safety and Health (NIOSH) Manual of Analytical Methods (NMAM)*, *Fourth Edition.*⁷ Additional sampling and analytical procedures were taken from Internal Operating Procedures (IOP) developed by the ATC Chemistry Team and from Standard Operating Procedures (SOP) developed by the U.S. Army Center for Health Promotion and Preventive Medicine (USACHPPM) Directorate of Laboratory Services (DLS). Where necessary, the test methods were adapted to reflect application to the unique testing of ordnance detonation in the LOTC.

The following sections identify and briefly describe the test methods used to measure each compound or group of compounds. Additional information regarding the operation of the LOTC and the test methods used is presented in Reference 2. EPA-approved methods were used by the laboratories that provided sampling and analysis data.

TABLE 3 EMISSION TEST METHODS USED

| Compound | Test Method | |
|--|--|--|
| СО | 40 CFR 60, Appendix A, EPA Method 10 – Determination of Carbon Monoxide Emissions from Stationary Sources [sampling and analysis] | |
| CO ₂ | 40 CFR 60, Appendix A, EPA Method 3A - Determination of Oxygen and Carbon Dioxide Concentrations in Emissions from Stationary Sources (Instrumental Analyzer Procedure [sampling and analysis] | |
| NO_X | 40 CFR 60, Appendix A, EPA Method 7E - Determination of Nitrogen Oxides Emissions from Stationary Sources (Instrumental Analyzer Procedure) [sampling and analysis] | |
| SO_2 | 40 CFR 60, Appendix A, EPA Method 6C - Determination of Sulfur Dioxide Emissions from Stationary Sources (Instrumental Analyzer Procedure) [sampling and analysis] | |
| SF ₆ , NH ₃ , CO, CO ₂ , and CH ₄ | 40 CFR 63, Appendix A, EPA Proposed Method 320 - Measurement of Vapor Phase Organic and Inorganic Emissions By Extractive Fourier Transform Infrared (FTIR) Spectroscopy [sampling and analysis] | |
| TSP | 40 CFR 60, Appendix A, EPA Method 17 - Determination of Particulate Emissions from Stationary Sources (in-Stack Filtration Method) [sampling] | |
| | 40 CFR 50, Appendix B - Reference Method for the Determination of Suspended Particulate Matter as in the Atmosphere (High Volume Method) [analysis] | |
| PM-10 and PM-2.5 | 40 CFR 50, Appendix J - Reference Method for the Determination of Particulate Matter as PM-10 in the Atmosphere [sampling and analysis] | |
| Metals Metal sample was obtained from TSP sample [sampling] | | |
| | 40 CFR 50, Appendix G - Reference Method for Determination of Lead in Suspended Particulate Matter Collected from Ambient Air [analysis] | |

| Compound | Test Method |
|----------------------------|--|
| VOC | EPA Compendium Method TO-14A - Determination of Volatile Organic Compounds (VOCs) in Ambient Air Using Specially Prepared Canisters with Subsequent Analysis by Gas Chromatography [sampling and analysis] |
| РАН | EPA Compendium Method TO-13A - Determination of Polycyclic Aromatic Hydrocarbons (PAHs) in Ambient Air Using Gas Chromatography/Mass Spectrometry (GC/MS) [sampling and analysis] |
| | USA CHPPM SOP SV 72.4 [analysis] |
| SVOC | EPA Compendium Method TO-13A - Determination of Polycyclic Aromatic Hydrocarbons (PAHs) in Ambient Air Using Gas Chromatography/Mass Spectrometry (GC/MS) [sampling] |
| | SW-846 Method 8270M - Semi-Volatile Organic Compounds by Gas Chromatography/ Mass Spectrometry (GC/MS) [analysis] |
| Dioxins and Furans | EPA Compendium Method TO-9A - Determination of Polychlorinated, Polybrominated, and Brominated/Chlorinated Dibenzo-p-Dioxins and Dibenzofurans in Ambient Air [sampling and analysis] |
| Acid Gases | NIOSH Method 7903 - Inorganic Acids [sampling and analysis] |
| | ATC IOP 372 - Anion Analysis Using the AS 14 Column [analysis] |
| Hydrogen | NIOSH Method 7904 - Aerosol and Gas Cyanides [sampling] |
| Cyanide | ATC IOP No. 231 - Analysis of Cyanide and Sulfide by Ion Chromatography [analysis] |
| Aldehydes and Carbonyls | EPA Compendium Method TO-11A - Determination of Formaldehyde in Ambient Air Using Adsorbent Cartridge Followed by High Performance Liquid Chromatography (HPLC) [sampling and analysis] |
| | ATC IOP No. 227, Revision 3.0 - Analysis of Formaldehyde and Other Carbonyl Compounds by High Performance Liquid Chromatography. [analysis] |
| Hydrocarbons and n-Hexane | 40 CFR 60, Appendix A, EPA Method 18 - Measurement of Gaseous Organic Compound Emissions by Gas Chromatography [sampling and analysis] |
| SF ₆ | 40 CFR 60, Appendix A, EPA Method 18 - Measurement of Gaseous Organic Compound Emissions by Gas Chromatography [sampling] |
| | Project specific IOP, IR and gas chromatography with electron capture detection (GC/ECD) [analysis] |
| Energetic Materials | EPA Compendium Method TO-13A - Determination of Polycyclic Aromatic Hydrocarbons (PAHs) in Ambient Air Using Gas Chromatography/Mass Spectrometry (GC/MS) [sampling] |
| | USACHPPM-DLS SOP CAD 26.2 - Procedure for Analysis of Explosives in Ambient Air [analysis] |
| Total Carbon | NIOSH Method 5040 - <i>Diesel Particulate Matter as Elemental Carbon</i> [sampling and analysis] |
| Hexavalent Chromium | OSHA Method ID-215 - Hexavalent Chromium in Workplace Atmospheres [sampling and analysis] |

2.1 Carbon Monoxide, Carbon Dioxide, Nitrogen Oxides, and Sulfur Dioxide

Real-time concentrations of CO, CO_2 , NO_X , and SO_2 that resulted from the detonation of ordnance in the LOTC were measured using a continuous emissions measurement system (CEMS). The final test report¹ states that the compounds were measured using EPA reference methods as described in the following paragraph. Although the report does include results from some quality control checks, due to a lack of calibration data, conformance with the stated methodologies actually could not be confirmed.

CO sampling and analysis was conducted in accordance with 40 CFR Part 60, Appendix A, Method 10 - Determination of Carbon Monoxide Emissions from Stationary Sources, with a Thermo Environmental Instruments, Inc. (TEI) nondispersive infrared analyzer. CO₂ sampling and analysis was conducted in accordance with 40 CFR Part 60, Appendix A, Method 3A - Determination of Oxygen and Carbon Dioxide Concentrations in Emissions from Stationary Sources, with a TEI nondispersive infrared analyzer. NO_X sampling and analysis was conducted in accordance with 40 CFR Part 60, Appendix A, Method 7E - Determination of Nitrogen Oxides Emissions from Stationary Sources, with a TEI model No. 10AR chemiluminescent analyzer. SO₂ sampling and analysis was conducted in accordance with 40 CFR Part 60, Appendix A, Method 6C - Determination of Sulfur Dioxide Emissions from Stationary Sources, with a TEI 43H pulsed fluorescent analyzer.

Each analyzer channel sampled at approximately 4 liters per minute (lpm) through Teflon-lined polyethylene tubing with Mine Safety Appliance (MSA) ultra type H cartridge filters used to control particulate. An inline Balston filter was used to trap condensation. Due to the ambient condition of the samples, no heated probes or sample lines were used. Data was collected four times per second.

2.2 Sulfur Hexafluoride, Ammonia, Methane, Carbon Monoxide, and Carbon Dioxide

Extractive Fourier Transform Infrared (FTIR) spectroscopy was used to measure real-time concentrations of the SF₆ tracer gas and NH₃, CO, CO₂, and CH₄, as well as to identify qualitatively the presence of other permanent gas species using protocols modified from procedures in 40 CFR 63, Appendix A, EPA Proposed Method 320 - *Measurement of Vapor Phase Organic and Inorganic Emissions By Extractive Fourier Transform Infrared (FTIR) Spectroscopy*. Samples were drawn continuously from the LOTC at a nominal rate of 0.5 cubic feet per minute (ft³/min) using a Teflon-lined polyethylene tube and an Alcatel Pascal[®] rotary vane-type pump, for a minimum of 20 minutes. The FTIR spectra were collected and stored at a minimum of every 20 seconds.

2.3 Total Suspended Particulate

The TSP concentration that resulted from the detonation of ordnance in the LOTC was determined using modified procedures based on 40 CFR 60, Appendix A, Reference Method 17 - Determination of Particulate Emissions from Stationary Sources (In-Stack Filtration Method). A Graseby® Model PS-1 ambient air sampling module equipped with a quartz fiber filter was connected by a flex line to the Method 17 metering system. The metering system extracted test chamber air through a Whatman® filter at a nominal rate of 1ft³/min for a target minimum sampling time of 20 minutes. The TSP concentration was computed, in accordance with 40 CFR 50, Appendix B - Reference Method for the Determination of Suspended Particulate Matter as in the Atmosphere, by dividing the mass of TSP collected by the volume of air sampled, corrected to standard conditions.

2.4 Particulate Matter with an Aerodynamic Diameter Less than or Equal to 10 or 2.5 Microns

The PM-10 and PM-2.5 concentrations that resulted from the detonation of ordnance in the LOTC were determined using a sampling and analysis procedure based on 40 CFR 50, Appendix J - Reference Method for the Determination of Particulate Matter as PM-10 in the Atmosphere. An ambient

sampler with a Graseby Model No. G246B dichotomous PM-10 size selective inlet head was used to collect the PM-10 and smaller particulate at a sampling rate of 16.7 lpm for 20 minutes. The PM-10 fraction was then passed through a virtual impactor to separate the particulate larger than PM-2.5 from PM-2.5. The PM-10 and PM-2.5 concentrations were computed by dividing the mass of PM-10 and PM-2.5 collected by the respective volume of air sampled, corrected to standard conditions.

2.5 Metals

Metal concentrations that resulted from the detonation of ordnance in the LOTC were determined using particulate matter from the TSP samples collected as described in Section 2.3. The analyses followed the procedures of 40 CFR 50, Appendix G - *Reference Method for the Determination of Lead in Suspended Particulate Matter Collected from Ambient Air.* The method was modified to use inductively coupled plasma (ICP) spectroscopy for analysis and to include additional metals. Concentrations of target metals were computed by dividing the mass of metal collected by the volume of air sampled, corrected to standard conditions.

2.6 Volatile Organic Compounds

VOC concentrations that resulted from the detonation of ordnance in the LOTC were determined using procedures from the *Second Supplement to Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air*, Method TO-14A - *Determination of Volatile Organic Compounds (VOCs) in Ambient Air Using Specially-Prepared Canisters with Subsequent Analysis by Gas Chromatography*. Air samples were collected for 20 minutes using a Scientific Instruments Specialists, Inc., pressurized canister sampler and 6-liter silicon-lined SUMMA canister that was certified VOC-free prior to use. An analysis by gas chromatography/mass spectrometry in full scan mode (GC/MS/SCAN) followed. Unknown compounds were tentatively identified using computerized mass spectral matching techniques.

2.7 Semivolatile Organic Compounds and Polycyclic Aromatic Hydrocarbons

SVOC and PAH concentrations that resulted from the detonation of ordnance in the LOTC were determined based on procedures found in Method TO-13A - *Determination of Polycyclic Aromatic Hydrocarbons (PAHs) in Ambient Air Using Gas Chromatography/Mass Spectrometry (GC/MS)*. During each run, samples were collected using a PS-1 sampler equipped with a quartz fiber filter and XAD-2 adsorbent resin cartridges. The PS-1 sampling module was used to sample for 20 minutes at 1 ft³/min. Following sampling, the filters and resin cartridges underwent solvent extraction and analysis by GC/MS analysis in selected ion monitoring (SIM) mode. The sample extracts were reanalyzed in scan mode following modified procedures in SW-846 Method 8270 - *Semivolatile Organic Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)*. Modifications are described in the USACHPPM SOP SV 72.4. Unknown compounds were tentatively identified using computerized mass spectral matching techniques.

2.8 Dioxin and Furan Compounds

Polychlorinated dibenzodioxin (PCDD) and polychlorinated dibenzofuran (PCDF) compound concentrations that resulted from the detonation of ordnance in the LOTC were determined based on procedures found in Method TO-9A - *Determination of Polychlorinated, Polybrominated, and Brominated/Chlorinated Dibenzo-p-Dioxins and Dibenzofurans in Ambient Air.* Samples were collected using modified PS-1 samplers with quartz fiber particulate filters and polyurethane foam (PUF) packed cartridges. The PS-1 sampling module was connected to a stack source airflow metering system with a flex line and used to sample for 20 minutes at a nominal rate of 1 ft³/min. After sampling, the filters and

cartridges underwent extraction and analysis by GC/High Resolution Mass Spectroscopy according to Method TO-9A.

2.9 Acid Gases

Acid gas concentrations that resulted from the detonation of ordnance in the LOTC were determined based on procedures found in NIOSH Method 7903 - *Inorganic Acids*. Acid gases were collected by using silica gel sorbent tubes (ORBO-53) and a sampling pump. Sample gases were passed through the sorbent tubes for 20 minutes at a rate of 0.5 lpm. Analysis was performed according to NIOSH Method 7903 - *Inorganic Acids*, by ion chromatography as detailed in ATC IOP No. 372 - *Anion Analysis Using the AS14 Column*.

2.10 Hydrogen Cyanide

HCN concentrations that resulted from the detonation of ordnance in the LOTC were determined using NIOSH Method 7904 - *Aerosol and Gas Cyanides*, but utilizing modified sampling and analytical procedures. The sample gas was drawn through a 0.8 micrometer PVC filter and an impinger containing 0.1 normal potassium hydroxide (KOH) for 20 minutes at a rate of 0.5 lpm. The impinger solution and extracted filter were analyzed by ion chromatography in accordance with NIOSH method 7904 and ATC IOP No. 231 - *Analysis of Cyanide and Sulfide by Ion Chromatography*.

2.11 Aldehydes and Carbonyl Compounds

Aldehyde and carbonyl compound concentrations that resulted from the detonation of ordnance in the LOTC were determined using EPA Compendium Method TO-11A - *Determination of Formaldehyde in Ambient Air Using Adsorbent Cartridge Followed by High Performance Liquid Chromatography (HPLC)*, but utilizing modified sampling and analytical procedures. Dinitrophenylhydrazine (DNPH) laden cartridge tubes were used as a direct probe to trap and derivatize aldehyde and carbonyl compounds. Sample gases were extracted for 20 minutes (nominal) from the test chamber through ozone denuders at a rate of 1 to 2 lpm using a diaphragm pump, manifold, critical orifice, and needle valve. Analysis was by reverse-phase HPLC with ultraviolet (UV) detection using Compendium Method TO-11A modified as detailed in ATC IOP 227 Rev. 3.0 - *Analysis of Formaldehyde and Other Carbonyl Compounds by High Performance Liquid Chromatography*.

2.12 Hydrocarbons, n-Hexane, and Sulfur Hexafluoride

Hydrocarbon, n-hexane, and SF₆ concentrations that resulted from the detonation of ordnance in the LOTC were sampled using 40 CFR 60, Appendix A, EPA Method 18 - *Measurement of Gaseous Organic Compound Emissions by Gas Chromatography*. The samples were collected in Tedlar[®] bags at a nominal rate of 1 lpm for 20 minutes. Analysis for the light hydrocarbons was by gas chromatography with a flame ionization detector (GC/FID) in accordance with 40 CFR 60, Appendix A, EPA Method 18. SF₆ was analyzed by GC/MS or by gas chromatography with electron capture detection (GC/ECD). This measurement served as a backup to the FTIR measurement of SF₆.

2.13 Energetic Materials

Energetic compound concentrations that resulted from the detonation of ordnance in the LOTC were determined based on procedures found in EPA Compendium Method TO-13A - *Determination of Polycyclic Aromatic Hydrocarbons (PAHs) in Ambient Air Using Gas Chromatography/Mass Spectrometry (GC/MS)*. Samples were collected using a Graseby Model PS-1 sampling module equipped with a quartz fiber particulate filter in series with an XAD-2 adsorbent resin cartridge. The PS-1 sampling module was connected by flex line to a stack source airflow metering system and was used to

sample for 20 minutes at approximately 1 ft³/min. Analyses followed procedures outlined in USACHPPM, Directorate of Laboratory Sciences, Chromatographic Analysis Division - *Procedure for Analysis of Explosives in Ambient Air*, Standard Operating Procedure CAD 26.2.

This method specifically address the extraction and analysis of XAD-2 resin used to sample energetic compounds from air. The energetic compounds were extracted from the XAD-2 resin using isoamyl acetate and analyzed by gas chromatography with electron capture detection/mass spectrometry (GC/ECD/MS). A surrogate compound (3,4-dinitrotoluene) was added to the XAD-2 resin prior to sampling. The quartz fiber filter used in conjunction with the XAD-2 resin was extracted and analyzed separately to characterize the energetic materials present as particulate solids. The filter was placed in a vial containing isoamyl acetate immediately after testing to preserve the sample and prevent sublimation of any of the analytes. No surrogate compound was used in the analysis of the filter portion of the sample.

2.14 Total Carbon

Total carbon concentrations that resulted from the detonation of ordnance in the LOTC were determined based on the procedures in NIOSH Method 5040 - *Diesel Particulate Matter as Elemental Carbon*. Samples were collected at a rate of 2.0 lpm using an APEX Instruments Model No. 623 meter console. A dry gas meter contained in the console was used to measure the sample volume. Flow rates were verified using an electronic flow meter. The particulate sample was collected on a 37-mm heat treated quartz fiber filter and was analyzed using a thermo-optical analyzer to differentiate between organic and elemental carbon.

2.15 Hexavalent Chromium

The hexavalent chromium concentrations that resulted from the detonation of ordnance in the LOTC were determined based on the procedures in OSHA Method ID-215 - *Hexavalent Chromium in Workplace Atmospheres*. Since particulate results from previous weapons firing showed that a large fraction of the total particulate produced was less than 2.5 microns in size, the method was modified to utilize a 0.8-µm pore size polyvinyl chloride (PVC) filter instead of a 5-µm PVC filter. Samples were collected at a rate of 2.0 lpm using an APEX Instruments Model No. 623 meter console. A dry gas meter contained in the console was used to measure the sample volume. Flow rates were verified using an electronic flow meter. The filter media was extracted and analyzed using ion chromatography.

3.0 TEST DATA ANALYSIS AND RATING

3.1 EPA Guidance Regarding Test Data Quality Ratings

Prior to inclusion of emission factors in AP-42, the reliability of the underlying emission test data must be appraised in accordance with the rating system specified in Reference 4. Under this rating system, test data are assigned a rating from A to D, where an "A" rating is assigned to the highest quality data. The criteria used to assign a specific data quality rating are summarized below.

- A Tests are performed by using an EPA reference test method, or when not applicable, a sound methodology. Tests are reported in enough detail for adequate validation and raw data are provided that can be used to duplicate the emission results presented in the report.
- **B** Tests are performed by a generally sound methodology, but lacking enough detail for adequate validation. Data are insufficient to completely duplicate the emission result presented in the report.

- C Tests are based on an unproven or new methodology, or are lacking a significant amount of background information.
- **D** Tests are based on a generally unacceptable method, but the method may provide an order-of-magnitude value for the source.

Four specific criteria are identified in Reference 4 for consideration to assist in the assignment of a test data quality rating. These four criteria are:

- 1. <u>Source operation</u>. If the manner in which the source was operated is well documented in the report and the source was operating within typical parameters during the test, an "A" rating should be assigned. If the report stated parameters that were typical, but lacked detailed information, a "B" rating should be assigned. If there is reason to believe the operation was not typical, a "C" or "D" rating should be assigned.
- 2. Test methods and sampling procedures. In developing the ratings, the estimated accuracy and precision of the test method as well as the adequacy of the documentation should be considered. In general, if a current EPA reference test method, appropriate for the source, was followed, the rating should be higher ("A" or "B"). If other methods were used, an assessment should be made of their validity. If it is judged that the method was likely to be inaccurate or biased, a lower rating ("C" or "D") should be given. A complete report should indicate whether any procedures deviated from standard methods and explain any deviations. If deviations were reported, an evaluation should be made of whether these were likely to influence the test results.
- 3. Process information. During testing, many variations in the process can occur without warning and sometimes without being noticed. Such variations can induce wide deviations in sampling results. If a large variation between test run results cannot be explained by information contained in the site final test report or from test reports of other sources, the data are suspect and should be given a lower rating or excluded. However, it should be recognized that a process may have highly variable emissions and a lower rating may not be appropriate solely on the basis of wide deviations in sampling results.
- 4. Analysis and calculations. Ideally, final test reports should contain original raw data sheets and other documentation such as gas parameters (dry cubic feet per minute, oxygen percentage), calculation sheets, or example calculations describing how the calculated emission results were obtained. If there are data sheets, the nomenclature and equations used should be compared to those specified by EPA to establish equivalency. The depth of review of the calculations should be dictated by the reviewers' confidence in the ability and conscientiousness of the tester, based on such factors as consistency of results and completeness of other areas of the final test report. Reports may indicate that raw data sheets were available, but were not included. If the final test report is of high quality based on the other criteria, the quality rating should not be lowered due to a lack of data sheets.

An overall test data quality rating should be assigned based upon the ratings assigned for each of the four criteria.

3.2 Analysis of Test Data

Data included in the final test report,¹ the detailed test plan,² and in the analytical data supporting the test report³ were rated in accordance with the rating system described above. Results for each of the four criteria described above are presented in the following sections.

3.2.1 Source Operations

The manner by which the ordnance were deployed (i.e., used) is documented in the final test report. With the exception of the adaptation discussed below that were made to facilitate testing in the LOTC, each ordnance that was tested was deployed in a manner similar to that which would occur in the field. DODICs K145 and M032, the M18A1 Antipersonnel Mine and the 1-Pound Demolition Block charge, respectively, were detonated using an M6 blasting cap. M6 blasting caps were also detonated during the background runs for these ordnance so that emissions associated with the blasting caps could be subtracted from emissions associated with the blasting caps and the ordnance. DODIC G911, the MK3A2 Offensive Hand Grenade, was placed in a special cup and the pin was removed. The grenade was pulled from the cup using a cable, the safety handle was actuated, and detonation in its tactical scenario occurred. Finally, DODIC L709, the M25 Target-Hit Simulator, was detonated by electrical pulse from a blasting machine. The tests appear to have replicated typical ordnance operating parameters; consequently, the test data should be assigned an "A" rating based on this criterion.

3.2.2 Test Methods and Sampling Procedures

The test methods and sampling procedures were evaluated as being appropriate and consistent with EPA test methods or sound methodology. Except as noted below, no problems of any significance were identified; consequently, the test data should be assigned an "A" rating based on this criterion.

3.2.2.1 CEMS-Measured Data

CEMS raw data were provided for the tests³, but not for the pre- and post-test quality control (QC) activities. CEMS QC measures specified under 40 CFR 60, Appendix A, EPA Methods 3A, 7E, and 6C were not followed. Furthermore, none of the calibration gas certifications were supplied. There was no evidence of bias in the data; however, based on the issues noted above, the test data for the CEMS-measured compounds (i.e., CO, CO₂, NO_x, and SO₂) should be assigned a "B" rating based on this criterion.

3.2.2.2 Compounds Sampled or Analyzed Using More than One Test Method or Analytical Method

Thirty-four compounds were either sampled or analyzed using two methods; these compounds are identified in Table 4. For each of these compounds, emission factors were calculated based upon the data measured using the more appropriate test or analytical method; data obtained using the less appropriate method were ignored. The more appropriate method was identified by reviewing the methods and the target compound lists associated with each method. If a specific compound appeared on the target compound list for one method but not the other, the method targeting the compound was selected. If a specific compound appeared on the target compound lists for both methods, the method judged to provide the most accurate data was selected.

If a compound was analyzed using both the SW8270M (SVOC) and TO-14A (VOC) methods, the TO-14A method analysis was judged to be more accurate and was therefore selected. For compounds analyzed using both the SW8270M (SVOC) and TO-13A (PAH) methods, the TO-13A value was selected because TO-13A allowed quantification of compounds at lower levels than SW8270M. For compounds sampled using a CEMS and FTIR spectroscopy, the CEMS measurements were judged to be

more accurate. For compounds analyzed using both SW8270M (SVOC) and SOP CAD 26.2 (energetic materials), the SOP CAD 26.2 data were selected because SOP CAD 26.2 allows detection of compounds at lower levels than SW8270M. If a compound analyzed using both the TO-11A (aldehydes) and TO-14A (VOC) methods, the TO-11A method analysis was judged to be more accurate and was therefore selected. For compounds analyzed using both Method 18 (hydrocarbons) and TO-14A (VOC), and for which the compounds appeared on both target compound lists, the Method 18 analysis was judged to be more accurate and was selected. For compounds analyzed using both Method 18 (hydrocarbons) and FTIR spectroscopy, the Method 18 analysis was judged to be more accurate and was selected.

Occasionally, the compound measurement from the less accurate method was chosen because the compound was present in the method blank of the preferred method. In addition, if a compound sampled under two methods had poor precision between test runs for the sampling method that would have been more accurate under normal circumstances, data from the other method were used to calculate emission factors. These cases are also noted in the footnotes to Table 4.

TABLE 4 COMPOUNDS MEASURED USING MORE THAN ONE TEST OR ANALYTICAL METHOD

| Compound | Selected Method | Other Method Employed |
|-----------------------------------|-------------------------------------|------------------------------|
| Carbon dioxide | 40 CFR 60, Method 3A (CEMS) | 40 CFR 63, Method 320 (FTIR) |
| Carbon monoxide | 40 CFR 60, Method 10 (CEMS) | 40 CFR 63, Method 320 (FTIR) |
| Methane | 40 CFR 60, Method 18 (Hydrocarbons) | 40 CFR 63, Method 320 (FTIR) |
| Acenaphthene | TO-13A (PAH) | SW8270M (SVOC) |
| Acenaphthylene | TO-13A (PAH) | SW 8270M (SVOC) |
| Acrolein | TO-11A (Aldehydes) | TO-14A (VOC) |
| Anthracene | TO-13A (PAH) | SW8270M (SVOC) |
| Benzo[a]anthracene ^a | TO-13A (PAH) | SW8270M (SVOC) |
| Benzo[b]fluoranthene | TO-13A (PAH) | SW8270M (SVOC) |
| Benzo[k]fluoranthene ^b | TO-13A (PAH) | SW8270M (SVOC) |
| Benzo[g,h,i]perylene ^c | TO-13A (PAH) | SW8270M (SVOC) |
| Benzo[a]pyrene | TO-13A (PAH) | SW8270M (SVOC) |
| 1,3-Butadiene | 40 CFR 60 Method 18 (Hydrocarbons) | TO-14A (VOC) |
| Chrysene ^d | TO-13A (PAH) | SW8270M (SVOC) |
| Dibenz[a,h]anthracene | TO-13A (PAH) | SW8270M (SVOC) |
| 1,2-Dichlorobenzene | TO-14A (VOC) | SW8270M (SVOC) |
| 1,3-Dichlorobenzene | TO-14A (VOC) | SW8270M (SVOC) |
| 1,4-Dichlorobenzene | TO-14A (VOC) | SW8270M (SVOC) |
| 2,4-Dinitrotoluene | SOP CAD 26.2 (Energetics) | SW8270M (SVOC) |
| 2,6-Dinitrotoluene | SOP CAD 26.2 (Energetics) | SW8270M (SVOC) |
| Fluoranthene ^e | TO-13A (PAH) | SW8270M (SVOC) |

TABLE 4 (cont.)

| Compound | Selected Method | Other Method Employed |
|-------------------------------------|------------------------------------|-----------------------|
| Fluorene | TO-13A (PAH) | SW8270M (SVOC) |
| Hexachlorobutadiene | TO-14A (VOC) | SW8270M (SVOC) |
| Hexachlorethane | TO-14A (VOC) | SW8270M (SVOC) |
| Hexane | 40 CFR 60 Method 18 (Hydrocarbons) | TO-14A (VOC) |
| Indeno[1,2,3-cd]pyrene | TO-13A (PAH) | SW8270M (SVOC) |
| Naphthalene ^f | TO-13A (PAH) | SW8270M (SVOC) |
| Nitrobenzene | SOP CAD 26.2 (Energetics) | SW8270M (SVOC) |
| Phenanthrene ^g | TO-13A (PAH) | SW8270M (SVOC) |
| Propylene | 40 CFR 60 Method 18 (Hydrocarbons) | TO-14A (VOC) |
| Pyrene ^h | TO-13A (PAH) | SW8270M (SVOC) |
| 1,2,4-Trichlorobenzene ⁱ | TO-14A (VOC) | SW8270M (SVOC) |
| Acetone | TO-11A (Aldehydes) | TO-14A (VOC) |
| Pentane | 40 CFR 60 Method 18 (Hydrocarbons) | TO-14A (VOC) |

- ^a SW8270M data were used to develop the Benzo[a]anthracene emission factor for DODIC M032 because the relative percent difference between test samples as measured using the TO-13A method was greater than 100 percent, indicating poor precision.
- b SW8270M data were used to develop the Benzo[k]fluoranthene emission factor for DODIC M032 because Benzo[k]fluoranthene was detected in the lab blank at a concentration greater than 20 percent using the TO-13A method.
- ^c SW8270M data were used to develop the Benzo[g,h,i]perylene emission factor for DODIC K145 because Benzo[g,h,i]perylene was detected in the lab blank at a concentration greater than 20 percent using the TO-13A method.
- ^d SW8270M data were used to develop the Chrysene emission factor for DODIC M032 because the relative percent difference between test samples as measured using the TO-13A method was greater than 100 percent, indicating poor precision.
- ^e SW8270M data were used to develop the Fluoranthene emission factor for DODIC M032 because the relative percent difference between test samples as measured using the TO-13A method was greater than 100 percent, indicating poor precision.
- SW8270M data were used to develop the Naphthalene emission factor for DODIC L709 because Naphthalene was detected in the lab blank at a concentration greater than 20 percent using the TO-13A method.
- SW8270M data were used to develop the Phenanthrene emission factor for DODIC L709 because Phenanthrene was detected in the lab blank at a concentration greater than 20 percent using the TO-13A method.
- SW8270M data were used to develop the Pyrene emission factor for DODIC L709 because Pyrene was detected in the lab blank at a concentration greater than 20 percent using the TO-13A method. SW8270M data were also used to develop the Pyrene emission factor for DODIC M032 because the relative percent difference between test samples as measured using the TO-13A method was greater than 100 percent, indicating poor precision.
- SW8270M data were used to develop the 1,2,4-Trichlorobenzene for DODICs L709 and M032 because the relative percent difference between test samples as measured using the TO-14A method was greater than 100 percent, indicating poor precision.
- 3.2.2.4 Tentatively Identified Compounds

During the analysis of the SVOC data, the 20 highest nontarget "peaks" were tentatively identified using computerized mass spectral matching techniques. Similarly, during the analysis of the VOC data, the 20 highest nontarget "peaks" were tentatively identified. Emission factors were developed for these tentatively identified compounds (TICs) if all of the following criteria were met.

- 1. The TIC corresponded to a unique compound (e.g., ethylbenzene). Emission factors were not developed if the TIC corresponded to a class of compounds (e.g., unknown alcohol).
- 2. The TIC was not identified using another analysis method that provided higher confidence data. Emission factors were developed based upon the higher confidence analysis method if such data were available.
- 3. The TIC was not present in the method blank. Emission factors were not developed if the TIC was found in the corresponding method blank.

The number of SVOC that were tentatively identified as unique compounds, that were not identified using a higher confidence method, and that were not present in the method blank varied from one to six compounds per ordnance. Similarly, the number of VOC that were tentatively identified as unique compounds, that were not identified using a higher confidence method, and that were not present in the method blank varied from one to eleven compounds per ordnance. Emission factors were developed for all of these TICs, but because of the uncertainty in the true identity of the TICs, the test data were assigned a "C" rating.

3.2.3 Process Information

Ordnance are manufactured to tight tolerances and are expected to deploy in a very repeatable fashion. Consequently, the test data should be assigned an "A" rating based upon this criterion. However, large relative percent differences (i.e., greater than 100 percent) between test runs were noted for several compounds. Specific instances in which these differences were noted are identified in Table 5. The equation below illustrates calculation of relative percent difference:

$$relative\ percent\ difference = \frac{test\ 1\ concentration - test\ 2\ concentration}{average\ of\ test\ 1\ and\ test\ 2\ concentrations} \times 100\%$$

TABLE 5 COMPOUNDS FOR WHICH LARGE RELATIVE PERCENT DIFFERENCES WERE NOTED BETWEEN TEST RUNS

| Compound | Applicable DODIC |
|------------------------------------|------------------|
| 1,2-Dichloroethane | G911 |
| Ethylbenzene | L709 |
| bis(2-Ethylhexyl)phthalate | L709 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | K145, M032 |
| Methyl methacrylate | K145 |
| Nickel | G911 |
| 2,3,4,7,8-Pentachlorodibenzofuran | M032 |
| 1,1,1-Trichloroethane | L709 |

TABLE 5 (cont.)

| Compound | Applicable DODIC | |
|-----------------------|------------------|--|
| Vinyl chloride | M032 | |
| o-Xylene | L709 | |
| 4-Ethyltoluene | L709 | |
| Heptane | K145 | |
| RDX | L709 | |
| 2,4,6-Trinitrotoluene | G911, M032 | |

Due to the large relative percent differences between test runs, the test data specifically identified in Table 5 were assigned a "C" rating. The remainder of the data should be assigned an "A" rating based on this criterion.

3.2.4 Analysis and Calculations

The test report, detailed test plan, and analytical data supporting the test report were reviewed to determine whether they contained all of the original raw data, other documentation, and example calculations. Although the test report did not contain raw field data, the data were made available upon request. The test report also lacked certain calibration data. However, the missing information was judged insufficient to result in a downgrade of the test data quality rating.

The raw data and sample calculations presented in the final test report, detailed test plan and analytical data supporting the test report were reviewed to determine if the emission factors presented in the report could be duplicated. Where differences were found between the emission factors calculated using the Excel spreadsheets and those presented in the test report, an examination was made to determine the reason for the differences.

Several minor errors were noted in the calculation of the emission factors within the test report, particularly with respect to correcting the sample volumes to standard conditions and the incorporation of analytical detection limits into the emission factors (see Section 4.2 for a discussion of the methodology). However, the emission factors presented in AP-42 are based upon the corrected spreadsheets. Based upon the raw data, other documentation, and the Excel spreadsheet calculations, the test data should be assigned an "A" rating.

Emission factors developed for compounds present in the method blank at levels of 20 percent to 50 percent of test values were assumed to be biased high. Five compounds met this criterion. Specifically, methane met the criterion for DODIC G911; 1,2,3,4,6,7,8-heptachlorodibenzofuran met the criterion for DODICs M032 and K145; 1,2,3,4,6,7,8,9-octachlorodibenzofuran met the criterion for DODICs M032 and K145; and 1,2,3,4,6,7,8-heptachlorodibenzo-p-dioxin and 1,2,3,4,6,7,8,9-octachlorodibenzo-p-dioxin met the criterion for DODIC L709. For these compounds, the test data were assigned a "B" rating.

When compounds were found in the method blank at levels greater than 50 percent of test values, the data were assumed to be suspect and no emission factors were developed. Three compounds met this criterion. Specifically, 1,2,3,4,7,8-hexachlorodibenzofuran met the criterion for DODICs G911, M032, and K145; and 1,2,3,4,6,7,8,9-octachlorodibenzofuran and 1,2,3,4,6,7,8-heptachlorodibenzofuran met this criterion for DODIC L709.

3.3 Test Data Quality Ratings

Upon completing the analysis described in the preceding section of this document, the test data quality ratings assigned as a result of the four criteria were reviewed. This review led to a downgrading of some of the test data from an "A" rating to either a "B" rating or a "C" rating. Table 6 identifies the data quality ratings for all compounds that did not receive an "A" rating.

TABLE 6 DOWNGRADED DATA QUALITY RATINGS

| Compound | Data Quality Rating | Applicable DODIC |
|--|------------------------|------------------|
| Carbon dioxide | В | All DODICs |
| Carbon monoxide | В | All DODICs |
| Oxides of nitrogen | В | All DODICs |
| Sulfur dioxide | В | All DODICs |
| Acetophenone | С | K145 |
| 1,2-Dichloroethane | С | G911 |
| Ethylbenzene | С | L709 |
| bis(2-Ethylhexyl)phthalate | С | L709 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | С | L709 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | В | K145, M032 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | С | K145, M032 |
| Methyl methacrylate | С | K145 |
| Nickel | С | G911 |
| 1,2,3,4,6,7,8,9-Octachlorodibenzofuran | В | K145, M032 |
| 1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin | В | L709 |
| 2,3,4,7,8-Pentachlorodibenzofuran | С | M032 |
| 1,1,1-Trichloroethane | С | L709 |
| Vinyl chloride | С | M032 |
| o-Xylene | С | L709 |
| Benzene, (epoxyethyl)-,R- | С | K145 |
| Benzene, (2-nitroethenyl)- | С | K145 |
| Benzene, propyl- | С | K145 |
| Benzofuran | С | K145 |
| Benzoic Acid | С | K145, L709 |
| 1,3-Butadiyne | С | G911, K145 |
| 2-Butenal, (E)- | С | G911 |
| 3-Buten-2-one | С | G911 |
| 1-Buten-3-yne | С | G911 |

TABLE 6 (cont.)

| Compound | Data Quality Rating | Applicable DODIC |
|-------------------------------------|------------------------|------------------|
| Cyclohexanone | С | K145 |
| Docosane | С | L709, M032 |
| 4-Ethyltoluene | С | L709 |
| Furan, 2,5-dihydro- | С | K145 |
| Furan, 2,3-dihydro- | С | K145 |
| Furan, tetrahydro- | С | L709 |
| Heptane | С | K145 |
| 1-Heptene | С | G911 |
| Hexacosane | С | L709 |
| Hexadecane | С | M032 |
| Hexanedioic acid, bis(2-ethylhexyl) | С | K145, M032 |
| Hexatriacontane | С | G911 |
| 1-Hexene | С | G911, K145 |
| Icosane | С | M032 |
| Methane | В | G911 |
| Methane, nitro- | С | G911, K145 |
| Methenamine | С | L709 |
| Phenylethyne | С | K145 |
| 2-Propenal, 2-methyl | С | G911, K145 |
| RDX | С | L709 |
| Silane, difluorodimethyl- | С | M032 |
| 2,4,6-Trinitrotoluene | С | G911, M032 |

4.0 EMISSION FACTOR CALCULATIONS

The methodologies and procedures that were used to develop emission factors from the test data are described in this section. A similar approach was used to calculate emission factors for TSP, PM-10, PM-2.5, metals, SVOC, PAH, and energetic materials. The calculation steps that were performed for each sample and each test run are summarized below.

- 1. The sample volumes associated with the background and test runs were corrected to EPA Stationary Source Program standard temperature and pressure (i.e., 68°F and 29.92 inches of mercury).
- 2. For compounds for which more than one test sample was obtained, analytical detection limits were incorporated into the test data.

- The background compound concentration was calculated by dividing the mass of compound detected during the background run by the background run sample volume corrected to standard conditions.
- 4. The test compound concentration was calculated by dividing the mass of compound detected during the test run by the test run sample volume corrected to standard conditions.
- 5. A background-corrected concentration was calculated by subtracting the background concentration from the test concentration.
- 6. A dilution-corrected concentration was calculated by dividing the background-corrected concentration by the test dilution correction factor.
- 7. The mass of compound released during the test run was calculated by multiplying the dilution-corrected concentration by the volume of the LOTC.
- 8. Emission factors for each sample and test run were calculated by dividing the mass of compound released by the number of ordnance detonated during the test run or by the NEW detonated during the test run, as appropriate.
- 9. Average emission factors were calculated for each compound.

Because concentration data (i.e., milligrams per cubic meter [mg/m³], parts per million by volume [ppmv], or parts per billion by volume [ppbv]) were recorded for VOC, hydrocarbons, FTIR-measured compounds, and CEMS-measured compounds, it was not necessary to calculate a corrected sample volume as described in step 1 or to calculate background and test concentrations as described in steps 3 and 4. Detection limits were applied directly to test compound concentrations of VOC, hydrocarbons, FTIR-measured compounds, and CEMS-measured compounds, as described in step 2. Where present, ppmv and ppbv values were converted to mg/m³. Emission factors for VOC and hydrocarbons were then estimated in accordance with steps 5 through 9 described above. FTIR- and CEMS-measured compound emission factors were estimated in accordance with steps 5 through 9 described above, except that step 6, the dilution correction, was not required. FTIR- and CEMS-measured concentrations were determined from initial steady-state points that occurred prior to significant dilution.

Test and background concentrations of acid gases, hydrogen cyanide, aldehyde, dioxin/furan, and total carbon were provided by the laboratory. The laboratory calculated concentrations based upon the sample volumes as measured at actual conditions. Concentration data received from the laboratory were corrected for volume at EPA Stationary Source Program standard conditions and analytical detection limits were incorporated as described in step 2. Emission factors for acid gases, hydrogen cyanide, aldehydes, and dioxin/furan compounds were then estimated in accordance with steps 5 through 9 described above.

The following sections describe the emission factor calculation steps in more detail. Sections 4.1 through 4.9 discuss the calculations involved with the completion of the nine basic steps listed above.

4.1 Correction of Sample Volume to Standard Conditions

Because sample volumes were initially reported at actual conditions, calculations to correct the sample volumes to the EPA Stationary Source Program standard temperature of 68°F and pressure of 29.92 inches of Hg (20°C and 760 mm Hg) were performed. To obtain the sample volume at EPA Stationary Source Program standard conditions (V_{stat}), the sample volume at actual conditions (V_{act}) was multiplied by the ratio of pressure at actual conditions (P_{act}) to temperature at actual conditions (P_{act}) and by the ratio of temperature at Stationary Source Program standard conditions (P_{stat}). This calculation is illustrated by the following equation:

$$V_{stat} = V_{act} \times \frac{P_{act}}{T_{act}} \times \frac{T_{stat}}{P_{stat}}$$

4.2 Incorporation of Analytical Detection-Limits to the Test Data

In most cases, more than one test sample was obtained for a specific compound (i.e., more than one test run was conducted). When multiple samples were obtained for the same compound, a comparison was made of all the sample data collected. Based upon the results of the comparison, the following adjustments were made to the test data:

- 1. If all of the samples indicated that a compound was "not detected," the sample data were not adjusted.
- 2. If all of the samples indicated that a compound was detected, the sample data were not adjusted.
- 3. If one or more of the samples indicated that a compound was detected and one or more of the samples indicated that a compound was not detected, the "not detected" values were replaced with a value equal to one half of the compound's analytical detection limit. The assumption inherent to this adjustment was that the measured presence of a compound in one or more samples was indicative of the compound's presence in all samples. The analytical detection limits for each sample were obtained from the analytical data supporting the test report.

4.3 Determination of Background Concentration

For TSP, PM-10, PM-2.5, metals, hexavalent chromium, SVOC, and PAH, the background compound concentration (BC) was calculated by dividing the mass of compound detected during the background run (Bkgd mass) by the background run sample volume corrected to EPA Stationary Source Program standard conditions (Bkgd V_{stat}). This calculation is illustrated by the following equation:

$$BC = \frac{Bkgd \ mass}{Bkgd \ V_{stat}}$$

The background compound concentrations of acid gases, hydrogen cyanide, aldehydes, dioxin/furan compounds, and total carbon provided by the lab were calculated using the sample volume at actual conditions (V_{act}). In order to correct the concentrations to EPA Stationary Source Program standard conditions (V_{stat}), the uncorrected concentrations (UC) were multiplied by the ratio of volume at actual conditions to volume at standard conditions. This calculation is illustrated by the following equation:

$$BC = UC \times \frac{V_{act}}{V_{stat}}$$

FTIR- and CEMS-measured compounds were not sampled for during the background runs. However, background data were recorded for each test run between the time the FTIR and CEMS began sampling and the time of detonation. The only FTIR- or CEMS-measured compound for which significant background concentrations were recorded was CO₂. The CO₂ background concentration was assumed to equal a representative value over the sampling period. Background concentrations for the other FTIR- and CEMS-measured compounds were assumed to be zero.

4.4 Determination of Test Compound Concentration

For TSP, PM-10, PM-2.5, metals, SVOC, PAH, and energetic materials, the test compound concentration (TC) was calculated by dividing the mass of compound detected during the test run (Test

mass) by the test run sample volume corrected to EPA Stationary Source Program standard conditions (Test V_{stat}). This calculation is illustrated by the following equation:

$$TC = \frac{Test \, mass}{Test \, V_{stat}}$$

Test compound concentrations of acid gases, hydrogen cyanide, aldehydes, dioxin/furan compounds, and total carbon provided by the lab were calculated using the sample volume at actual conditions. To correct the concentrations to EPA Stationary Source Program standard conditions, the uncorrected concentrations (UC) were multiplied by the ratio of the volume at actual conditions (V_{act}) to the volume at EPA Stationary Source Program standard conditions (V_{stat}). This calculation is illustrated by the following equation:

$$TC = UC \times \frac{V_{act}}{V_{stat}}$$

For VOC, hydrocarbons, FTIR-measured compounds, and CEMS-measured compounds, sample volume correction was not required. VOC and hydrocarbon test concentration values were used directly. Test concentrations of CEMS- and FTIR-measured compounds were determined based upon the initial steady-state values reached after detonation. The time at which a compound's concentration reached an initial steady state was determined by visually inspecting plots of the concentration data versus time. The initial steady state corresponds to the time at which a compound was well mixed within the LOTC, but before the compound's concentration began to decrease at a steady rate due to dilution.

4.5 Determination of Background-Corrected Concentration

For all compounds, the calculation of the background-corrected concentration (BCC) was dependent on whether the background (BC) and test concentrations (TC) were detected and whether they were less than, equal to, or greater than one another. The procedures used to calculate the background-corrected concentration for each sampling train and compound are described below and are displayed graphically in Figure 1.

- 1. If the test concentration was not detected (ND), the background-corrected concentration equaled ND
- 2. If the test concentration was detected and the background concentration was not detected, the background-corrected concentration equaled the test concentration.
- 3. If the test and background concentrations were detected and the test concentration was less than or equal to the background concentration, the background-corrected concentration equaled 0.
- 4. If the test and background concentrations were detected and the background concentration was less than the test concentration, the background concentration, was subtracted from the test concentration. This calculation is illustrated by the following equation:

$$BCC = TC - BC$$

Blasting caps were used in the detonation of DODICs M032 and K145, but are not used when the weapons are tactically detonated. To quantify emissions associated with the blasting caps and thereby

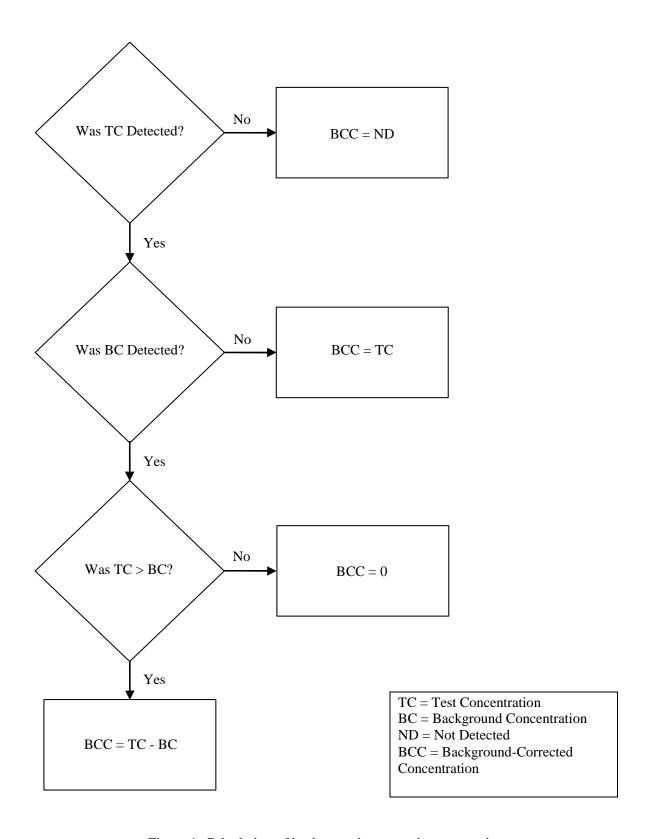


Figure 1 Calculation of background-corrected concentration.

allow the emissions to be subtracted from the resulting emission factors, the blasting caps were detonated during the background runs. As a result, it was necessary to correct these background concentrations for dilution. This was accomplished by multiplying the background concentration by the ratio of test dilution correction factor (DCF_{test}) to background dilution correction factor (DCF_{background}) prior to completing either step 3 or step 4 as described above. The calculation of the background-corrected concentration for those DODICs that were tested with a blasting cap is illustrated by the following equation:

$$BCC = TC - BC \times \frac{DCF_{test}}{DCF_{background}}$$

4.6 Determination of Dilution-Corrected Concentration

The dilution-corrected concentration (DCC) was calculated by dividing the background-corrected concentration by the applicable test dilution correction factor. This calculation is illustrated by the following equation:

$$DCC = \frac{BCC}{DCF_{test}}$$

With the exception of CEMS- and FTIR-measured compounds, dilution correction factors were obtained from the test contractor.³ Because the test concentrations of CEMS-and FTIR-measured compounds were obtained prior to dilution, it was not necessary to calculate a dilution corrected concentration for these compounds.

4.7 Determination of Mass of Compound Released

The mass of compound released was calculated by multiplying the dilution-corrected concentration by the volume of the LOTC. This calculation is illustrated by the following equation:

$$Mass\ compound\ released = DCC\ x\ LOTC\ volume$$

4.8 Determination of Emission Factors

Once the mass of compound released was calculated, two emission factors were developed for each sample and for each test run: the mass of compound released per item (i.e., per single ordnance) and the mass of compound released per pound NEW. The NEW for all ordnance were determined from Reference 1.

4.9 Determination of Average Emission Factors

Steps 1 through 8, as described in Sections 4.1 through 4.8, are applicable to individual samples within individual test runs. The final step in the emission factor calculation process was to calculate average emission factors for each compound in terms of mass released per item and mass released per pound NEW. The average emission factors for each compound were calculated as the arithmetic mean of the individual samples associated with the compound. If all samples indicated that the compound was not detected (ND), then the average emission factor was assigned a value of ND. [Note: The minimum detection levels associated with the compounds that were not detected are presented in Appendix A.] Total dioxin/furan emission factors ere calculated by summing the average emission factors of all dioxin/furan compounds.

Samples of energetic materials were collected using both filters and resin packed cartridges. However, the filter and resin were analyzed separately. Because the filter and resin packed cartridge were arranged in series, average emission factors calculated for the filter were added to average emission factors calculated for resin packed cartridge to determine total emission factors.

5.0 EMISSION FACTOR RATINGS

The emission factors were appraised in accordance with the rating system specified in Reference 4. Under this rating system, emission factors are assigned a rating from A to E, where an "A" rating is assigned to the highest quality factors. The criteria used to assign a specific emission factor rating are summarized below.

- **A** Excellent. The emission factor was developed primarily from A- and B-rated source test data taken from many randomly chosen facilities in the industry population. The source category population was sufficiently specific to minimize variability.
- **B** Above average. The emission factor was developed primarily from A- or B-rated test data from a moderate number of facilities. Although no specific bias was evident, it was not clear if the facilities tested represented a random sample of the industry. As with the "A" rating, the source category population was sufficiently specific to minimize variability.
- C <u>Average</u>. The emission factor was developed primarily from A-, B-, and/or C-rated test data from a reasonable number of facilities. Although no specific bias was evident, it was not clear if the facilities tested represented a random sample of the industry. As with the "A" rating, the source category population was sufficiently specific to minimize variability.
- **D** <u>Below average</u>. The emission factor was developed primarily from A-, B-, and C-rated test data from a small number of facilities, and there may have been reason to suspect that these facilities did not represent a random sample of the industry. There also may have been evidence of variability within the source category population.
- E <u>Poor</u>. The emission factor was developed from C- and D-rated test data from a very limited number of facilities, and there may have been reason to suspect that the facilities tested did not represent a random sample of the industry. There also may have been evidence of variability within the source category population.

Two analyses were conducted to assign ratings to the ordnance emission factors. First, an analysis was conducted on an ordnance-specific basis. Second, an analysis was conducted using all available ordnance emission factor data. The second analysis was conducted to determine whether a sufficient correlation existed between emission factors for different but similar ordnance to allow the number of test data points to be increased to the point that higher emission factor ratings could be assigned than were possible when using the ordnance-specific approach. Both analyses are described below.

5.1 Emission Factor Ratings Assigned – Based on Ordnance-Specific Test Data

As previously described, emission factor ratings are dependent upon the test data quality, the number of test data points, the amount of variability present within a source category population, and the randomness of the source category sample. The following test data facts pertain to these rating criteria:

- 1. As described in Section 3 of this Background Document, the ordnance test data was primarily rated A or B. The test data for a few compounds was rated C.
- 2. Only two tests were conducted per ordnance.

- 3. Ordnance are manufactured to very tight tolerance levels so there is little variability within a specific type of ordnance.
- 4. There was no evidence that suggested the tested items within each type of ordnance were specially selected.

Emission factor ratings were assigned based upon these facts. The rationale used to accept or reject specific emission factor ratings follow.

- A: Rejected. The number of test data points was deemed to be insufficient to assign an A emission factor rating.
- B: Rejected. The number of test data points was deemed to be insufficient to assign a B emission factor rating.
- C: Accepted for most ordnance. The emission factors were developed using A- and B-rated test data, there is little variability among items, and there was no evidence that suggested the tested items were specially selected. Because of the limited number of data points, a C rating was deemed appropriate for this set of circumstances.
- D: Accepted for some ordnance. The emission factors were developed using C-rated test data, there is little variability among items, and there was no evidence that suggested the tested items were specially selected. Because of the limited number of data points, a D rating was deemed appropriate for this set of circumstances.
- E: Rejected. The ordnance described in this report were developed primarily using A- and B-rated test data rather than C- or D-rated data, there is little variability among items, and there was no evidence that suggested the tested items were specially selected. Therefore, an E emission factor rating was deemed inappropriate.

5.2 Emission Factor Ratings Assigned – Based on All Available Test Data

The proceeding sections of this Background Document concern the emission measurement methods, data analysis, and calculations used to develop emission factors for specific ordnance. However, USAEC's ordnance emission factor development program includes more than 200 ordnance that have been tested under more than 25 separate test series. Because many of these ordnance are similar in size and/or chemical composition, a statistical analysis was conducted to assess the similarity of the emission factors developed for similar ordnance. The results of this analysis were used to reevaluate the emission factor ratings assigned on an ordnance-specific basis.

USAEC characterized individual ordnance as falling into one of 17 separate categories, depending upon the size and/or chemical composition of the ordnance. The ordnance and their respective categories are identified in Table 7 along with a comment field describing the number of data points.

TABLE 7 ORDNANCE CATEGORIZATION FOR EMISSION FACTOR CORRELATION ASSESSMENT

| Category | DODIC | Ordnance Description | Test Series | Comment |
|------------|-------------------|---|----------------|-----------------|
| CS | G963 ^a | M73A CS Riot Control Agent Hand Grenade | DPG VI | Data not yet |
| | K765 ^a | CS Riot Control Agent Capsule | DPG VI | available |
| | G900 | TH3 AN-M14 Incendiary Grenade | EO5 | |
| | G911 | MK3A2 Offensive Hand Grenade | EO2 | |
| | G911 | MK3A2 Offensive Hand Grenade | EO6 | |
| | K010 | M4 Field Incendiary Burster | EO5 | |
| | K145 | M18A1 Antipersonnel Mine | EO2 | |
| | M023 | M112 Demolition Block Charge | EO1 | |
| | M030 | 1/4-Pound Demolition Block Charge | EO1 | |
| | M030 | 1/4-Pound Demolition Block Charge | EO3 | |
| | M031 | 1/2-Pound Demolition Block Charge | DPG IVA | |
| | M032 | 1-Pound Demolition Block Charge | EO2 | |
| | M032 | 1-Pound Demolition Block Charge | EO3 | |
| | M130 ^a | M6 Electric Blasting Cap | DPG VII | |
| | M130 ^a | M6 Electric Blasting Cap | EO7 | |
| | M131 ^a | M7 Blasting Cap | EO7 | |
| Demolition | M241 | M10 High Explosive Universal Destructor | DPG IVB | 20+ data points |
| | M456 | PETN Type 1 Detonating Cord | DPG IVA | • |
| | M500 ^a | M21 REEF Line Cutter | FP10 | |
| | M591 | M1 Military Dynamite Demolition Block Charge | EO1 | |
| | M913 | M58A3 Linear Demolition Charge | EO3 | |
| | ML05 ^a | MK24 High Explosive Cutter | EO11 | |
| | ML09 | Linear Demolition Charge, Shaped 20 gr/ft | DPG IVA | |
| | ML15 | Linear Demolition Charge, Shaped 225 gr/ft | DPG IVA | |
| | ML47 ^a | M11 Blasting Cap | EO7 | |
| | MM50 ^a | M221 Shaped Charge | EO11 | |
| | MN02 ^a | M12 Blasting Cap | EO7 | |
| | MN03 ^a | M13 Blasting Cap | EO7 | |
| | MN06 ^a | M14 Blasting Cap | EO7 | |
| | MN07 ^a | M15 Delay Blasting Cap | EO11 | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|--------------|-------------------|---|----------------|------------------|
| | MN08 ^a | M81 Igniter | EO9 | |
| Demolition, | MN68 ^a | M151 Booster Demolition Charge | DPG VIII | 20 . 1-4 |
| continued | None | PAX-11, Granular Powder Burn | EO4 | 20+ data points |
| | None | PAX-11, Molded Pellet Detonation | EO4 | |
| | G878 ^a | M228 Practice Hand Grenade Fuse | DPG VI | |
| | K051 | M604 Anti-Tank Practice Mine Fuze | EO6 | |
| | N278 ^a | M564 MTSQ Fuze | EO11 | |
| Fuze | N285 ^a | M577 Fuze | EO9 | <10 data points |
| Tuze | N286 ^a | M582 Fuze | EO7 | <10 data points |
| | N335 | M557 Point Detonating Fuze | EO5 | |
| | N340 | M739A1 Point Detonating Fuze | EO5 | |
| | N464 ^a | M732 Fuze | EO9 | |
| | G881 | M67 Fragmentation Grenade | EO1 | |
| Grenade | G978 | M82 Smoke Simulant Screening Grenade Launcher | DPG V | <10 data points |
| | G982 | Terephthalic Acid Smoke Hand Grenade | DPG V | <u>,</u> |
| | GG09 ^a | M84 Non-Lethal Stun Hand Grenade | EO12 | |
| | B535 | M583A1 40-mm White Star Parachute Cartridge | DPG IVB | |
| | B536 | M585 40-mm White Star Cluster Cartridge | DPG IVB | |
| | B627 | M83A3, M83A2, & M83A1 60-mm Illuminating Cartridge with Fuze | DPG V | |
| | D505 | M485A2 155-mm Illumination Round (projectile) | DPG I | |
| | L305 | M195 Green Star Parachute Signal Flare | DPG I | |
| Illumination | L306 | M158 Red Star Cluster Signal Illumination | DPG II | 20 L data points |
| Humination | L307 | M159 White Star Cluster Signal Illumination | DPG II | 20+ data points |
| | L311 | M126A1 Red Star Parachute Signal Flare | DPG II | |
| | L312 | M127A1 White Star Parachute Signal Flare | DPG I | |
| | L314 | M125A1 Green Star Cluster Signal Flare | DPG I | |
| | L367 ^a | M22 Anti-Tank, Guided Missile, and Rocket Launching Simulator | DPG VI | |
| | L410 ^a | M206 Aircraft Countermeasure Flare | DPG VI | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|------------|--------------------|---|----------------|---------------------------------------|
| Inert | HA11 ^a | Rocket, 2.75-inch Flechette with M255A1 Warhead | DPG VIII | Data not yet available |
| | C511 | M490 105-mm Target Practice Tracer Cartridge (tracer) | EO6 | |
| Large | C784 ^a | M831 120-mm Target Practice Tracer Cartridge | EO12 | Only 1 data point yet available |
| | C785 ^a | M865 120-mm Target Practice Discarding Sabot Tracer Cartridge | EO12 | avanaore |
| 3.6.11 | BA11 ^a | M1001 40-mm HVCC Cartridge | EO12 | Data not yet |
| Medium | BA15 ^a | M769 60-mm FRP Cartridge | EO12 | available |
| | A652 | M220 20-mm TP-T Cartridge | FP9 | |
| | A940 | M910 25-mm Target Practice Discarding Sabot Tracer Cartridge | FP8 | |
| | A976 | M793 25-mm Target Practice Tracer Cartridge | FP8 | 10+ data points |
| Medium-FP | B129 ^a | M789 30 mm CTG | EO9 | |
| | B519 | M781 40-mm Practice Cartridge | FP2 | |
| | B584 | M918 40-mm Practice Cartridge | FP2 | |
| | B505 ^a | M662 40-mm Red Star Parachute Cartridge | DPG VIII | |
| Mine | K042 ^a | M88 Volcano Practice Canister Mine | DPG VIII | Data not yet available |
| Mortar | CA03 | XM929 120 mm Smoke Cartridge with M7334A1 Fuze | DPG V | Only 1 data point |
| | B542 | M430 40-mm High Explosive Dual Purpose (HEDP) Cartridge (projectile) | EO3 | |
| | B571 | M383 40-mm High Explosive Cartridge (projectile) | EO3 | |
| Projectile | B632 | M49A4 60-mm High Explosive Cartridge (projectile) | EO3 | 10 doto = sinte |
| | B642 | M720 60-mm High Explosive Cartridge (projectile) | EO6 | 10+ data points |
| | BZ-13 ^a | M888 60-mm Cartridge with M935 PD Fuze | EO11 | |
| | C995 | M136 AT4 Recoilless Rifle, 84-mm Cartridge (projectile) | EO3 | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|------------|-------------------|---|----------------|------------------|
| | H557 | M72A3 66-mm High Explosive Antitank Rocket (warhead) | EO1 | |
| Projectile | H708 ^a | M73 35-mm Subcaliber Practice Rocket | DPG VIII | 10 - data mainta |
| continued | None ^a | PAX-21, 60-mm Mortar | EO8 | 10+ data points |
| | PJ02 | FIM-92A Stinger-Basic Guided Missile (warhead) | EO6 | |
| | B642 | M720 60-mm High Explosive Cartridge (propelling charge) | FP4 | |
| | B653 ^a | M766 60-mm Short Range Practice Mortar Cartridge | FP10 | |
| | C226 | M301A3 81-mm Illuminating Cartridge (propelling charge) | FP4 | |
| | C379 | M934 120-mm High Explosive Cartridge (Zone 1 - propelling charge) | FP8 | |
| | C511 | M490 105-mm Target Practice Tracer Cartridge (propelling charge) | FP5 | |
| | C784 | M831 120-mm Target Practice Tracer Cartridge (propelling charge) | FP5 | |
| | C785 | M865 120-mm Target Practice Discarding Sabot Tracer Cartridge (propelling charge) | FP5 | |
| Dan 11- 11 | C868 | M821 81-mm High Explosive Cartridge (propelling charge) | FP4 | 20. 4.4 |
| Propellant | C876 | M880 81-mm Target Practice Short Range Cartridge (propelling charge) | FP4 | 20+ data points |
| | CA09 | M931 120-mm Full Range Practice Cartridge (Zone 1 - propelling charge) | FP8 | |
| | CA09 | M931 120-mm Full Range Practice Cartridge (Zone 4 - propelling charge) | FP8 | |
| | D533 | M119A2 155-mm Propelling Charge (Zone 7) | FP5 | |
| | D540 | M3 155-mm Propelling Charge (Zone 3, M199 Cannon) | FP1 | |
| | D540 | M3 155-mm Propelling Charge (Zone 3, M199 Cannon) | FP5 | |
| | D540 | M3 155-mm Propelling Charge (Zone 3, M284 Cannon) | FP1 | |
| | D540 | M3 155-mm Propelling Charge (Zone 5, M199 Cannon) | FP1 | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|-----------------------|-------------------|---|----------------|-----------------|
| | D540 | M3A1 155-mm Propelling Charge (Zone 3, M199 Cannon) | FP1 | |
| | D540 | M3A1 155-mm Propelling Charge (Zone 3, M284 Cannon) | FP1 | |
| | D541 | M4A2 155-mm Propelling Charge (Zone 7) | FP5 | |
| Dropollant | H557 | M72A3 66-mm High Explosive Antitank Rocket (propelling rocket) | FP7 | |
| Propellant, continued | M174 ^a | MK209 Impulse Cartridge | FP10 | 20+ data points |
| | M842 ^a | M1 Squib | EO7 | |
| | M842 ^a | M79 Igniter | EO9 | |
| | MD73 ^a | M796 Impulse Cartridge | FP10 | |
| | PJ02 | FIM-92A Stinger-Basic Guided Missile (flight motor) | FP7 | |
| | PJ02 | FIM-92A Stinger-Basic Guided Missile (launch motor) | EO5 | |
| | H975 ^a | M274 2.75-inch Signature Smoke with H872 Warhead | DPG VIII | |
| | L366 | M74A1 Projectile Air Burst Simulator | DPG IVB | |
| | L495 | M49A1 Surface Trip Flare | DPG II | |
| | L508 ^a | M72 Red Railroad Warning Fuse | DPG VI | |
| | L592 | TOW Blast Simulator | DPG V | |
| | L594 | M115A2 Ground Burst Simulator | DPG I | |
| | L595 ^a | M9 Liquid Projectile Air Burst Simulator | EO12 | |
| | L596 | M110 Flash Artillery Simulator | DPG I | |
| Pyrotechnic | L598 | M117 Flash Booby Trap Simulator | DPG I | 10+ data points |
| · | L599 | M118 Illuminating Booby Trap Simulator | DPG II | |
| | L600 | M119 Whistling Booby Trap Simulator | DPG II | |
| | L601 | M116A1 Hand Grenade Simulator | DPG I | |
| | L602 | M21 Artillery Flash Simulator | DPG IVB | |
| | L709 | M25 Target Hit Simulator | EO2 | |
| | L709 | XM25 Target Hit Simulator | DPG V | |
| | L720 | M26 Target Kill Simulator | EO6 | |
| | M327 ^a | Coupling Base Firing Device | FP10 | |
| | M448 ^a | M2 Percussion Detonator | EO11 | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|------------------------|-------------------|---|----------------|-----------------|
| | M626 ^a | M1 Push Igniter | EO9 | |
| | M627 ^a | M5 Pressure Release Igniter | FP10 | |
| | M630 | M1 Pull Igniter | DPG V | |
| Pyrotechnic, continued | M630 ^a | M1 Pull Igniter | EO9 | 10+ data points |
| Continued | M670 ^a | M700 Blasting Fuse | EO11 | |
| | M766 ^a | M60 Igniter | EO9 | |
| | ML03 ^a | M142 Firing Device | EO11 | |
| | H459 | Rocket, 2.75-inch Flechette, MK40 Mod 3 Motor (propelling rocket) | FP7 | |
| | H557 | M72A3 66-mm High Explosive Antitank Rocket (propelling rocket) | FP4 | |
| Rocket/Missile | H708 | M73 35-mm Subcaliber Practice Rocket Motor | FP9 | <10 data points |
| | H974 | Rocket, 2.75-inch M267 Practice Warhead, MK66 Mod 3 Motor (propelling rocket) | FP7 | |
| | A010 | M220 10 Gage Blank/Subcaliber Salute Cartridge | FP9 | |
| | A011 | 12 Gage #00 Shot Cartridge | FP9 | |
| | A017 ^a | 12 Gage #9 Shot Cartridge | FP10 | |
| | A059 | M855 5.56-mm Ball Cartridge (fired from the M16A1 Rifle) | FP3 | |
| | A059 | M855 5.56-mm Ball Cartridge (fired from the M16A2 Rifle) | FP3 | |
| | A059 | M855 5.56-mm Ball Cartridge (No-Lead) | FP4 | |
| Small Arm-FP | A063 | M856 5.56-mm Tracer Cartridge | FP3 | 20+ data points |
| | A065 | M862 5.56-mm Practice Ball Cartridge | FP3 | |
| | A066 | M193 5.56-mm Ball Cartridge | FP6 | |
| | A068 | M196 5.56-mm Tracer Cartridge | FP6 | |
| | A080 | M200 5.56-mm Blank Cartridge | FP3 | |
| | A086 | .22 Caliber Long Rifle Ball Cartridge | FP4 | |
| | A106 | .22 Caliber Standard Velocity Long Rifle Ball Cartridge | FP4 | |
| | A111 | M82 7.62-mm Blank Cartridge | FP3 | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|---------------|-------------------|---|----------------|-----------------|
| | A131 | M62 7.62-mm Tracer Cartridge | FP6 | |
| | A136 | M118 7.62-mm Ball Match Cartridge | FP6 | |
| | A143 | M80 7.62-mm Ball Cartridge | FP3 | |
| | A171 | M852 7.62-mm Ball Match Cartridge | FP6 | |
| | A182 | M1 .30 Caliber Ball Cartridge | FP6 | |
| | A212 | M2 .30 Caliber Ball Cartridge | FP6 | |
| | A218 | M25 .30 Caliber Tracer Cartridge | FP9 | |
| | A247 | M72 .30 Caliber Ball Match Cartridge | FP6 | |
| | A363 | M882 9-mm Ball Cartridge | FP3 | |
| Small Arm-FP, | A365 | M181 14.5-mm Trainer-Spotter Cartridge with 3-sec Delay (Artillery) | DPG V | 20+ data points |
| continued | A366 ^a | M182 14.5-mm Cartridge | EO11 | _ |
| | A400 | M41 .38 Caliber Special Ball Cartridge | FP9 | |
| | A403 | .38 Caliber Special Blank Cartridge | FP9 | |
| | A475 | M1911 .45 Caliber Ball Cartridge | FP3 | |
| | A518 | M903 .50 Caliber SLAP Cartridge | FP9 | |
| | A518 ^a | M962 .50 Caliber SLAP/T | FP10 | |
| | A525 | M2 .50 Caliber Armor Piercing Cartridge | FP8 | |
| | A557 | M17 .50 Caliber Tracer Cartridge | FP3 | |
| | A557 | M33 .50 Caliber Ball Cartridge | FP3 | |
| | A598 | M1A1 .50 Caliber Blank Cartridge | FP3 | |
| | C870 ^a | M819 (IUK) 81-mm Red Phosphorous Smoke | DPG VII | |
| | G815 ^a | Red Phosphorous Smoke Screening Grenade Launcher (UK) | DPG VII | |
| | G930 | Hexachloroethane Smoke Grenade | DPG V | |
| Smoke | G940 | M18 Green Smoke Hand Grenade | DPG III | |
| | G945 | M18 Yellow Smoke Hand Grenade | DPG III | 10+ data points |
| | G950 | M18 Red Smoke Hand Grenade | DPG III | |
| | G950 | M18 Red Smoke Hand Grenade (new formulation) | DPG V | |
| | G955 | M18 Violet Smoke Hand Grenade | DPG III | |
| | G955 | M18 Violet Smoke Hand Grenade (new formulation) | DPG V | |

TABLE 7 (cont.)

| Category | DODIC | Ordnance Description | Test Series | Comment |
|-----------|-------------------|------------------------------------|----------------|-----------------|
| Smoke, | K866 ^a | ABC-M5 HC Ground Smoke Pot (MILES) | DPG VII | 10+ data points |
| continued | K867 ^a | M4A2 Floating Smoke Pot | DPG VII | 10+ data points |

^a Although testing may have been completed, emission factors for this ordnance have not yet been analyzed for inclusion in AP-42; therefore, these data were not included when the data correlation was assessed.

Within each of the 17 ordnance categories identified by USAEC, emission factors for each compound were compared. To allow the comparison of emission factors for ordnance with similar constituents but significant differences in net explosive weight, the comparison was made using the normalized emission factor units of mass of compound released per pound NEW. Based upon information provided by EPA, 8 the following procedures were used to assess the data correlation:

- 1. The relative standard deviation, defined as the standard deviation divided by the mean, was calculated for each compound within each ordnance category.
- 2. If the relative standard deviation was less than 1.0, the evaluated emission factors were considered to demonstrate good correlation. As such, the rating for these emission factors <u>could</u> be elevated to a maximum of an A, depending on the number of data points within the evaluated ordnance category.
- 3. If the relative standard deviation was between 1.0 and 2.0, the evaluated emission factors were considered to demonstrate fair correlation. As such, the rating for these emission factors <u>could</u> be elevated to a maximum of a B, depending on the number of data points within the evaluated ordnance category.
- 4. If the relative standard deviation was greater than 2.0, the evaluated emission factors were considered to demonstrate poor correlation. As such, the emission factor rating could not be elevated, regardless of the amount of data available.

A poor correlation between emission factors was not necessarily construed as being indicative of poor test data. Rather, a poor correlation was more likely to indicate that the ordnance included in the category were not as similar in nature as anticipated by USAEC when the ordnance categories were defined.

In addition to assessing the data correlation, an assessment was made of the number of test data points available within each of the 17 ordnance categories. Because each ordnance test consisted of two test data points (i.e., two test runs per ordnance or two independent sampling trains were used during an ordnance test), the number of test data points available in each of the ordnance categories varied from 2 to 68. Based upon information provided by EPA, the following assumptions were used to assess whether sufficient category-specific test data points were available to justify elevating the emission factor ratings based on ordnance-specific data only:

- 1. If 20 or more data points were available, the emission factor rating could be elevated to a maximum of an A, provided that the data also demonstrated a good correlation.
- 2. If at least 10 but less than 20 data points were available, the emission factor rating could be elevated to a maximum of a B, provided that the data also demonstrated a good correlation.
- 3. If less than 10 data points were available, the emission factor rating could not be elevated, regardless of the data correlation.

- 4. If the data demonstrated a fair correlation and 20 or more data points were available, the emission factor rating could be elevated to a maximum of a B.
- 5. If the data demonstrated a fair correlation and at least 10 but less than 20 data points were available, the emission factor rating could be elevated to a maximum of a C.

Using the criteria specified above, the emission factor ratings assigned to ordnance in each of the 17 ordnance categories were reevaluated. This evaluation indicated that some of the emission factor ratings associated with ordnance included in eight categories could be elevated from a C or D rating to an A or B rating. These eight categories are:

- 1. Demolition
- 2. Illumination
- 3. Medium Firing Point
- 4. Projectiles
- 5. Propellants
- 6. Pyrotechnics
- 7. Small Arms Firing Point
- 8. Smokes

A final assessment was made as to the emission factor rating assigned based on ordnance-specific test data only. If the original emission factor data rating assigned was a C, then the emission factor rating was elevated to an A or B, as appropriate, based upon the data for the whole ordnance category. If the original emission factor data rating assigned was a D, then the emission factor rating was elevated to a B or C, as appropriate, based upon the data for the whole ordnance category. The analysis is documented in an Excel spreadsheet that is located on the EPA website at: http://www.epa.gov/ttn/chief/ap42/index.html.

Within the current test series, DODICS G911, K145 and M032 were included in the Demolition category, which included more than 20 test data points. DODIC L709 was included in the Pyrotechnics category, which included more than 10 test data points. As a result, some emission factor ratings associated with each of these ordnance were elevated. The emission factor ratings assigned are presented in Appendix A.

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- 7. National Institute for Occupational Safety and Health (NIOSH) Manual of Analytical Methods (NMAM), Fourth Edition, National Institute for Occupational Safety and Health, Cincinnati, OH, 2003.
- 8. Information regarding the relationship between emission factor data correlation, the number of data points available, and the resulting emission factor rating assigned supplied upon request by Mr. Ron Myers, Measurement Policy Group, Office of Air Quality Planning and Standards, U.S. Environmental Protection Agency, Research Triangle Park, NC, June 2006.

APPENDIX A

COMPOUNDS ANALYZED AND EMISSION FACTORS DEVELOPED FOR ORDNANCE INCLUDED IN EXPLODING ORDNANCE EMISSION STUDY PHASE II SERIES 2 TESTING AT ABERDEEN PROVING GROUND, MARYLAND

TABLE A1 COMPOUNDS ANALYZED AND EMISSION FACTORS DEVELOPED FOR DODIC G911, MK3A2 OFFENSIVE HAND GRENADE

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|---|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| (| Carbon Dioxide, Criteria Pollutants, Me | thane, and Total | Suspended Parti | culate |
| 124-38-9 | Carbon dioxide ^f | 6.9 E-01 | 1.4 | |
| 630-08-0 | Carbon monoxide ^g | 1.5 E-02 | 3.0 E-02 | |
| 7439-92-1 | Lead | 3.3 E-04 | 6.7 E-04 | |
| 74-82-8 | Methane ^g | 2.7 E-04 | 5.5 E-04 | |
| | Oxides of nitrogen ^f | 7.1 E-03 | 1.4 E-02 | |
| | PM-2.5 ^g | 2.0 E-02 | 4.0 E-02 | |
| | PM-10 | 3.5 E-02 | 7.0 E-02 | |
| 7446-09-5 | Sulfur dioxide | ND | ND | 5.3 E-01 |
| 12789-66-1 | Total suspended particulate | 3.9 E-02 | 7.9 E-02 | |
| | Hazardous Air Pollutant | s and Toxic Che | micals | |
| 83-32-9 | Acenaphtheneg | 8.7 E-08 | 1.8 E-07 | |
| 208-96-8 | Acenaphthyleneg | 1.1 E-06 | 2.3 E-06 | |
| 75-07-0 | Acetaldehyde ^g | 7.2 E-05 | 1.5 E-04 | |
| 75-05-8 | Acetonitrile ^g | 2.1 E-05 | 4.2 E-05 | |
| 107-02-8 | Acrolein | ND | ND | 1.9 E-01 |
| 107-13-1 | Acrylonitrile ^g | 3.9 E-06 | 7.8 E-06 | |
| 107-05-1 | Allyl chloride | ND | ND | 8.9 E-04 |
| 7429-90-5 | Aluminum ^g | 3.1 E-04 | 6.2 E-04 | |
| 7664-41-7 | Ammonia | ND | ND | 2.1 |
| 120-12-7 | Anthraceneg | 1.5 E-07 | 3.0 E-07 | |
| 7440-36-0 | Antimony | ND | ND | 2.1 E-02 |
| 7440-38-2 | Arsenic | ND | ND | 2.1 E-02 |
| 7440-39-3 | Barium | 3.3 E-04 | 6.6 E-04 | |
| 71-43-2 | Benzene ^g | 3.7 E-05 | 7.5 E-05 | |
| 56-55-3 | Benzo[a]anthracene | 1.1 E-07 | 2.2 E-07 | |
| 205-99-2 | Benzo[b]fluoranthene | 8.0 E-08 | 1.6 E-07 | |
| 207-08-9 | Benzo[k]fluoranthene | 6.6 E-08 | 1.3 E-07 | |
| 191-24-2 | Benzo[g,h,i]perylene | 8.8 E-08 | 1.8 E-07 | |
| 50-32-8 | Benzo[a]pyrene | 8.1 E-08 | 1.6 E-07 | |
| 192-97-2 | Benzo[e]pyrene | 1.2 E-07 | 2.4 E-07 | |
| 100-44-7 | Benzyl chloride | ND | ND | 2.6 E-04 |
| 7440-41-7 | Beryllium | ND | ND | 2.1 E-02 |

TABLE A1 (cont.)

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|--------------------------------|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 75-25-2 | Bromoform | ND | ND | 5.3 E-04 |
| 74-83-9 | Bromomethane | ND | ND | 3.6 E-04 |
| 101-55-3 | 4-Bromophenylphenylether | ND | ND | 3.1 E-02 |
| 106-99-0 | 1,3-Butadiene ^g | ND | ND | 4.5 E-02 |
| 75-65-0 | t-Butyl alcohol | ND | ND | 4.0 E-04 |
| 85-68-7 | Butylbenzylphthalatef | 1.7 E-06 | 3.4 E-06 | |
| 123-72-8 | Butyraldehyde | ND | ND | 1.9 E-01 |
| 7440-43-9 | Cadmium | 1.8 E-05 | 3.6 E-05 | |
| 75-15-0 | Carbon disulfide | 1.7 E-06 | 3.4 E-06 | |
| 56-23-5 | Carbon tetrachloride | ND | ND | 3.2 E-04 |
| 106-47-8 | 4-Chloroaniline | ND | ND | 3.1 E-02 |
| 108-90-7 | Chlorobenzene | ND | ND | 2.4 E-04 |
| 75-45-6 | Chlorodifluoromethane | ND | ND | 2.5 E-04 |
| 75-00-3 | Chloroethaneg | ND | ND | 6.2 E-04 |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | ND | 1.6 E-02 |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | ND | 1.6 E-02 |
| 67-66-3 | Chloroform | ND | ND | 4.0 E-04 |
| 74-87-3 | Chloromethane | 1.5 E-07 | 3.0 E-07 | |
| 91-58-7 | 2-Chloronaphthalene | ND | ND | 1.6 E-02 |
| 95-57-8 | 2-Chlorophenol | ND | ND | 1.6 E-02 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | ND | ND | 1.6 E-02 |
| 7440-47-3 | Chromium ^g | 2.8 E-05 | 5.6 E-05 | |
| 18540-29-9 | Hexavalent chromium | 8.2 E-07 | 1.7 E-06 | |
| 218-01-9 | Chrysene | 1.3 E-07 | 2.6 E-07 | |
| 7440-48-4 | Cobalt ^g | ND | ND | 2.1 E-02 |
| 7440-50-8 | Copper ^g | 2.3 E-04 | 4.6 E-04 | |
| 106-44-5 | p-Cresol | ND | ND | 1.6 E-02 |
| 4170-30-3 | Crotonaldehyde | ND | ND | 1.9 E-01 |
| 98-82-8 | Cumene ^g | ND | ND | 2.0 E-04 |
| 57-12-5 | Particulate cyanide | ND | ND | 1.7 E-02 |
| 53-70-3 | Dibenz[a,h]anthracene | 3.0 E-08 | 6.1 E-08 | |
| 132-64-9 | Dibenzofuran | ND | ND | 1.6 E-02 |
| 106-93-4 | 1,2-Dibromoethane | ND | ND | 3.9 E-04 |
| 84-74-2 | Dibutyl phthalate ^f | 3.2 E-06 | 6.4 E-06 | |

TABLE A1 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 95-50-1 | 1,2-Dichlorobenzene | ND | ND | 4.9 E-04 |
| 541-73-1 | 1,3-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 106-46-7 | 1,4-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | ND | 1.6 E-02 |
| 75-27-4 | Dichlorobromomethane | ND | ND | 4.1 E-04 |
| 75-71-8 | Dichlorodifluoromethaneg | 0 | 0 | |
| 75-34-3 | 1,1-Dichloroethane | ND | ND | 3.7 E-04 |
| 107-06-2 | 1,2-Dichloroethane ⁱ | 3.7 E-07 | 7.4 E-07 | |
| 75-43-4 | Dichlorofluoromethane | ND | ND | 2.1 E-04 |
| 120-83-2 | 2,4-Dichlorophenol | ND | ND | 1.6 E-02 |
| 78-87-5 | 1,2-Dichloropropane | ND | ND | 3.3 E-04 |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | ND | 2.3 E-04 |
| 76-14-2 | Dichlorotetrafluoroethane | ND | ND | 5.0 E-04 |
| 105-67-9 | 2,4-Dimethylphenol | ND | ND | 1.6 E-02 |
| 131-11-3 | Dimethyl phthalate | ND | ND | 1.6 E-02 |
| 99-65-0 | 1,3-Dinitrobenzene | ND | ND | 6.5 E-03 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | ND | ND | 1.6 E-02 |
| 51-28-5 | 2,4-Dinitrophenol | ND | ND | 1.6 E-02 |
| 121-14-2 | 2,4-Dinitrotoluene | 1.4 E-05 | 2.8 E-05 | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | ND | 1.3 E-03 |
| 123-91-1 | 1,4-Dioxane | ND | ND | 3.3 E-04 |
| | Total dioxin/furan compounds | 2.9 E-10 | 5.9 E-10 | |
| 140-88-5 | Ethyl acrylate | ND | ND | 1.7 E-04 |
| 100-41-4 | Ethylbenzene ^g | 8.5 E-07 | 1.7 E-06 | |
| 74-85-1 | Ethylene ^g | 3.0 E-04 | 6.1 E-04 | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate ^f | 6.5 E-06 | 1.3 E-05 | |
| 206-44-0 | Fluorantheneg | 3.2 E-07 | 6.4 E-07 | |
| 86-73-7 | Fluorene ^g | 2.0 E-07 | 4.0 E-07 | |
| 50-00-0 | Formaldehyde | 4.3 E-05 | 8.6 E-05 | |
| 76-13-1 | Freon 113 ^g | ND | ND | 5.5 E-04 |
| 35822-46-9 | 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 2.1 E-11 | 4.2 E-11 | |
| 67562-39-4 | 1,2,3,4,6,7,8- Heptachlorodibenzofuran | 2.8 E-12 | 5.6 E-12 | |

TABLE A1 (cont.)

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|---|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 55673-89-7 | 1,2,3,4,7,8,9- Heptachlorodibenzofuran | ND | ND | 3.4 E-09 |
| 118-74-1 | Hexachlorobenzene | ND | ND | 1.6 E-02 |
| 87-68-3 | Hexachlorobutadiene | ND | ND | 7.6 E-04 |
| 77-47-4 | Hexachlorocyclopentadiene | ND | ND | 1.6 E-02 |
| 39227-28-6 | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | ND | ND | 3.8 E-09 |
| 57653-85-7 | 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 5.4 E-13 | 1.1 E-12 | |
| 19408-74-3 | 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 8.5 E-13 | 1.7 E-12 | |
| 57117-44-9 | 1,2,3,6,7,8-Hexachlorodibenzofuran | ND | ND | 2.2 E-09 |
| 72918-21-9 | 1,2,3,7,8,9-Hexachlorodibenzofuran | ND | ND | 3.2 E-09 |
| 60851-34-5 | 2,3,4,6,7,8-Hexachlorodibenzofuran | ND | ND | 2.9 E-09 |
| 67-72-1 | Hexachloroethane | ND | ND | 5.9 E-04 |
| 110-54-3 | Hexane | ND | ND | 7.2 E-02 |
| 7647-01-0 | Hydrochloric acid | ND | ND | 2.1 E-01 |
| 74-90-8 | Hydrogen cyanide | 2.6 E-04 | 5.2 E-04 | |
| 7664-39-3 | Hydrogen fluoride | ND | ND | 2.1 E-01 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 4.1 E-08 | 8.4 E-08 | |
| 78-59-1 | Isophorone | ND | ND | 1.6 E-02 |
| 7439-92-1 | Lead | 3.3 E-04 | 6.7 E-04 | |
| 7439-96-5 | Manganese ^g | 6.2 E-05 | 1.3 E-04 | |
| 96-33-3 | Methyl acrylate | ND | ND | 2.5 E-04 |
| 74-95-3 | Methylene bromide | ND | ND | 5.1 E-04 |
| 75-09-2 | Methylene chloride ^g | 9.4 E-07 | 1.9 E-06 | |
| 78-93-3 | Methyl ethyl ketone ^f | 1.6 E-06 | 3.2 E-06 | |
| 74-88-4 | Methyl iodide | ND | ND | 3.5 E-04 |
| 108-10-1 | Methyl isobutyl ketone | ND | ND | 3.3 E-04 |
| 80-62-6 | Methyl methacrylate | ND | ND | 4.6 E-04 |
| 91-57-6 | 2-Methylnaphthalene | ND | ND | 1.6 E-02 |
| 95-48-7 | 2-Methylphenol | ND | ND | 1.6 E-02 |
| 1634-04-4 | Methyl tert-butyl ether | ND | ND | 3.7 E-04 |
| 91-20-3 | Naphthalene ^g | 2.2 E-06 | 4.4 E-06 | |
| 7440-02-0 | Nickel ^h | 1.5 E-05 | 3.0 E-05 | |

TABLE A1 (cont.)

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|--|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 7697-37-2 | Nitric acid ^g | 1.1 E-04 | 2.2 E-04 | |
| 100-01-6 | 4-Nitroaniline | ND | ND | 1.6 E-02 |
| 98-95-3 | Nitrobenzene | ND | ND | 6.5 E-04 |
| 55-63-0 | Nitroglycerin | ND | ND | 6.5 E-04 |
| 88-75-5 | 2-Nitrophenol | ND | ND | 1.6 E-02 |
| 100-02-7 | 4-Nitrophenol | ND | ND | 1.6 E-02 |
| 62-75-9 | n-Nitrosodimethylamine | ND | ND | 1.6 E-02 |
| 86-30-6 | n-Nitrosodiphenylamine | ND | ND | 1.6 E-02 |
| 621-64-7 | n-Nitroso-di-n-propylamine | ND | ND | 1.6 E-02 |
| 3268-87-9 | 1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin | 2.6 E-10 | 5.2 E-10 | |
| 39001-02-0 | 1,2,3,4,6,7,8,9- Octachlorodibenzofuran | 8.7 E-12 | 1.8 E-11 | |
| 40321-76-4 | 1,2,3,7,8-Pentachlorodibenzo-p- dioxin | ND | ND | 4.0 E-09 |
| 57117-41-6 | 1,2,3,7,8-Pentachlorodibenzofuran | ND | ND | 3.3 E-09 |
| 57117-31-4 | 2,3,4,7,8-Pentachlorodibenzofuran | ND | ND | 1.8 E-09 |
| 87-86-5 | Pentachlorophenol | ND | ND | 1.6 E-02 |
| 85-01-8 | Phenanthrene ^g | 7.1 E-07 | 1.4 E-06 | |
| 108-95-2 | Phenol | 1.7 E-06 | 3.5 E-06 | |
| 123-38-6 | Propionaldehyde | ND | ND | 1.9 E-01 |
| 115-07-1 | Propylene ^g | 7.5 E-05 | 1.5 E-04 | |
| 129-00-0 | Pyrene ^g | 4.8 E-07 | 9.6 E-07 | |
| 7782-49-2 | Selenium | ND | ND | 2.1 E-02 |
| 7440-22-4 | Silver | ND | ND | 2.1 E-02 |
| 100-42-5 | Styrene | ND | ND | 3.0 E-04 |
| 7664-93-9 | Sulfuric acid | 3.7 E-05 | 7.5 E-05 | |
| 1746-01-6 | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | ND | ND | 2.4 E-09 |
| 51207-31-9 | 2,3,7,8-Tetrachlorodibenzofuran | 1.3 E-12 | 2.6 E-12 | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | ND | 3.5 E-04 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | ND | 3.5 E-04 |
| 127-18-4 | Tetrachloroethylene | ND | ND | 4.1 E-04 |
| 7440-28-0 | Thallium | ND | ND | 2.1 E-02 |
| 108-88-3 | Toluene ^g | 8.4 E-06 | 1.7 E-05 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | ND | 1.3 E-03 |

TABLE A1 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|-----------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 71-55-6 | 1,1,1-Trichloroethane | ND | ND | 4.4 E-04 |
| 79-00-5 | 1,1,2-Trichloroethane | ND | ND | 2.8 E-04 |
| 75-69-4 | Trichlorofluoromethane | ND | ND | 3.4 E-04 |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | ND | 1.6 E-02 |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | ND | 1.6 E-02 |
| 96-18-4 | 1,2,3-Trichloropropane | ND | ND | 4.9 E-04 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 4.4 E-07 | 8.9 E-07 | |
| 540-84-1 | 2,2,4-Trimethylpentane ^g | ND | ND | 1.9 E-04 |
| 7440-62-2 | Vanadium | 3.9 E-03 | 8.0 E-03 | |
| 108-05-4 | Vinyl acetate | ND | ND | 3.9 E-04 |
| 75-01-4 | Vinyl chloride | ND | ND | 2.6 E-04 |
| 75-35-4 | Vinylidene chloride | ND | ND | 2.1 E-04 |
| 106-42-3, 108-38-3 | m-Xylene, p-Xylene ^g | 7.7 E-07 | 1.6 E-06 | |
| 95-47-6 | o-Xylene ^g | 2.3 E-07 | 4.7 E-07 | |
| 7440-66-6 | Zinc ^g | 6.6 E-03 | 1.3 E-02 | |
| | Other Po | llutants | | |
| 67-64-1 | Acetone | 1.3 E-05 | 2.6 E-05 | |
| 74-86-2 | Acetylene ^g | 1.8 E-04 | 3.7 E-04 | |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene ^g | 4.7 E-07 | 9.4 E-07 | |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 3.3 E-07 | 6.7 E-07 | |
| 100-52-7 | Benzaldehyde | ND | ND | 1.9 E-01 |
| 100-51-6 | Benzyl alcohol | ND | ND | 1.6 E-02 |
| 108-86-1 | Bromobenzene | ND | ND | 3.9 E-04 |
| 460-12-8 | 1,3-Butadiyne ⁱ | 1.8 E-06 | 3.7 E-06 | |
| 106-97-8 | Butane | ND | ND | 5.3 E-02 |
| 123-73-9 | 2-Butenal, (E)- ⁱ | 1.5 E-06 | 3.1 E-06 | |
| 106-98-9 | 1-Butene | 1.9 E-05 | 3.8 E-05 | |
| 590-18-1 | cis-Butene | ND | ND | 4.7 E-02 |
| 624-64-6 | trans-Butene ^g | ND | ND | 4.7 E-02 |
| 78-94-4 | 3-Buten-2-one ⁱ | 3.6 E-06 | 7.2 E-06 | |
| 689-97-4 | 1-Buten-3-yne ⁱ | 1.3 E-06 | 2.7 E-06 | |
| 107-00-6 | 1-Butyne | ND | ND | 4.5 E-02 |
| 503-17-3 | 2-Butyne | ND | ND | 4.5 E-02 |
| 7440-70-2 | Calcium ^f | 4.8 E-04 | 9.6 E-04 | |

TABLE A1 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|----------------------------------|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 7440-44-0 | Total carbon ^g | 6.4 E-03 | 1.3 E-02 | |
| 39638-32-9 | bis(2-Chloroisopropyl)ether | ND | ND | 1.6 E-02 |
| 59-50-7 | 4-Chloro-3-methylphenol | ND | ND | 1.6 E-02 |
| 124-48-1 | Dibromochloromethane | ND | ND | 5.2 E-04 |
| 156-59-2 | cis-1,2-Dichloroethene | ND | ND | 2.0 E-04 |
| 156-60-5 | trans-1,2-Dichloroethene | ND | ND | 4.8 E-04 |
| 84-66-2 | Diethylphthalate | ND | ND | 1.6 E-02 |
| 5779-94-2 | 2,5-Dimethylbenzaldehyde | 0 | 0 | |
| 117-84-0 | Di-n-octylphthalate ^g | ND | ND | 1.6 E-02 |
| 74-84-0 | Ethane ^g | 1.8 E-05 | 3.7 E-05 | |
| 141-78-6 | Ethyl acetate | ND | ND | 1.4 E-03 |
| 97-63-2 | Ethyl methacrylate | ND | ND | 3.3 E-04 |
| 622-96-8 | 4-Ethyltoluene | 0 | 0 | |
| 142-82-5 | Heptane | 6.6 E-07 | 1.3 E-06 | |
| 592-76-7 | 1-Heptene ⁱ | 6.1 E-06 | 1.2 E-05 | |
| 66-25-1 | Hexaldehyde | ND | ND | 1.9 E-01 |
| 591-78-6 | 2-Hexanone | ND | ND | 2.9 E-04 |
| 630-06-8 | Hexatriacontane ⁱ | 7.5 E-06 | 1.5 E-05 | |
| 2691-41-0 | HMX | ND | ND | 1.3 E-03 |
| 592-41-6 | 1-Hexene ⁱ | 1.3 E-05 | 2.6 E-05 | |
| 10035-10-6 | Hydrogen bromide | ND | ND | 2.0 E-01 |
| 75-28-5 | Isobutane | ND | ND | 4.8 E-02 |
| 590-86-3 | Isovaleraldehyde | ND | ND | 1.9 E-01 |
| 7439-95-4 | Magnesium ^g | 2.9 E-03 | 5.9 E-03 | |
| 75-52-5 | Methane, nitro-h | 2.2 E-06 | 4.5 E-06 | |
| 98-83-9 | alpha-Methylstyrene | ND | ND | 2.9 E-04 |
| 88-74-7 | 2-Nitroaniline | ND | ND | 1.6 E-02 |
| 99-09-2 | 3-Nitroaniline | ND | ND | 1.6 E-02 |
| 88-72-2 | 2-Nitrotoluene | ND | ND | 6.5 E-04 |
| 99-08-1 | 3-Nitrotoluene | ND | ND | 6.5 E-04 |
| 99-99-0 | 4-Nitrotoluene | ND | ND | 6.5 E-04 |
| 111-65-9 | Octane | 6.6 E-07 | 1.3 E-06 | |
| 109-66-0 | Pentane | ND | ND | 6.0 E-02 |
| 7664-38-2 | Phosphoric acid | ND | ND | 2.0 E-01 |

TABLE A1 (cont.)

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|-------------------------------------|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 74-98-6 | Propane | ND | ND | 3.7 E-02 |
| 78-85-3 | 2-Propenal, 2-methyl- ⁱ | 2.0 E-06 | 4.1 E-06 | |
| 115-11-7 | 1-Propene, 2-methyl- | 1.3 E-05 | 2.6 E-05 | |
| 74-99-7 | Propyne ^g | 1.3 E-05 | 2.6 E-05 | |
| 121-82-4 | RDX | 7.7 E-06 | 1.6 E-05 | |
| 7440-24-6 | Strontium | 2.9 E-03 | 5.9 E-03 | |
| 479-45-8 | Tetryl | ND | ND | 6.5 E-04 |
| 7440-32-6 | Titanium | 1.8 E-05 | 3.6 E-05 | |
| 1334-78-7 | o,m,p-Tolualdehyde | ND | ND | 5.6 E-01 |
| 25323-89-1 | Trichloroethane | ND | ND | 3.0 E-04 |
| 108-67-8 | 1,3,5-Trimethylbenzene ^g | 0 | 0 | |
| 99-35-4 | 1,3,5-Trinitrobenzene ^g | 4.2 E-06 | 8.5 E-06 | |
| 118-96-7 | 2,4,6-Trinitrotoluene ⁱ | 2.7 E-05 | 5.5 E-05 | |
| 110-62-3 | Valeraldehyde | ND | ND | 1.9 E-01 |

^a CASRN = Chemical Abstracts Service Registry Number.

^b ND = nondetected.

^c Emission factors rated C unless otherwise noted.

^d NEW = net explosive weight. The NEW for this ordnance is 4.94 E-01 pounds per item.

^e Data provided for compounds that were not detected.

Emission factor rated A because of correlation with emission factors for similar ordnance and number of test data points.

g Emission factor rated B because of correlation with emission factors for similar ordnance and number of test data points.

^h Emission factor based upon C-rated test data, but because of correlation with emission factors for similar ordnance and number of data points the factor was upgraded from a D rating to a C rating.

ⁱ Emission factor rated D because the factor is based upon C-rated test data.

TABLE A2 COMPOUNDS ANALYZED AND EMISSION FACTORS DEVELOPED FOR DODIC K145, M18A1 ANTIPERSONNEL MINE

| | | Emission | Factor ^{b,c} | Detection Limit |
|--------------------|---|------------------|-------------------------------|---|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Based Concentration mg/m ^{3,e} |
| (| Carbon Dioxide, Criteria Pollutants, Me | thane, and Total | Suspended Parti | culate |
| 124-38-9 | Carbon dioxide ^f | 1.6 | 1.0 | |
| 630-08-0 | Carbon monoxide ^g | 2.0 E-02 | 1.3 E-02 | |
| 7439-92-1 | Lead | 5.7 E-05 | 3.8 E-05 | |
| 74-82-8 | Methane ^g | 3.8 E-04 | 2.5 E-04 | |
| | Oxides of nitrogen ^f | 1.8 E-02 | 1.2 E-02 | |
| | PM-2.5 ^g | 2.6 E-02 | 1.7 E-02 | |
| | PM-10 | 4.9 E-02 | 3.2 E-02 | |
| 7446-09-5 | Sulfur dioxide | 9.1 E-05 | 6.1 E-05 | |
| 12789-66-1 | Total suspended particulate | 5.4 E-02 | 3.6 E-02 | |
| | Hazardous Air Pollutant | s and Toxic Che | micals | |
| 83-32-9 | Acenaphtheneg | 6.3 E-08 | 4.2 E-08 | |
| 208-96-8 | Acenaphthyleneg | 1.6 E-06 | 1.1 E-06 | |
| 75-07-0 | Acetaldehyde ^g | 5.4 E-05 | 3.6 E-05 | |
| 75-05-8 | Acetonitrile ^g | 4.5 E-05 | 3.0 E-05 | |
| 98-86-2 | Acetophenonei | 2.6 E-06 | 1.8 E-06 | |
| 107-02-8 | Acrolein | ND | ND | 1.8 E-01 |
| 107-13-1 | Acrylonitrileg | 9.7 E-06 | 6.5 E-06 | |
| 107-05-1 | Allyl chloride | ND | ND | 8.9 E-04 |
| 7429-90-5 | Aluminum ^g | 6.3 E-04 | 4.2 E-04 | |
| 7664-41-7 | Ammonia | ND | ND | 2.1 |
| 120-12-7 | Anthracene ^g | 2.0 E-07 | 1.3 E-07 | |
| 7440-36-0 | Antimony | ND | ND | 2.0 E-02 |
| 7440-38-2 | Arsenic | ND | ND | 2.0 E-02 |
| 7440-39-3 | Barium | 2.4 E-04 | 1.6 E-04 | |
| 71-43-2 | Benzene ^g | 1.3 E-04 | 8.9 E-05 | |
| 56-55-3 | Benzo[a]anthracene | 5.1 E-08 | 3.4 E-08 | |
| 205-99-2 | Benzo[b]fluoranthene | 2.6 E-08 | 1.7 E-08 | |
| 207-08-9 | Benzo[k]fluoranthene | 2.6 E-08 | 1.7 E-08 | |
| 191-24-2 | Benzo[g,h,i]perylene | ND | ND | 1.6 E-02 |
| 50-32-8 | Benzo[a]pyrene | 1.6 E-08 | 1.0 E-08 | |
| 192-97-2 | Benzo[e]pyrene | 2.5 E-08 | 1.7 E-08 | |
| 100-44-7 | Benzyl chloride | 3.3 E-06 | 2.2 E-06 | |
| 7440-41-7 | Beryllium | ND | ND | 2.0 E-02 |

TABLE A2 (cont.)

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|--------------------------------|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 75-25-2 | Bromoform | ND | ND | 5.3 E-04 |
| 74-83-9 | Bromomethane | ND | ND | 3.6 E-04 |
| 101-55-3 | 4-Bromophenylphenylether | ND | ND | 3.2 E-02 |
| 106-99-0 | 1,3-Butadiene ^g | ND | ND | 4.5 E-02 |
| 75-65-0 | t-Butyl alcohol | ND | ND | 4.0 E-04 |
| 85-68-7 | Butylbenzylphthalate | 2.7 E-06 | 1.8 E-06 | |
| 123-72-8 | Butyraldehyde | ND | ND | 1.8 E-01 |
| 7440-43-9 | Cadmium | 1.1 E-04 | 7.4 E-05 | |
| 75-15-0 | Carbon disulfide | ND | ND | 3.2 E-04 |
| 56-23-5 | Carbon tetrachloride | ND | ND | 3.2 E-04 |
| 106-47-8 | 4-Chloroaniline | ND | ND | 3.2 E-02 |
| 108-90-7 | Chlorobenzene | 3.3 E-07 | 2.2 E-07 | |
| 75-45-6 | Chlorodifluoromethane | ND | ND | 2.5 E-04 |
| 75-00-3 | Chloroethane | ND | ND | 6.2 E-04 |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | ND | 1.6 E-02 |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | ND | 1.6 E-02 |
| 67-66-3 | Chloroform | ND | ND | 4.0 E-04 |
| 74-87-3 | Chloromethane ^g | 1.3 E-06 | 8.7 E-07 | |
| 91-58-7 | 2-Chloronaphthalene | ND | ND | 1.6 E-02 |
| 95-57-8 | 2-Chlorophenol | ND | ND | 1.6 E-02 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | ND | ND | 1.6 E-02 |
| 7440-47-3 | Chromium ^g | 1.0 E-04 | 6.7 E-05 | |
| 18540-29-9 | Hexavalent chromium | 2.9 E-06 | 2.0 E-06 | |
| 218-01-9 | Chrysene | 6.7 E-08 | 4.5 E-08 | |
| 7440-48-4 | Cobalt ^g | ND | ND | 2.0 E-02 |
| 7440-50-8 | Copper ^g | 1.3 E-04 | 8.8 E-05 | |
| 106-44-5 | p-Cresol | ND | ND | 1.6 E-02 |
| 4170-30-3 | Crotonaldehyde | ND | ND | 1.8 E-01 |
| 98-82-8 | Cumene ^g | 1.8 E-06 | 1.2 E-06 | |
| 57-12-5 | Particulate cyanide | ND | ND | 1.9 E-02 |
| 53-70-3 | Dibenz[a,h]anthracene | ND | ND | 1.6 E-05 |
| 132-64-9 | Dibenzofuran | ND | ND | 1.6 E-02 |
| 106-93-4 | 1,2-Dibromoethane | ND | ND | 3.9 E-04 |
| 84-74-2 | Dibutyl phthalate ^f | 6.6 E-06 | 4.4 E-06 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | ND | 4.9 E-04 |

TABLE A2 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 541-73-1 | 1,3-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 106-46-7 | 1,4-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | ND | 1.6 E-02 |
| 75-27-4 | Dichlorobromomethane | ND | ND | 4.1 E-04 |
| 75-71-8 | Dichlorodifluoromethaneg | 5.3 E-09 | 3.5 E-09 | |
| 75-34-3 | 1,1-Dichloroethane | ND | ND | 3.7 E-04 |
| 107-06-2 | 1,2-Dichloroethane | 3.4 E-05 | 2.3 E-05 | |
| 75-43-4 | Dichlorofluoromethane | ND | ND | 2.1 E-04 |
| 120-83-2 | 2,4-Dichlorophenol | ND | ND | 1.6 E-02 |
| 78-87-5 | 1,2-Dichloropropane | ND | ND | 3.3 E-04 |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | ND | 2.3 E-04 |
| 76-14-2 | Dichlorotetrafluoroethane | ND | ND | 5.0 E-04 |
| 105-67-9 | 2,4-Dimethylphenol | ND | ND | 1.6 E-02 |
| 131-11-3 | Dimethyl phthalate | ND | ND | 1.6 E-02 |
| 99-65-0 | 1,3-Dinitrobenzene | ND | ND | 6.8 E-04 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | ND | ND | 1.6 E-02 |
| 51-28-5 | 2,4-Dinitrophenol | ND | ND | 1.6 E-02 |
| 121-14-2 | 2,4-Dinitrotoluene | 7.3 E-07 | 4.9 E-07 | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | ND | 3.4 E-03 |
| 123-91-1 | 1,4-Dioxane | ND | ND | 3.3 E-04 |
| | Total dioxin/furan compounds | 2.5 E-10 | 1.7 E-10 | |
| 140-88-5 | Ethyl acrylate | ND | ND | 1.7 E-04 |
| 100-41-4 | Ethylbenzene ^g | 1.0 E-05 | 6.7 E-06 | |
| 74-85-1 | Ethylene ^g | 1.2 E-04 | 8.3 E-05 | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate ^f | 6.3 E-06 | 4.2 E-06 | |
| 206-44-0 | Fluorantheneg | 4.0 E-07 | 2.6 E-07 | |
| 86-73-7 | Fluoreneg | 5.8 E-07 | 3.9 E-07 | |
| 50-00-0 | Formaldehyde | 1.7 E-04 | 1.1 E-04 | |
| 76-13-1 | Freon 113 ^g | ND | ND | 5.5 E-04 |
| 35822-46-9 | 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 2.3 E-11 | 1.5 E-11 | |
| 67562-39-4 | 1,2,3,4,6,7,8- Heptachlorodibenzofuran | 2.9 E-12 | 1.9 E-12 | |
| 55673-89-7 | 1,2,3,4,7,8,9- Heptachlorodibenzofuran | 3.9 E-13 | 2.6 E-13 | |

TABLE A2 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 118-74-1 | Hexachlorobenzene | ND | ND | 1.6 E-02 |
| 87-68-3 | Hexachlorobutadiene | ND | ND | 7.6 E-04 |
| 77-47-4 | Hexachlorocyclopentadiene | ND | ND | 1.6 E-02 |
| 39227-28-6 | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | ND | ND | 2.9 E-09 |
| 57653-85-7 | 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 8.2 E-13 | 5.5 E-13 | |
| 19408-74-3 | 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 8.8 E-13 | 5.9 E-13 | |
| 57117-44-9 | 1,2,3,6,7,8-Hexachlorodibenzofuran ⁱ | 3.6 E-13 | 2.4 E-13 | |
| 72918-21-9 | 1,2,3,7,8,9-Hexachlorodibenzofuran | ND | ND | 2.8 E-09 |
| 60851-34-5 | 2,3,4,6,7,8-Hexachlorodibenzofuran | ND | ND | 2.6 E-09 |
| 67-72-1 | Hexachloroethane | ND | ND | 5.9 E-04 |
| 110-54-3 | Hexane | ND | ND | 7.2 E-02 |
| 7647-01-0 | Hydrochloric acid | ND | ND | 2.3 E-01 |
| 74-90-8 | Hydrogen cyanide | 3.6 E-04 | 2.4 E-04 | |
| 7664-39-3 | Hydrogen fluoride | ND | ND | 2.3 E-01 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 7.4 E-09 | 4.9 E-09 | |
| 78-59-1 | Isophorone | ND | ND | 1.6 E-02 |
| 7439-92-1 | Lead | 5.7 E-05 | 3.8 E-05 | |
| 7439-96-5 | Manganese ^g | 7.3 E-05 | 4.9 E-05 | |
| 96-33-3 | Methyl acrylate | ND | ND | 2.5 E-04 |
| 74-95-3 | Methylene bromide | ND | ND | 5.1 E-04 |
| 75-09-2 | Methylene chloride ^g | 3.8 E-07 | 2.5 E-07 | |
| 78-93-3 | Methyl ethyl ketone ^f | 1.5 E-05 | 9.7 E-06 | |
| 74-88-4 | Methyl iodide | ND | ND | 3.5 E-04 |
| 108-10-1 | Methyl isobutyl ketone | ND | ND | 3.3 E-04 |
| 80-62-6 | Methyl methacrylate ⁱ | 1.7 E-07 | 1.1 E-07 | |
| 91-57-6 | 2-Methylnaphthalene | ND | ND | 1.6 E-02 |
| 95-48-7 | 2-Methylphenol | 1.4 E-06 | 9.6 E-07 | |
| 1634-04-4 | Methyl tert-butyl ether | ND | ND | 3.7 E-04 |
| 91-20-3 | Naphthaleneg | 5.4 E-06 | 3.6 E-06 | |
| 7440-02-0 | Nickel ^g | ND | ND | 2.0 E-02 |
| 7697-37-2 | Nitric acid ^g | 8.5 E-04 | 5.6 E-04 | |
| 100-01-6 | 4-Nitroaniline | ND | ND | 1.6 E-02 |

TABLE A2 (cont.)

| | | Emission | Factor ^{b,c} | Minimum | |
|--------------------|--|-------------|-------------------------------|-------------------------------------|--|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} | |
| 98-95-3 | Nitrobenzene | ND | ND | 6.8 E-03 | |
| 55-63-0 | Nitroglycerin | ND | ND | 6.8 E-04 | |
| 88-75-5 | 2-Nitrophenol | 5.5 E-06 | 3.7 E-06 | | |
| 100-02-7 | 4-Nitrophenol | 7.3 E-06 | 4.8 E-06 | | |
| 62-75-9 | n-Nitrosodimethylamine | ND | ND | 1.6 E-02 | |
| 86-30-6 | n-Nitrosodiphenylamine | ND | ND | 1.6 E-02 | |
| 621-64-7 | n-Nitroso-di-n-propylamine | ND | ND | 1.6 E-02 | |
| 3268-87-9 | 1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin | 2.2 E-10 | 1.4 E-10 | | |
| 39001-02-0 | 1,2,3,4,6,7,8,9- Octachlorodibenzofuran | 5.7 E-12 | 3.8 E-12 | | |
| 40321-76-4 | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | ND | ND | 1.5 E-09 | |
| 57117-41-6 | 1,2,3,7,8-Pentachlorodibenzofuran | 4.5 E-13 | 3.0 E-13 | | |
| 57117-31-4 | 2,3,4,7,8-Pentachlorodibenzofuran | 8.4 E-13 | 5.6 E-13 | | |
| 87-86-5 | Pentachlorophenol | ND | ND | 1.6 E-02 | |
| 85-01-8 | Phenanthrene ^g | 1.8 E-06 | 1.2 E-06 | | |
| 108-95-2 | Phenol | 3.9 E-06 | 2.6 E-06 | | |
| 123-38-6 | Propionaldehyde | ND | ND | 1.8 E-01 | |
| 115-07-1 | Propylene ^g | 2.2 E-05 | 1.5 E-05 | | |
| 129-00-0 | Pyrene ^g | 4.1 E-07 | 2.7 E-07 | | |
| 7782-49-2 | Selenium | ND | ND | 2.0 E-02 | |
| 7440-22-4 | Silver | ND | ND | 2.0 E-02 | |
| 100-42-5 | Styrene | ND | ND | 3.0 E-04 | |
| 7664-93-9 | Sulfuric acid | ND | ND | 2.2 E-01 | |
| 1746-01-6 | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | ND | ND | 2.1 E-09 | |
| 51207-31-9 | 2,3,7,8-Tetrachlorodibenzofuran | 1.3 E-12 | 8.6 E-13 | | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | ND | 3.5 E-04 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | ND | 3.5 E-04 | |
| 127-18-4 | Tetrachloroethylene | ND | ND | 4.1 E-04 | |
| 7440-28-0 | Thallium | ND | ND | 0 | |
| 108-88-3 | Toluene ^g | 2.0 E-05 | 1.3 E-05 | | |
| 120-82-1 | 1,2,4-Trichlorobenzene ^g | ND | ND | 1.3 E-03 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | ND | 4.4 E-04 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | ND | 2.8 E-04 | |

TABLE A2 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|-----------------------|---------------------------------|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 75-69-4 | Trichlorofluoromethane | 3.7 E-09 | 2.4 E-09 | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | ND | 1.6 E-02 |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | ND | 1.6 E-02 |
| 96-18-4 | 1,2,3-Trichloropropane | ND | ND | 4.9 E-04 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 7.2 E-06 | 4.8 E-06 | |
| 540-84-1 | 2,2,4-Trimethylpentane | ND | ND | 1.9 E-04 |
| 7440-62-2 | Vanadium | 5.4 E-04 | 3.6 E-04 | |
| 108-05-4 | Vinyl acetate | ND | ND | 3.9 E-04 |
| 75-01-4 | Vinyl chloride | 3.0 E-06 | 2.0 E-06 | |
| 75-35-4 | Vinylidene chloride | ND | ND | 2.1 E-04 |
| 106-42-3, 108-38-3 | m-Xylene, p-Xylene ^g | 4.0 E-06 | 2.6 E-06 | |
| 95-47-6 | o-Xylene ^g | 1.9 E-06 | 1.3 E-06 | |
| 7440-66-6 | Zinc ^g | 8.2 E-04 | 5.5 E-04 | |
| | Other Po | llutants | | |
| 67-64-1 | Acetone | 1.6 E-05 | 1.1 E-05 | |
| 74-86-2 | Acetylene ^g | 1.8 E-04 | 1.2 E-04 | |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | ND | ND | 6.8 E-04 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | ND | ND | 6.8 E-04 |
| 100-52-7 | Benzaldehyde | 1.4 E-04 | 9.6 E-05 | |
| 20780-53-4 | Benzene, (epoxyethyl)-, R-i | 3.0 E-05 | 2.0 E-05 | |
| 102-96-5 | Benzene,(2-nitroethenyl)-i | 2.4 E-05 | 1.6 E-05 | |
| 103-65-1 | Benzene, propyl-i | 1.8 E-06 | 1.2 E-06 | |
| 271-89-6 | Benzofuran ⁱ | 1.1 E-05 | 7.2 E-06 | |
| 65-85-0 | Benzoic acid ⁱ | 3.7 E-05 | 2.4 E-05 | |
| 100-51-6 | Benzyl alcohol | 1.7 E-06 | 1.2 E-06 | |
| 108-86-1 | Bromobenzene | ND | ND | 3.9 E-04 |
| 460-12-8 | 1,3-Butadiyne ⁱ | 1.8 E-06 | 1.2 E-06 | |
| 106-97-8 | Butane | ND | ND | 5.3 E-02 |
| 590-18-1 | cis-Butene | ND | ND | 4.7 E-02 |
| 624-64-6 | trans-Butene ^g | ND | ND | 4.7 E-02 |
| 107-00-6 | 1-Butyne | ND | ND | 4.5 E-02 |
| 503-17-3 | 2-Butyne | ND | ND | 4.5 E-02 |
| 7440-70-2 | Calcium ^f | 1.6 E-03 | 1.0 E-03 | |
| 7440-44-0 | Total carbon ^g | 6.9 E-03 | 4.6 E-03 | |

TABLE A2 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|--|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 39638-32-9 | bis(2-Chloroisopropyl)ether | ND | ND | 1.6 E-02 |
| 59-50-7 | 4-Chloro-3-methylphenol | ND | ND | 1.6 E-02 |
| 108-94-1 | Cyclohexanonei | 1.4 E-06 | 9.5 E-07 | |
| 124-48-1 | Dibromochloromethane | ND | ND | 5.2 E-04 |
| 156-59-2 | cis-1,2-Dichloroethene | ND | ND | 2.0 E-04 |
| 156-60-5 | trans-1,2-Dichloroethene | ND | ND | 4.8 E-04 |
| 84-66-2 | Diethylphthalate | ND | ND | 1.6 E-02 |
| 5779-94-2 | 2,5-Dimethylbenzaldehyde | ND | ND | 1.8 E-01 |
| 117-84-0 | Di-n-octylphthalate ^g | 1.6 E-06 | 1.1 E-06 | |
| 74-84-0 | Ethane ^g | 2.1 E-05 | 1.4 E-05 | |
| 141-78-6 | Ethyl acetate | ND | ND | 1.4 E-03 |
| 97-63-2 | Ethyl methacrylate | ND | ND | 3.3 E-04 |
| 622-96-8 | 4-Ethyltoluene | 3.3 E-06 | 2.2 E-06 | |
| 1191-99-7 | Furan, 2,3-dihydro-i | 1.0 E-06 | 6.9 E-07 | |
| 1708-29-8 | Furan, 2,5-dihydro-i | 1.0 E-06 | 6.9 E-07 | |
| 142-82-5 | Heptane ⁱ | 1.2 E-07 | 8.0 E-08 | |
| 66-25-1 | Hexaldehyde | ND | ND | 1.8 E-01 |
| 103-23-1 | Hexanedioic acid, bis(2-ethylhexyl) ester ⁱ | 3.1 E-05 | 2.1 E-05 | |
| 591-78-6 | 2-Hexanone | ND | ND | 2.9 E-04 |
| 592-41-6 | 1-Hexene ⁱ | 3.1 E-06 | 2.1 E-06 | |
| 2691-41-0 | HMX | ND | ND | 1.4 E-03 |
| 10035-10-6 | Hydrogen bromide | ND | ND | 2.2 E-01 |
| 75-28-5 | Isobutane | ND | ND | 4.8 E-02 |
| 590-86-3 | Isovaleraldehyde | ND | ND | 1.8 E-01 |
| 7439-95-4 | Magnesium ^g | 1.2 E-03 | 7.7 E-04 | |
| 75-52-5 | Methane, nitro-h | 2.7 E-06 | 1.8 E-06 | |
| 98-83-9 | alpha-Methylstyrene | ND | ND | 2.9 E-04 |
| 88-74-7 | 2-Nitroaniline | ND | ND | 1.6 E-02 |
| 99-09-2 | 3-Nitroaniline | ND | ND | 1.6 E-02 |
| 88-72-2 | 2-Nitrotoluene | ND | ND | 6.8 E-04 |
| 99-08-1 | 3-Nitrotoluene | ND | ND | 6.8 E-04 |
| 99-99-0 | 4-Nitrotoluene | ND | ND | 6.8 E-04 |
| 111-65-9 | Octane | 3.8 E-07 | 2.5 E-07 | |
| 109-66-0 | Pentane | ND | ND | 6.0 E-02 |

TABLE A2 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|-------------------------------------|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Compound | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 536-74-3 | Phenyl acetylene ⁱ | 1.6 E-05 | 1.0 E-05 | |
| 7664-38-2 | Phosphoric acid | ND | ND | 2.2 E-01 |
| 74-98-6 | Propane | ND | ND | 3.7 E-02 |
| 78-85-3 | 2-Propenal, 2-methyl- ⁱ | 1.0 E-06 | 6.8 E-07 | |
| 115-11-7 | 1-Propene, 2-methyl- | ND | ND | 4.7 E-02 |
| 74-99-7 | Propyne ^g | 1.1 E-05 | 7.3 E-06 | |
| 121-82-4 | RDX | 2.1 E-06 | 1.4 E-06 | |
| 7440-24-6 | Strontium | 1.2 E-03 | 7.7 E-04 | |
| 479-45-8 | Tetryl | ND | ND | 6.8 E-04 |
| 7440-32-6 | Titanium | 1.1 E-04 | 7.4 E-05 | |
| 1334-78-7 | o,m,p-Tolualdehyde | ND | ND | 5.5 E-01 |
| 25323-89-1 | Trichloroethane | ND | ND | 3.0 E-04 |
| 108-67-8 | 1,3,5-Trimethylbenzene ^g | 2.2 E-06 | 1.5 E-06 | |
| 99-35-4 | 1,3,5-Trinitrobenzene ^g | 7.5 E-07 | 5.0 E-07 | |
| 118-96-7 | 2,4,6-Trinitrotoluene | 1.0 E-06 | 6.7 E-07 | |
| 110-62-3 | Valeraldehyde | ND | ND | 1.8 E-01 |

^a CASRN = Chemical Abstracts Service Registry Number.

^b ND = nondetected.

^c Emission factors rated C unless otherwise noted.

^d NEW = net explosive weight. The NEW for this ordnance is 1.50 pounds per item.

^e Data provided for compounds that were not detected.

^f Emission factor rated A because of correlation with emission factors for similar ordnance and number of test data points.

g Emission factor rated B because of correlation with emission factors for similar ordnance and number of test data points.

^h Emission factor based upon C-rated test data, but because of correlation with emission factors for similar ordnance and number of data points the factor was upgraded from a D rating to a C rating.

ⁱ Emission factor rated D because the factor is based upon C-rated test data.

TABLE A3 COMPOUNDS ANALYZED AND EMISSION FACTORS DEVELOPED FOR DODIC L709, M25 TARGET-HIT SIMULATOR

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|--|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| (| Carbon Dioxide, Criteria Pollutants, Me | thane, and Total | Suspended Parti | culate |
| 124-38-9 | Carbon dioxide ^g | 3.8 E-02 | 7.5 E-01 | |
| 630-08-0 | Carbon monoxide ^g | 1.3 E-03 | 2.4 E-02 | |
| 7439-92-1 | Lead ^g | 1.9 E-05 | 3.8 E-04 | |
| 74-82-8 | Methane | 5.8 E-06 | 1.1 E-04 | |
| | Oxides of nitrogen ^g | 1.8 E-03 | 3.6 E-02 | |
| | PM-2.5 | 2.9 E-02 | 5.6 E-01 | |
| | PM-10 ^g | 3.4 E-02 | 6.7 E-01 | |
| 7446-09-5 | Sulfur dioxide | ND | ND | 5.3 E-01 |
| 12789-66-1 | Total suspended particulate ^g | 3.6 E-02 | 7.0 E-01 | |
| | Hazardous Air Pollutant | s and Toxic Che | micals | |
| 83-32-9 | Acenaphthene | ND | ND | 1.6 E-05 |
| 208-96-8 | Acenaphthylene | 0 | 0 | |
| 75-07-0 | Acetaldehyde | ND | ND | 1.9 E-01 |
| 75-05-8 | Acetonitrile ^g | 0 | 0 | |
| 107-02-8 | Acrolein | ND | ND | 1.9 E-01 |
| 107-13-1 | Acrylonitrile ^g | ND | ND | 6.0 E-04 |
| 107-05-1 | Allyl chloride | ND | ND | 8.9 E-04 |
| 7429-90-5 | Aluminum ^g | 1.8 E-05 | 3.5 E-04 | |
| 7664-41-7 | Ammonia | ND | ND | 2.1 |
| 120-12-7 | Anthracene | ND | ND | 1.6 E-05 |
| 7440-36-0 | Antimony | ND | ND | 1.5 E-02 |
| 7440-38-2 | Arsenic | ND | ND | 1.5 E-02 |
| 7440-39-3 | Barium | 4.8 E-04 | 9.3 E-03 | |
| 71-43-2 | Benzene | 6.8 E-07 | 1.3 E-05 | |
| 56-55-3 | Benzo[a]anthracene | ND | ND | 1.6 E-05 |
| 205-99-2 | Benzo[b]fluoranthene | ND | ND | 1.6 E-05 |
| 207-08-9 | Benzo[k]fluoranthene | ND | ND | 1.6 E-05 |
| 191-24-2 | Benzo[g,h,i]perylene | ND | ND | 1.6 E-05 |
| 50-32-8 | Benzo[a]pyrene | ND | ND | 1.6 E-05 |
| 192-97-2 | Benzo[e]pyrene | ND | ND | 1.6 E-05 |
| 100-44-7 | Benzyl chloride | ND | ND | 2.6 E-04 |
| 7440-41-7 | Beryllium | ND | ND | 1.5 E-02 |

TABLE A3 (cont.)

| | | Emission Factor ^{b,c} | Factor ^{b,c} | Minimum |
|--------------------|-----------------------------------|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 75-25-2 | Bromoform | ND | ND | 5.3 E-04 |
| 74-83-9 | Bromomethane | ND | ND | 3.6 E-04 |
| 101-55-3 | 4-Bromophenylphenylether | ND | ND | 3.2 E-02 |
| 106-99-0 | 1,3-Butadiene | ND | ND | 4.5 E-02 |
| 75-65-0 | t-Butyl alcohol | ND | ND | 4.0 E-04 |
| 85-68-7 | Butylbenzylphthalate ^g | ND | ND | 1.6 E-02 |
| 123-72-8 | Butyraldehyde | ND | ND | 1.9 E-01 |
| 7440-43-9 | Cadmium | ND | ND | 1.5 E-02 |
| 75-15-0 | Carbon disulfide ^g | 1.7 E-06 | 3.2 E-05 | |
| 56-23-5 | Carbon tetrachloride | ND | ND | 3.2 E-04 |
| 106-47-8 | 4-Chloroaniline | ND | ND | 3.2 E-02 |
| 108-90-7 | Chlorobenzene | ND | ND | 2.4 E-04 |
| 75-45-6 | Chlorodifluoromethane | ND | ND | 2.5 E-04 |
| 75-00-3 | Chloroethane | ND | ND | 6.2 E-04 |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | ND | 1.6 E-02 |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | ND | 1.6 E-02 |
| 67-66-3 | Chloroform | ND | ND | 4.0 E-04 |
| 74-87-3 | Chloromethane | 1.8 E-08 | 3.6 E-07 | |
| 91-58-7 | 2-Chloronaphthalene | ND | ND | 1.6 E-02 |
| 95-57-8 | 2-Chlorophenol | ND | ND | 1.6 E-02 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | ND | ND | 1.6 E-02 |
| 7440-47-3 | Chromium ^g | ND | ND | 1.5 E-02 |
| 18540-29-9 | Hexavalent chromium | 1.8 E-07 | 3.5 E-06 | |
| 218-01-9 | Chrysene | ND | ND | 1.6 E-05 |
| 7440-48-4 | Cobalt | ND | ND | 1.5 E-02 |
| 7440-50-8 | Copper | 5.6 E-07 | 1.1 E-05 | |
| 106-44-5 | p-Cresol | ND | ND | 1.6 E-02 |
| 4170-30-3 | Crotonaldehyde | ND | ND | 1.9 E-01 |
| 98-82-8 | Cumene | ND | ND | 2.0 E-04 |
| 57-12-5 | Particulate cyanide | ND | ND | 1.6 E-02 |
| 53-70-3 | Dibenz[a,h]anthracene | ND | ND | 1.6 E-05 |
| 132-64-9 | Dibenzofuran | ND | ND | 1.6 E-02 |
| 106-93-4 | 1,2-Dibromoethane | ND | ND | 3.9 E-04 |
| 84-74-2 | Dibutyl phthalate | ND | ND | 1.6 E-02 |

TABLE A3 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 95-50-1 | 1,2-Dichlorobenzene | ND | ND | 4.9 E-04 |
| 541-73-1 | 1,3-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 106-46-7 | 1,4-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | ND | 1.6 E-02 |
| 75-27-4 | Dichlorobromomethane | ND | ND | 4.1 E-04 |
| 75-71-8 | Dichlorodifluoromethane | 0 | 0 | |
| 75-34-3 | 1,1-Dichloroethane | ND | ND | 3.7 E-04 |
| 107-06-2 | 1,2-Dichloroethane | ND | ND | 2.1 E-04 |
| 75-43-4 | Dichlorofluoromethane | ND | ND | 2.1 E-04 |
| 120-83-2 | 2,4-Dichlorophenol | ND | ND | 1.6 E-02 |
| 78-87-5 | 1,2-Dichloropropane | ND | ND | 3.3 E-04 |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | ND | 2.3 E-04 |
| 76-14-2 | Dichlorotetrafluoroethane | ND | ND | 5.0 E-04 |
| 105-67-9 | 2,4-Dimethylphenol | ND | ND | 1.6 E-02 |
| 131-11-3 | Dimethyl phthalate | ND | ND | 1.6 E-02 |
| 99-65-0 | 1,3-Dinitrobenzene | ND | ND | 6.8 E-04 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | ND | ND | 1.6 E-02 |
| 51-28-5 | 2,4-Dinitrophenol | ND | ND | 1.6 E-02 |
| 121-14-2 | 2,4-Dinitrotoluene | ND | ND | 6.8 E-04 |
| 606-20-2 | 2,6-Dinitrotoluene | ND | ND | 6.8 E-04 |
| 123-91-1 | 1,4-Dioxane | ND | ND | 3.3 E-04 |
| | Total dioxin/furan compounds | 4.1 E-12 | 7.9 E-11 | |
| 140-88-5 | Ethyl acrylate | ND | ND | 1.7 E-04 |
| 100-41-4 | Ethylbenzene ^h | 9.1 E-08 | 1.8 E-06 | |
| 74-85-1 | Ethylene ^g | ND | ND | 2.3 E-02 |
| 117-81-7 | bis(2-Ethylhexyl)phthalate ^h | 5.2 E-06 | 1.0 E-04 | |
| 206-44-0 | Fluoranthene | 4.0 E-09 | 7.8 E-08 | |
| 86-73-7 | Fluorene | 1.6 E-10 | 3.1 E-09 | |
| 50-00-0 | Formaldehyde | ND | ND | 1.9 E-01 |
| 76-13-1 | Freon 113 | ND | ND | 5.5 E-04 |
| 35822-46-9 | 1,2,3,4,6,7,8- Heptachlorodibenzo-p-dioxin | 4.4 E-13 | 8.7 E-12 | |
| 55673-89-7 | 1,2,3,4,7,8,9- Heptachlorodibenzofuran | ND | ND | 2.6 E-09 |
| 118-74-1 | Hexachlorobenzene | ND | ND | 1.6 E-02 |

TABLE A3 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|--|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 87-68-3 | Hexachlorobutadiene | ND | ND | 7.6 E-04 |
| 77-47-4 | Hexachlorocyclopentadiene | ND | ND | 1.6 E-02 |
| 39227-28-6 | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | ND | ND | 3.2 E-09 |
| 57653-85-7 | 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | ND | ND | 3.3 E-09 |
| 19408-74-3 | 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | ND | ND | 3.3 E-09 |
| 70648-26-9 | 1,2,3,4,7,8-Hexachlorodibenzofuran | ND | ND | 1.9 E-09 |
| 57117-44-9 | 1,2,3,6,7,8-Hexachlorodibenzofuran | ND | ND | 1.9 E-09 |
| 72918-21-9 | 1,2,3,7,8,9-Hexachlorodibenzofuran | ND | ND | 2.5 E-09 |
| 60851-34-5 | 2,3,4,6,7,8-Hexachlorodibenzofuran | ND | ND | 2.5 E-09 |
| 67-72-1 | Hexachloroethane | ND | ND | 5.9 E-04 |
| 110-54-3 | Hexane | ND | ND | 7.2 E-02 |
| 7647-01-0 | Hydrochloric acid | 5.7 E-05 | 1.1 E-03 | |
| 74-90-8 | Hydrogen cyanide | ND | ND | 1.7 E-02 |
| 7664-39-3 | Hydrogen fluoride | ND | ND | 2.2 E-01 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | ND | ND | 1.6 E-05 |
| 78-59-1 | Isophorone | ND | ND | 1.6 E-02 |
| 7439-92-1 | Lead | 1.9 E-05 | 3.8 E-04 | |
| 7439-96-5 | Manganese ^g | 2.4 E-06 | 4.7 E-05 | |
| 96-33-3 | Methyl acrylate | ND | ND | 2.5 E-04 |
| 74-95-3 | Methylene bromide | ND | ND | 5.1 E-04 |
| 75-09-2 | Methylene chloride | 0 | 0 | |
| 78-93-3 | Methyl ethyl ketone | ND | ND | 8.1 E-04 |
| 74-88-4 | Methyl iodide | ND | ND | 3.5 E-04 |
| 108-10-1 | Methyl isobutyl ketone | ND | ND | 3.3 E-04 |
| 80-62-6 | Methyl methacrylate | ND | ND | 4.6 E-04 |
| 91-57-6 | 2-Methylnaphthalene | ND | ND | 1.6 E-02 |
| 95-48-7 | 2-Methylphenol | ND | ND | 1.6 E-02 |
| 1634-04-4 | Methyl tert-butyl ether | ND | ND | 3.7 E-04 |
| 91-20-3 | Naphthalene | ND | ND | 1.6 E-02 |
| 7440-02-0 | Nickel ^g | ND | ND | 1.5 E-02 |
| 7697-37-2 | Nitric acid | 3.8 E-06 | 7.3 E-05 | |
| 100-01-6 | 4-Nitroaniline | ND | ND | 1.6 E-02 |

TABLE A3 (cont.)

| | | Emission Factor ^{b,c} | | Minimum |
|--------------------|--|--------------------------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 98-95-3 | Nitrobenzene | ND | ND | 1.4 E-03 |
| 55-63-0 | Nitroglycerin | ND | ND | 6.8 E-04 |
| 88-75-5 | 2-Nitrophenol | ND | ND | 1.6 E-02 |
| 100-02-7 | 4-Nitrophenol | ND | ND | 1.6 E-02 |
| 62-75-9 | n-Nitrosodimethylamine | ND | ND | 1.6 E-02 |
| 86-30-6 | n-Nitrosodiphenylamine | ND | ND | 1.6 E-02 |
| 621-64-7 | n-Nitroso-di-n-propylamine | ND | ND | 1.6 E-02 |
| 3268-87-9 | 1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin | 4.1 E-12 | 7.9 E-11 | |
| 40321-76-4 | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | ND | ND | 3.8 E-09 |
| 57117-41-6 | 1,2,3,7,8-Pentachlorodibenzofuran | ND | ND | 2.9 E-09 |
| 57117-31-4 | 2,3,4,7,8-Pentachlorodibenzofuran | ND | ND | 1.7 E-09 |
| 87-86-5 | Pentachlorophenol | ND | ND | 1.6 E-02 |
| 85-01-8 | Phenanthrene | ND | ND | 3.2 E-02 |
| 108-95-2 | Phenol | ND | ND | 1.6 E-02 |
| 123-38-6 | Propionaldehyde | ND | ND | 1.9 E-01 |
| 115-07-1 | Propylene ^g | ND | ND | 3.5 E-02 |
| 129-00-0 | Pyrene | ND | ND | 1.6 E-02 |
| 7782-49-2 | Selenium | ND | ND | 1.5 E-02 |
| 7440-22-4 | Silver | ND | ND | 1.5 E-02 |
| 100-42-5 | Styrene | ND | ND | 3.0 E-04 |
| 7664-93-9 | Sulfuric acid | 1.6 E-04 | 3.1 E-03 | |
| 1746-01-6 | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | ND | ND | 2.3 E-09 |
| 51207-31-9 | 2,3,7,8-Tetrachlorodibenzofuran | ND | ND | 2.5 E-09 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | ND | 3.5 E-04 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | ND | 3.5 E-04 |
| 127-18-4 | Tetrachloroethylene | ND | ND | 4.1 E-04 |
| 7440-28-0 | Thallium | ND | ND | 1.5 E-02 |
| 108-88-3 | Toluene | 2.0 E-07 | 3.9 E-06 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | ND | 1.6 E-02 |
| 71-55-6 | 1,1,1-Trichloroethane ^h | 1.2 E-07 | 2.3 E-06 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | ND | 2.8 E-04 |
| 75-69-4 | Trichlorofluoromethane | 0 | 0 | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | ND | 1.6 E-02 |

TABLE A3 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|-----------------------|-------------------------------|-------------|-------------------------------|--|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | ND | 1.6 E-02 |
| 96-18-4 | 1,2,3-Trichloropropane | ND | ND | 4.9 E-04 |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | ND | 2.5 E-04 |
| 540-84-1 | 2,2,4-Trimethylpentane | ND | ND | 1.9 E-04 |
| 7440-62-2 | Vanadium | 1.6 E-05 | 3.2 E-04 | |
| 108-05-4 | Vinyl acetate | ND | ND | 3.9 E-04 |
| 75-01-4 | Vinyl chloride | ND | ND | 2.6 E-04 |
| 75-35-4 | Vinylidene chloride | ND | ND | 2.1 E-04 |
| 106-42-3, 108-38-3 | m-Xylene, p-Xylene | 1.9 E-07 | 3.8 E-06 | |
| 95-47-6 | o-Xylene ^h | 1.4 E-07 | 2.7 E-06 | |
| 7440-66-6 | Zinc ^f | ND | ND | 1.2 E-01 |
| | Other Po | llutants | | |
| 67-64-1 | Acetone ^g | 1.1 E-07 | 2.2 E-06 | |
| 74-86-2 | Acetylene | 3.2 E-06 | 6.3 E-05 | |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | ND | ND | 6.8 E-04 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | ND | ND | 6.8 E-04 |
| 100-52-7 | Benzaldehyde | ND | ND | 1.9 E-01 |
| 65-85-0 | Benzoic acid ^h | 7.1 E-06 | 1.4 E-04 | |
| 100-51-6 | Benzyl alcohol | ND | ND | 1.6 E-02 |
| 108-86-1 | Bromobenzene | ND | ND | 3.9 E-04 |
| 106-97-8 | Butane | ND | ND | 5.3 E-02 |
| 590-18-1 | cis-Butene | ND | ND | 4.7 E-02 |
| 624-64-6 | trans-Butene ^g | ND | ND | 4.7 E-02 |
| 107-00-6 | 1-Butyne | ND | ND | 4.5 E-02 |
| 503-17-3 | 2-Butyne | ND | ND | 4.5 E-02 |
| 7440-70-2 | Calcium | ND | ND | 1.5 E-01 |
| 7440-44-0 | Total carbon | 1.0 E-03 | 2.0 E-02 | |
| 39638-32-9 | bis(2-Chloroisopropyl)ether | ND | ND | 1.6 E-02 |
| 59-50-7 | 4-Chloro-3-methylphenol | ND | ND | 1.6 E-02 |
| 124-48-1 | Dibromochloromethane | ND | ND | 5.2 E-04 |
| 156-59-2 | cis-1,2-Dichloroethene | ND | ND | 2.0 E-04 |
| 156-60-5 | trans-1,2-Dichloroethene | ND | ND | 4.8 E-04 |
| 84-66-2 | Diethylphthalate ^g | ND | ND | 1.6 E-02 |
| 5779-94-2 | 2,5-Dimethylbenzaldehyde | ND | ND | 1.9 E-01 |

TABLE A3 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|-----------------------------|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 117-84-0 | Di-n-octylphthalate | ND | ND | 1.6 E-02 |
| 629-97-0 | Docosane ^h | 2.7 E-06 | 5.3 E-05 | |
| 74-84-0 | Ethane ^g | ND | ND | 2.5 E-02 |
| 141-78-6 | Ethyl acetate | ND | ND | 1.4 E-03 |
| 97-63-2 | Ethyl methacrylate | ND | ND | 3.3 E-04 |
| 622-96-8 | 4-Ethyltoluene ^h | 1.0 E-07 | 2.0 E-06 | |
| 109-99-9 | Furan, tetrahydro-h | 5.2 E-07 | 1.0 E-05 | |
| 142-82-5 | Heptane | ND | ND | 2.1 E-04 |
| 630-01-3 | Hexacosane ^h | 2.8 E-06 | 5.5 E-05 | |
| 66-25-1 | Hexaldehyde | ND | ND | 1.9 E-01 |
| 591-78-6 | 2-Hexanone | ND | ND | 2.9 E-04 |
| 2691-41-0 | HMX | ND | ND | 1.4 E-03 |
| 10035-10-6 | Hydrogen bromide | ND | ND | 2.1 E-01 |
| 75-28-5 | Isobutane | ND | ND | 4.8 E-02 |
| 590-86-3 | Isovaleraldehyde | ND | ND | 1.9 E-01 |
| 98-83-9 | alpha-Methylstyrene | ND | ND | 2.9 E-04 |
| 7439-95-4 | Magnesium ^g | 6.7 E-03 | 1.3 E-01 | |
| 100-97-0 | Methenamine ^h | 1.8 E-06 | 3.4 E-05 | |
| 88-74-7 | 2-Nitroaniline | ND | ND | 1.6 E-02 |
| 99-09-2 | 3-Nitroaniline | ND | ND | 1.6 E-02 |
| 88-72-2 | 2-Nitrotoluene | ND | ND | 6.8 E-04 |
| 99-08-1 | 3-Nitrotoluene | ND | ND | 6.8 E-04 |
| 99-99-0 | 4-Nitrotoluene | ND | ND | 6.8 E-04 |
| 111-65-9 | Octane | ND | ND | 1.9 E-04 |
| 109-66-0 | Pentane | ND | ND | 6.0 E-02 |
| 7664-38-2 | Phosphoric acid | ND | ND | 2.1 E-01 |
| 74-98-6 | Propane | ND | ND | 3.7 E-02 |
| 115-11-7 | 1-Propene, 2-methyl- | ND | ND | 4.7 E-02 |
| 74-99-7 | Propyne | ND | ND | 3.3 E-02 |
| 121-82-4 | RDX ^h | 1.2 E-07 | 2.3 E-06 | |
| 7440-24-6 | Strontium | 6.7 E-03 | 1.3 E-01 | |
| 479-45-8 | Tetryl | ND | ND | 6.8 E-04 |
| 7440-32-6 | Titanium | ND | ND | 1.5 E-02 |
| 1334-78-7 | o,m,p-Tolualdehyde | ND | ND | 5.6 E-01 |

TABLE A3 (cont.)

| | | Emission Factor ^{b,c} | | Minimum | |
|--------------------|------------------------|--------------------------------|-------------------------------|--|--|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} | |
| 25323-89-1 | Trichloroethane | ND | ND | 3.0 E-04 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | ND | 2.5 E-04 | |
| 99-35-4 | 1,3,5-Trinitrobenzene | 1.4 E-07 | 2.7 E-06 | | |
| 118-96-7 | 2,4,6-Trinitrotoluene | ND | ND | 6.8 E-04 | |
| 110-62-3 | Valeraldehyde | ND | ND | 1.9 E-01 | |

^a CASRN = Chemical Abstracts Service Registry Number.

^b ND = nondetected.

^c Emission factors rated C unless otherwise noted.

^d NEW = net explosive weight. The NEW for this ordnance is 5.11 E-02 pounds per item.

^e Data provided for compounds that were not detected.

Emission factor rated A because of correlation with emission factors for similar ordnance and number of test data points.

^g Emission factor rated B because of correlation with emission factors for similar ordnance and number of test data points.

h Emission factor rated D because the factor is based upon C-rated test data.

TABLE A4 COMPOUNDS ANALYZED AND EMISSION FACTORS DEVELOPED FOR DODIC M032, 1-POUND DEMOLITION BLOCK CHARGE

| | | Emission Factor ^{b,c} | | Minimum | |
|------------|---|--------------------------------|-------------------------------|-------------------------------------|--|
| CASRNª | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} | |
| (| Carbon Dioxide, Criteria Pollutants, Methane, and Total Suspended Particulate | | | | |
| 124-38-9 | Carbon dioxide ^f | 1.2 | 1.2 | | |
| 630-08-0 | Carbon monoxide ^g | 5.3 E-03 | 5.3 E-03 | | |
| 7439-92-1 | Lead | 2.0 E-04 | 2.0 E-04 | | |
| | Oxides of nitrogen ^f | 1.4 E-02 | 1.4 E-02 | | |
| | PM-2.5 ^g | 1.4 E-02 | 1.4 E-02 | | |
| | PM-10 | 2.5 E-02 | 2.5 E-02 | | |
| 7446-09-5 | Sulfur dioxide | ND | ND | 5.3 E-01 | |
| 12789-66-1 | Total suspended particulate | 3.3 E-02 | 3.3 E-02 | | |
| | Hazardous Air Pollutant | s and Toxic Che | micals | | |
| 83-32-9 | Acenaphtheneg | 9.2 E-09 | 9.2 E-09 | | |
| 208-96-8 | Acenaphthyleneg | 1.0 E-07 | 1.0 E-07 | | |
| 75-07-0 | Acetaldehyde ^g | ND | ND | 1.7 E-01 | |
| 75-05-8 | Acetonitrile ^g | 0 | 0 | | |
| 107-02-8 | Acrolein | ND | ND | 1.7 E-01 | |
| 107-13-1 | Acrylonitrile ^g | 3.1 E-07 | 3.1 E-07 | | |
| 107-05-1 | Allyl chloride | ND | ND | 8.9 E-04 | |
| 7429-90-5 | Aluminum ^g | 9.1 E-04 | 9.1 E-04 | | |
| 7664-41-7 | Ammonia | ND | ND | 2.1 | |
| 120-12-7 | Anthraceneg | 1.2 E-08 | 1.2 E-08 | | |
| 7440-36-0 | Antimony | ND | ND | 2.2 E-02 | |
| 7440-38-2 | Arsenic | ND | ND | 2.2 E-02 | |
| 7440-39-3 | Barium | 6.2 E-04 | 6.2 E-04 | | |
| 71-43-2 | Benzene ^g | 1.7 E-07 | 1.7 E-07 | | |
| 56-55-3 | Benzo[a]anthracene | ND | ND | 1.6 E-02 | |
| 205-99-2 | Benzo[b]fluoranthene | ND | ND | 1.6 E-05 | |
| 207-08-9 | Benzo[k]fluoranthene | ND | ND | 1.6 E-02 | |
| 191-24-2 | Benzo[g,h,i]perylene | ND | ND | 1.6 E-05 | |
| 50-32-8 | Benzo[a]pyrene | ND | ND | 1.6 E-05 | |
| 192-97-2 | Benzo[e]pyrene | 0 | 0 | | |
| 100-44-7 | Benzyl chloride | ND | ND | 2.6 E-04 | |
| 7440-41-7 | Beryllium | ND | ND | 2.2 E-02 | |
| 75-25-2 | Bromoform | ND | ND | 5.3 E-04 | |

TABLE A4 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|--------------------------------|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 74-83-9 | Bromomethane | ND | ND | 3.6 E-04 |
| 101-55-3 | 4-Bromophenylphenylether | ND | ND | 3.2 E-02 |
| 75-65-0 | t-Butyl alcohol | ND | ND | 4.0 E-04 |
| 85-68-7 | Butylbenzylphthalatef | 1.7 E-06 | 1.7 E-06 | |
| 123-72-8 | Butyraldehyde | ND | ND | 1.7 E-01 |
| 7440-43-9 | Cadmium | ND | ND | 2.2 E-02 |
| 75-15-0 | Carbon disulfide | ND | ND | 3.2 E-04 |
| 56-23-5 | Carbon tetrachloride | ND | ND | 3.2 E-04 |
| 106-47-8 | 4-Chloroaniline | ND | ND | 3.2 E-02 |
| 108-90-7 | Chlorobenzene | ND | ND | 2.4 E-04 |
| 75-45-6 | Chlorodifluoromethane | ND | ND | 2.5 E-04 |
| 75-00-3 | Chloroethane | ND | ND | 6.2 E-04 |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | ND | 1.6 E-02 |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | ND | 1.6 E-02 |
| 67-66-3 | Chloroform | ND | ND | 4.0 E-04 |
| 74-87-3 | Chloromethaneg | 1.1 E-07 | 1.1 E-07 | |
| 91-58-7 | 2-Chloronaphthalene | ND | ND | 1.6 E-02 |
| 95-57-8 | 2-Chlorophenol | ND | ND | 1.6 E-02 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | ND | ND | 1.6 E-02 |
| 7440-47-3 | Chromium ^g | 8.7 E-06 | 8.7 E-06 | |
| 18540-29-9 | Hexavalent chromium | 2.1 E-06 | 2.1 E-06 | |
| 218-01-9 | Chrysene | ND | ND | 1.6 E-02 |
| 7440-48-4 | Cobalt ^g | ND | ND | 2.2 E-02 |
| 7440-50-8 | Copper ^g | 5.3 E-04 | 5.3 E-04 | |
| 106-44-5 | p-Cresol | ND | ND | 1.6 E-02 |
| 4170-30-3 | Crotonaldehyde | ND | ND | 1.7 E-01 |
| 98-82-8 | Cumene ^g | ND | ND | 2.0 E-04 |
| 57-12-5 | Particulate cyanide | ND | ND | 1.6 E-02 |
| 53-70-3 | Dibenz[a,h]anthracene | ND | ND | 1.6 E-05 |
| 132-64-9 | Dibenzofuran | ND | ND | 1.6 E-02 |
| 106-93-4 | 1,2-Dibromoethane | ND | ND | 3.9 E-04 |
| 84-74-2 | Dibutyl phthalate ^f | 2.9 E-06 | 2.9 E-06 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | ND | 4.9 E-04 |
| 541-73-1 | 1,3-Dichlorobenzene | ND | ND | 4.3 E-04 |

TABLE A4 (cont.)

| | | Emission | Factor ^{b,c} | Minimum Detection Level mg/m ^{3,e} |
|--------------------|---|-------------|-------------------------------|---|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | ND | 4.3 E-04 |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | ND | 1.6 E-02 |
| 75-27-4 | Dichlorobromomethane | ND | ND | 4.1 E-04 |
| 75-71-8 | Dichlorodifluoromethaneg | 1.0 E-09 | 1.0 E-09 | |
| 75-34-3 | 1,1-Dichloroethane | ND | ND | 3.7 E-04 |
| 107-06-2 | 1,2-Dichloroethane | ND | ND | 2.1 E-04 |
| 75-43-4 | Dichlorofluoromethane | ND | ND | 2.1 E-04 |
| 120-83-2 | 2,4-Dichlorophenol | ND | ND | 1.6 E-02 |
| 78-87-5 | 1,2-Dichloropropane | ND | ND | 3.3 E-04 |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | ND | 2.3 E-04 |
| 76-14-2 | Dichlorotetrafluoroethane | ND | ND | 5.0 E-04 |
| 105-67-9 | 2,4-Dimethylphenol | ND | ND | 1.6 E-02 |
| 131-11-3 | Dimethyl phthalate | ND | ND | 1.6 E-02 |
| 99-65-0 | 1,3-Dinitrobenzene | ND | ND | 6.7 E-04 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | ND | ND | 1.6 E-02 |
| 51-28-5 | 2,4-Dinitrophenol | ND | ND | 1.6 E-02 |
| 121-14-2 | 2,4-Dinitrotoluene | 1.5 E-06 | 1.5 E-06 | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | ND | 6.7 E-04 |
| 123-91-1 | 1,4-Dioxane | ND | ND | 3.3 E-04 |
| | Total dioxin/furan compounds | 2.5 E-10 | 2.5 E-10 | |
| 140-88-5 | Ethyl acrylate | ND | ND | 1.7 E-04 |
| 100-41-4 | Ethylbenzene ^g | ND | ND | 3.1 E-04 |
| 117-81-7 | bis(2-Ethylhexyl)phthalate ^f | 9.9 E-06 | 9.9 E-06 | |
| 206-44-0 | Fluorantheneg | ND | ND | 1.6 E-02 |
| 86-73-7 | Fluoreneg | 2.1 E-08 | 2.1 E-08 | |
| 50-00-0 | Formaldehyde | 5.8 E-05 | 5.8 E-05 | |
| 76-13-1 | Freon 113 ^g | ND | ND | 5.5 E-04 |
| 35822-46-9 | 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 1.5 E-11 | 1.5 E-11 | |
| 67562-39-4 | 1,2,3,4,6,7,8- Heptachlorodibenzofuran | 1.8 E-12 | 1.8 E-12 | |
| 55673-89-7 | 1,2,3,4,7,8,9- Heptachlorodibenzofuran | 5.5 E-13 | 5.5 E-13 | |
| 118-74-1 | Hexachlorobenzene | ND | ND | 1.6 E-02 |
| 87-68-3 | Hexachlorobutadiene | ND | ND | 7.6 E-04 |

TABLE A4 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 77-47-4 | Hexachlorocyclopentadiene | ND | ND | 1.6 E-02 |
| 39227-28-6 | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | ND | ND | 3.9 E-09 |
| 57653-85-7 | 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | ND | ND | 4.0 E-09 |
| 19408-74-3 | 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | ND | ND | 4.0 E-09 |
| 57117-44-9 | 1,2,3,6,7,8-Hexachlorodibenzofuran | 4.4 E-13 | 4.4 E-13 | |
| 72918-21-9 | 1,2,3,7,8,9-Hexachlorodibenzofuran ^h | ND | ND | 3.1 E-09 |
| 60851-34-5 | 2,3,4,6,7,8-Hexachlorodibenzofuran | 5.4 E-13 | 5.4 E-13 | |
| 67-72-1 | Hexachloroethane | ND | ND | 5.9 E-04 |
| 7647-01-0 | Hydrochloric acid | ND | ND | 1.7 E-01 |
| 74-90-8 | Hydrogen cyanide | 4.4 E-05 | 4.4 E-05 | |
| 7664-39-3 | Hydrogen fluoride | ND | ND | 1.8 E-01 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | ND | ND | 1.6 E-05 |
| 78-59-1 | Isophorone | ND | ND | 1.6 E-02 |
| 7439-92-1 | Lead | 2.0 E-04 | 2.0 E-04 | |
| 7439-96-5 | Manganese ^g | 4.3 E-05 | 4.3 E-05 | |
| 96-33-3 | Methyl acrylate | ND | ND | 2.5 E-04 |
| 74-95-3 | Methylene bromide | ND | ND | 5.1 E-04 |
| 75-09-2 | Methylene chloride ^g | 0 | 0 | |
| 78-93-3 | Methyl ethyl ketone ^f | ND | ND | 8.1 E-04 |
| 74-88-4 | Methyl iodide | ND | ND | 3.5 E-04 |
| 108-10-1 | Methyl isobutyl ketone | ND | ND | 3.3 E-04 |
| 80-62-6 | Methyl methacrylate | ND | ND | 4.6 E-04 |
| 91-57-6 | 2-Methylnaphthalene | ND | ND | 1.6 E-02 |
| 95-48-7 | 2-Methylphenol | ND | ND | 1.6 E-02 |
| 1634-04-4 | Methyl tert-butyl ether | ND | ND | 3.7 E-04 |
| 91-20-3 | Naphthaleneg | 2.6 E-07 | 2.6 E-07 | |
| 7440-02-0 | Nickel ^g | ND | ND | 2.2 E-02 |
| 7697-37-2 | Nitric acid ^g | 4.5 E-04 | 4.5 E-04 | |
| 100-01-6 | 4-Nitroaniline | ND | ND | 1.6 E-02 |
| 98-95-3 | Nitrobenzene | ND | ND | 6.7 E-04 |
| 55-63-0 | Nitroglycerin | ND | ND | 6.7 E-04 |
| 88-75-5 | 2-Nitrophenol | ND | ND | 1.6 E-02 |

TABLE A4 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|--|-------------|-------------------------------|--|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 100-02-7 | 4-Nitrophenol | ND | ND | 1.6 E-02 |
| 62-75-9 | n-Nitrosodimethylamine | ND | ND | 1.6 E-02 |
| 86-30-6 | n-Nitrosodiphenylamine | ND | ND | 1.6 E-02 |
| 621-64-7 | n-Nitroso-di-n-propylamine | ND | ND | 1.6 E-02 |
| 3268-87-9 | 1,2,3,4,6,7,8,9-Octachlorodibenzo-p-dioxin | 2.2 E-10 | 2.2 E-10 | |
| 39001-02-0 | 1,2,3,4,6,7,8,9- Octachlorodibenzofuran | 3.4 E-12 | 3.4 E-12 | |
| 40321-76-4 | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | ND | ND | 4.2 E-09 |
| 57117-41-6 | 1,2,3,7,8-Pentachlorodibenzofuran | ND | ND | 3.4 E-09 |
| 57117-31-4 | 2,3,4,7,8-Pentachlorodibenzofuran ^h | 7.0 E-13 | 7.0 E-13 | |
| 87-86-5 | Pentachlorophenol | ND | ND | 1.6 E-02 |
| 85-01-8 | Phenanthrene ^g | 1.3 E-07 | 1.3 E-07 | |
| 108-95-2 | Phenol | ND | ND | 1.6 E-02 |
| 123-38-6 | Propionaldehyde | ND | ND | 1.7 E-01 |
| 129-00-0 | Pyrene | ND | ND | 1.6 E-02 |
| 7782-49-2 | Selenium | 3.8 E-06 | 3.8 E-06 | |
| 7440-22-4 | Silver | ND | ND | 2.2 E-02 |
| 100-42-5 | Styrene | ND | ND | 3.0 E-04 |
| 7664-93-9 | Sulfuric acid | ND | ND | 1.7 E-01 |
| 1746-01-6 | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | ND | ND | 2.7 E-09 |
| 51207-31-9 | 2,3,7,8-Tetrachlorodibenzofuran | 8.6 E-13 | 8.6 E-13 | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | ND | 3.5 E-04 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | ND | 3.5 E-04 |
| 127-18-4 | Tetrachloroethylene | ND | ND | 4.1 E-04 |
| 7440-28-0 | Thallium | ND | ND | 2.2 E-02 |
| 108-88-3 | Toluene ^g | ND | ND | 2.7 E-04 |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | ND | 1.6 E-02 |
| 71-55-6 | 1,1,1-Trichloroethane | ND | ND | 4.4 E-04 |
| 79-00-5 | 1,1,2-Trichloroethane | ND | ND | 2.8 E-04 |
| 75-69-4 | Trichlorofluoromethane | 5.8 E-10 | 5.8 E-10 | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | ND | 1.6 E-02 |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | ND | 1.6 E-02 |
| 96-18-4 | 1,2,3-Trichloropropane | ND | ND | 4.9 E-04 |

TABLE A4 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|-----------------------|---|-------------|-------------------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item | lb per lb NEW ^d | Detection Level mg/m ^{3,e} |
| 95-63-6 | 1,2,4-Trimethylbenzene ^g | 0 | | |
| 540-84-1 | 2,2,4-Trimethylpentane ^g | ND | ND | 1.9 E-04 |
| 7440-62-2 | Vanadium | 2.6 E-04 | 2.6 E-04 | |
| 108-05-4 | Vinyl acetate | ND | ND | 3.9 E-04 |
| 75-01-4 | Vinyl chloride ^h | 8.0 E-08 | 8.0 E-08 | |
| 75-35-4 | Vinylidene chloride | ND | ND | 2.1 E-04 |
| 106-42-3, 108-38-3 | m-Xylene, p-Xylene ^g | ND | ND | 4.0 E-04 |
| 95-47-6 | o-Xylene ^g | ND | ND | 4.9 E-04 |
| 7440-66-6 | Zinc ^g | 3.3 E-04 | 3.3 E-04 | |
| | Other Po | llutants | | |
| 67-64-1 | Acetone | 0 | 0 | |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene ^g | ND | ND | 6.7 E-04 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | ND | ND | 6.7 E-04 |
| 100-52-7 | Benzaldehyde | ND | ND | 1.7 E-01 |
| 100-51-6 | Benzyl alcohol | ND | ND | 1.6 E-02 |
| 108-86-1 | Bromobenzene | ND | ND | 3.9 E-04 |
| 7440-70-2 | Calcium ^f | 3.5 E-04 | 3.5 E-04 | |
| 7440-44-0 | Total carbon ^g | 8.4 E-03 | 8.4 E-03 | |
| 39638-32-9 | bis(2-Chloroisopropyl)ether | ND | ND | 1.6 E-02 |
| 59-50-7 | 4-Chloro-3-methylphenol | ND | ND | 1.6 E-02 |
| 124-48-1 | Dibromochloromethane | ND | ND | 5.2 E-04 |
| 156-59-2 | cis-1,2-Dichloroethene | ND | ND | 2.0 E-04 |
| 156-60-5 | trans-1,2-Dichloroethene | ND | ND | 4.8 E-04 |
| 84-66-2 | Diethylphthalate | ND | ND | 1.6 E-02 |
| 5779-94-2 | 2,5-Dimethylbenzaldehyde | ND | ND | 1.7 E-01 |
| 117-84-0 | Di-n-octylphthalate ^g | ND | ND | 1.6 E-02 |
| 629-97-0 | Docosane ^h | 1.8 E-06 | 1.8 E-06 | |
| 141-78-6 | Ethyl acetate | ND | ND | 1.4 E-03 |
| 97-63-2 | Ethyl methacrylate | ND | ND | 3.3 E-04 |
| 622-96-8 | 4-Ethyltoluene ^g | ND | ND | 3.0 E-04 |
| 142-82-5 | Heptane | ND | ND | 2.1 E-04 |
| 544-76-3 | Hexadecane ^h | 1.6 E-06 | 1.6 E-06 | |
| 66-25-1 | Hexaldehyde | ND | ND | 1.7 E-01 |

TABLE A4 (cont.)

| | | Emission | Factor ^{b,c} | Minimum |
|--------------------|--|--|-----------------------|-------------------------------------|
| CASRN ^a | Pollutant | lb per item lb per lb NEW ^d | | Detection Level mg/m ^{3,e} |
| 103-23-1 | Hexanedioic acid, bis(2-ethylhexyl) ester ^h | 5.2 E-06 | 5.2 E-06 | |
| 591-78-6 | 2-Hexanone | ND | ND | 2.9 E-04 |
| 2691-41-0 | HMX | ND | ND | 1.3 E-03 |
| 10035-10-6 | Hydrogen bromide | ND | ND | 1.7 E-01 |
| 112-95-8 | Icosane ^h | 4.2 E-06 | 4.2 E-06 | |
| 590-86-3 | Isovaleraldehyde | ND | ND | 1.7 E-01 |
| 7439-95-4 | Magnesium ^g | 3.8 E-04 | 3.8 E-04 | |
| 98-83-9 | alpha-Methylstyrene | ND | ND | 2.9 E-04 |
| 88-74-7 | 2-Nitroaniline | ND | ND | 1.6 E-02 |
| 99-09-2 | 3-Nitroaniline | ND | ND | 1.6 E-02 |
| 88-72-2 | 2-Nitrotoluene | ND | ND | 6.7 E-04 |
| 99-08-1 | 3-Nitrotoluene | ND | ND | 6.7 E-04 |
| 99-99-0 | 4-Nitrotoluene | ND | ND | 6.7 E-04 |
| 111-65-9 | Octane | ND | ND | 1.9 E-04 |
| 7664-38-2 | Phosphoric acid | ND | ND | 1.7 E-01 |
| 121-82-4 | RDX | 2.4 E-06 | 2.4 E-06 | |
| 353-66-2 | Silane, difluorodimethyl-h | 2.1 E-06 | 2.1 E-06 | |
| 7440-24-6 | Strontium | 3.8 E-04 | 3.8 E-04 | |
| 479-45-8 | Tetryl | ND | ND | 6.7 E-04 |
| 7440-32-6 | Titanium | ND | ND | 0.02167542 |
| 1334-78-7 | o,m,p-Tolualdehyde | ND | ND | 5.3 E-01 |
| 25323-89-1 | Trichloroethane | ND | ND | 3.0 E-04 |
| 108-67-8 | 1,3,5-Trimethylbenzene ^g | ND | ND | 2.5 E-04 |
| 99-35-4 | 1,3,5-Trinitrobenzene ^g | 8.3 E-06 | 8.3 E-06 | |
| 118-96-7 | 2,4,6-Trinitrotoluene ^h | 1.3 E-07 | 1.3 E-07 | |
| 110-62-3 | Valeraldehyde | ND | ND | 1.7 E-01 |

^a CASRN = Chemical Abstracts Service Registry Number.

b ND = nondetected.

^c Emission factors rated C unless otherwise noted.

d NEW = net explosive weight. The NEW for this ordnance is 1.00 pound per item.

^e Data provided for compounds that were not detected.

f Emission factor rated A because of correlation with emission factors for similar ordnance and number of test data points.

g Emission factor rated B because of correlation with emission factors for similar ordnance and number of test data points. $^{\rm h}$ Emission factor rated D because the factor is based upon C-rated test data.

APPENDIX B

NEW AP-42 SECTIONS FOR ORDNANCE INCLUDED IN EXPLODING ORDNANCE EMISSION STUDY PHASE II SERIES 2 TESTING AT ABERDEEN PROVING GROUND, MARYLAND

Electronic versions of the new AP-42 sections for ordnance included in Exploding Ordnance Emission Study Phase II Series 2 testing at Aberdeen Proving Ground, Maryland, are located on the EPA website at: http://www.epa.gov/ttn/chief/ap42/index.html.

APPENDIX C

PUBLIC COMMENTS ON AP-42 SECTIONS FOR ORDNANCE INCLUDED IN PHASE II SERIES 2 TESTING AT ABERDEEN PROVING GROUND, MARYLAND

No public comments were received on the draft version of this document. However, the document has been modified since the publication of the draft to incorporate revised emission factor ratings for numerous pollutants. The methodology used to revise the emission factor ratings is presented in Section 5 of this document, while the revised emission factor ratings are presented in Appendix A.