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CONTROLLED BURN AIR SAMPLING TECHNICAL REPORT

FINAL
August 2001

Prepared for

Directorate of Safety, Health and Environment
U.S. Army Garrison
Aberdeen Proving Ground, Maryland

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Docket No. 40-8838-ML

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ACRONYMS AND ABBREVIATIONS

| | |
|---------------------|---|
| ALI | Annual Limit on Intake |
| ANL | Argonne National Laboratory |
| APG | Aberdeen Proving Ground |
| ATC | Aberdeen Test Center |
| CFR | Code of Federal Regulations |
| DAAMS | Depot Area Air Monitoring System |
| DAC | Derived Air Concentration |
| DDT | Dichlorodiphenyltrichloroethane |
| DU | Depleted Uranium |
| ECBC | Edgewood Chemical Biological Center |
| EPA | U.S. Environmental Protection Agency |
| GC/ECD | Gas Chromatography/Electron Capture Detector |
| GC/FID | Gas Chromatography/Flame Ionization Detector |
| GC/FPD | Gas Chromatography/Flame Photometric Detector |
| GC/MS | Gas Chromatograph/Mass Spectrometer |
| HPLC | High Performance Liquid Chromatography |
| HQ | Hazard Quotient |
| ICP | Inductively Coupled Plasma |
| IRP | Installation Restoration Program |
| K-40 | Potassium-40 |
| MDE | Maryland Department of the Environment |
| µg | microgram |
| µg/m ³ | microgram per cubic meter |
| m ³ /min | cubic meters per minute |
| mph | miles per hour |
| OSHA | Occupational Safety and Health Administration |
| Pb-212 | Lead-212 |
| PCB | Polychlorinated Biphenyl |
| PCi | pico-Curies |
| pCi/m ³ | pico-Curies per cubic meter |
| PEL | Permissible Exposure Limit |
| ppb | parts per billion |
| PUF | Polyurethane Foam |
| Ra-223 | Radium-223 |
| Ra-224 | Radium-224 |
| RBC | Risk-Based Concentration |
| RfD | Reference Dose |
| RF-RBC | Range Fire Risk-Based Concentration |
| TAP | Toxic Air Pollutant |
| TCE | Trichloroethene |
| TIC | Tentatively Identified Compound |
| TLV | Threshold Limit Value |
| VOC | Volatile Organic Compound |
| U-235 | Uranium-235 |
| UXO | Unexploded Ordnance |

1.0 INTRODUCTION

The Aberdeen Proving Ground (APG) Installation Restoration Program (IRP) identified the need to conduct a limited number of controlled burns at test ranges in the Aberdeen and Edgewood Areas of APG to support sampling of air emissions produced by range fires. The purpose of the study was to generate, to the extent possible, quantifiable emissions representative of fires occurring at APG test ranges to allow a screening assessment of potential human health impacts. This Technical Report documents the controlled burn air emissions sampling conducted during three events, and provides human health risk screening of the analytical data obtained.

APG lies on the northwestern shore of the Chesapeake Bay in southern Harford County and southeastern Baltimore County, Maryland (Figure 1). Major geographical areas bordering APG include the Chesapeake Bay and its tributaries; Gunpowder Falls State Park; the Crane Power Plant; and the towns of Bel Air, Edgewood, Joppatowne/Magnolia, Aberdeen, and various smaller residential areas. APG occupies 72,500 acres of land and water, of which approximately 75 percent are range areas. The Bush River divides the Installation into two noncontiguous areas commonly referred to as the Aberdeen Area and the Edgewood Area.

Since 1917, the Edgewood Area has been the center for the research, development, testing and manufacturing of military-related chemicals and agents. Activities at the Edgewood Area have included laboratory research and development, field testing, and pilot- and production-scale manufacture of chemical warfare agents. Chemical warfare materiel, hazardous wastes, and low-level radiological wastes have been stored at the Edgewood Area. The Aberdeen Area was established as the Ordnance Proving Ground in 1917, and has been the site of intense research and development, large-scale testing of munitions, weapons, and materiel, and a training school for ordnance officers and enlisted specialists.

2.0 BACKGROUND

Ordnance firing, other test activities, residual white phosphorus in subsurface soils, and lightning strikes occasionally cause accidental fires in the test range areas of APG. These accidental fires may occur under unfavorable weather conditions (e.g., meteorological conditions may create wind directions that transport range fire smoke plumes toward residential communities), and the amount of vegetative fuel and acreage burned cannot always be controlled during these unplanned burns.

APG's long history of weapons testing and past disposal practices caused members of the public to express concerns that contaminants accumulated in surface soils and vegetation could be transported in smoke plumes produced by such fires. The potential transport and deposition of contaminants via the smoke plume and the associated health risks were of greatest concern to the public. Sources of contamination could include residues in and on vegetative matter and surface soils from previous weapons testing and disposal of hazardous substances; chemicals released from burning of uncontaminated vegetation; and detonation or rupture of unexploded ordnance (UXO).

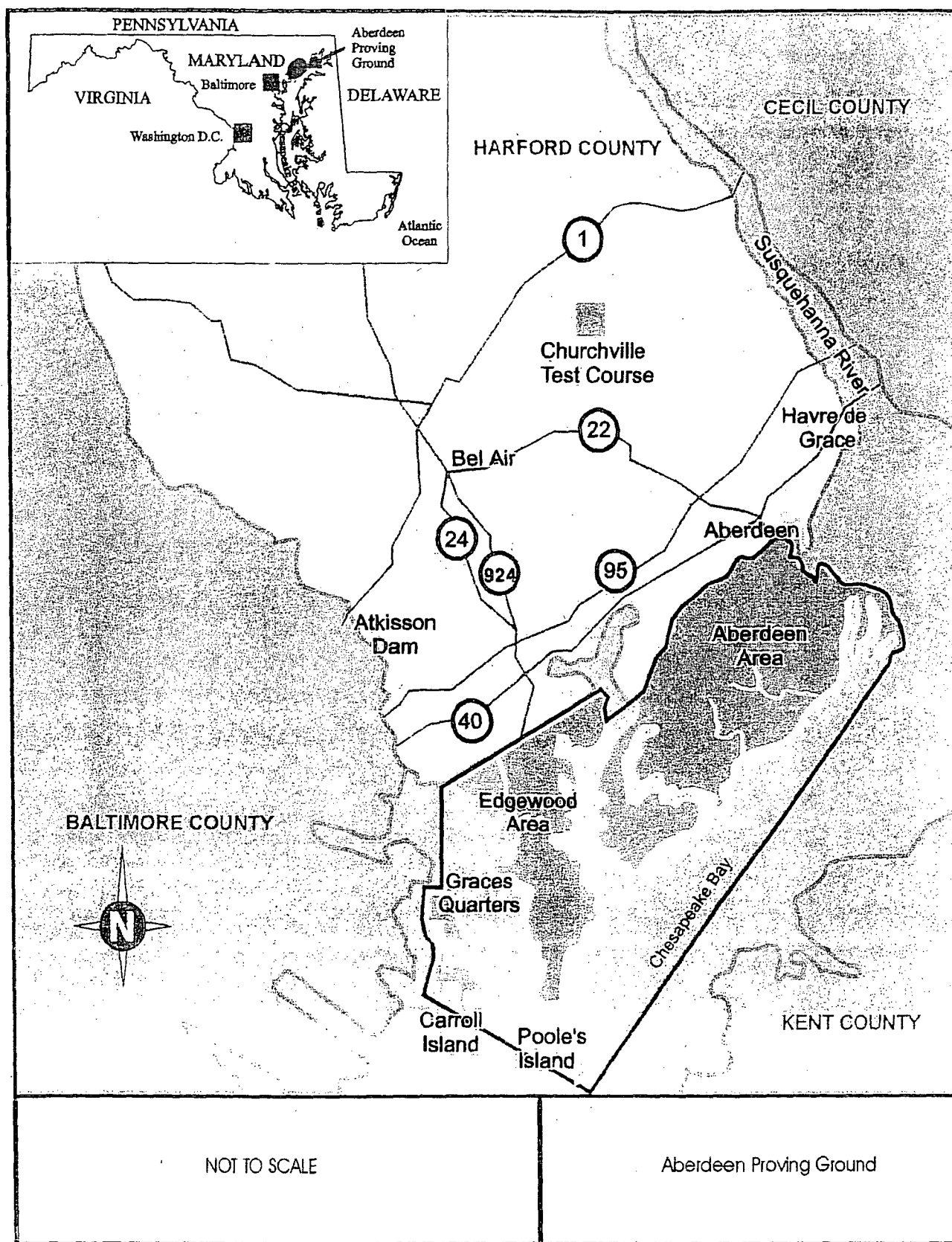


Figure 1. Location of Aberdeen Proving Ground

3.0 OTHER STUDIES

3.1 Argonne National Laboratory 1998 Report

The Environmental Assessment Division of Argonne National Laboratory (ANL) completed a study in 1998 in response to the public concerns. The study, entitled "Potential Human Health Impacts from Range Fires at Aberdeen Proving Ground, Maryland," used atmospheric dispersion models to evaluate potential human health impacts resulting from exposure to contaminants resulting from range fires. The screening study focused specifically on five contaminants considered most likely to be present in surface soils and vegetation as a result of past activities at APG, two chemical agents, and two naturally-occurring compounds released during burning of uncontaminated vegetation. The contaminants, selected with input from APG personnel and a citizens advisory committee, were lead, arsenic, depleted uranium (DU), trichloroethene (TCE), dichlorodiphenyltrichloroethane (DDT), vinyl acetate, 2-furaldehyde, and mustard and phosgene in UXO. The modeling results were compared to U.S. Environmental Protection Agency (EPA) action levels.

The study concluded that range fires at APG do not pose a significant health risk to APG workers or the surrounding populations. Use of conservative assumptions in the study provided an upper bound estimate of potential risk. The study recommended future efforts be directed at fire management and control to reduce the occurrence and duration of range fires. The IRP elected to conduct a series of controlled burns for data collection purposes in response to on-going public concerns relating to range fires and potential risk to human health.

3.2 Argonne National Laboratory 2000 Report

The original report prepared by ANL in 1998 was modified in October 2000 to include actual air emissions data collected during the J-Field controlled burn conducted in April 2000. The updated report incorporated measured contaminant levels in vegetation samples taken from the Toxic Burn Pit area of J-Field. The data was used in the FIREPLUME computer model to calculate estimated ground-level contaminant concentrations during a range fire. The study then estimated exposure levels using conservative assumptions to evaluate impacts to human health. The model-predicted concentrations were one to two orders of magnitude greater than the field measured concentrations due to the use of conservative assumptions. The study concluded that the risk of adverse health effects from mobilization of contaminants as a result of range fires is extremely small. The study again recommended that future efforts be directed at range management to reduce the number of unplanned fires. The range management efforts could effectively include controlled burns.

3.3 Environmental Protection Agency Data Collection

Lockheed Martin, under contract to the EPA through the Response, Engineering, and Analytical Contract, collected air samples for analysis during two O-Field burn attempts and the J-Field controlled burn. Samples were collected for analysis of dioxins, metals, polynuclear hydrocarbons, inorganic acids, volatile organic compounds (VOCs), and chemical agents. Particulate monitoring was also conducted using an MIE DataRAM. The three trip reports for these sampling events are included in Appendix A for reference. The data was not incorporated into the evaluation performed as part of this report.

4.0 CONTROLLED BURN STUDY PARAMETERS

The controlled burn study parameters were developed in close coordination with the U.S. Army Aberdeen Test Center (ATC), the agency responsible for range management and control at APG. Meteorological and range control personnel supported the development of the study parameters. In addition, close coordination with the APG Fire Department and Safety personnel provided guidance for developing sampling protocols and selecting range sites for the controlled burns. Input from the citizens advisory committee was solicited regarding potential controlled burn locations.

4.1 *Meteorological Conditions*

The controlled burns for air emissions sampling were conducted under specific meteorological conditions to minimize potential impacts to civilian communities and to facilitate data collection. Wind directions were selected to minimize travel of the plumes toward populated areas. In general, the controlled burns were initiated with north-northeast or west-southwest wind directions. Controlled burns were initiated only with wind speeds of 15 miles per hour (mph) or less. Greater wind speeds would have resulted in difficulty in controlling and extinguishing the fires, as well as a reduction in the sampling period. Atmospheric stability Class D or Class C conditions were selected as burn parameters to obtain the most rapid return of range fire smoke to ground level and limited atmospheric dispersion. The site-specific burn plans developed for each controlled burn location provide specific details and procedures.

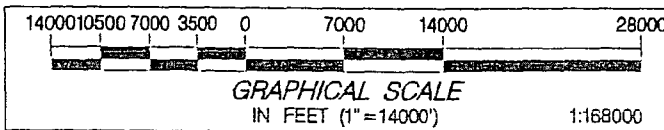
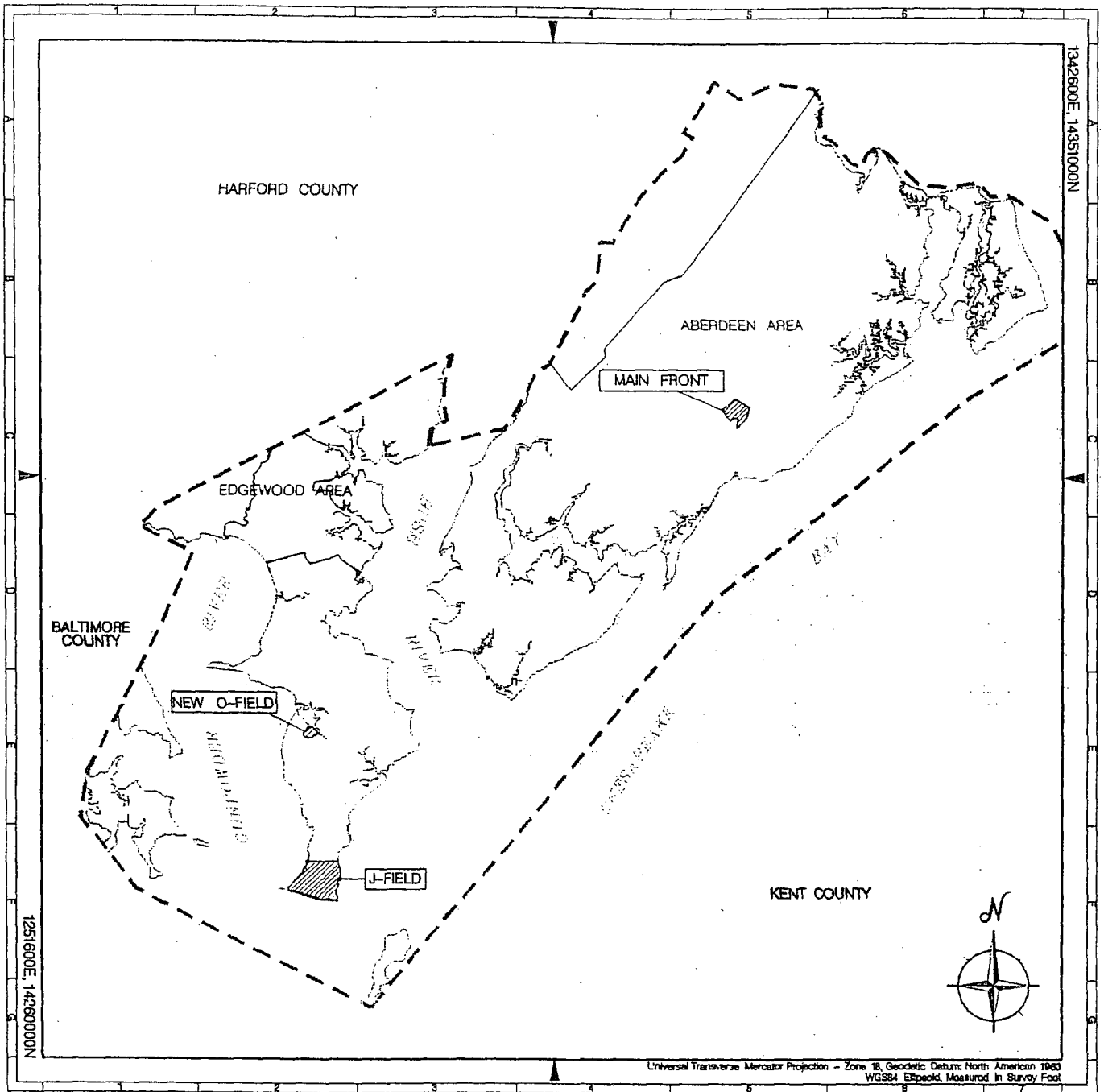
4.2 *Test Range Selection*

Selection of range areas with the most potential surface soil contamination provided a "worst case" scenario for the controlled burn sampling events. With input from IRP, ATC, Fire Department, and Safety personnel, and the citizens advisory committee, three range areas (Figure 2) were selected for controlled burns and air emissions sampling:

- Main Front range in APG Aberdeen Area — selected as representative of test ranges with the highest potential DU contamination and other toxic compounds
- J-Field in APG Edgewood Area — selected as representative of worst-case air emissions due to historical testing and disposal activities, and based on soil contamination data collected as part of the IRP
- New O-Field in APG Edgewood Area — selected as representative of worst-case air emissions due to historical testing and disposal activities, and based on contamination data collected as part of the IRP

4.3 *Sampling Locations and Analytes*

For each controlled burn, monitoring involved the collection of both upwind and downwind air samples. Upwind samples were collected during each burn to measure the level of potential contaminants in ambient air. Downwind sample locations were placed at specified distances from the fire ignition point to capture air samples within the smoke plumes upon return to ground level.



LEGEND

- Aberdeen Proving Ground
- Water
- Study Area Boundary
- Installation Boundary
- Range Area Selected For Controlled Burn

| | | | |
|--|---------------------------|--------------------------------------|--------------|
| 6700 Alexander Bell Drive Columbia, MD 21046 | | (800) 727-6677 www.genphysics.com | |
| TITLE: RANGE AREAS SELECTED FOR CONTROLLED BURNS | | | |
| CARTOGRAPHER: M. BROOKS | APPROVED BY: K. THORPE | DATE: 11-07-00 | FIGURE: 2 |

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Sampling collection and analysis methods are listed in Table 1. Equipment setup is illustrated in photographs provided in Appendix B.

Table 1. APG Controlled Burn Sample Collection and Analysis Methods

| Analyte | Sampling Method | Equipment | Analytical Method |
|------------------------------------|-----------------|---|--|
| Volatile Organic Compounds | USEPA TO-14 | Summa Canister – 3-hour sampling valve | GC/MS GC/FID |
| Explosives | TO-4 Modified | High-Volume Sampler Glass Fiber Filter and PUF | HPLC |
| Pesticides/PCBs | TO-4 | High-Volume Sampler Glass Fiber Filter and PUF | GC/ECD |
| Inorganics | 6010 Modified | High-Volume Samplers (2) Quartz Filter | ICP |
| Chemical Agents | DAAMS | Calibrated Pump and DAAMS Tubes | ECBC Analytical Method |
| Gross Alpha/Beta and Gamma Spectra | Quartz Filter | High-Volume Sampler Quartz Filter | EPA 900.0 (Modified) EPA 901.1 (Modified) |

DAAMS – Depot Area Air Monitoring System
ECBC – Edgewood Chemical Biological Center
GC/ECD – Gas Chromatograph/Electron Capture Detector
GC/FID – Gas Chromatograph/Flame Ionization Detector
PCB – Polychlorinated Biphenyl

PUF – Polyurethane Foam
GC/MS – Gas Chromatograph/Mass Spectrometer
HPLC – High Performance Liquid Chromatography
ICP – Inductively Coupled Plasma
EPA – Environmental Protection Agency

The EPA Method TO-14 is designed for sampling and analysis of volatile organic compounds (VOCs) in ambient air as collected in summa canisters or other specially prepared canisters. A sample of ambient air is drawn through a sampling train, comprised of components that regulate the rate and duration of sampling, into a pre-evacuated, passivated canister. The VOCs are separated by gas chromatography and measured by a mass spectrometer or by multi-detector techniques. Analysis of VOCs included reporting of up to 10 tentatively identified compounds (TICs).

Method TO-4 is a procedure for determination of a variety of organochlorine pesticides and polychlorinated biphenyls (PCBs) in ambient air. Method TO-4 utilizes a modified high volume sampler consisting of a glass fiber filter with a polyurethane foam (PUF) backup adsorbent cartridge used to sample ambient air at a rate of approximately 200 – 280 liters per minute. Flow rates for the high volume samplers are calculated during the calibration process prior to each sampling event. The high volume particulate sampler operates at an average flow rate of approximately 1.2 cubic meters per minute (m^3/min); the average flow rate for the high volume PUF sampler is approximately 0.2 m^3/min . The PCBs and pesticides are recovered by Soxhlet extraction and analyzed using gas chromatography with electron capture detection (GC/ECD). Samples collected using TO-4 (modified) are analyzed for explosives using high performance liquid chromatography (HPLC).

The EPA method 6010 utilizes Inductively Coupled Plasma (ICP) instrumentation with a high-temperature source for metals analysis. The samples are collected using a high-volume sampler and quartz filter media. The sample is prepared for ICP analysis by digesting the quartz filter in nitric acid.

The Depot Area Air Monitoring System (DAAMS) sampling method for chemical agents requires air flow through a solid sorbent tube at a controlled flow rate and a measured time period. The sampler flow rate is calibrated prior to the sampling event. The DAAMS tubes are analyzed by the

U.S. Army Edgewood Chemical Biological Center (ECBC) laboratory using thermal desorption techniques to strip the analytes into a gas chromatography/flame photometric detector (GC/FPD) analytical system.

For radioactivity, the analysis included gross alpha analysis for radionuclides that emit alpha particles, gross beta analysis for radionuclides that emit beta particles, and gamma ray spectroscopy for radionuclides that emit gamma rays. Of the methods employed, only gamma spectroscopy is capable of identifying the specific radionuclides and the amount of radioactive material present (in pico-Curies (pCi)) from that radionuclide. Gross alpha and gross beta analyses provide only the amount of radioactivity (pCi).

The specific analytes included in each sampling and analysis method are provided for reference in Appendix C.

4.4 Quality Assurance

The ambient air sampling of these short-term events (the controlled burns conducted at APG) is considered representative of fires in fields where these burns occurred. These events may not be representative of all fires, but can be considered "worst case" for evaluation of release of potential contaminants in vegetation at sites with soil contamination. The sampling efforts incorporated numerous quality assurance methods to provide the best possible results.

- Equipment calibration was performed prior to each sampling event to provide accuracy in field measurements. Field instruments were calibrated according to manufacturers' specifications, and the calibration results were recorded.
- Use of high-volume sampling equipment, as appropriate, reduced errors potentially associated with low sample volumes, and achieved lower detection limits.
- Filter or media blanks for each sampling method (except the summa canister) were sent to the laboratory for analysis to detect filter or media contamination unrelated to the range fire sampling. The summa canisters were cleaned and evacuated by the analytical laboratory.
- Generators providing power supply for the sampling equipment were placed downwind or cross-wind from the sampling points to prevent interferences.
- Vehicles used to transport personnel and sampling equipment were parked downwind of the sampling equipment or removed from the sampling location.
- Samples were collected at an upwind location during each range fire sampling event to allow evaluation of ambient concentrations of detected analytes.

Sampling locations were dictated by the availability of established roads and by explosive fragment hazard distances. No sampling points were selected in off-road locations due to UXO safety considerations.

5.0 CONTROLLED BURN EVENTS AND RESULTS

Three controlled burn events were conducted from April 1999 through April 2000. Burn events were conducted in the J-Field and New O-Field ranges of the Edgewood Area, and in the Main

Front range of the Aberdeen Area of APG. A second burn event in the Main Front area was attempted in April 2001.

5.1 Main Front Range Controlled Burn – April 1999

A controlled burn was conducted on 28 April 1999, at the Main Front range in the Aberdeen Area of APG (Figure 3). Samples were collected at three sites downwind of the fire, and at one upwind site to serve as a background location. Downwind sampling sites SP1, SP2, and SP3 were located southwest of the burn site at distances of approximately 1000, 2000, and 3000 meters, respectively.

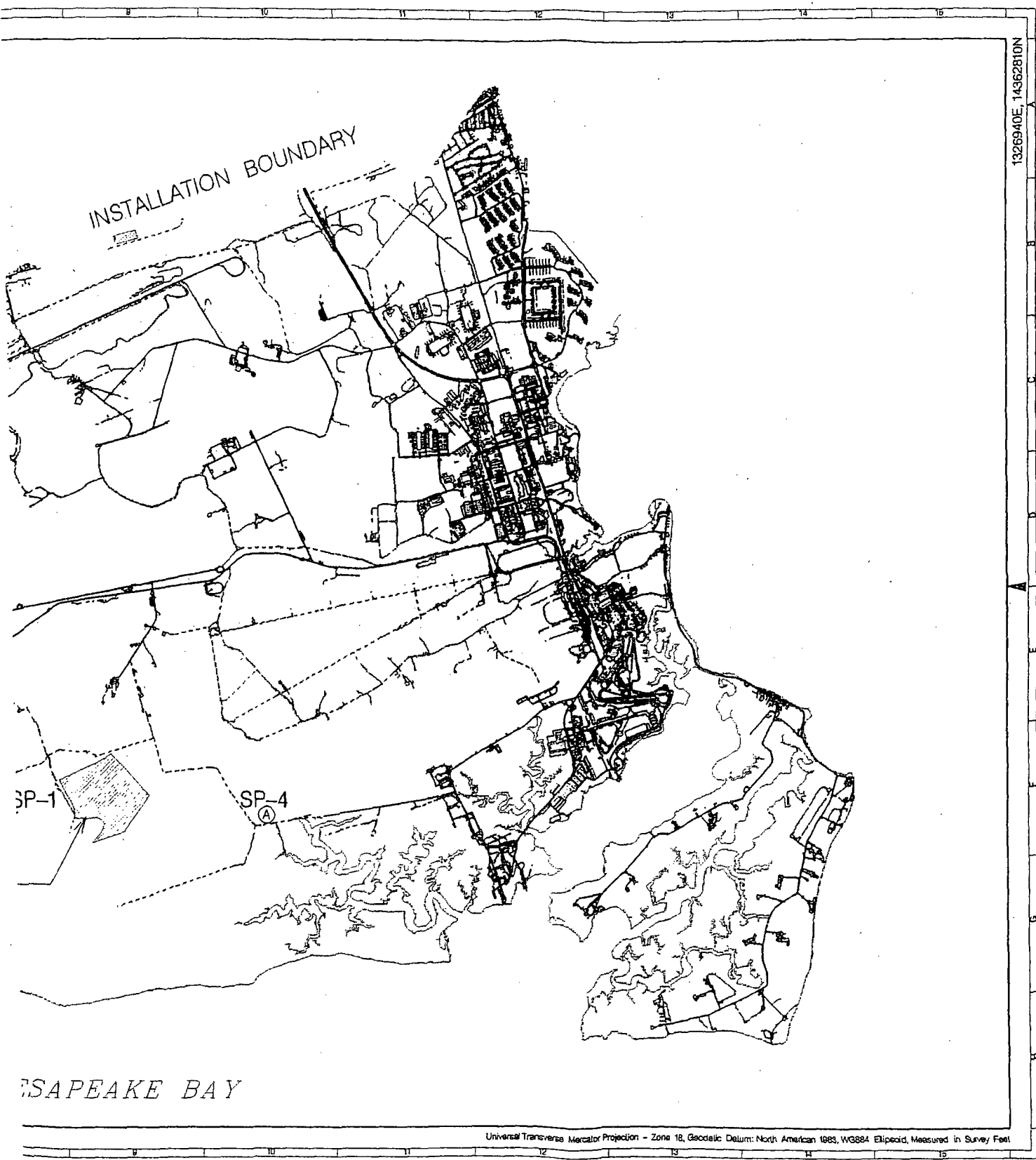
With favorable meteorological conditions forecasted by the ATC Meteorological Office, the fire ignition by the APG Fire Department occurred at approximately 1500 hours. The sampling duration was approximately four hours.

Meteorological data collected during the controlled burn show that the wind direction shifted widely during the course of the burn. The forecasted wind direction was from the northeast (i.e., blowing toward the Edgewood Area and down the Chesapeake Bay); the average winds during the sampling event were from the southeast. The shifting wind direction resulted in a reduction of the burn area and intensity of the fire, causing less smoke to be produced. The variable wind direction resulted in exposure of the upwind sampling point (SP4) to smoke during a portion of the sampling period. Photographs taken during the Main Front burn event are presented in Appendix D-1.

Detections of several analytes were reported for the 1999 Main Front controlled burn event:

- Several VOCs were detected, including:
 - acetone
 - nonane
 - toluene
 - decane
 - methylene chloride
 - xylene.
- Analysis of the PUF media yielded detection of 2,2',3,4,5'-pentachlorobiphenyl and 2,4',5-trichlorobiphenyl in the SP3 sample, and 2,4',5-trichlorobiphenyl in the SP4 sample. These PCBs were detected at a concentration of approximately 1 ppb.
- One pesticide (dieldrin) was detected by the PUF samplers at all downwind sampling locations (SP-1, -2, and -3) in the parts per trillion range.
- Numerous metals were detected, but not at levels exceeding blank concentrations.

Chemical agents and explosives were not detected at the downwind or upwind sampling locations during the Main Front controlled burn. Additionally, no specific radionuclides were detected above the minimum detection activity, the analytical error, or the blank radionuclide activity level. Appendix E-1: Tables E-1 through E-10 contain the results from the analysis of the air samples collected during the Main Front controlled burn.



1500 5000 10000
GRAPHICAL SCALE
(1" = 5000')
1:60000

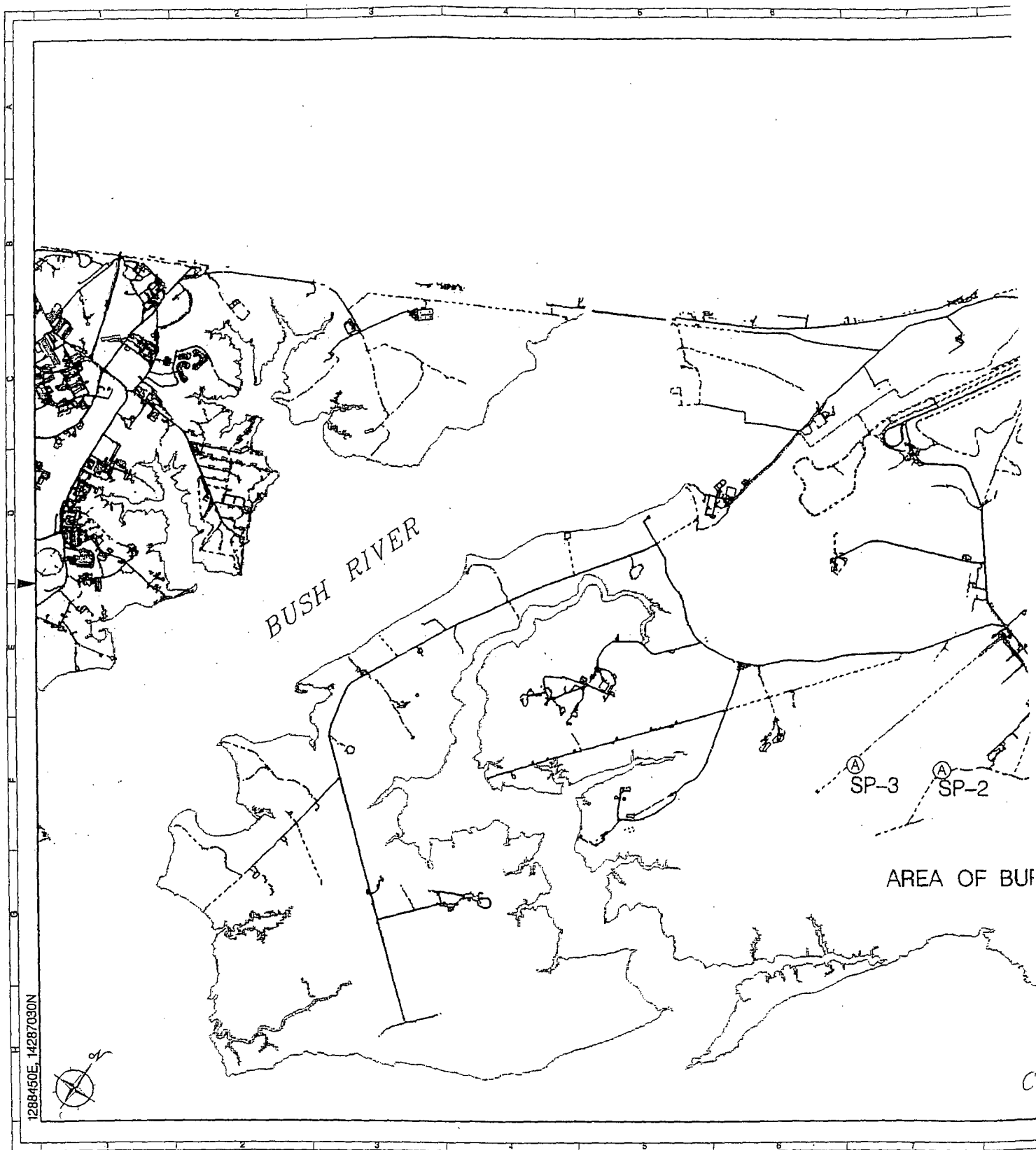
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MAIN FRONT CONTROLLED BURN AREA AND
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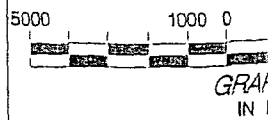


L E G E N D

— Paved Roads
 - - - Unpaved Roads

▬ Structures
 — Coastline

▬ Controlled Burn Area
 (A) Sampling Location



5.2 New O-Field Controlled Burn – December 1999

The New O-Field controlled burn occurred on 3 December 1999 in the Edgewood Area of APG (Figures 4 and 5). Downwind sampling locations SP1, SP2, and SP3 were located 335, 1200, and 2300 meters, respectively, from the burn area. Sampling point SP1 was located at Watson Creek, SP2 at Ricketts Point Road, and SP3 at Briery Point on the Bush River shoreline. Due to a slight variation in wind direction during the burn ($238^{\circ} \pm 20^{\circ}$), the SP2 sampling location was re-positioned within the smoke plume. The upwind sampling point SP4 was located on the Gunpowder River shoreline, approximately 500 meters from the burn location.

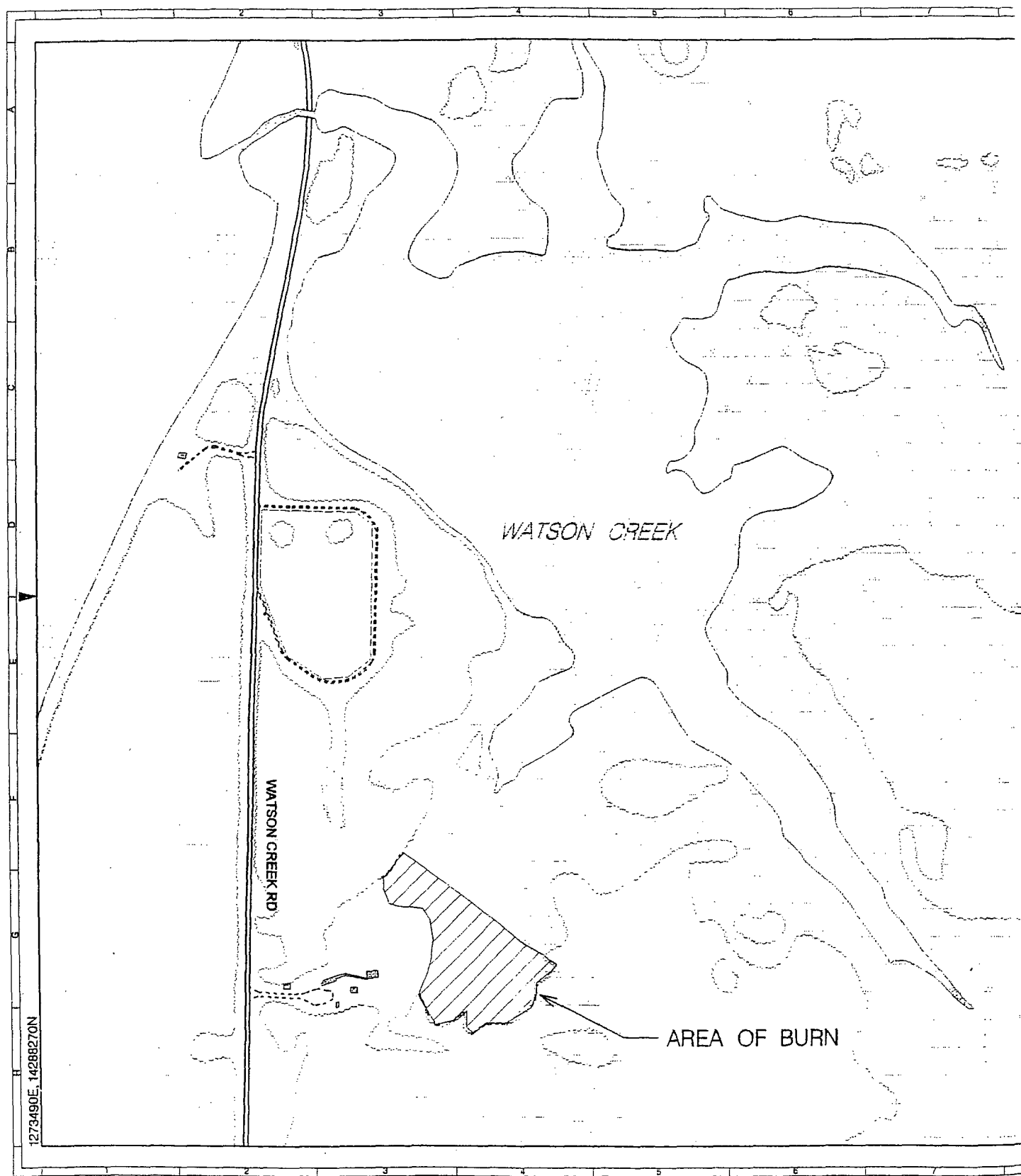
The meteorological forecast provided by the ATC Meteorological Office indicated winds speeds of less than 15 mph, a southwest wind direction, and Class D stability. Given the favorable forecast, the fire was ignited by the APG Fire Department at approximately 1530 hours. The sampling duration was approximately four hours.

The meteorological data collected during the controlled burn show that the wind direction generally remained from the southwest, with only slight variation during the course of the burn. However, reduced wind speeds, coupled with wet conditions in New O-Field, limited the size of the burn area during this event. Given that the area and the intensity of the burn were much less than anticipated, a reduced amount of smoke was produced from the fire for the air sampling event. Photographs of the area following the controlled burn are included in Appendix C-2.

Samples collected were analyzed for chemical agents, explosives, VOCs, PCBs, pesticides, inorganics, and radionuclides; detections of several analytes were reported:

- Several VOCs were detected in the ppb range, including:
 - acetone
 - benzene
 - benzonitrile
 - carbon disulfide
 - dedecene
 - hexane
 - methylene chloride
 - toluene
 - xylene
- 2,2',3,4,5-pentachlorobiphenyl was detected at sampling locations SP1, SP2, and SP3 in concentrations ranging from 0.0004 to 0.0011 micrograms per cubic meter ($\mu\text{g}/\text{m}^3$).
2,2',3,4,4'-tetrachlorobiphenyl was detected in the SP4 (upwind) sampling location at a concentration of 0.0020 $\mu\text{g}/\text{m}^3$.
- Numerous metals were detected in the ppb range in samples collected both upwind and downwind of the burn area.

Chemical agents, explosive compounds, and pesticides were not detected at the downwind or upwind sampling locations during the New O-Field burn. Appendix E-2: Tables E-11 through E-17 contain the results from the analysis of the air samples collected during the controlled burn at New O-Field.



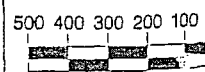
L E G E N D

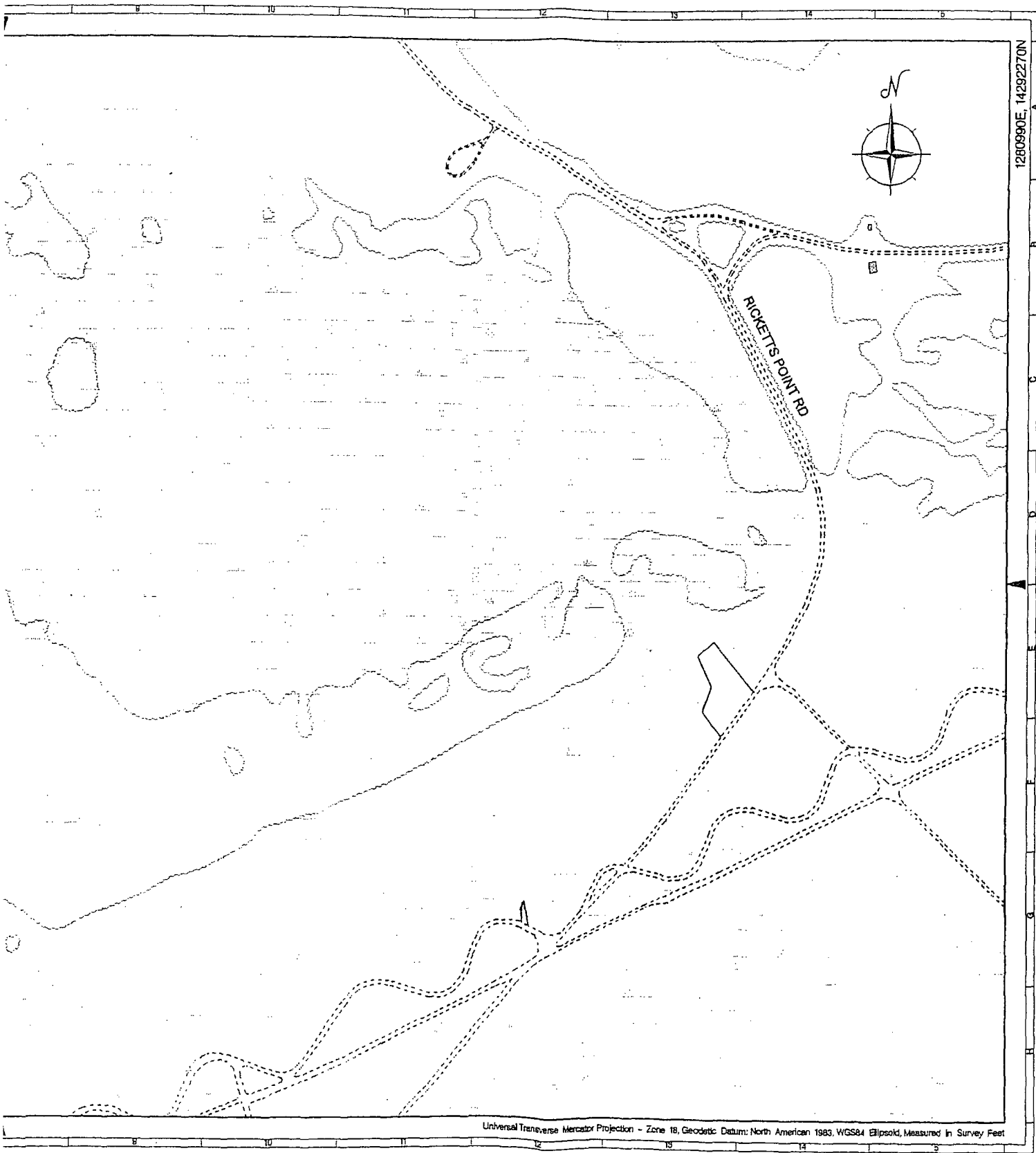
Structure
Paved Road

Unpaved Road
Treeline

Wetland
Water

Fence
Controlled Burn Area





Universal Transverse Mercator Projection - Zone 18, Geodetic Datum: North American 1983, WGS84 Ellipsoid, Measured in Survey Feet



GRAPHICAL SCALE
FEET (1" = 500')

1:6000

TITLE:

NEW O-FIELD CONTROLLED BURN AREA DECEMBER 1999



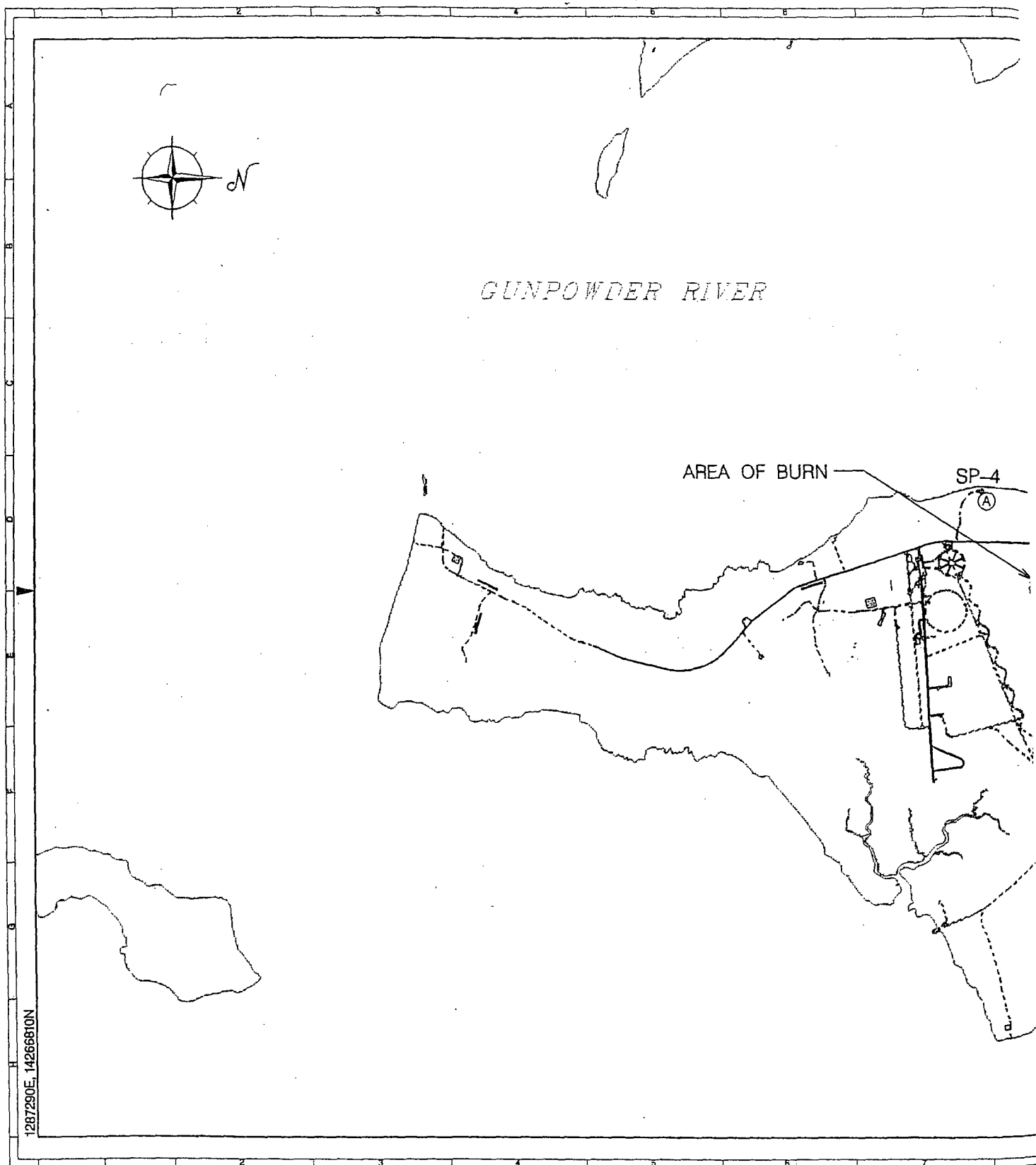
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
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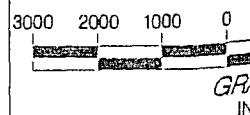
APPROVED BY:
K. THORPE

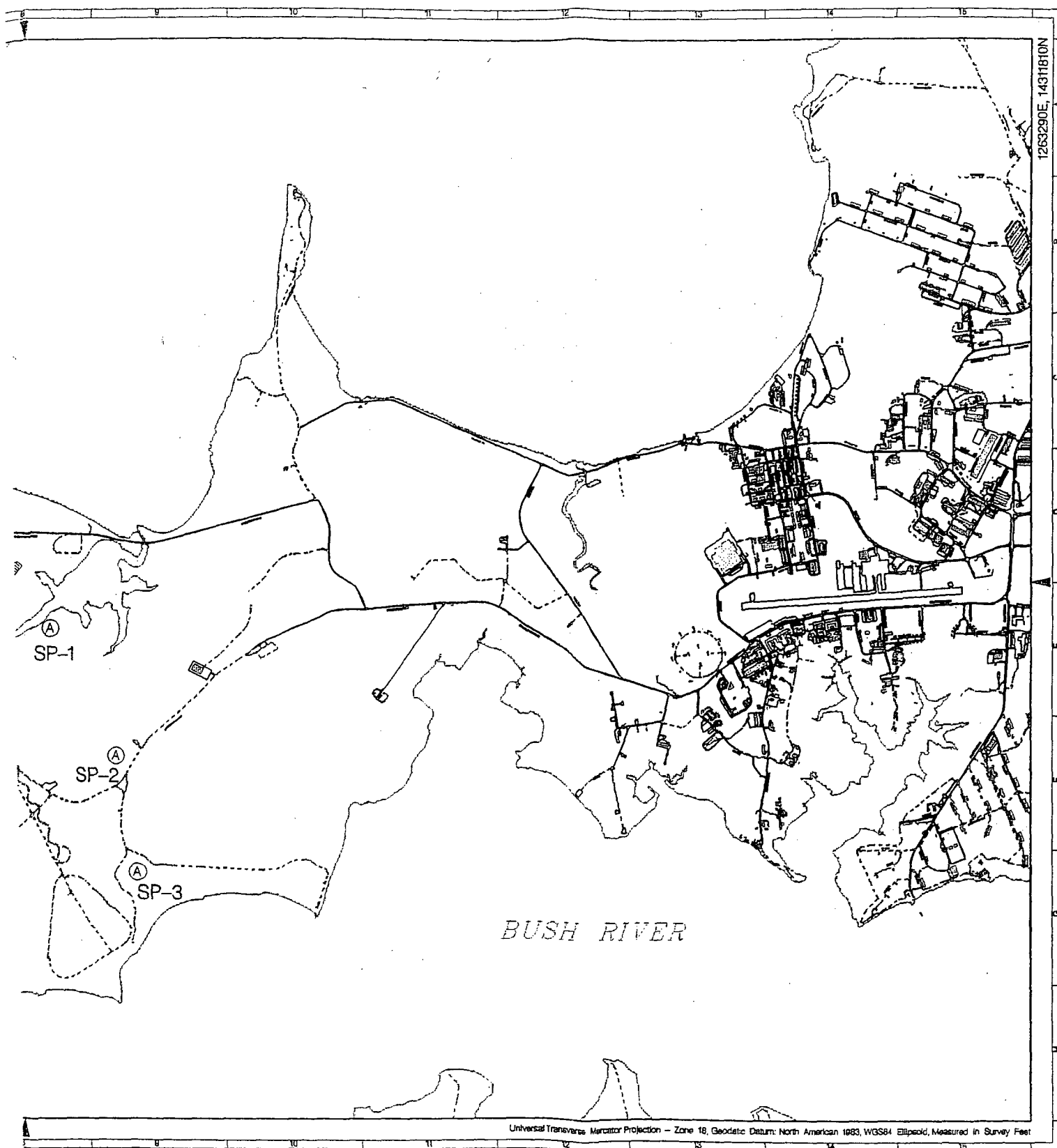
FIGURE:
4



L E G E N D

- Paved Road
- - - - - Unpaved Road
-  Structure
- Coastline
- (A) Sampling Location
-  Controlled Burn Area





1500 3000 6000

GRAPHICAL SCALE

1" = 3000'

1:36000

TITLE:

NEW O-FIELD CONTROLLED BURN SAMPLING LOCATIONS DECEMBER 1999



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DATE:
11-07-00

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K. THORPE

FIGURE:
5

5.3 J-Field Controlled Burn – April 2000

The J-Field controlled burn occurred on 6 April 2000 in the Edgewood Area of APG (Figure 6). The burn area extended over the southeast portion of J-Field, covering both marsh and forest environments. Robbins Point Road and the Bush River served respectively as the northern and eastern firebreaks.

Air samples were collected at two monitoring locations during the J-Field controlled burn: one located downwind of the fire to capture smoke constituents (SP1), and one located upwind of the fire (SP4). The downwind sampling location was northeast of the burn area, along the end of Robbins Point Road on the shore of the Bush River. The SP1 sampling location was approximately 10 meters from the northernmost edge of the burned area. The upwind sampling location (SP4) was located on the Gunpowder River shoreline at the end of Ricketts Point Road, approximately 500 meters from the fire location. Collection of additional downwind samples did not occur due to the logistics of staging samplers at offshore locations in the Bush River.

The ATC Meteorological Office provided a favorable forecast for wind speeds of less than 15 mph from the southwest, and atmospheric stability Class D conditions. The APG Fire Department

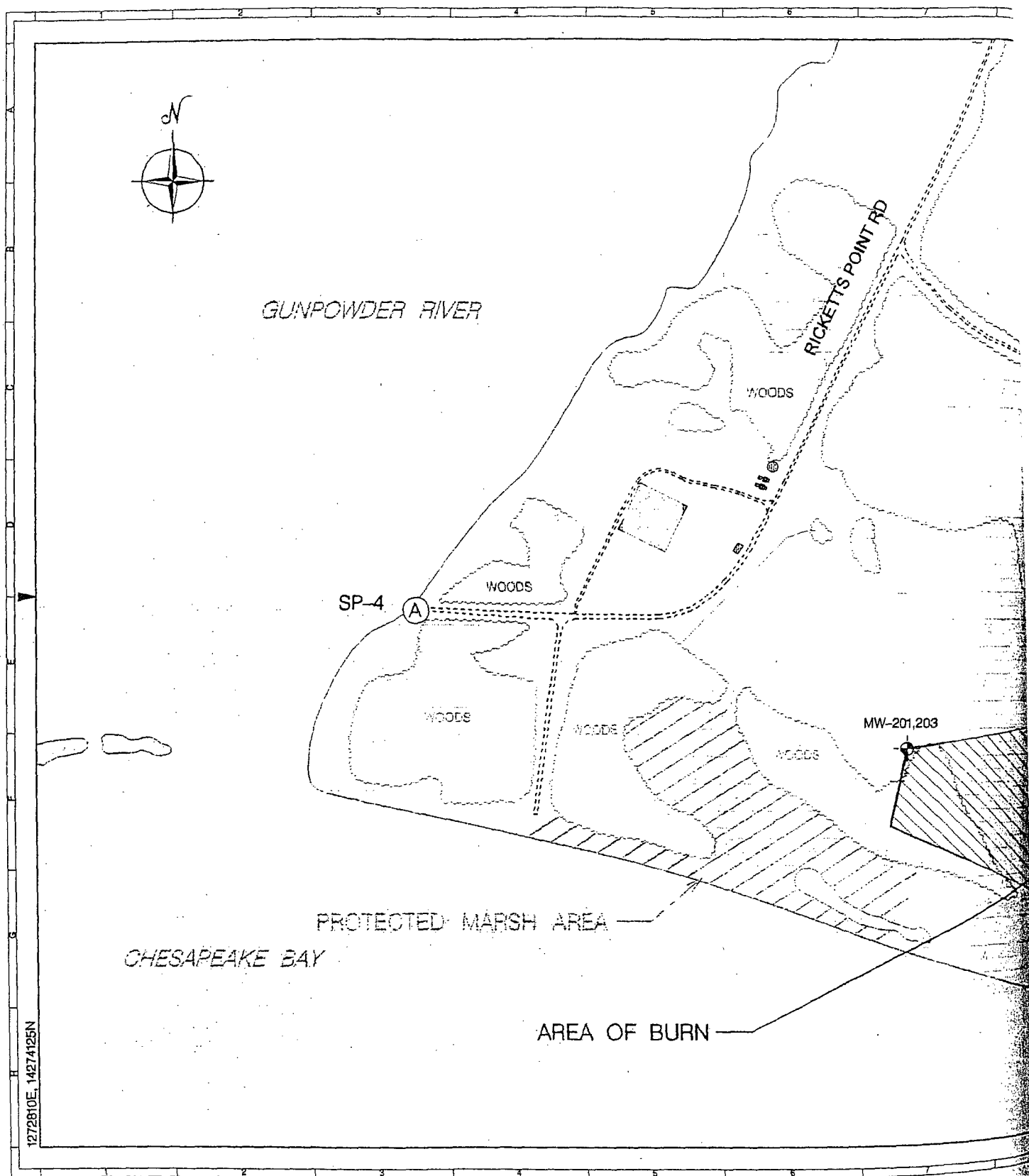
initiated the controlled burn at approximately 1725 hours. The sampling duration was approximately three hours.

Meteorological data collected during the controlled burn period indicated stable wind directions from the southwest, with only slight variations. Wind gusts of up to 15 mph were recorded by an on-site weather station. Wind speeds, coupled with dry conditions and adequate vegetative fuel, sustained the fire during the J-Field controlled burn. A visible smoke plume extended from the burn area in a northeasterly direction.

Photographs taken during and following the J-Field controlled burn are presented in Appendix D-3. The J-Field controlled burn revealed a significant amount of surface waste and debris throughout the burned area, indicating disposal had previously occurred in the area. A separate removal action was conducted in May 2000 to remove the surface debris, including ordnance-related items.

Sampling was performed for chemical agents, explosives, VOCs, PCBs, pesticides, inorganics, and radionuclides; detections of several analytes were reported:

- Several VOCs were detected in the ppb range at the upwind location (SP4):
 - acetic acid
 - acetone
 - hydrocarbon compound (no identification from the TIC library)



L E G E N D

Structure

Unpaved Road

Water

Wetland

Topography

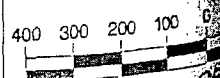
Treeline

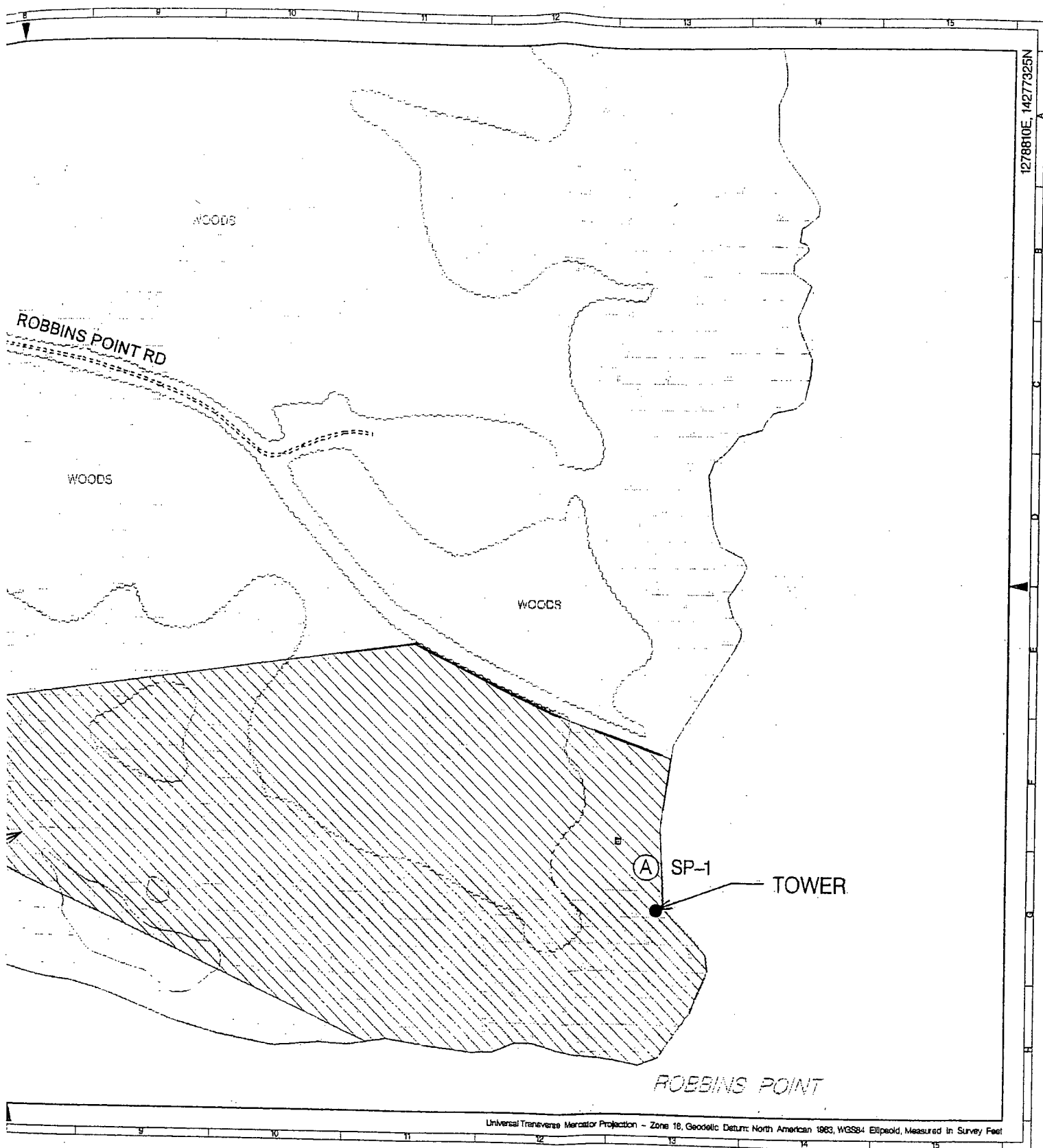
Fence

Monitoring Well

Controlled Burn Area

(A) Sampling Location





200 400 800

GRAPHICAL SCALE
FEET (1" = 400')

1:4800

TITLE:

J-FIELD CONTROLLED BURN AREA AND SAMPLING LOCATIONS APRIL 2000



6700 Alexander Bell Drive (800) 727-6677
Columbia, MD 21046 www.genphysics.com

CARTOGRAPHER:

M. BROOKS

DATE:

11-07-00

APPROVED BY:

K. THORPE

FIGURE:

6

- Several VOCs were also detected in the ppb range at the downwind sampling location (SP1):
 - acetaldehyde
 - acetone
 - acetonitrile
 - ethylhexanol
 - furan
 - furfural
 - methylester acetic acid
 - methylfuran
 - methylpropene
 - hydrocarbon compound (no identification from the TIC library)
- An isolated pesticide detection (heptachlor) was reported in the ppb range at the upwind sampling location (SP4). No pesticides were detected at the downwind sampling location (SP1).
- Two explosive-related compounds (2-amino-4,6-dinitrotoluene and 4,-amino-2,6-dinitrotoluene) were detected in the ppb range at the downwind sampling location. No explosive-related compounds were detected at the upwind sampling location.
- Numerous metals were detected in the ppb range in samples collected both upwind and downwind of the burn area.
- U-235 was reported as detected in the downwind sampling location.

No chemical agents or PCBs were detected in samples collected either upwind or downwind of the burn area. Appendix E-3: Tables E-18 through E-24 contain the results from the analysis of the air samples collected during the J-Field controlled burn.

5.4 Main Front Controlled Burn Attempt – April 2001

A second controlled burn in the Main Front Range was planned in an area where testing of DU weapons has occurred. Immediately following the successful completion of the April 2000 J-Field controlled burn, coordination resumed for the second Main Front controlled burn. Wind directions under which the controlled burn could be conducted were northeast or southwest. However, given the active testing schedule and other limitations (wind direction and greening vegetation), the controlled burn could not be accomplished in the spring, and was delayed until fall.

Coordination resumed in late fall when the vegetation was determined by the APG Fire Department officials to be sufficiently dried to provide adequate fuel and a successful burn. Once again, the active testing schedule and unfavorable meteorological conditions prevented successful completion of the controlled burn. Coordination again resumed in spring of 2001. Under favorable wind conditions, the controlled burn was attempted in the Main Front Range on 6 April 2001. However, light precipitation and the wet condition of the underlying vegetative fuel prevented successful ignition of the burn area.

Evaluation of the selected burn area by Fire Department personnel indicated that a successful controlled burn was unlikely, given the wet conditions and reduced available fuel volume as a result of previous unplanned burns. Active test schedules were projected by ATC for the selected area. Given these limitations, completion of a second burn in the Main Front Range is not feasible.

6.0 RISK ANALYSIS

The evaluation of risk involves comparison of air sampling data collected from the controlled burn events to available human health screening criteria, and calculations to evaluate potential risk associated with exposure to range fire smoke via the inhalation pathway.

6.1 Risk-Based Screening Criteria

To provide a screening level evaluation of potential human health impacts from range fire smoke, concentrations of contaminants detected above quantitation limits are compared to the Maryland Toxic Air Pollutant (TAP) Screening Levels and EPA Region III Risk-Based Concentrations (RBCs). The TAP Screening Levels and RBCs for inhalation are more conservative than other screening criteria such as the Occupational Safety and Health Administration (OSHA) Permissible Exposure Limits (PELs). The PELs establish workplace exposure limits for inhalation by healthy workers, generally based on an 8- or 10-hour workday in a 40-hour workweek. The available TAP Screening Levels and RBCs used in this evaluation are lower values than the PELs for a given chemical. The TAP Screening Levels and RBCs provide consideration of the general population in evaluating exposure and associated risk.

6.1.1 Maryland Toxic Air Pollutant Screening Levels

The Maryland TAP regulations were promulgated in September 1988 to protect the public from TAP emissions from stationary sources of air pollution. The Maryland Department of the Environment (MDE) maintains a list of screening levels for over 1700 compounds. These TAP Screening Levels are tools used to predict whether emissions from a source will unreasonably endanger public health. Emissions from a pollutant source are compared to benchmark concentrations known as "screening levels" which are considered safe or sufficiently conservative that no one would be endangered by that level of exposure. The TAP Screening Levels are included in Appendix E, Tables E-1 through E-24, for the compounds detected in range fire samples collected as part of the controlled burn project.

6.1.2 EPA Region III Ambient Air Risk-Based Concentrations

The RBCs were developed originally for use in the EPA Region III Superfund Program. The primary use of RBCs is for chemical screening during baseline risk assessments. The RBCs combine toxicity factors with "standard" exposure scenarios to provide a numerical estimation of the concentration that relates to a specified risk level. The inhalation RBCs for ambient air presented in Tables E-1 through E-24 (Appendix E) are based on an increased lifetime cancer risk of 1×10^{-6} for carcinogens (i.e., one in one million), or a hazard quotient of 0.1 for non-carcinogens. The exposure factors used in the calculation of the ambient air RBCs are conservative, and are based on residential exposure to contaminants (i.e., 350 days per year).

Noncarcinogenic effects are evaluated by calculating the ratio of a site-specific exposure level for a specified time period to a reference dose (RfD). The RfD for a specific chemical is an estimate of the daily exposure level, with consideration of sensitive populations, that is not expected to cause adverse health effects over the course of a lifetime. The calculated ratio is known as the hazard quotient (HQ). Unlike cancer risk estimates, HQs are not expressed as a probability. An HQ of less than one indicates that exposures are not likely to be associated with adverse noncarcinogenic effects. As the hazard quotient approaches or exceeds 10, the likelihood of adverse effects is increased to the point where action to reduce human exposure should be considered (although the

magnitude of the uncertainty factors used to derive the RfD should also be considered). Because of the uncertainties involved with these estimates, values between one and 10 may be of concern, particularly when additional significant risk factors are present. Since RfDs do not have equal accuracy or precision and they are not based on the same severity of toxic effects, evaluation of hazard indices (i.e., the sum of two or more HQ values for multiple substances and/or multiple exposure pathways) should take into account the uncertainties associated with chemical-specific RfDs. Using this approach, contaminants can then be excluded when they contribute an HQ of less than 0.1 (for noncarcinogens).

6.1.3 Radiological Parameters

Air samples collected were analyzed for gross alpha and gross beta activity, and specific radionuclides by gamma spectroscopy. Results were evaluated against upwind (background) concentrations as well as blank analysis results. Further evaluation was on the basis of Title 10 Code of Federal Regulations (CFR), Part 20, *Standards for Protection Against Radiation*, Appendix B – Table 2, *Annual Limits on Intake (ALIs) and Derived Air Concentrations (DACs) of Radionuclides for Effluent Concentrations*. Table 2 of Appendix B of 10 CFR Part 20 provides concentration limits for radionuclides in airborne effluents released to the general public.

Main Front

Gross alpha and beta radioactivity were detected in the blank and samples. Gross alpha results were not statistically different between the blank and samples. Gross beta results for SP3 and SP4 were also not statistically different from the blank result.

Although gross beta activity was detected in the SP1 and SP2 samples at levels statistically different from the blank, the concentrations of radioactivity are less than 30 times the most restrictive limit for radioactivity per 10 CFR Part 20, Appendix B, Table 2 for unidentified radionuclides. Gamma ray spectroscopy identified the presence of only Potassium-40 (K-40), Lead-212 (Pb-212), Radium-223 (Ra-223), and Uranium-235 (U-235). All of these radionuclides are naturally occurring and were detected with amounts so small that they could not be quantified as statistically significant above the background for the detector used by the gamma ray spectroscopy system. These radionuclides were detected in the background spectrum for the instrument and are therefore considered as not detected.

The levels of airborne radioactivity detected during the Main Front controlled burn sampling event could not be distinguished from ambient concentrations, and do not pose an increased health risk.

New O-Field

Gross alpha and beta radioactivity were detected in the blank and the samples. Gross alpha results were not statistically different between the blank and sample results. Gross beta results were not significant between SP1 and the blank. The gross beta results for SP2, SP3, and SP4, although statistically different from the blank, are present at concentrations less than 10 times the most restrictive limit for radioactivity in air per 10 CFR Part 20, Appendix B, Table 2, for unidentified radionuclides.

Gamma spectroscopy identified the presence of only K-40, Pb-212, Radium-224 (Ra-224), and U-235. All of these radionuclides are naturally occurring, were detected in the blank, and were detected at levels too low to be quantified as statistically significant above background for the detector utilized for the analysis. The levels of radioactivity measured in air samples collected during the New O-Field controlled burn could not be distinguished from ambient concentrations and do not pose an increased health risk.

J-Field

Gross alpha and beta activity detected was not statistically different between the blank and sample results, and could not be distinguished from ambient concentrations.

Gamma spectroscopy identified the presence of the naturally occurring radionuclides K-40, Pb-212, and U-235 at levels too low to be quantified as statistically different from background for the detector used. Thus, these radionuclides were considered not detected. Uranium-235 was reported as detected in the downwind sampling location (SP1) at 0.0005 pico-Curies per cubic meter (pCi/m^3), less than one percent of the most restrictive limit for U-235 in air as per 10 CFR Part 20, Appendix B, Table 2 (i.e., $0.06 \text{ pCi}/\text{m}^3$). On this basis, the detected levels of U-235 are not considered to pose a health risk.

6.2 Results of Risk-Based Screening

Several analytes detected in the controlled burn sampling events conducted at APG occurred at levels exceeding either the Maryland TAP Screening Levels or the EPA Region III ambient air RBCs. The analytes exceeding these criteria are highlighted in the data tables (Appendix E, Tables E-1 through E-24) and included in Table 2. Analytes for which screening levels are not available are not further evaluated.

Table 2. Calculated Range Fire RBCs

| Analyte | Noncarcinogenic RF-RBC ($\mu\text{g}/\text{m}^3$) | Carcinogenic RF-RBC ($\mu\text{g}/\text{m}^3$) | Maximum Reported Concentration ($\mu\text{g}/\text{m}^3$) ^a |
|----------------------------|---|--|--|
| VOLATILE ORGANICS | | | |
| Acetaldehyde | --- | 170.1 | 3.98 |
| Benzene | --- | 46.2 | 19.9 |
| Furan | 777 | --- | 8.58 |
| Methylene Chloride | N/A | 798 | 25.25 |
| Trimethylbenzene | 1302 | --- | 12.54 (Upwind) |
| | | | |
| PCBS | | | |
| 2,4',5-Trichlorobiphenyl | --- | 0.651 | 0.0110 |
| | | | |
| PESTICIDES | | | |
| Dieldrin | --- | 0.0819 | 0.0030 |
| Heptachlor | --- | 0.294 | 0.0020 (Upwind) |
| | | | |
| EXPLOSIVES | | | |
| 2-Amino-4,6-dinitrotoluene | 46.2 | --- | 0.4570 |
| | | | |
| INORGANICS | | | |
| Aluminum | 777 | --- | 51.19 |
| Arsenic | --- | 0.0861 | 0.0147 |
| Cadmium | --- | 0.208 | 0.0036 |
| Manganese | 10.92 | --- | 0.5476 |

^a Maximum reported concentration is the maximum concentration detected based on three burn events.

6.3 Risk Calculations

Risk calculations were performed to further assess potential human health impacts from airborne range fire contaminant concentrations that exceeded the screening levels. The Maryland TAP Screening Levels assume emissions from a stationary source, indicating frequent emissions and associated exposure. The assumptions used in determining the EPA Region III RBCs for evaluating a residential exposure to contaminants in ambient air are overly conservative for evaluating potential human health impacts due to infrequent exposure to range fire smoke. Therefore, the default exposure parameters used in the RBC calculations are modified to reflect a more realistic scenario for exposure to smoke from infrequent range fires (Table 2). The revised exposure parameters are then employed in back-calculating a revised risk-based concentration for the chemicals detected during range fire smoke sampling at concentrations in excess of the risk-based screening criteria. The calculated concentration represents the upper bound of the risk levels established by EPA as acceptable: for carcinogens, increased lifetime cancer risk of 1 in 1,000,000 (1×10^{-6}); for non-carcinogens, a hazard quotient of 0.1.

Data evaluated by ANL in preparation of the "Potential Human Health Impacts from Range Fires at Aberdeen Proving Ground, Maryland" report indicate that, from the period of 1992 – 1997, an average of 80 fires occurred per year at APG. Of those fires, 84 percent involved areas less than 5 acres in size. Fires that burned 25 acres or more constituted only 2 percent of the fires during that period, although some fire reports did not include an estimate of the area burned. ANL used the assumption that five 25-acre or larger fires occurred per year. The ANL report also indicated that the average duration of range fires at APG is approximately one hour. The use of a helicopter with "Bambi bucket" to drop water directly onto the burning areas allows the fires to be extinguished in a short time. These factors were used to develop conservative exposure duration and frequency parameters for calculating revised risk-based concentrations.

The approach used in this report for determining the range fire RBCs (RF-RBCs) is based on modification of the EPA Region III RBCs to reflect a conservative frequency for the exposure of the general population to range fire smoke. The EPA Region III RBCs used for screening purposes assume a residential exposure to airborne contaminants from an ongoing source, with a frequency of 350 days per year. The RF-RBCs are derived on the basis of exposure to 10 range fires per year, with the assumption that wind direction would control exposure. The residential EPA Region III RBCs assume an exposure basis of 24 hours per day. For range fires, that basis is reduced to 4 hours per event. The conservative exposure duration assumes that the receptor would be exposed to smoke from 10 of 80 range fires occurring per year for a maximum duration of 4 hours per fire. The calculated RF-RBCs (presented in $\mu\text{g}/\text{m}^3$) are compared (Table 2) to maximum concentrations detected (also presented in $\mu\text{g}/\text{m}^3$) in the controlled burn sampling events.

7.0 CONCLUSIONS

Air emissions sampling was conducted during the course of three controlled burns at APG. To assess the potential impacts to human health resulting from exposure to smoke from range fires at APG, the analytical results obtained from the sampling events were compared against EPA Region III RBCs and Maryland TAP Screening Levels. Thirteen analytes were reported at concentrations exceeding at least one of the two screening criteria. To further screen the data, revised RF-RBCs were calculated using parameters conservatively considered representative of exposure of residential receptors to range fire smoke at APG. The conservative RF-RBCs were calculated based on the assumption that a receptor is exposed to smoke from 10 range fires per year at APG, for a

duration of four hours each. The resulting RF-RBCs were then compared to the maximum reported concentrations for the 13 analytes in any of the controlled burn sampling events.

The risk analysis presented in this report does not indicate significant impacts to human health resulting from range fires at APG. The risk analysis assumes that the data collected during the controlled burn events are representative of "typical" range fires occurring at APG.

8.0 REFERENCES

- Argonne National Laboratory, 2000. *Potential Health Impacts from Range Fires at Aberdeen Proving Ground, Maryland, Draft_01*. Prepared by Argonne National Laboratory, Argonne, Illinois for the U.S. Army Garrison, Aberdeen Proving Ground, MD. Draft, October 2000.
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- U.S. EPA, 2000. *Risk-Based Concentration Table*. United States Environmental Protection Agency, Region III, Superfund Technical Support Section, Philadelphia, PA. 5 October 2000.
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APPENDIX A

LOCKHEED MARTIN RANGE FIRE SAMPLING TRIP REPORTS

APPENDIX A-1

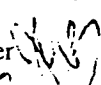
O-FIELD TRIP REPORT
(JULY 2000)

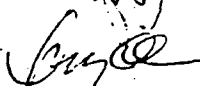
Lockheed Martin Technology Services Group
Environmental Services REAC
20 Woodbridge Avenue, Building 209 Annex Edison, NJ 08837-3679
Telephone 732-321-4200 Facsimile 732-494-4021

LOCKHEED MARTIN 

DATE: July 5, 2000

TO: David Mickunas, U.S. EPA/ERTC Work Assignment Manager

THROUGH: Jeff Bradstreet, REAC Air Group Leader 

FROM: Amy DuBois, REAC Task Leader 

SUBJECT: AIR MONITORING AND SAMPLING AT THE AIR MONITORING SAMPLING, ANALYSIS, AND MODELING SUPPORT, AND UNDERWATER SURVEY ACTIVITIES SITE, ABERDEEN PROVING GROUND, ABERDEEN, MD, WORK ASSIGNMENT #0-110 - TRIP REPORT - O-FIELD

BACKGROUND

The United States Environmental Protection Agency/Environmental Response Team Center (U.S. EPA/ERTC) issued Work Assignment Number 0-110 to Lockheed Martin under the Response, Engineering, and Analytical Contract (REAC) to provide air monitoring and air sampling during two controlled burns in the Edgewood Area of Aberdeen Proving Ground (APG). One burn was to be conducted at O-Field and one at J-Field.

Ordnance firing, ongoing test activities, and lightning strikes occasionally cause accidental fires in the test range areas at APG. Because of APG's long history of weapons testing and disposal practices, there is concern that contaminants have accumulated in the surface soils and vegetation at these locations and could be transported in the smoke plumes produced by such fires, posing a health risk to exposed individuals on and off the installation.

The scope of work for this work assignment included air sampling for dioxins, metals, polynuclear aromatic hydrocarbons (PAHs), inorganic acids, volatile organic compounds (VOCs) and chemical warfare agents (CWAs). Particulate monitoring was conducted utilizing an MIE DataRAM at each location.

OBSERVATIONS AND ACTIVITIES

REAC personnel mobilized to APG on December 3, 1999. Air sampling and monitoring were conducted at 5 downwind and 2 upwind locations (Figure 1).

VOC sampling and analysis was conducted following EPA Method TO-14A: *Determination of Volatile Organic Compounds in Ambient Air Using SUMMA Passivated Canister Sampling and Gas Chromatographic Mass Spectrometric (GC/MS) Analysis*. A sampling orifice was connected to each SUMMA canister to control the flow at 15 cubic centimeters per minute (cc/min). A solenoid valve was then connected to the SUMMA orifice. A trip wire was attached to each solenoid valve to trigger the solenoid to open just before personnel exited the downwind area.

PAH sampling and analysis was conducted following National Institute for Occupational Safety and Health (NIOSH) Method # 5515: *Polynuclear Aromatic Hydrocarbons*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (2 Liters per minute (L/min)) through a sampling train containing a teflon

prefilter cassette and an XAD-2 sorbent tube. The pumps were programmed for a delayed start with a 4-hour sampling period.

Sampling and analysis for inorganic acids was conducted following NIOSH Method # 7903: *Acids, Inorganic*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (250 cc/min) through a sampling train containing a silica gel sorbent tube. The pumps were programmed for a delayed start with a 4-hour sampling period.

Sampling and analysis for dioxins was conducted following modified U.S. EPA Method TO9A, *Determination of Polychlorinated, Polybrominated and Brominated/Chlorinated Dibenzo-p-Dioxins and Dibenzofurans in Ambient Air*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (3 L/min) through a sampling train containing a polyurethane foam (PUF) plug and quartz filter. The pumps were programmed for a delayed start with a 4-hour sampling period. PUF glassware, plugs, and quartz filters were cleaned and certified by Southwest Research Institute in San Antonio, Texas prior to use.

Sampling and analysis for metals was conducted following modified NIOSH Method # 7300: *Elements (ICP)*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (3 L/min) through a sampling train containing a mixed cellulose ester filter cassette. The pumps were programmed for a delayed start with a 4-hour sampling period.

Samples were collected for CWAs utilizing a personal sampling pump (SKC) to draw a measured volume of air (100 cc/min) through a sampling train containing two Depot Area Air Monitoring System (DAAMS) sorbent tubes in a dual-sampling manifold. The CWAs analyzed for included: Sarin (GB), Soman (GD), Mustard (HD), and VX. The pumps were programmed for a delayed start with a 4-hour sampling period. Tubes and analysis were provided by Soldiers Biological and Chemical Command (SBC COM).

Air monitoring for total particulates was performed utilizing an MIE DataRAM portable real-time aerosol monitor. Concentration data was logged every 10 seconds for the duration of the burn.

APG personnel positioned support poles, at each of the five downwind locations, prior to REAC's mobilization to the site. Due to the heavy equipment required to position the poles, and the potential for unexploded ordinance in the marsh/brush area downwind of the proposed burn area, the support poles were positioned on solid ground along the edge of the marsh off Ricketts Point Road. Two nights before the scheduled burn, a spontaneous fire burned the marsh area between Watsons Creek and Ricketts Point Road right up to the support poles. The support poles were used to hold the sampling devices 15 feet above the ground, this positioned the samplers in the plume but out of the potential burn path of the fire. The collection of sampling devices was hoisted up the support pole after setting the timers on the individual pumps. The trip wire for each SUMMA canister allowed the solenoid valve for each SUMMA to be triggered from ground level. Each SUMMA was triggered just before sampling personnel left the potential burn area for a safe zone upwind. When all personnel were out of the area, the APG Fire Department initiated the burn.

RESULTS

VOCs: Benzene and toluene were the only target VOCs detected in any of the samples. The detected concentrations of these two compounds were between 0.4 and 0.6 parts per billion volume (ppbv). These concentrations should be regarded as not detected because 0.6 ppbv each of benzene and toluene were detected in the trip blank. For complete analytical results for VOCs, see the Analytical Report in Appendix A.

PAHs: No PAHs were detected in any of the samples. For complete analytical results for PAHs see the Analytical Report in Appendix B.

Inorganic Acids: No inorganic acids were detected in any of the samples. For complete analytical results for inorganic acids see the Analytical Report in Appendix B.

Dioxins/Furans: A summary of dioxins/furans results can be found in Table 1. The method blank contained OCDD, 1234678-HpCDF, and OCDF; none of the sample results for these compounds were greater than five times the concentration detected in the method blank. The results for each of those compounds should be regarded as not detected. The trip blanks contained 123678-HxCDD, 1234678-HpCDD, OCDD, 12378-PeCDF, 1234678-HpCDF, and OCDF. None of the samples contained concentrations of 123678-HxCDD, 1234678-HpCDD, or 12378-PeCDF exceeding five times the concentrations detected in the trip blank; the results for these compounds should be regarded as not detected. The field blank contained 12378-PeCDD. None of the samples contained 12378-PeCDD at concentrations greater than five times the detected field blank concentration. The results for 12378-PeCDD should be regarded as not detected. The total dioxins/furans detected at each location after adjusting for the compounds regarded as not detected are as follows: O-2(not detected), O-3(not detected), O-4(0.0491 picograms per cubic meter (pg/m³)), O-5(0.705 pg/m³), O-UW1(not detected), and O-UW2(not detected). For complete analytical results for dioxins/furans, see the Analytical Report in Appendix B.

Metals: A summary of metals results can be found in Table 2. The tin concentration detected in sample 28050 should be regarded as estimated because the acceptable quality control (QC) limits for the percent recovery of the blank spike (BS) and blank spike duplicate (BSD) were exceeded. All other concentrations should be regarded as not detected because they were each less than 5 times the lot blank concentration. For complete analytical results for metals, see the Analytical Report in Appendix B.

CWAs: No chemical warfare agents were detected in any of the samples. CWA results are provided by SBC COM, see Appendix C.

Particulates: Particulates results are shown in Figures 2 through 8. The overall maximum concentration of 54.9 micrograms per cubic meter (µg/m³) was detected at location O-UW1.

Meteorological data: Windroses representing local wind speed and wind direction during the burn period are provided in Appendix D. The data was collected at H-Field using a 10-meter tower, and at Poverty Island using a 5-meter tower. Winds at Poverty Island were predominantly out of the southwest, but were light and variable. H-Field recorded stronger winds at the 10-meter level, predominantly out of the south southwest.

Analysis for VOCs and PAHs were provided by REAC, Edison, NJ. Analysis for dioxins/furans, inorganic acids, and metals were provided by Southwest Research Institute, San Antonio, TX. Analysis for CWAs was provided by SBC COM, APG, MD.

FUTURE ACTIVITIES

Due to light winds and the wet marsh, the proposed burn area did not burn, only the area near the fire initiation line ignited. The sampling devices were too far away to capture the plume from the small burned area. A second controlled burn will be conducted at O-Field when the conditions are more favorable. After the O-Field burn is completed, the J-Field burn will be initiated. There are no eagle nesting restrictions affecting the J-Field burn.

Table 1
Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
Summary of Dioxins/Furans Sampling Results - O-Field Controlled Burn - December 3, 2000

| Sample Number Sample Location | 28080 O-1(Field Blank) | 28081 O-2 | 28082 O-3 | 28083 O-4 | 28084 O-5 | 28085 O-UW1 | 28086 O-UW2 | 28088 Trip Blank | 28089 Trip Blank |
|-------------------------------------|---------------------------|---------------|---------------|-------------------|-------------------|-------------------|-------------------|---------------------|---------------------|
| Adjusted concentration ¹ | pg | pg | pg | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ |
| 1,2,3,7,8-PeCDD ⁴ | 4.35 | U | U | 6.9 | U | 7.85 | 5.85 | U | U |
| 1,2,3,6,7,8-HxCDD ³ | U | U | U | U | U | U | 1.02 | U | 0.862 |
| 1,2,3,4,6,7,8-HpCDD ³ | U | U | 0.192 | U | U | 0.277 | U | U | 0.0574 |
| OCDD ² | U | 0.0381 | 0.0548 | U | U | 0.0918 | 0.0306 | 0.0172 | 0.019 |
| 2,3,7,8-TCDF | U | U | U | U | 0.526 | U | U | U | U |
| 1,2,3,7,8-PeCDF ³ | U | 0.3845 | U | 0.3895 | 0.351 | U | 0.52 | 0.374 | 0.209 |
| 1,2,3,4,7,8-HxCDF | U | U | U | U | 0.179 | U | U | U | U |
| 1,2,3,7,8,9-HxCDF | U | U | U | 0.0491 | U | U | U | U | U |
| 1,2,3,4,6,7,8-HpCDF ² | U | U | U | U | U | U | 0.146 | 0.0712 | U |
| OCDF ² | 0.0115 | 0.0215 | 0.0263 | 0.0201 | U | 0.0511 | 0.0203 | 0.0113 | U |
| Total | 4.3615 | 0.4441 | 0.2731 | 7.3587 | 1.056 | 8.2699 | 7.5869 | 0.4737 | 1.1474 |

pg - picograms

pg/m³ - picograms per cubic meter

¹ Adjusted concentration - detected concentration multiplied by the toxicity equivalency factor (TEF) for each compound.

² The OCDD results for samples 28081, 28082, 28085, 28086, 28088, and 28089; the 1,2,3,4,6,7,8-HpCDF results for samples 28086 and 28088; and the OCDF results for samples 28080, 28081, 28082, 28083, 28085, 28086, and 28088 should be regarded as not detected because the concentrations in the samples were less than 5 times the concentration in the method blank.

³ The 1,2,3,6,7,8-HxCDD result for sample 28086; the 1,2,3,4,6,7,8-HpCDD result for samples 28082 and 28085; and the 1,2,3,7,8-PeCDF result for samples 28081, 28083, 28084, and 28086 should be regarded as not detected because the concentration in the sample is less than 5 times the concentration in the trip blank.

⁴ The 1,2,3,7,8-PeCDD results for samples 28083, 28085, and 28086 should be regarded as not detected because the concentration in the samples were less than 5 times the concentration in the field blank.

Table 2
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
Summary of Metals Sampling Results - O-Field Controlled Burn - December 3, 2000

| Sample Number | 28050 | 28051 | 28052 | 28053 | 28054 | 28055 | 28056 | 28057 | 28058 | 28059 |
|---------------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|-----------------------------|-----------------------------|-----------------------------|
| Location | O-1 | O-2 | O-3 | O-4 | O-5 | O-UW1 | O-UW2 | Field Blank | Trip Blank | Lot Blank |
| Parameter | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{m}^3$ | $\mu\text{g}/\text{filter}$ | $\mu\text{g}/\text{filter}$ | $\mu\text{g}/\text{filter}$ |
| Aluminum | U | 1.5 | 3.5 | 2.7 | 4.0 | 2.0 | 2.0 | 1.3 | 1.3 | 2.3 |
| Calcium | 8.8 | 8.3 | 9.0 | 9.0 | 9.7 | 10 | 9.0 | 5.4 | 5.5 | 5.5 |
| Chromium | 0.59 | 0.47 | 0.64 | 0.66 | 0.76 | 0.97 | 0.7 | 0.56 | 0.49 | 0.44 |
| Iron | 1.5 | 1.1 | 1.1 | 0.95 | 2.7 | 1.2 | 2.1 | 0.82 | 0.45 | 0.45 |
| Phosphorus | U | U | U | U | U | U | U | U | U | U |
| Sodium | 9.0 | 12.6 | 12.5 | 11.6 | 14 | 13.8 | 12 | 12.3 | 7.8 | 10 |
| Tin | 1.3 ¹ | U | U | U | U | U | U | U | U | U |
| Zinc | 0.21 | U | 0.25 | 0.23 | 0.33 | 0.31 | 0.16 | 0.11 | U | 0.12 |

¹Regard concentration as estimated, acceptable QC limits for the %Recovery of the BS and the BSD were exceeded.

All detected concentrations for all compounds in this table should be regarded as not detected because they are each less than 5*(Lot Blank Concentration).

$\mu\text{g}/\text{m}^3$ - micrograms per cubic meter

QC - Quality control

BS - Blank spike

BSD - Blank spike duplicate

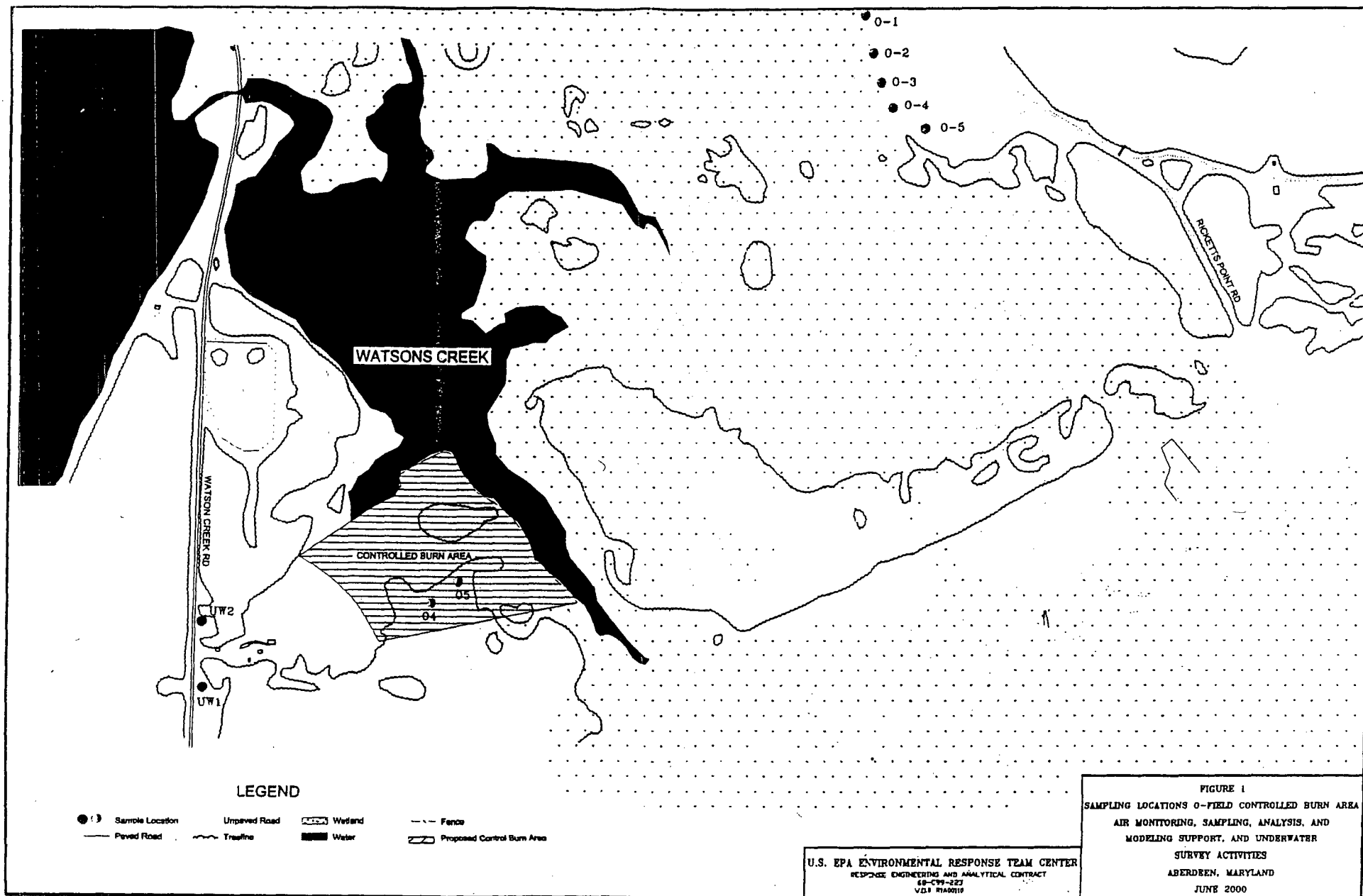


Figure 2
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-1

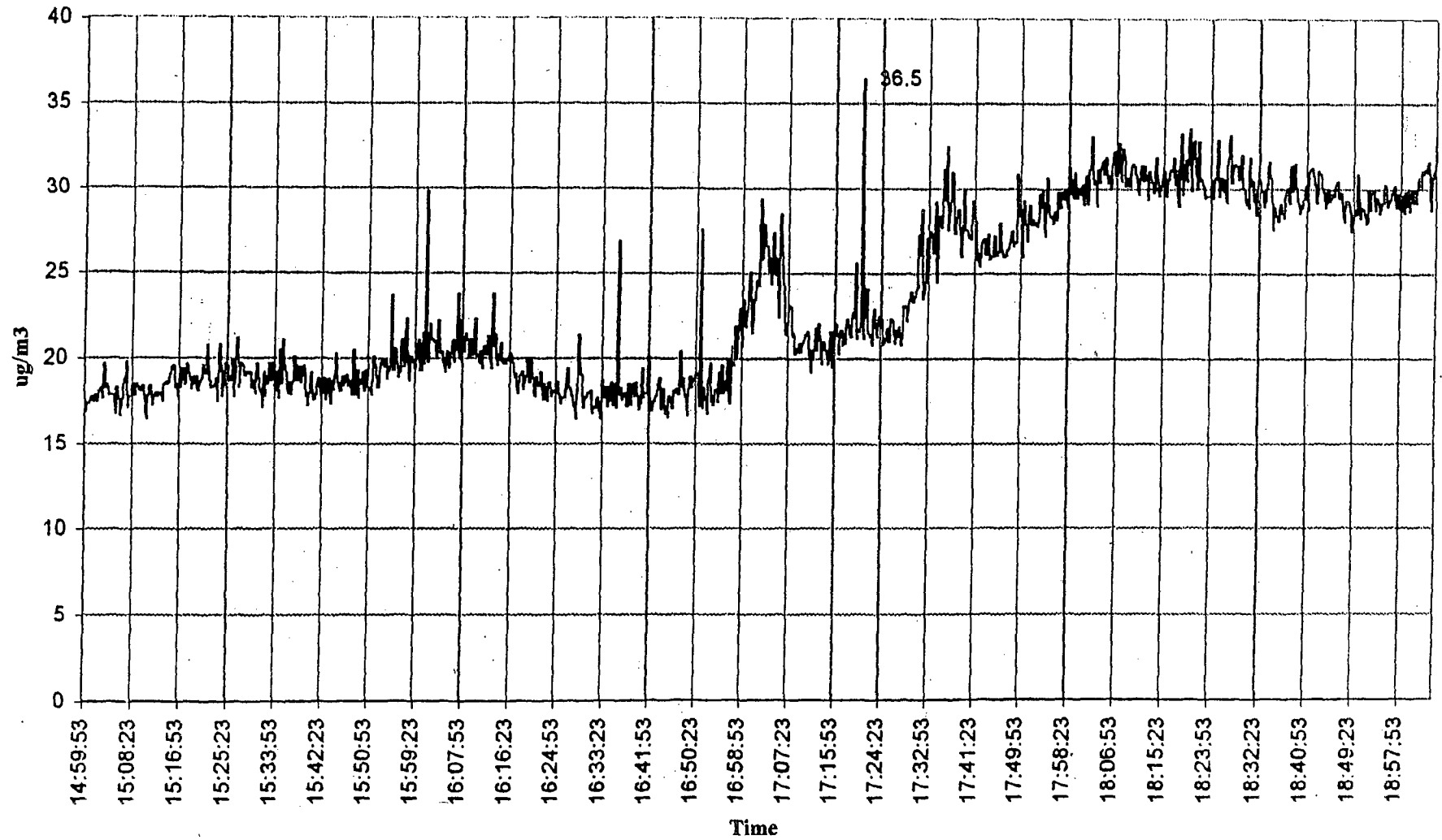


Figure 3
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-2

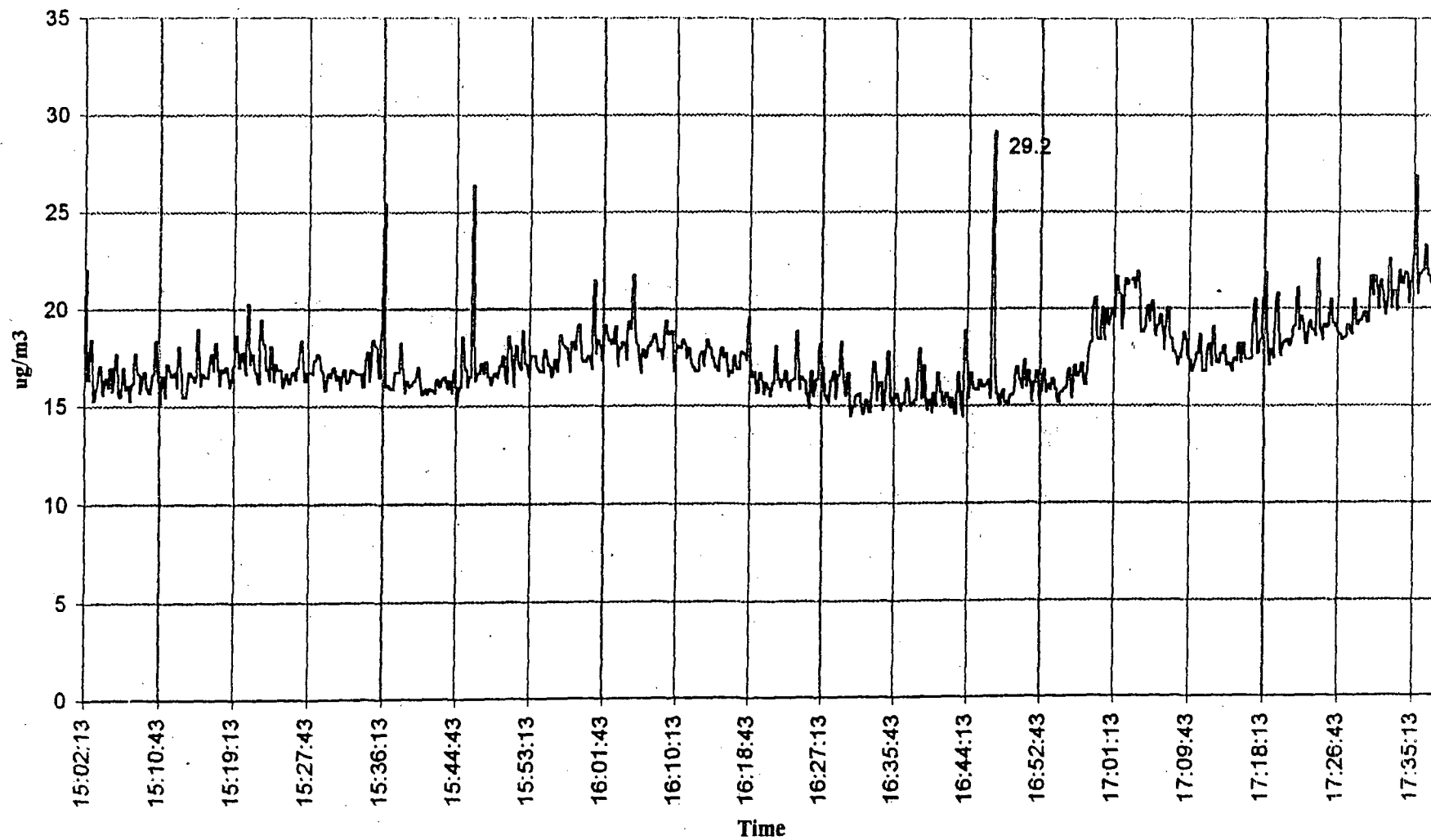


Figure 4
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-3

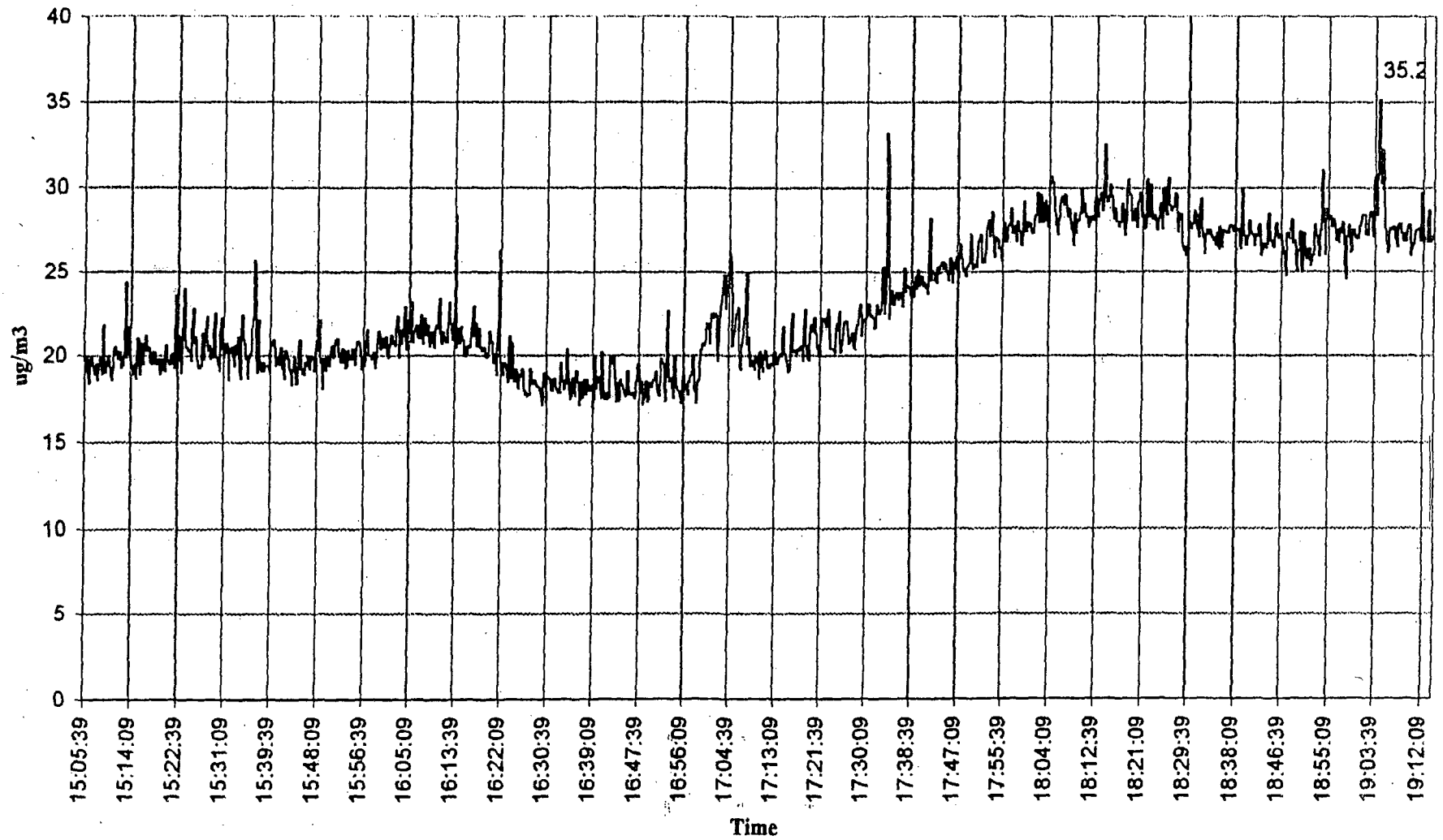


Figure 5
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-4

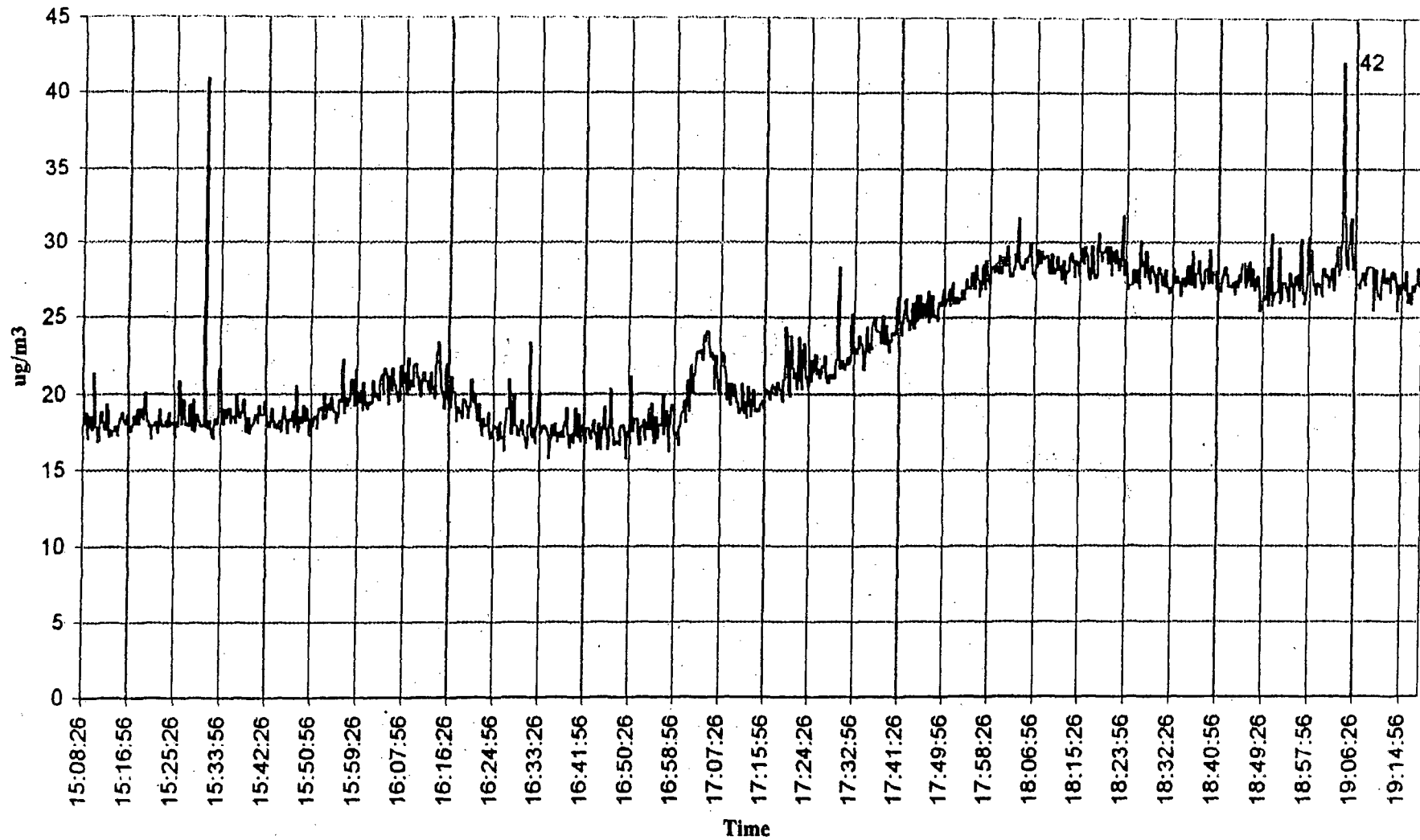


Figure 6
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-5

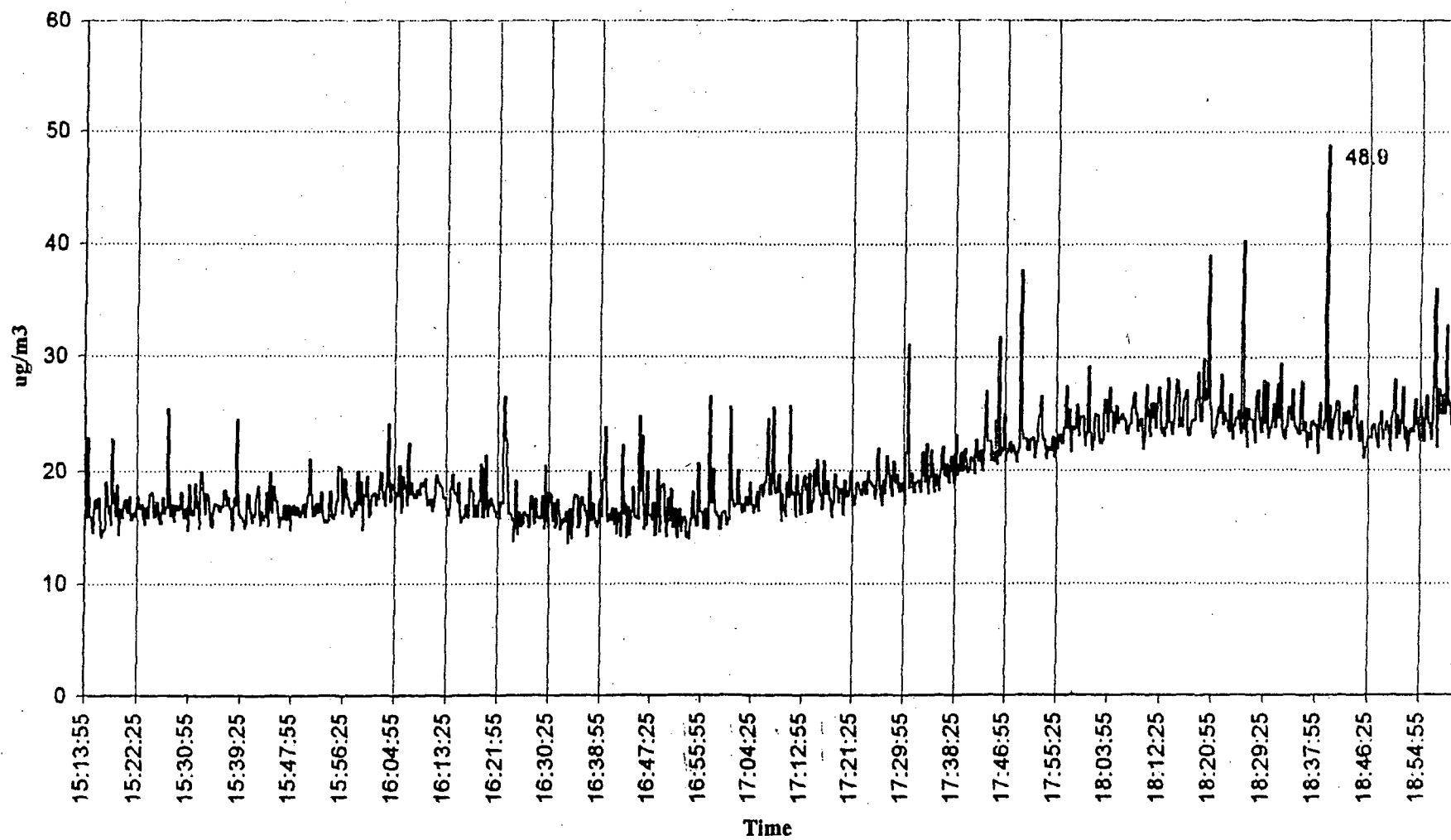


Figure 7
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-UW1

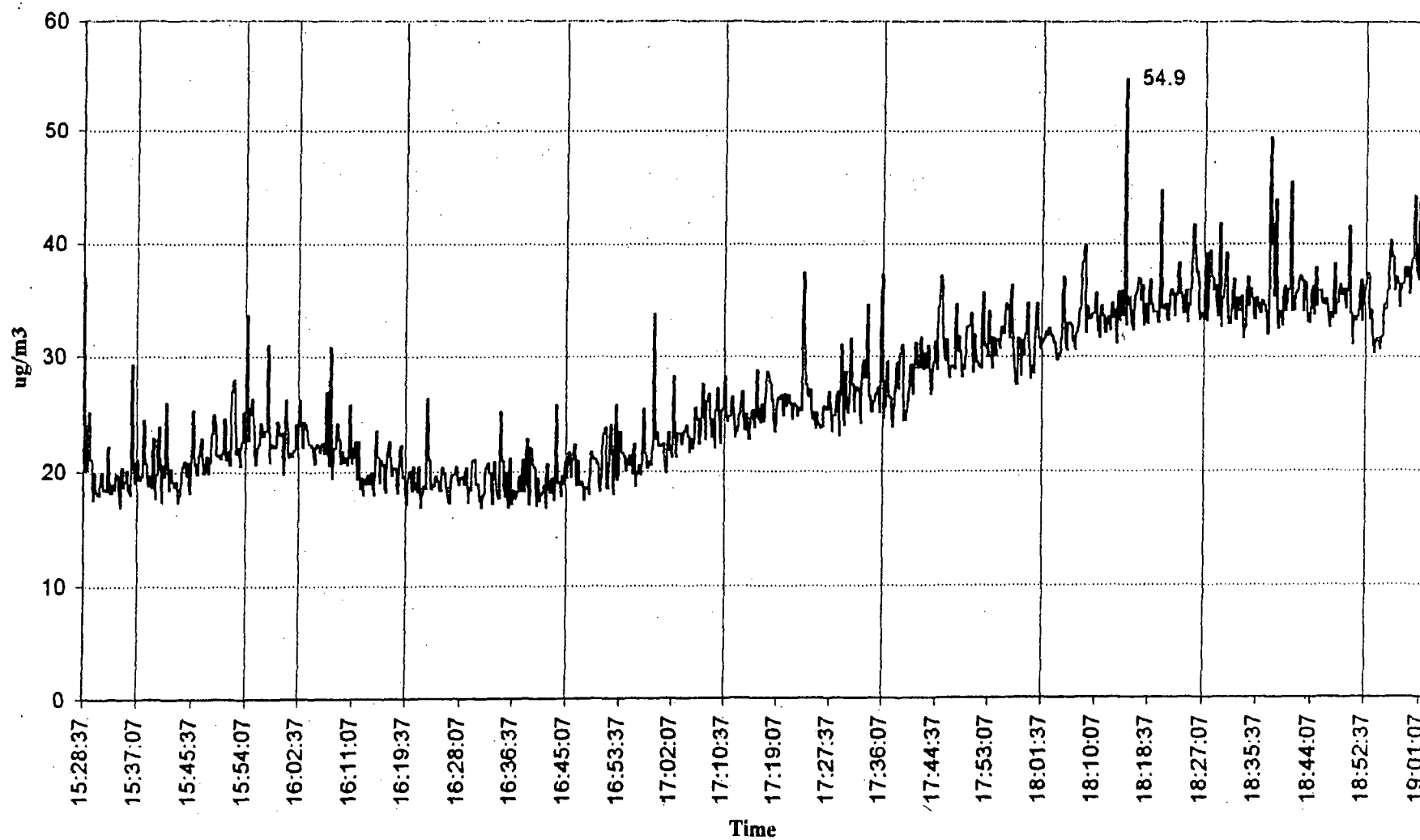
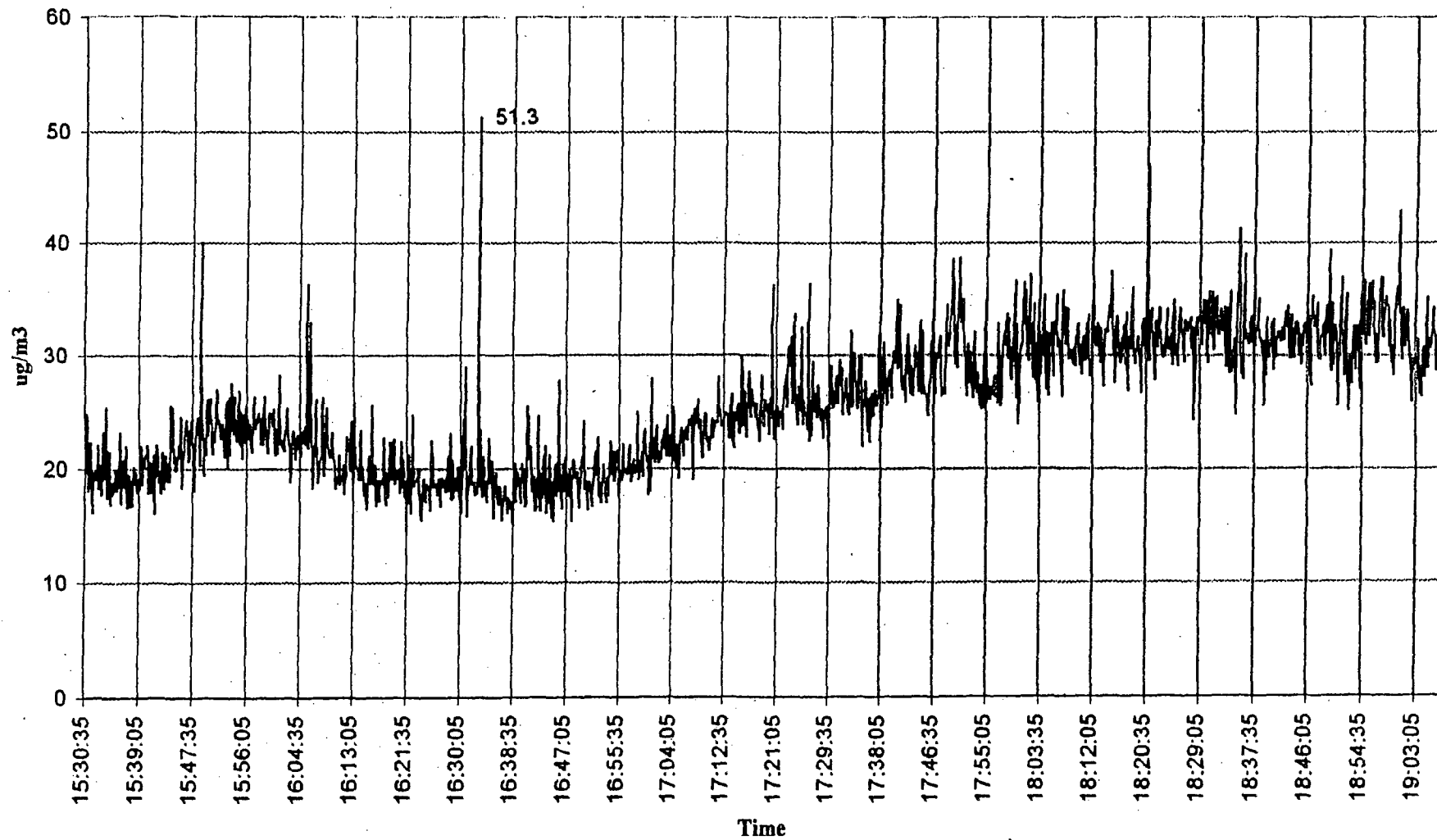


Figure 8
Aberdeen Proving Ground
O Field Burn Data - Particulates
December 3, 1999
Location: O-UW2



APPENDIX A
Analytical Report (VOCs)
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

ANALYTICAL REPORT

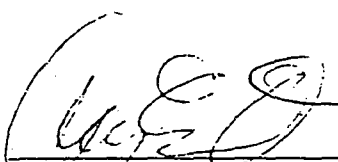
Prepared by
Lockheed Martin Technology Services Group

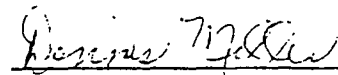
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
Aberdeen, MD

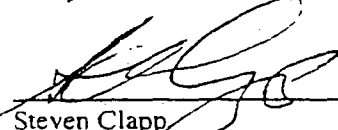
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EPA Work Assignment No:0-110
LOCKHEED MARTIN Work Order No. R1A00110
EPA Contract No. 68-C99-223

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1.0 INTRODUCTION

Summa canister samples were collected at the Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site in Aberdeen, MD on 03 December 1999. A total of eight (8) samples were collected in 6-liter passivated Summa canisters, the samples were transported back to the Environmental Response Team Center (ERTC) facility in Edison, New Jersey. These samples were analyzed by the Response Engineering and Analytical Contract (REAC) using gas chromatography/mass spectrometry (GC/MS) on 06, 07, and 08 December 1999.

2.0 GC/MS CANISTER PROCEDURES

2.1 Sample Pressurization

The Summa canisters used for sampling were cleaned by REAC using REAC Standard Operating Procedure (SOP) #1703 and were selected from batches certified clean by REAC. Before analysis, all canisters were pressurized. A pressurizing train was setup with a pressure gauge accurate to ± 0.1 pounds per square inch absolute (psia). The gauge and train were purged with nitrogen gas (Ultra High Pure grade) for 5 minutes. The train was then connected to the canister, an initial reading was taken. Nitrogen was added to all canister samples to bring the canister pressure to 3 times the initial reading, except 29007 trip blank.

| <u>Sample</u> | <u>Location</u> | <u>Initial Pressure (psia)</u> | <u>Final Pressure (psia)</u> |
|---------------|-----------------|------------------------------------|----------------------------------|
| 29007 | Trip Blank | 0.7 | 20.0 |
| 29000 | 01 | 8.2 | 24.6 |
| 29001 | 02 | 10.2 | 30.6 |
| 29002 | 03 | 9.2 | 27.6 |
| 29003 | 04 | 8.5 | 25.5 |
| 29004 | 05 | 9.8 | 27.4 |
| 29005 | UPW1 | 10.4 | 31.2 |
| 29006 | UPW2 | 8.3 | 24.9 |

2.2 Summa Canister Analysis

Samples were analyzed by cryogenic trapping of aliquots from Summa canisters via a canister using a Hewlett-Packard 5890 gas chromatography (GC) and 5971 A mass selective detector (MSD) running ChemStation software. Table 1 lists cryogenic trap and GC/MS conditions.

All canisters were attached to the Summa canister autosampler. Sample analysis began by cooling the first cryotrap, module -1 (M-1), to -160 degree Celsius ($^{\circ}\text{C}$). Once M-1 was cooled, a specified aliquot of sample or standard was cryotrapped. This aliquot was transferred to a Tenax trap, M-2, to eliminate most of the water, and then cryofocused at a third trap, M-3, before injection by direct heating.

2.3 Calibration and Sample Spiking

A twenty-five (25) compound standard was provided in compressed gas cylinder No ALM009519 by Scott Specialty Gases, Inc. This standard was diluted from 1 part per million volume (ppmv) to 20 parts per billion volume (ppbv) in a Silco canister. An initial calibration range was obtained by varying the volume of the 20 ppbv standard from 50 to 1250 milliliters (mL), equivalent to 1 nanoliter (nL) to 25 nL. Daily standards were obtained by analyzing the 20 ppbv standard at 500 mL (equivalent to 10 nL).

Bromochloromethane (BCM) and p-bromofluorobenzene (BFB) were added to both samples and standards. Both standards were provided in compressed gas cylinder No. ALM046281 by Scott Specialty Gases. BCM was used as an internal standard and BFB was used as a surrogate standard. This standard was diluted from 1 ppmv to 100 ppbv in a Silco canister. An aliquot of 100 mL (equivalent to 10 nL) was added to all standards and samples. To validate the mass spectrometer tuning, an aliquot of 70 mL (equivalent to 50 nanograms of BFB) was analyzed alone. Standard cylinder I.D. numbers, concentrations, and their quantitation ions are listed in Table 2.

2.4 Compound Identification/Quantitation

Contaminants in samples were identified and quantitated by the ChemStation software. This software was designed in order to tentatively identify and quantitate target compounds, using reconstructed and extracted ion chromatogram which were matched with retention time windows. The report format prints the identified compound mass spectra (both raw and background subtracted), quantitation, and qualifier ion chromatogram.

Target compound results are originally reported in nL. The limit of quantitation (LOQ) for all the target compounds is estimated to be 1 nL, being the lowest volume of standard on the calibration curve. The target compound results are calculated in ppbv using the following equation:

$$\text{Concentration (ppbv)} = \frac{\text{Quant Result (nL)} \times 1000}{\text{Undiluted Sample Volume (mL)}}$$

The quantitation limit was 4 ppbv.

Non-target compounds were identified by a library search of all peaks in a chromatogram. The library search report prints out the sample spectrum along with the ten best library matches and the three best library match spectra. These matches were used along with mass spectral interpretation techniques to tentatively identify the unknowns. Concentrations were calculated based on the total ion response of bromochloromethane in the daily standard. All compounds appearing in the method blank as well as other background compounds commonly found in Summa canister GC/MS analyses (siloxanes, carbon dioxide, etc.) were deleted from the sample results to provide a true listing of the compounds in the samples.

2.5 QA/QC

The following QA/QC procedures were performed for this analysis:

- The HP 5971A was tuned daily for perfluorotributylamine (PFTBA) to meet abundance criteria for p-bromofluorobenzene as listed in EPA Method 624. Tuning results are included in the QA/QC data section (Appendix B). The tune was adjusted when necessary.
- An initial calibration by automated injection from a Silco canister standard at 20 ppbv was performed on 25 September 1999. All compounds met the acceptance criteria of having relative standard deviations (RSD) of less than 25%, except chloroethane (29.03 %), 1,1,1-trichloroethane (25.71 %) and carbon tetrachloride (26.97 %).
- Continuing calibrations were performed on 06, 07, and 08 December 1999 to satisfy the 12 hour requirement. All compounds met the acceptance criteria of having relative percent difference (RPD) less than 25%.
- A surrogate standard of BFB was added to all standards and samples. Percent recoveries were calculated against daily standards, and are listed in Table 3. Recoveries should be within 70% to 130% for BFB.

- ▶ Method blanks were analyzed after continuing calibrations to ensure that the system was clean.
- ▶ A duplicate was analyzed on sample 29000 (01).
- ▶ One set of matrix spike and matrix spike duplicates (MS/MSD) was analyzed on sample 29006 (UPW2) by spiking the samples with 500 mL of the 20 ppbv standard. There is no specific recovery range established according to SOP # 1705.

3.0 RESULTS

Summa canister target and non-target results are listed in Tables 3 and 4, respectively. The recoveries for the MS/MSD are presented in Table 5. All results are reported in ppbv for Summa canister samples and blanks. The chain-of-custody is in Appendix A. The Summa canister data is in Appendix B.

In Appendix B, the Analysis Log is followed by the calibration package for each day of analysis. The calibration package includes the daily analysis log, canister pressurization log, BFB tune, and initial or continuing calibration quant report. The quant report lists the retention time, quantitation ion, peak area, and concentration in nL. Concentrations listed on this quant report are generated by using the average response factors of the initial calibration and the response factors of the continuing calibrations.

The following is a list of the QA/QC flags used in qualifying the results:

- A - Assumed volume for method blank.
- B - Concentration less than 3 times method blank value.
- C - Compound calibration relative standard deviation (RSD) >25% (concentrations calculated by average response factor only).
- E - Exceeds calibration range.
- J - Below 1.0 nL quantitation limit.
- U - Not detected.

4.0 DATA ASSESSMENT

The following summarizes the data validation performed on the air toxic analysis of 8 Summa canister air samples received at REAC on 12/6/99, chain of custody 03218, collected on 12/3/99 for the Air Monitoring, Sampling, Analysis and Modeling Support, and Underwater Survey Activities project, WA# 0-110.

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

The samples were treated with procedures consistent with those specified in SOP #1008.

The method blank of 12/7/99 contained 0.4 ppbv of benzene. The concentrations of benzene in samples 29004, 29005 and 29006 should be regarded as not detected.

The trip blank, sample 29007, contained 0.7 ppbv of trichlorofluoromethane, 0.6 ppbv of 1,1-dichloroethene, 0.8 ppbv methylene chloride, 0.6 ppbv trans 1,2-dichloroethylene, 0.6 ppbv 1,1-dichloroethane, 0.7 ppbv trichloroethane, 0.5 ppbv 1,2-dichloroethane, 0.6 ppbv benzene, 0.7 ppbv carbon tetrachloride, 0.6 ppbv trichloroethylene, 0.6 ppbv dibromomethane, 0.5 ppbv bromochloromethane, 0.6 ppbv of toluene and 0.6 ppbv tetrachloroethylene. The data are affected as follows:

The concentrations of benzene and toluene in samples 29000, 29001, 29003, 29004, 29005 and 29006 should be regarded as not detected.

The concentration of toluene in sample 29002 should be regarded as not detected.

The remainder of the data are not affected as the other analytes detected in the trip blank were not detected in the samples.

In the initial calibration of 9/25/99 the acceptable QC limits were exceeded for the percent relative standard deviation for 1,1,1-trichloroethane (26%) and carbon tetrachloride (27%). The data are not affected because these analytes were not detected in the associated samples.

TABLE 1 - GC/MS Instrument Conditions

A. Single Tube Desorber Conditions

| | |
|-------------------------|---------------|
| Cool Desorb Temperature | : 20°C |
| Cool Desorb Time | : 1 minute |
| Cool Desorb Flow | : 50 mL/min |
| Hot Desorb Temperature | : 240°C |
| Hot Desorb Time | : 10.0 minute |
| Hot Desorb Flow | : 50 mL/min |

(2) Preconcentrator Conditions:

| | |
|---------------------------------|---------------|
| M-1 Cryotrap Temperature | : -160°C |
| Internal Standard Trap Time | : 1.0 minute |
| Sample flow | : 150 mL/min |
| M-1 Cryotrap Desorb Temperature | : 20°C |
| M-2 Cryotrap Temperature | : -10°C |
| Transfer (M-1 to M-2) Time | : 4.5 minutes |
| M-2 Cryotrap Desorb Temperature | : 240°C |
| M-3 Cryotrap Temperature | : -160°C |
| Transfer (M-2 to M-3) Time | : 3.5 minutes |
| Injection Time | : 2.0 minutes |

C. GC/MS Conditions, Sample Analysis:

| | |
|---------------------|-----------------|
| Initial Temperature | : 40.0°C |
| Initial Time | : 6.0 minutes |
| Ramp Rate | : 8.0°C/min |
| Final Temperature | : 185.0°C |
| Final Time | : 11.4 minutes |
| Run Time | : 35.5 minutes |
| Mass Scan Range: | : 35 to 250 AMU |

Column: 0.32 mm x 60 meter Restek RTx-5, 1.50 um film thickness (Restek Corporation)

TABLE 2 - Air Toxic Standards (Concentrations and Quantitation Ions)

| <u>Compound</u> | <u>Cylinder</u> | <u>Conc. (ppmv)</u> | <u>Quant. Ion</u> |
|----------------------------|-----------------|---------------------|-------------------|
| chloromethane | ALM009519 | 0.98 | 50 |
| vinyl chloride | ALM009519 | 0.97 | 62 |
| chloroethane | ALM009519 | 1.00 | 64 |
| trichlorofluoromethane | ALM009519 | 1.04 | 101 |
| 1,1-dichloroethene | ALM009519 | 1.02 | 61 |
| dichloromethane | ALM009519 | 1.00 | 49 |
| trans-1,2-dichloroethene | ALM009519 | 1.00 | 61 |
| 1,1-dichloroethane | ALM009519 | 1.02 | 63 |
| trichloromethane | ALM009519 | 1.02 | 83 |
| 1,1,1-trichloroethane | ALM009519 | 1.01 | 97 |
| 1,2-dichloroethane | ALM009519 | 1.02 | 62 |
| benzene | ALM009519 | 1.00 | 78 |
| carbon tetrachloride | ALM009519 | 0.98 | 117 |
| trichloroethene | ALM009519 | 1.00 | 130 |
| dibromomethane | ALM009519 | 0.98 | 174 |
| bromodichloromethane | ALM009519 | 1.01 | 83 |
| toluene | ALM009519 | 1.01 | 91 |
| 1,1,2-trichloroethane | ALM009519 | 0.98 | 97 |
| tetrachloroethene | ALM009519 | 1.00 | 166 |
| ethylbenzene | ALM009519 | 1.01 | 91 |
| meta-xylene | ALM009519 | 1.02 | 91 |
| styrene | ALM009519 | 1.04 | 104 |
| ortho-xylene | ALM009519 | 1.04 | 91 |
| 1,1,2,2-tetrachloroethane | ALM009519 | 1.00 | 83 |
| 1,3,5-trimethylbenzene | ALM009519 | 1.05 | 120 |
| <u>Surrogate Standards</u> | | | |
| bromochloromethane | ALM046281 | 1.06 | 49 |
| p-bromofluorobenzene | ALM046281 | 1.06 | 95 |

TABLE 3 Air Toxic Target Compound Results for Summa Canister Samples
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
(concentrations in ppbv)

| Sample Number | Method | 29007 | 29000 | 29000 Rep | 29001 |
|--------------------------------|----------|------------|----------|-----------|----------|
| Sample Location | Blank | Trip Blank | 01 | 01 | 02 |
| Date Sampled | N/A | 12/03/99 | 12/03/99 | 12/03/99 | 12/03/99 |
| Date Analyzed | 12/07/99 | 12/06/99 | 12/06/99 | 12/06/99 | 12/06/99 |
| Data File | CET016 | ABS001 | ABS002 | ABS013 | ABS003 |
| Chloromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Vinyl Chloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| Chloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Trichlorofluoromethane | 4 U | 0.7 J | 4 U | 4 U | 4 U |
| 1,1-Dichloroethene | 4 U | 0.6 J | 4 U | 4 U | 4 U |
| Methylene Chloride | 4 U | 0.8 J | 4 U | 4 U | 4 U |
| trans-1,2-Dichloroethylene | 4 U | 0.6 J | 4 U | 4 U | 4 U |
| 1,1-Dichloroethane | 4 U | 0.6 J | 4 U | 4 U | 4 U |
| Trichloromethane | 4 U | 4.0 U | 4 U | 4 U | 4 U |
| 1,1,1-Trichloroethane | 4 U | 0.7 J | 4 U | 4 U | 4 U |
| 1,2-Dichloroethane | 4 U | 0.5 J | 4 U | 4 U | 4 U |
| Benzene | 4 U | 0.6 J | 0.4 J | 0.4 J | 0.4 J |
| Carbon Tetrachloride | 4 U | 0.7 J | 4 U | 4 U | 4 U |
| Trichloroethylene | 4 U | 0.6 J | 4 U | 4 U | 4 U |
| Dibromomethane | 4 U | 0.6 J | 4 U | 4 U | 4 U |
| Bromodichloromethane | 4 U | 0.5 J | 4 U | 4 U | 4 U |
| Toluene | 4 U | 0.6 J | 0.6 J | 0.6 J | 0.5 J |
| 1,1,2-Trichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Tetrachloroethylene | 4 U | 0.6 J | 4 U | 4 U | 4 U |
| Ethylbenzene | 4 U | 4 U | 4 U | 4 U | 4 U |
| m & p-Xylenes | 4 U | 4 U | 4 U | 4 U | 4 U |
| Styrene | 4 U | 4 U | 4 U | 4 U | 4 U |
| o-Xylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1,2,2-Tetrachloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,3,5-Trimethylbenzene | 4 U | 4 U | 4 U | 4 U | 4 U |
| p-Bromofluorobenzene (% Rec) | 117 | 116 | 123 | 125 | 125 |
| Pressurized Sample Volume (mL) | 250 | 250 | 750 | 750 | 750 |
| Initial Pressure (psia) | N/A | N/A | 8.2 | 10.4 | 10.2 |
| Final Pressure (psia) | N/A | N/A | 24.6 | 31.2 | 30.6 |
| Quantitation Limit (ppbv) | 4 | 4 | 4 | 4 | 4 |

A - Assumed volume for Blanks
 B - <3 times Method Blank value
 C - Compound Calibration >25% RSD
 D - Compound Calibration Check >25% RPD
 E - Concentration exceeded calibration limit (25nL)
 J - Below 1.00 nL Quantitation Limit
 U - Not Detected
 N/A - Not Applicable

Table 3 (cont.) Air Toxic Target Compound Results for Summa Canister Samples
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
(concentrations in ppbv)

| Sample Number | 29002 | 29003 | Method | 29004 | 29005 |
|--------------------------------|----------|----------|----------|----------|----------|
| Sample Location | 03 | 04 | Blank | 05 | UPW1 |
| Date Sampled | 12/03/99 | 12/03/99 | N/A | 12/03/99 | 12/03/99 |
| Date Analyzed | 12/06/99 | 12/06/99 | 12/07/99 | 12/07/99 | 12/07/99 |
| Data File | ABS004 | ABS005 | ABS007 | ABS011 | ABS012 |
| Chloromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Vinyl Chloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| Chloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Trichlorofluoromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1-Dichloroethene | 4 U | 4 U | 4 U | 4 U | 4 U |
| Methylene Chloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| trans-1,2-Dichloroethylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1-Dichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Trichloromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1,1-Trichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,2-Dichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Benzene | 4 U | 0.4 J | 0.4 J | 0.4 J | 0.4 J |
| Carbon Tetrachloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| Trichloroethylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| Dibromomethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Bromodichloromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Toluene | 0.6 J | 0.5 J | 4 U | 0.6 J | 0.5 J |
| 1,1,2-Trichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Tetrachloroethylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| Ethylbenzene | 4 U | 4 U | 4 U | 4 U | 4 U |
| m & p-Xylenes | 4 U | 4 U | 4 U | 4 U | 4 U |
| Styrene | 4 U | 4 U | 4 U | 4 U | 4 U |
| o-Xylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1,2,2-Tetrachloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,3,5-Trimethylbenzene | 4 U | 4 U | 4 U | 4 U | 4 U |
| p-Bromofluorobenzene (% Rec) | 130 | 128 | 117 | 126 | 129 |
| Pressurized Sample Volume (mL) | 750 | 750 | 250 | 750 | 750 |
| Initial Pressure (psia) | 9.2 | 8.5 | N/A | 9.8 | 10.4 |
| Final Pressure (psia) | 27.6 | 25.5 | N/A | 27.4 | 31.2 |
| Quantitation Limit (ppbv) | 4 | 4 | 4 | 4 | 4 |

A - Assumed volume for Blanks
B - <3 times Method Blank value
C - Compound Calibration >25% RSD
D - Compound Calibration Check >25% RPD
E - Concentration exceeded calibration limit (25nL)
J - Below 1.00 nL Quantitation Limit
U - Not Detected
N/A - Not Applicable

Table 3 (cont.) Air Toxic Target Compound Results for Summa Canister Samples
 WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
 (concentrations in ppbv)

| | |
|--------------------------------|----------|
| Sample Number | 29006 |
| Sample Location | UPW2 |
| Date Sampled | 12/03/99 |
| Date Analyzed | 12/07/99 |
| Data File | ABS014 |
| Chloromethane | 4 U |
| Vinyl Chloride | 4 U |
| Chloroethane | 4 U |
| Trichlorofluoromethane | 4 U |
| 1,1-Dichloroethene | 4 U |
| Methylene Chloride | 4 U |
| trans-1,2-Dichloroethylene | 4 U |
| 1,1-Dichloroethane | 4 U |
| Trichloromethane | 4 U |
| 1,1,1-Trichloroethane | 4 U |
| 1,2-Dichloroethane | 4 U |
| Benzene | 0.4 J |
| Carbon Tetrachloride | 4 U |
| Trichloroethylene | 4 U |
| Dibromomethane | 4 U |
| Bromodichloromethane | 4 U |
| Toluene | 0.5 J |
| 1,1,2-Trichloroethane | 4 U |
| Tetrachloroethylene | 4 U |
| Ethylbenzene | 4 U |
| m & p-Xylenes | 4 U |
| Styrene | 4 U |
| o-Xylene | 4 U |
| 1,1,2,2-Tetrachloroethane | 4 U |
| 1,3,5-Trimethylbenzene | 4 U |
| p-Bromofluorobenzene (% Rec) | 126 |
| Pressurized Sample Volume (mL) | 750 |
| Initial Pressure (psia) | 8.3 |
| Final Pressure (psia) | 24.9 |
| Quantitation Limit (ppbv) | 4 |

A - Assumed volume for Blanks
 B - <3 times Method Blank value
 C - Compound Calibration >25% RSD
 D - Compound Calibration Check >25% RPD
 E - Concentration exceeded calibration limit (25nL)
 J - Below 1.00 nL Quantitation Limit
 U - Not Detected
 N/A - Not Applicable

**Table 4 Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities**

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | Method | Reference Standard: | Bromochloromethane |
| Sample Location: | Blank | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 250 | Reference Std Volume (mL): | 500 |
| Date Sampled: | N/A | Reference Std Area: | 11910887 |
| Date Analyzed: | 12/07/99 | Initial Pressure (psig): | N/A |
| Data File: | CET016 | Final Pressure (psig): | N/A |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.114 | 589021 | 2 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|------------|-----------------------------|--------------------|
| Sample Number: | 29007 | Reference Standard: | Bromochloromethane |
| Sample Location: | Trip Blank | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 250 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 11910887 |
| Date Analyzed: | 12/06/99 | Initial Pressure (psig): | N/A |
| Data File: | ABS001 | Final Pressure (psig): | N/A |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.114 | 605224 | 2 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29000 | Reference Standard: | Bromochloromethane |
| Sample Location: | 01 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 11910887 |
| Date Analyzed: | 12/06/99 | Initial Pressure (psig): | 8.2 |
| Data File: | ABS002 | Final Pressure (psig): | 24.6 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.098 | 689263 | 2 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29001 | Reference Standard: | Bromochloromethane |
| Sample Location: | 02 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 11910887 |
| Date Analyzed: | 12/06/99 | Initial Pressure (psig): | 10.2 |
| Data File: | ABS003 | Final Pressure (psig): | 30.6 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.098 | 664275 | 2 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29002 | Reference Standard: | Bromochloromethane |
| Sample Location: | 03 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 11910887 |
| Date Analyzed: | 12/06/99 | Initial Pressure (psig): | 9.2 |
| Data File: | ABS004 | Final Pressure (psig): | 27.6 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.114 | 636487 | 2 |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities**

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29003 | Reference Standard: | Bromochloromethane |
| Sample Location: | 04 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 11910887 |
| Date Analyzed: | 12/06/99 | Initial Pressure (psig): | 8.5 |
| Data File: | ABS005 | Final Pressure (psig): | 25.5 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.122 | 629600 | 2 * |

* - Below 4 ppbv Limit of Quantitation

N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | Method | Reference Standard: | Bromochloromethane |
| Sample Location: | Blank | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 250 | Reference Std Volume (mL): | 500 |
| Date Sampled: | N/A | Reference Std Area: | 10549361 |
| Date Analyzed: | 12/07/99 | Initial Pressure (psig): | N/A |
| Data File: | ABS007 | Final Pressure (psig): | N/A |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.066 | 555849 | 2 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities**

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29004 | Reference Standard: | Bromochloromethane |
| Sample Location: | 05 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 10549361 |
| Date Analyzed: | 12/07/99 | Initial Pressure (psig): | 9.8 |
| Data File: | ABS011 | Final Pressure (psig): | 27.4 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.082 | 615240 | 2 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29005 | Reference Standard: | Bromochloromethane |
| Sample Location: | UPW1 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 10549361 |
| Date Analyzed: | 12/07/99 | Initial Pressure (psig): | 10.4 |
| Data File: | ABS012 | Final Pressure (psig): | 31.2 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.075 | 599625 | 2 * |

* - Below 4 ppbv Limit of Quantitation

N/A - Not Applicable

Table 4 (cont.) Air Toxic Non-target Compounds
Summa Canister Sample Results
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 29006 | Reference Standard: | Bromochloromethane |
| Sample Location: | UPW2 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 750 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 12/03/99 | Reference Std Area: | 10549361 |
| Date Analyzed: | 12/07/99 | Initial Pressure (psig): | 8.3 |
| Data File: | ABS014 | Final Pressure (psig): | 24.9 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|--------|----------------------|
| dichlorodifluoro-methane | 6.066 | 638029 | 3 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 5 Air Toxic MS/MSD Recovery Summary for Summa Canister Samples
 WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample Number | 29006 | 29006 MS | 29006 MSD | | | | |
|-------------------------------|----------|----------|-----------|----------|----------|----------|-----|
| Sample Location | UPW2 | UPW2 | UPW2 | | | | |
| Date Sampled | 12/03/99 | 12/03/99 | 12/03/99 | | | | |
| Date Analyzed | Spike | 12/07/99 | 12/08/99 | % | 12/08/99 | % | |
| Data File | Amount | ABS014 | ABS018 | Recovery | ABS019 | Recovery | RPD |
| Chloromethane | 9.8 | U | 10.13 | 103 | 10.03 | 102 | 1 |
| Vinyl Chloride | 9.7 | U | 10.05 | 104 | 9.87 | 102 | 2 |
| Chloroethane | 10.0 | U | 10.59 | 106 | 10.54 | 105 | 0.5 |
| Trichlorofluoromethane | 10.4 | U | 9.93 | 95 | 9.85 | 95 | 1 |
| 1,1-Dichloroethene | 10.2 | U | 10.27 | 101 | 10.15 | 100 | 1 |
| Methylene Chloride | 10.0 | U | 10.14 | 101 | 10.01 | 100 | 1 |
| trans-1,2-Dichloroethene | 10.0 | U | 10.18 | 102 | 10.08 | 101 | 1 |
| 1,1-Dichloroethane | 10.2 | U | 10.09 | 99 | 9.87 | 97 | 2 |
| Trichloromethane | 10.2 | U | 10.22 | 100 | 9.98 | 98 | 2 |
| 1,1,1-Trichloroethane | 10.1 | U | 9.67 | 96 | 9.25 | 92 | 4 |
| 1,2-Dichloroethane | 10.2 | U | 10.04 | 98 | 9.76 | 96 | 3 |
| Benzene | 10.0 | 0.11 | 10.02 | 99 | 9.79 | 97 | 2 |
| Carbon Tetrachloride | 9.8 | U | 9.49 | 97 | 9.40 | 96 | 1 |
| Trichloroethylene | 10.0 | U | 10.05 | 101 | 9.84 | 98 | 2 |
| Dibromomethane | 9.8 | U | 10.13 | 103 | 9.96 | 102 | 2 |
| Bromodichloromethane | 10.1 | U | 10.36 | 103 | 10.08 | 100 | 3 |
| Toluene | 10.1 | 0.13 | 10.33 | 101 | 10.34 | 101 | 0.1 |
| 1,1,2-Trichloroethane | 9.8 | U | 10.77 | 110 | 10.67 | 109 | 1 |
| Tetrachloroethylene | 10.0 | U | 10.09 | 101 | 9.81 | 98 | 3 |
| Ethylbenzene | 10.1 | U | 11.20 | 111 | 11.06 | 110 | 1 |
| meta & para-Xylenes | 10.2 | U | 11.01 | 108 | 11.05 | 108 | 0.4 |
| Styrene | 10.4 | U | 10.63 | 102 | 10.75 | 103 | 1 |
| ortho-Xylene | 10.4 | U | 11.15 | 107 | 11.04 | 106 | 1 |
| 1,1,2,2-Tetrachloroethane | 10.0 | U | 10.33 | 103 | 10.24 | 102 | 1 |
| 1,3,5-trimethylbenzene | 10.5 | U | 9.69 | 92 | 9.51 | 91 | 2 |
| p-Bromofluorobenzene (% Rec.) | N/A | 126 | 104 | N/A | 103 | N/A | N/A |

N/A - Not Applicable

APPENDIX A

CHAIN-OF-CUSTODY

**Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities,
Aberdeen, MD**

Sampled on 3 December 1999

WA #: 0-110

REAC, EL NJ

(998) 321- J

EPA Contract 68-04-0022 ^{APB}
099-213

CHAIN OF CUSTODY RECORD

Project Name: ARC Farm Support

Project Number: LA00110

RFW Contact: John Johnson Phone: 321-4248
LM

No: 03218

SHEET NO. 1 OF 1

1:0699-

Sample Identification

VOCs

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | VOLUND (VOCs) | | |
|--------------------------------|------------|-------------------|--------|----------------|--------------|------------------------|---------------|---|--|
| 643 | 29000 | 02 | A | 12/3/99 | 1 | Summa/none | 3.6 | ✓ | |
| 644 | 29001 | 02 | | | 1 | | 3.6 | ✓ | |
| 645 | 29002 | 03 | | | | | 3.6 | ✓ | |
| 646 | 29003 | 04 | | | | | 3.6 | ✓ | |
| 647 | 29004 | 05 | | | | | 3.6 | ✓ | |
| 648 | 29005 | UPW1 | | | | | 3.6 | ✓ | |
| 649 | 29006 | UPW2 | | | | | 3.6 | ✓ | |
| 650 | 29007 | TRIP blank | ✓ | ✓ | ✓ | ✓ | 0 | ✓ | |
| <div>0022</div> <div>APB</div> | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

Matrix:

SD - Sediment PW - Potable Water S - Soil
DS - Drum Solids GW - Groundwater W - Water
DL - Drum Liquids SW - Surface Water O - Oil
X - Other SL - Sludge A - Air

Special Instructions:

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF CUSTODY #

| Items/Reason | Relinquished By | Date | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|------------------|-----------------|---------|------------------|---------|--------|--------------|------------------|---------|--------------|---------|---------|
| Special Analysis | Maple | 12/4/99 | David Hargreaves | 12/4/99 | 9:50am | All Analyses | David Hargreaves | 12/6/99 | John Johnson | 12/6/99 | 10:15am |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |

APPENDIX B
Analytical Report (PAH, Inorganic Acids, Metals, and Dioxins/Furans)
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

ANALYTICAL REPORT

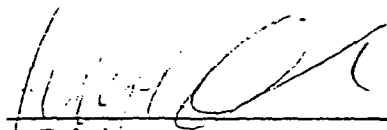
Prepared by
Lockheed Martin

Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
Aberdeen Proving Ground, Aberdeen, MD

February 2000

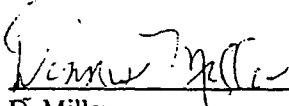
EPA Work Assignment No. 0-110
LOCKHEED MARTIN Work Order No. R1A00110
EPA Contract No. 68-C99-223

Submitted to
D. Mickunas
EPA-ERTC

 2/14/00

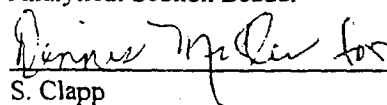
A. Dubois
Task Leader Date

Analysis by:
REAC
SWRI

 2/15/2000

D. Miller
Analytical Section Leader Date

Prepared by:
G. Karustis

 2/15/2000

S. Clapp
Program Manager Date

Reviewed by:
D. Killeen

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Appendices will be furnished on request.

Introduction

REAC in response to WA # 0-110, provided analytical support for environmental samples collected from Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities, located in Aberdeen Proving Ground, Aberdeen, MD as described in the following table. The support also included QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

The samples were treated with procedures consistent with those described in SOP # 1008 and are summarized in the following table:

| COC # | Number of Samples | Sampling Date | Date Received | Matrix | Analysis | Laboratory |
|-------|-------------------|---------------|---------------|--------|-----------------|------------|
| 03215 | 9 | 12/3/99 | 12/6/99 | Air | Dioxin | SWRI* |
| 03217 | 10 | 12/3/99 | 12/6/99 | Air | NIOSH 5515 | REAC |
| 03132 | 10 | 12/3/99 | 12/6/99 | Air | Inorganic Acids | SWRI* |
| 03133 | 10 | 12/3/99 | 12/6/99 | Air | Metals | SWRI* |

* SWRI denotes Southwest Research Institute

Case Narrative

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

PAH in Air Package I 475

The data were examined and were found to be satisfactory.

Metals in Air Package J 012

The lot blank contained 2.3 µg/ filter aluminum, 5.5 µg/ filter calcium, 0.44 µg/ filter chromium, 0.45 µg/ filter iron, 10.0 µg/ filter sodium and 0.12 µg/ filter zinc. The data are affected as follows:

The results for calcium, chromium, iron, sodium and zinc in sample 28050 should be regarded as not detected because the concentration of analyte was less than five times that of the lot blank.

The results for aluminum, calcium, chromium, iron and sodium in samples 28051 and 28058 (the trip blank) should be regarded as not detected because the concentration of analyte was less than five times that of the lot blank.

The results for aluminum, calcium, chromium, iron, sodium and zinc in samples 28052, 28053, 28054, 28055, 28056 and 28057 (field blank) should be regarded as not detected because the concentration of analyte was less than five times that of the lot blank.

The acceptable QC limits for the percent recovery were exceeded in the laboratory control sample for phosphorous (73%), tin (63%) and zirconium (62%). The concentrations of these metals in samples 28050, 28051, 28052, 28053, 28054, 28055, 28056, 28057, 28058 and 28059 should be regarded as estimated.

The acceptable QC limits for the percent recovery were exceeded in the blank spike for phosphorous (52%), tin (36%) and zirconium (37%) and in the blank spike duplicate for the same metals (51%, 34% and 35%, respectively). The concentrations of these metals in samples 28050, 28051, 28052, 28053, 28054, 28055, 28056, 28057, 28058 and 28059 should be regarded as estimated.

Dioxins in Air Package J 015

The samples were received at 12° C by the subcontract laboratory.

The method blank contained 38.2 pg OCDD, 12.4 pg 1234678-HpCDF and 24.3 pg OCDF. The data are affected as follows:

| <u>Sample ID</u> | <u>Analyte</u> | <u>The data should be regarded as</u> |
|------------------|--------------------------|---------------------------------------|
| 28080 | OCDF | Not detected |
| 28081 | OCDD, OCDF | Not detected |
| 28082 | OCDD, OCDF | Not detected |
| 28083 | OCDF | Not detected |
| 28085 | OCDD, OCDF | Not detected |
| 28086 | OCDD, OCDF 1234678-HpCDF | Not detected |
| 28088 | OCDD, OCDF 1234678-HpCDF | Not detected |
| 28089 | OCDD | Not detected |

The values in the above samples are regarded as not detected because they are less than five times the mass found in the method blank.

The trip blank, 28088, contained 17.2 pg OCDD, 7.48 pg 12378-PeCDF, 7.12 pg 1234678-HpCDF and 11.3 pg OCDF. The trip blank, 28089, contained 8.62 pg 123678-HxCDD, 5.74 pg 1234678-HpCDD, 19.0 pg OCDD and 4.18 pg 12378-PeCDF. The data are affected as follows:

| <u>Sample ID</u> | <u>Analyte</u> | <u>The data should be regarded as</u> |
|------------------|----------------|---------------------------------------|
| 28081 | 12378-PeCDF | Not detected |
| 28082 | 1234678-HpCDD | Not detected |
| 28083 | 12378-PeCDF | Not detected |
| 28084 | 12378-PeCDF | Not detected |

| | | |
|-------|---------------------------|--------------|
| 28085 | 1234678-HpCDD | Not detected |
| 28086 | 12378-PeCDF, 123678-HxCDD | Not detected |

The values in the above samples are regarded as not detected because they are less than five times the mass found in the trip blank.

Samples 28083, 28085 and 28086 had masses of 12378-PeCDD that were less than five times that found in the field blank. The values of 12378-PeCDD for these samples should be regarded as not detected.

In the ending calibration verification standard of 12/11/99 (9:06), the acceptable percent difference QC limits were exceeded for 12378-PeCDD (34%), 123478-HxCDD (20.7%), ¹³C-12378-PeCDF (62%), ¹³C-12378-PeCDD (64%), and ¹³C-OCDD (87%). The subcontract laboratory used the average relative response factor calculated from the two continuing calibrations bracketing samples, method blank 12/7/99, 28080 and 28081. Only estimated values or EMPC values were reported in the samples. The data are not affected.

The acceptable QC limits were exceeded for the percent recoveries of several internal standards. The internal standards in question and the samples and analytes involved are summarized as follows:

| <u>Sample ID</u> | <u>Internal standard</u> | <u>Effect</u> |
|------------------|-------------------------------|---|
| 28086 / | ¹³ C-2378-TCDF | The data are not affected |
| 28088 | ¹³ C-2378-TCDF | The data are not affected |
| | ¹³ C-1234678-HpCDF | The data for 1234678-HpCDF should be regarded as estimated. |
| 28089 | ¹³ C-2378-TCDF | The data are not affected |
| | ¹³ C-1234678-HpCDF | The data are not affected |
| Blank Spike | ¹³ C-2378-TCDF | The data for 2378-TCDF should be regarded as estimated |

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The data were examined and were found to be satisfactory.

Summary of Abbreviations

| | |
|----------------|---|
| AA | Atomic Absorption |
| B | The analyte was found in the blank |
| BFB | Bromofluorobenzene |
| C | Centigrade |
| D | (Surrogate Table) this value is from a diluted sample and was not calculated (Result Table) this result was obtained from a diluted sample |
| Dioxin | denotes Polychlorinated Dibenzo-p-dioxins and Polychlorinated Dibenzofurans and/or PCDD and PCDF |
| CLP | Contract Laboratory Protocol |
| COC | Chain of Custody |
| CONC | Concentration |
| CRDL | Contract Required Detection Limit |
| CRQL | Contract Required Quantitation Limit |
| DFTPP | Decafluorotriphenylphosphine |
| DL | Detection Limit |
| E | The value is greater than the highest linear standard and is estimated |
| EMPC | Estimated maximum possible concentration |
| ICAP | Inductively Coupled Argon Plasma |
| ISTD | Internal Standard |
| J | The value is below the method detection limit and is estimated |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| MDL | Method Detection Limit |
| MI | Matrix Interference |
| MS | Matrix Spike |
| MSD | Matrix Spike Duplicate |
| MW | Molecular Weight |
| NA | either Not Applicable or Not Available |
| NC | Not Calculated |
| NR | Not Requested |
| NS | Not Spiked |
| % D | Percent Difference |
| % REC | Percent Recovery |
| PPB | Parts per billion |
| PPBV | Parts per billion by volume |
| PPMV | Parts per million by volume |
| PQL | Practical Quantitation Limit |
| QA/QC | Quality Assurance/Quality Control |
| QL | Quantitation Limit |
| RPD | Relative Percent Difference |
| RSD | Relative Standard Deviation |
| SIM | Selected Ion Monitoring |
| TCLP | Toxic Characteristics Leaching Procedure |
| U | Denotes not detected |
| W | Weathered analyte; the results should be regarded as estimated |
| m ³ | cubic meter kg kilogram µg microgram |
| L | liter g gram pg picogram |
| mL | milliliter mg milligram ng nanogram |
| µL | microliter |
| * | denotes a value that exceeds the acceptable QC limit |
| | Abbreviations that are specific to a particular table are explained in footnotes on that table |

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Analytical Procedure for PAH in Air (XAD-2 Tubes)

XAD-2 Tube Preparation

The XAD-2 tubes were analyzed for polycyclic aromatic hydrocarbons (PAH) using modified NIOSH Method 5515. The front, back and filter portions of the tubes were analyzed separately by extracting them with 2.0 mL methylene chloride. A preweighed filter was also collected with these tubes and this filter was extracted with 4.0 mL methylene chloride. One mL of extract was spiked with 20 μ L of a 2000 ppm XAD internal standards solution consisting of naphthalene- d_8 , acenaphthene- d_{10} , phenanthrene- d_{10} , chrysene- d_{12} , and perylene- d_{12} , resulting in a 40.0 ppm concentration and analyzed.

GC/MS Analysis

An HP 6890 MSD, equipped with a 6890 autosampler and controlled by a personal computer equipped with HP-Enviroquant software was used to analyze the samples.

The instrument conditions were:

| | |
|--|--|
| Column | Restek Rtx-5 (cross bonded SE-54) 30 meter x 0.25 mm ID, 0.50 μ m film thickness. |
| Flow Rate | 1 mL/min, EPC enabled |
| Injection Temperature | 280° C |
| Transfer Temperature | 280° C |
| Source Temperature & Analyzer Temperature | Controlled by thermal transfer of heat from Transfer Line 280° C |
| Temperature Program | 70° C for 0.5 min 30° C/min to 295° C hold for 8 minutes 30° C/min to 315° C; hold for 7 min |
| Pulsed Splitless Injection | Pressure Pulse = 16 psi for 1.0 min, then normal flow 8:1 Split Ratio |
| Injection Volume | 1 μ L |
| | Must use 4 mm ID single gooseneck liners packed with 10 mm plug of silanized and conditioned glass wool |

The GC/MS was calibrated using 6 PAH standards at 10, 25, 50, 75, 100 and 150 ppm. Before analysis each day the system was tuned with 50-ng decafluorotriphenylphosphine (DFTPP) and passed a continuing calibration check by analyzing a 50 μ g/mL daily standard. The QC limit for the initial calibration is %RSD less than 30 and %D less than 25 for the daily check. Sample quantification is based on the average response factor of the calibration curve or the response factor of the daily calibration check.

The XAD-2 tube PAH results are listed in Table 1.1. Tentatively identified compounds are listed in Table 1.2. The following equations were used to calculate the analyte - total $\mu\text{g}/\text{sample}$:

$$\mu\text{g}/\text{sample} = C_u \times V \times DE = \frac{A_u \times C_{is} \times V \times DE}{A_{is} \times RRF}$$

where

- C_u = Concentration of the analyte ($\mu\text{g}/\text{mL}$)
- V = Extraction Volume (mL)
- DE = Desorption Efficiency = $100/(\% \text{ Recovery})$
- A_u = Area of the analyte
- C_{is} = Concentration of the internal standard ($\mu\text{g}/\text{mL}$)
- A_{is} = Area of the internal standard

The Relative Response Factor, RRF, is calculated from the calibration standard mixture using

$$RRF = \frac{A_u \times C_{is}}{A_{is} \times C_u}$$

where

- RRF = Relative Response Factor (unitless)
- A_u = Area of Analyte in the standard mixture
- C_{is} = Concentration of Internal Standard in the standard mixture ($\mu\text{g}/\text{mL}$)
- A_{is} = Area of Internal Standard in the standard mixture
- C_u = Concentration of Analyte in the standard mixture ($\mu\text{g}/\text{mL}$)

The concentration of the analyte in mg/m^3 and ppbv (parts per billion by volume) is calculated using the following:

$$\text{mg}/\text{m}^3 = \frac{(\text{Total } \mu\text{g front} + \text{Total } \mu\text{g back})}{\text{Liters Sampled}}$$

$$\text{ppbv} = \frac{\text{mg}/\text{m}^3 \times 24.45 \times 1000}{MW}$$

where MW is the molecular weight of the analyte

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LINEAR SCAN COMPOUND AND ION LIST FOR PAH/XAD TUBES

| <u>Compound</u> | <u>Quant Ion</u> | <u>Secondary Ions</u> |
|-----------------------------------|------------------|-----------------------|
| Naphthalene-d ₈ (IS) | 136 | 108 |
| Naphthalene | 128 | 127, 129 |
| 2-Methylnaphthalene | 142 | 141, 115 |
| 1-Methylnaphthalene | 142 | 141, 115 |
| Biphenyl | 154 | 153, 152 |
| 2,6-Dimethylnaphthalene | 156 | 141, 128 |
| 2-Fluorobiphenyl (SURR) | 172 | 171, 173 |
| Acenaphthene-d ₈ (IS) | 164 | 162 |
| Acenaphthylene | 152 | 151, 153 |
| Acenaphthene | 153 | 152, 151 |
| Dibenzofuran | 168 | 139 |
| Fluorene | 166 | 167, 165 |
| Phenanthrene-d ₁₀ (IS) | 188 | 189 |
| Phenanthrene | 178 | 179, 176 |
| Anthracene | 178 | 179, 176 |
| Carbazole | 167 | 166, 168 |
| Fluoranthene | 202 | 101, 200 |
| Pyrene | 202 | 101, 200 |
| Terphenyl-d ₁₄ (SURR) | 244 | 243 |
| Chrysene-d ₁₂ (IS) | 240 | 236 |
| Benzo(a)anthracene | 228 | 226, 229 |
| Chrysene | 228 | 226, 229 |
| Perylene-d ₁₂ (IS) | 264 | 260 |
| Benzo(b)fluoranthene | 252 | 250, 126 |
| Benzo(k)fluoranthene | 252 | 250, 126 |
| Benzo(e)pyrene | 252 | 250, 126 |
| Benzo(a)pyrene | 252 | 250, 126 |
| Indeno(1,2,3-cd)pyrene | 276 | 138, 277 |
| Dibenzo(a,h)anthracene | 278 | 139, 278 |
| Benzo(g,h,i)perylene | 276 | 277, 138 |

Analytical Procedure for Metals in Air

The subcontract laboratory determined the metal concentration in the samples by analyzing them according to NIOSH method 7300. The results of the analysis are listed in Table 1.3.

Analytical Procedure for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

The subcontract laboratory determined the concentration of polychlorinated dibenzodioxins and polychlorinated dibenzofurans in the samples by analyzing them according to USEPA Method 8290. The results of the analysis are listed in Table 1.4.

Analytical Procedure for Inorganic Acids in Air

The subcontract laboratory determined the concentration of inorganic acids in the samples by analyzing them according to NIOSH Method 7903. The results of the analysis for the soil samples are listed in Table 1.5.

Table 1.1 Results of the Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample No. | 28079 | | 28070 | | 28071 | | 28072 | | 28073 | |
|-------------------------|-------------|-----------|---------------|-------------|---------------|-------------|---------------|-------------|---------------|-------------|
| Sampling Location | Lot Blank | | 0-1 | | 0-2 | | 0-3 | | 0-4 | |
| Volume (L) | 0 | | 474.6 | | 458 | | 462 | | 460 | |
| Compound Name | Conc. µg | MDL µg | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv |
| Naphthalene | U | 8.6 | U | 3.5 | U | 3.6 | U | 3.6 | U | 3.6 |
| 2-Methylnaphthalene | U | 9.1 | U | 3.3 | U | 3.4 | U | 3.4 | U | 3.4 |
| 1-Methylnaphthalene | U | 9.0 | U | 3.2 | U | 3.4 | U | 3.3 | U | 3.4 |
| Biphenyl | U | 9.2 | U | 3.1 | U | 3.2 | U | 3.2 | U | 3.2 |
| 2,6-Dimethylnaphthalene | U | 9.3 | U | 3.1 | U | 3.2 | U | 3.1 | U | 3.2 |
| Acenaphthylene | U | 9.2 | U | 3.1 | U | 3.2 | U | 3.2 | U | 3.2 |
| Acenaphthene | U | 9.0 | U | 3.0 | U | 3.1 | U | 3.1 | U | 3.1 |
| Dibenzofuran | U | 9.0 | U | 2.7 | U | 2.8 | U | 2.8 | U | 2.8 |
| Fluorene | U | 9.1 | U | 2.8 | U | 2.9 | U | 2.9 | U | 2.9 |
| Phenanthrene | U | 9.2 | U | 2.7 | U | 2.8 | U | 2.7 | U | 2.8 |
| Anthracene | U | 8.9 | U | 2.6 | U | 2.7 | U | 2.6 | U | 2.7 |
| Carbazole | U | 9.7 | U | 3.0 | U | 3.1 | U | 3.1 | U | 3.1 |
| Fluoranthene | U | 9.2 | U | 2.4 | U | 2.4 | U | 2.4 | U | 2.4 |
| Pyrene | U | 9.2 | U | 2.3 | U | 2.4 | U | 2.4 | U | 2.4 |
| Benzo(a)anthracene | U | 9.2 | U | 2.1 | U | 2.2 | U | 2.1 | U | 2.2 |
| Chrysene | U | 8.9 | U | 2.0 | U | 2.1 | U | 2.1 | U | 2.1 |
| Benzo(b)fluoranthene | U | 9.6 | U | 2.0 | U | 2.0 | U | 2.0 | U | 2.0 |
| Benzo(k)fluoranthene | U | 9.3 | U | 1.9 | U | 2.0 | U | 1.9 | U | 2.0 |
| Benzo(e)pyrene | U | 9.5 | U | 1.9 | U | 2.0 | U | 2.0 | U | 2.0 |
| Benzo(a)pyrene | U | 9.6 | U | 2.0 | U | 2.0 | U | 2.0 | U | 2.0 |
| Indeno(1,2,3-cd)pyrene | U | 10 | U | 1.9 | U | 1.9 | U | 1.9 | U | 1.9 |
| Dibenzo(a,h)anthracene | U | 10 | U | 1.9 | U | 1.9 | U | 1.9 | U | 1.9 |
| Benzo(g,h,i)perylene | U | 10 | U | 1.9 | U | 2.0 | U | 2.0 | U | 2.0 |

Table 1.1 (cont.) Results of the Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample No. | 28074 | | 28075 | | 28076 | | 28077 | | 28078 | |
|-------------------------|---------------|-------------|---------------|-------------|---------------|-------------|-------------|-----------|-------------|-----------|
| Sampling Location | 0-5 | | 0-UW1 | | 0-UW2 | | Field Blank | | Trip Blank | |
| Volume (L) | 462 | | 424 | | 419.3 | | 0 | | 0 | |
| Compound Name | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. µg | MDL µg | Conc. µg | MDL µg |
| Naphthalene | U | 3.6 | U | 3.9 | U | 3.9 | U | 8.6 | U | 8.6 |
| 2-Methylnaphthalene | U | 3.4 | U | 3.7 | U | 3.7 | U | 9.1 | U | 9.1 |
| 1-Methylnaphthalene | U | 3.3 | U | 3.6 | U | 3.7 | U | 9.0 | U | 9.0 |
| Biphenyl | U | 3.2 | U | 3.4 | U | 3.5 | U | 9.2 | U | 9.2 |
| 2,6-Dimethylnaphthalene | U | 3.1 | U | 3.4 | U | 3.5 | U | 9.3 | U | 9.3 |
| Acenaphthylene | U | 3.2 | U | 3.5 | U | 3.5 | U | 9.2 | U | 9.2 |
| Acenaphthene | U | 3.1 | U | 3.4 | U | 3.4 | U | 9.0 | U | 9.0 |
| Dibenzofuran | U | 2.8 | U | 3.1 | U | 3.1 | U | 9.0 | U | 9.0 |
| Fluorene | U | 2.9 | U | 3.2 | U | 3.2 | U | 9.1 | U | 9.1 |
| Phenanthrene | U | 2.7 | U | 3.0 | U | 3.0 | U | 9.2 | U | 9.2 |
| Anthracene | U | 2.6 | U | 2.9 | U | 2.9 | U | 8.9 | U | 8.9 |
| Carbazole | U | 3.1 | U | 3.3 | U | 3.4 | U | 9.7 | U | 9.7 |
| Fluoranthene | U | 2.4 | U | 2.6 | U | 2.7 | U | 9.2 | U | 9.2 |
| Pyrene | U | 2.4 | U | 2.6 | U | 2.6 | U | 9.2 | U | 9.2 |
| Benzo(a)anthracene | U | 2.1 | U | 2.3 | U | 2.4 | U | 9.2 | U | 9.2 |
| Chrysene | U | 2.1 | U | 2.3 | U | 2.3 | U | 8.9 | U | 8.9 |
| Benzo(b)fluoranthene | U | 2.0 | U | 2.2 | U | 2.2 | U | 9.6 | U | 9.6 |
| Benzo(k)fluoranthene | U | 1.9 | U | 2.1 | U | 2.1 | U | 9.3 | U | 9.3 |
| Benzo(e)pyrene | U | 2.0 | U | 2.2 | U | 2.2 | U | 9.5 | U | 9.5 |
| Benzo(a)pyrene | U | 2.0 | U | 2.2 | U | 2.2 | U | 9.6 | U | 9.6 |
| Indeno(1,2,3-cd)pyrene | U | 1.9 | U | 2.1 | U | 2.1 | U | 10 | U | 10 |
| Dibenzo(a,h)anthracene | U | 1.9 | U | 2.1 | U | 2.1 | U | 10 | U | 10 |
| Benzo(g,h,i)perylene | U | 2.0 | U | 2.1 | U | 2.2 | U | 10 | U | 10 |

Table 1.2 Results of the TIC for PAHs in Air
 WA # 0-110 Air Monitoring, Sampling,
 Analysis, and Modeling Support,
 and Underwater Survey Activities

| Sample ID | Compound Identification |
|-------------------|-------------------------|
| 28079 Lot Blank | No TICs were found |
| 28070 | No TICs were found |
| 28071 | No TICs were found |
| 28072 | No TICs were found |
| 28073 | No TICs were found |
| 28074 | No TICs were found |
| 28075 | No TICs were found |
| 28076 | No TICs were found |
| 28077 Field Blank | No TICs were found |
| 28078 Trip Blank | No TICs were found |

| Client ID | PBW | | 28050 | | 28051 | | 28052 | | 28053 | | 28054 | |
|----------------|------------|-----------|---------------|--------------|---------------|--------------|---------------|--------------|---------------|--------------|---------------|--------------|
| Location | - | | 0-1 | | 0-2 | | 0-3 | | 0-4 | | 0-5 | |
| Air Volume (L) | 0 | | 678 | | 698.5 | | 713 | | 669.9 | | 693 | |
| Parameter | Conc µg | MDL µg | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ |
| Aluminum | U | 1.0 | U | 1.5 | 1.5 | 1.4 | 3.5 | 1.4 | 2.7 | 1.5 | 4.0 | 1.4 |
| Arsenic | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Beryllium | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Cadmium | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Calcium | U | 1.0 | 8.8 | 1.5 | 8.3 | 1.4 | 9.0 | 1.4 | 9.0 | 1.5 | 9.7 | 1.4 |
| Chromium | U | 0.1 | 0.59 | 0.15 | 0.47 | 0.14 | 0.64 | 0.14 | 0.66 | 0.15 | 0.76 | 0.14 |
| Cobalt | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Copper | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Iron | U | 0.4 | 1.5 | 0.59 | 1.1 | 0.57 | 1.1 | 0.56 | 0.95 | 0.60 | 2.7 | 0.56 |
| Lead | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Lithium | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Magnesium | U | 1.0 | U | 1.5 | U | 1.4 | U | 1.4 | U | 1.5 | U | 1.4 |
| Manganese | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Molybdenum | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Nickel | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Phosphorus | U | 0.4 | U | 0.59 | U | 0.57 | U | 0.56 | U | 0.60 | U | 0.56 |
| Platinum | U | 1.0 | U | 1.5 | U | 1.4 | U | 1.4 | U | 1.5 | U | 1.4 |
| Selenium | U | 0.2 | U | 0.29 | U | 0.29 | U | 0.28 | U | 0.30 | U | 0.29 |
| Silver | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Sodium | U | 6.0 | 9.0 | 8.8 | 12.6 | 8.6 | 12.5 | 8.4 | 11.6 | 9.0 | 14 | 8.7 |
| Tellurium | U | 1.0 | U | 1.5 | U | 1.4 | U | 1.4 | U | 1.5 | U | 1.4 |
| Thallium | U | 0.4 | U | 0.59 | U | 0.57 | U | 0.56 | U | 0.60 | U | 0.56 |
| Tin | U | 0.2 | 1.3 | 0.29 | U | 0.29 | U | 0.28 | U | 0.30 | U | 0.29 |
| Titanium | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Vanadium | U | 0.1 | U | 0.15 | U | 0.14 | U | 0.14 | U | 0.15 | U | 0.14 |
| Yttrium | U | 0.2 | U | 0.29 | U | 0.29 | U | 0.28 | U | 0.30 | U | 0.29 |
| Zinc | U | 0.1 | 0.21 | 0.15 | U | 0.14 | 0.25 | 0.14 | 0.23 | 0.15 | 0.33 | 0.14 |
| Zirconium | U | 0.2 | U | 0.29 | U | 0.29 | U | 0.28 | U | 0.30 | U | 0.29 |

Table 1.3 (cont.) Results of the Analysis for Metals in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Client ID | 28055 | | 28056 | | 28057 | | 28058 | | 28059 | |
|----------------|---------------|--------------|---------------|--------------|-------------------|------------------|-------------------|------------------|-------------------|------------------|
| Location | 0-UW1 | | 0-UW2 | | Field Blank | | Trip Blank | | Lot Blank | |
| Air Volume (L) | 636 | | 648 | | 0 | | 0 | | 0 | |
| Parameter | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/filter | MDL µg/filter | Conc µg/filter | MDL µg/filter | Conc µg/filter | MDL µg/filter |
| Aluminum | 2.0 | 1.6 | 2.0 | 1.5 | 1.3 | 1.0 | 1.3 | 1.0 | 2.3 | 1.0 |
| Arsenic | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Beryllium | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Cadmium | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Calcium | 10 | 1.6 | 9.0 | 1.5 | 5.4 | 1.0 | 5.5 | 1.0 | 5.5 | 1.0 |
| Chromium | 0.97 | 0.16 | 0.7 | 0.15 | 0.56 | 0.10 | 0.49 | 0.10 | 0.44 | 0.10 |
| Cobalt | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Copper | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Iron | 1.2 | 0.63 | 2.1 | 0.62 | 0.82 | 0.4 | 0.45 | 0.4 | 0.45 | 0.4 |
| Lead | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Lithium | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Magnesium | U | 1.6 | U | 1.5 | U | 1.0 | U | 1.0 | U | 1.0 |
| Manganese | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Molybdenum | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Nickel | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Phosphorus | U | 0.63 | U | 0.62 | U | 0.4 | U | 0.4 | U | 0.4 |
| Platinum | U | 1.6 | U | 1.5 | U | 1.0 | U | 1.0 | U | 1.0 |
| Selenium | U | 0.31 | U | 0.31 | U | 0.20 | U | 0.20 | U | 0.20 |
| Silver | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Sodium | 13.8 | 9.4 | 12 | 9.3 | 12.3 | 6.0 | 7.8 | 6.0 | 10 | 6.0 |
| Tellurium | U | 1.6 | U | 1.5 | U | 1.0 | U | 1.0 | U | 1.0 |
| Thallium | U | 0.63 | U | 0.62 | U | 0.4 | U | 0.4 | U | 0.4 |
| Tin | U | 0.31 | U | 0.31 | U | 0.20 | U | 0.20 | U | 0.20 |
| Titanium | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Vanadium | U | 0.16 | U | 0.15 | U | 0.10 | U | 0.10 | U | 0.10 |
| Yttrium | U | 0.31 | U | 0.31 | U | 0.20 | U | 0.20 | U | 0.20 |
| Zinc | 0.31 | 0.16 | 0.16 | 0.15 | 0.11 | 0.10 | U | 0.10 | 0.12 | 0.10 |
| Zirconium | U | 0.31 | U | 0.31 | U | 0.20 | U | 0.20 | U | 0.20 |

Table 1.4 Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID Location | Blank 12/07/99 | | | | 28081 0-2 | | | | |
|-----------------------------|-------------------|------------|-----------|-----------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-------|
| Matrix Volume of Air (L) | Air 0 | | | | Air 687 | | | | |
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | TEF |
| 2,3,7,8-TCDD | U | 3.34 | 10.0 | 0 | U | 2.85 | 14.6 | 0 | 1 |
| 1,2,3,7,8-PeCDD | U | 8.20 | 10.0 | 0 | U | 13.0 | 14.6 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDD | U | 3.36 | 25.0 | 0 | U | 1.40 | 36.4 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDD | U | 10.2 | 25.0 | 0 | U | 7.39 | 36.4 | 0 | 0.1 |
| 1,2,3,7,8,9-HxCDD | U | 2.34 | 25.0 | 0 | U | 4.60 | 36.4 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | U | 6.82 | 25.0 | 0 | U | 9.26 | 36.4 | 0 | 0.01 |
| OCDD | 38.2 J | | 50.0 | 0.0382 | 38.1 J | | 72.8 | 0.0381 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | U | | | | U | | | | |
| Total Hexa-Dioxins | U | | | | U | | | | |
| Total Hepta-Dioxins | U | | | | U | | | | |
| 2,3,7,8-TCDF | U | 5.34 | 10.0 | 0 | U | 5.68 | 14.6 | 0 | 0.1 |
| 1,2,3,7,8-PeCDF | U | 5.96 | 10.0 | 0 | 7.69 J | | 14.6 | 0.3845 | 0.05 |
| 2,3,4,7,8-PeCDF | U | 2.82 | 10.0 | 0 | U | 1.80 | 14.6 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDF | U | 1.74 | 25.0 | 0 | U | 2.91 | 36.4 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDF | U | 6.98 | 25.0 | 0 | U | 6.35 | 36.4 | 0 | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | 1.40 | 25.0 | 0 | U | 1.05 | 36.4 | 0 | 0.1 |
| 2,3,4,6,7,8-HxCDF | U | 2.76 | 25.0 | 0 | U | 1.02 | 36.4 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | 12.4 J | | 25.0 | 0.124 | U | 17.5 | 36.4 | 0 | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | 3.38 | 25.0 | 0 | U | 4.1 | 36.4 | 0 | 0.01 |
| OCDF | 24.3 J | | 50.0 | 0.0243 | 21.5 J | | 72.8 | 0.0215 | 0.001 |
| Total Tetra-Furans | U | | | | U | | | | |
| Total Penta-Furans | U | | | | 13.9 | | | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | U | | | | U | | | | |
| Total | | | | 0.1865 | | | | 0.4441 | |

Table 1.4 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 28080 | | | | 28082 | | | | |
|---------------------|-------------------|------------|-----------|-----------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-------|
| Location | 0-1 (Field Blank) | | | | 0-3 | | | | |
| Matrix | Air | | | | Air | | | | |
| Volume of Air (L) | 0 | | | | 513.3 | | | | |
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | TEF |
| 2,3,7,8-TCDD | U | 1.44 | 10.0 | 0 | U | 5.73 | 19.5 | 0 | 1 |
| 1,2,3,7,8-PeCDD | 8.7 J | | 10.0 | 4.35 | U | 16.3 | 19.5 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDD | U | 0.740 | 25.0 | 0 | U | 1.68 | 48.7 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDD | U | 5.46 | 25.0 | 0 | U | 15.2 | 48.7 | 0 | 0.1 |
| 1,2,3,7,8,9-HxCDD | U | 2.18 | 25.0 | 0 | U | 1.40 | 48.7 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | U | 4.20 | 25.0 | 0 | 19.2 J | | 48.7 | 0.192 | 0.01 |
| OCDD | U | 20.3 | 50.0 | 0 | 54.8 J | | 97.4 | 0.0548 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | 8.70 | | | | U | | | | |
| Total Hexa-Dioxins | U | | | | U | | | | |
| Total Hepta-Dioxins | U | | | | 19.2 | | | | |
| 2,3,7,8-TCDF | U | 2.16 | 10.0 | 0 | U | 7.75 | 19.5 | 0 | 0.1 |
| 1,2,3,7,8-PeCDF | U | 7.46 | 10.0 | 0 | U | 11.8 | 19.5 | 0 | 0.05 |
| 2,3,4,7,8-PeCDF | U | 1.16 | 10.0 | 0 | U | 4.32 | 19.5 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDF | U | 1.54 | 25.0 | 0 | U | 3.00 | 48.7 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDF | U | 3.00 | 25.0 | 0 | U | 5.65 | 48.7 | 0 | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | 0.640 | 25.0 | 0 | U | 1.36 | 48.7 | 0 | 0.1 |
| 2,3,4,6,7,8-HxCDF | U | 0.920 | 25.0 | 0 | U | 3.27 | 48.7 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | U | 8.80 | 25.0 | 0 | U | 19.1 | 48.7 | 0 | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | 1.14 | 25.0 | 0 | U | 2.18 | 48.7 | 0 | 0.01 |
| OCDF | 11.5 J | | 50.0 | 0.0115 | 26.3 J | | 97.4 | 0.0263 | 0.001 |
| Total Tetra-Furans | | | | | | | | | |
| Total Penta-Furans | U | | | | U | | | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | U | | | | U | | | | |
| Total | | | | 4.3615 | | | | 0.2731 | |

Table 1.4 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 28083 | | | | | 28084 | | | | |
|---------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-------|--|
| Location | 0-4 | | | | | 0-5 | | | | |
| Matrix | Air | | | | | Air | | | | |
| Volume of Air (L) | 693 | | | | | 646.8 | | | | |
| Analyte | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | TEF | |
| 2,3,7,8-TCDD | U | 3.46 | 14.4 | 0 | U | 4.76 | 15.5 | 0 | 1 | |
| 1,2,3,7,8-PeCDD | 13.8 J | | 14.4 | 6.9 | U | 9.89 | 15.5 | 0 | 0.5 | |
| 1,2,3,4,7,8-HxCDD | U | 0.895 | 36.1 | 0 | U | 2.10 | 38.7 | 0 | 0.1 | |
| 1,2,3,6,7,8-HxCDD | U | 8.63 | 36.1 | 0 | U | 7.17 | 38.7 | 0 | 0.1 | |
| 1,2,3,7,8,9-HxCDD | U | 2.51 | 36.1 | 0 | U | 1.42 | 38.7 | 0 | 0.1 | |
| 1,2,3,4,6,7,8-HpCDD | U | 14.9 | 36.1 | 0 | U | 6.00 | 38.7 | 0 | 0.01 | |
| OCDD | U | 32.6 | 72.2 | 0 | U | 23.0 | 77.3 | 0 | 0.001 | |
| Total Tetra-Dioxins | U | | | | U | | | | | |
| Total Penta-Dioxins | 13.8 | | | | U | | | | | |
| Total Hexa-Dioxins | U | | | | U | | | | | |
| Total Hepta-Dioxins | U | | | | U | | | | | |
| 2,3,7,8-TCDF | U | 7.33 | 14.4 | 0 | 5.26 J | | 15.5 | 0.526 | 0.1 | |
| 1,2,3,7,8-PeCDF | 7.79 J | | 14.4 | 0.3895 | 7.02 J | | 15.5 | 0.351 | 0.05 | |
| 2,3,4,7,8-PeCDF | U | 1.41 | 14.4 | 0 | U | 3.31 | 15.5 | 0 | 0.5 | |
| 1,2,3,4,7,8-HxCDF | U | 2.48 | 36.1 | 0 | 1.79 J | | 38.7 | 0.179 | 0.1 | |
| 1,2,3,6,7,8-HxCDF | U | 8.20 | 36.1 | 0 | U | 6.71 | 38.7 | 0 | 0.1 | |
| 1,2,3,7,8,9-HxCDF | 0.491 J | | 36.1 | 0.0491 | U | 0.866 | 38.7 | 0 | 0.1 | |
| 2,3,4,6,7,8-HxCDF | U | 1.70 | 36.1 | 0 | U | 1.08 | 38.7 | 0 | 0.1 | |
| 1,2,3,4,6,7,8-HpCDF | U | 16.1 | 36.1 | 0 | U | 15.2 | 38.7 | 0 | 0.01 | |
| 1,2,3,4,7,8,9-HpCDF | U | 2.77 | 36.1 | 0 | U | 1.08 | 38.7 | 0 | 0.01 | |
| OCDF | 20.1 J | | 72.2 | 0.0201 | U | 28.0 | 77.3 | 0 | 0.001 | |
| Total Tetra-Furans | U | | | | 7.64 | | | | | |
| Total Penta-Furans | 7.79 | | | | 7.02 | | | | | |
| Total Hexa-Furans | 0.491 | | | | 1.79 | | | | | |
| Total Hepta-Furans | U | | | | U | | | | | |
| Total | | | | 7.3587 | | | | 1.056 | | |

Table 1.4 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID Location | 28085 10-UW1 | | | | 28086 10-UW2 | | | | |
|-----------------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-------|
| Matrix Volume of Air (L) | Air 406 | | | | Air 612 | | | | |
| Analyte | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | TEF |
| 2,3,7,8-TCDD | U | 6.06 | 24.6 | 0 | U | 2.45 | 16.3 | 0 | 1 |
| 1,2,3,7,8-PeCDD | 15.7 J | | 24.6 | 7.85 | 11.7 J | | 16.3 | 5.85 | 0.5 |
| 1,2,3,4,7,8-HxCDD | U | 3.74 | 61.6 | 0 | U | 1.37 | 40.8 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDD | U | 13.0 | 61.6 | 0 | 10.2 J | | 40.8 | 1.02 | 0.1 |
| 1,2,3,7,8,9-HxCDD | U | 4.04 | 61.6 | 0 | U | 1.14 | 40.8 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | 27.7 J | | 61.6 | 0.277 | U | 14.5 | 40.8 | 0 | 0.01 |
| OCDD | 91.8 J | | 123.0 | 0.0918 | 30.6 J | | 81.7 | 0.0306 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | 15.7 | | | | 11.7 | | | | |
| Total Hexa-Dioxins | U | | | | 10.2 | | | | |
| Total Hepta-Dioxins | 39.6 | | | | 4.05 | | | | |
| 2,3,7,8-TCDF | U | 7.59 | 24.6 | 0 | U | 2.88 | 16.3 | 0 | 0.1 |
| 1,2,3,7,8-PeCDF | U | 15.0 | 24.6 | 0 | 10.4 J | | 16.3 | 0.52 | 0.05 |
| 2,3,4,7,8-PeCDF | U | 1.48 | 24.6 | 0 | U | 3.40 | 16.3 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDF | U | 2.96 | 61.6 | 0 | U | 1.08 | 40.8 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDF | U | 18.2 | 61.6 | 0 | U | 9.05 | 40.8 | 0 | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | 2.81 | 61.6 | 0 | U | 0.882 | 40.8 | 0 | 0.1 |
| 2,3,4,6,7,8-HxCDF | U | 3.15 | 61.6 | 0 | U | 1.05 | 40.8 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | U | 28.6 | 61.6 | 0 | 14.6 J | | 40.8 | 0.146 | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | 6.45 | 61.6 | 0 | U | 0.784 | 40.8 | 0 | 0.01 |
| OCDF | 51.1 J | | 123.0 | 0.0511 | 20.3 J | | 81.7 | 0.0203 | 0.001 |
| Total Tetra-Furans | U | | | | U | | | | |
| Total Penta-Furans | U | | | | 10.4 | | | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | U | | | | 14.6 | | | | |
| Total | | | | 8.2699 | | | | 7.5869 | |

Table 1.4 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 28088 | | | | 28089 | | | | |
|---------------------|--------------|------------|-----------|-----------------------|--------------|------------|-----------|-----------------------|-------|
| Location | Trip Blank | | | | Trip Blank | | | | |
| Matrix | Air | | | | Air | | | | |
| Volume of Air (L) | 0 | | | | 0 | | | | |
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | TEF |
| 2,3,7,8-TCDD | U | 3.24 | 10.0 | 0 | U | 2.72 | 10.0 | 0 | 1 |
| 1,2,3,7,8-PeCDD | U | 7.86 | 10.0 | 0 | U | 6.78 | 10.0 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDD | U | 0.880 | 25.0 | 0 | U | 1.16 | 25.0 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDD | U | 5.90 | 25.0 | 0 | 8.62 J | | 25.0 | 0.862 | 0.1 |
| 1,2,3,7,8,9-HxCDD | U | 0.920 | 25.0 | 0 | U | 1.18 | 25.0 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | U | 6.26 | 25.0 | 0 | 5.74 J | | 25.0 | 0.0574 | 0.01 |
| OCDD | 17.2 J | | 50.0 | 0.0172 | 19.0 J | | 50.0 | 0.019 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | U | | | | U | | | | |
| Total Hexa-Dioxins | U | | | | 8.62 | | | | |
| Total Hepta-Dioxins | U | | | | 5.74 | | | | |
| 2,3,7,8-TCDF | U | 1.72 | 10.0 | 0 | U | 2.90 | 10.0 | 0 | 0.1 |
| 1,2,3,7,8-PeCDF | 7.48 J | | 10.0 | 0.374 | 4.18 J | | 10.0 | 0.209 | 0.05 |
| 2,3,4,7,8-PeCDF | U | 0.960 | 10.0 | 0 | U | 0.620 | 10.0 | 0 | 0.5 |
| 1,2,3,4,7,8-HxCDF | U | 1.04 | 25.0 | 0 | U | 0.900 | 25.0 | 0 | 0.1 |
| 1,2,3,6,7,8-HxCDF | U | 4.46 | 25.0 | 0 | U | 4.78 | 25.0 | 0 | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | 0.420 | 25.0 | 0 | U | 0.540 | 25.0 | 0 | 0.1 |
| 2,3,4,6,7,8-HxCDF | U | 0.220 | 25.0 | 0 | U | 0.940 | 25.0 | 0 | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | 7.12 J | | 25.0 | 0.0712 | U | 8.96 | 25.0 | 0 | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | 0.740 | 25.0 | 0 | U | 1.72 | 25.0 | 0 | 0.01 |
| OCDF | 11.3 J | | 50.0 | 0.0113 | U | 15.3 | 50.0 | 0 | 0.001 |
| Total Tetra-Furans | U | | | | 1.28 | | | | |
| Total Penta-Furans | 11.1 | | | | 4.18 | | | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | 7.12 | | | | U | | | | |
| Total | | | | 0.4737 | | | | 1.1474 | |

Table 1.5 Results of the Analysis for Inorganic Acids in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 28090 | | 28091 | | 28092 | | 28060 | | 28061 | |
|-------------------|-------------|-----------|------------|-----------|------------|-----------|---------------------------|--------------------------|---------------------------|--------------------------|
| Location | Field Blank | | Trip blank | | Lot Blank | | 0-1 | | 0-2 | |
| Air Volume (L): | 0 | | 0 | | 0 | | 59.2 | | 41.2 | |
| Analyte | Conc mg | MDL mg | Conc mg | MDL mg | Conc mg | MDL mg | Conc mg/m ³ | MDL mg/m ³ | Conc mg/m ³ | MDL mg/m ³ |
| Hydrobromic acid | U | 0.0011 | U | 0.0011 | U | 0.0011 | U | 0.0171 | U | 0.0246 |
| Hydrochloric acid | U | 0.001 | U | 0.001 | U | 0.001 | U | 0.0174 | U | 0.0250 |
| Hydrofluoric acid | U | 0.001 | U | 0.001 | U | 0.001 | U | 0.0178 | U | 0.0256 |
| Nitric acid | U | 0.0045 | U | 0.0045 | U | 0.0045 | U | 0.0760 | U | 0.1092 |
| Phosphoric acid | U | 0.0032 | U | 0.0032 | U | 0.0032 | U | 0.0523 | U | 0.0752 |
| Sulfuric acid | U | 0.001 | U | 0.001 | U | 0.001 | U | 0.0172 | U | 0.0248 |

| Sample ID | 28062 | | 28063 | | 28064 | | 28065 | | 28066 | |
|-------------------|---------------------------|--------------------------|---------------------------|--------------------------|---------------------------|--------------------------|---------------------------|--------------------------|---------------------------|--------------------------|
| Location | 0-3 | | 0-4 | | 0-5 | | 0-UW1 | | 0-UW2 | |
| Air Volume (L): | 57.8 | | 52.0 | | 58.0 | | 53.0 | | 43.2 | |
| Analyte | Conc mg/m ³ | MDL mg/m ³ | Conc mg/m ³ | MDL mg/m ³ | Conc mg/m ³ | MDL mg/m ³ | Conc mg/m ³ | MDL mg/m ³ | Conc mg/m ³ | MDL mg/m ³ |
| Hydrobromic acid | U | 0.0175 | U | 0.0195 | U | 0.0174 | U | 0.0191 | U | 0.0234 |
| Hydrochloric acid | U | 0.0178 | U | 0.0198 | U | 0.0177 | U | 0.0194 | U | 0.0238 |
| Hydrofluoric acid | U | 0.0182 | U | 0.0203 | U | 0.0182 | U | 0.0199 | U | 0.0244 |
| Nitric acid | U | 0.0779 | U | 0.0865 | U | 0.0776 | U | 0.0849 | U | 0.1042 |
| Phosphoric acid | U | 0.0536 | U | 0.0596 | U | 0.0534 | U | 0.0584 | U | 0.0717 |
| Sulfuric acid | U | 0.0177 | U | 0.0196 | U | 0.0176 | U | 0.0193 | U | 0.0236 |

QA/QC for PAH in Air

Results of the BS/BSD Analysis for PAH in Air

A lot blank and a lot blank filter were chosen for the blank spike/blank spike duplicate (BS/BSD) analyses. The percent recoveries, for the lot blank, ranging from 78 to 97, are listed in Table 2.1. The relative percent differences, also listed in Table 2.1, ranged from 3 to 8. The percent recoveries, for the lot blank filter, ranging from 41 to 101, are also listed in Table 2.1. The relative percent differences, also listed in Table 2.1, ranged from 7 to 70. QC limits are not available for either the percent recoveries or the relative percent differences for this analysis.

Table 2.1 Results of BS/BSD Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

Sample ID: Lot Blank

| Compound | Spike Added µg | BS Rec. µg | % Rec. | BSD Rec. µg | % Rec. | RPD |
|-------------------------|----------------------|------------------|--------|-------------------|--------|-----|
| Naphthalene | 50 | 48.18 | 96 | 46.27 | 93 | 4 |
| 2-Methylnaphthalene | 50 | 47.78 | 96 | 45.73 | 91 | 5 |
| 1-Methylnaphthalene | 50 | 48.27 | 97 | 44.76 | 90 | 8 |
| Biphenyl | 50 | 47.72 | 95 | 45.26 | 91 | 5 |
| 2,6-Dimethylnaphthalene | 50 | 47.60 | 95 | 45.11 | 90 | 5 |
| Acenaphthylene | 50 | 47.86 | 96 | 46.05 | 92 | 4 |
| Acenaphthene | 50 | 48.20 | 96 | 46.41 | 93 | 4 |
| Dibenzofuran | 50 | 47.16 | 94 | 45.51 | 91 | 4 |
| Fluorene | 50 | 47.86 | 96 | 45.90 | 92 | 4 |
| Phenanthrene | 50 | 47.98 | 96 | 45.95 | 92 | 5 |
| Anthracene | 50 | 47.68 | 95 | 46.43 | 93 | 3 |
| Carbazole | 50 | 48.49 | 97 | 46.04 | 92 | 5 |
| Fluoranthene | 50 | 48.13 | 96 | 46.26 | 93 | 4 |
| Pyrene | 50 | 48.69 | 97 | 46.62 | 93 | 4 |
| Benzo(a)anthracene | 50 | 48.19 | 96 | 45.25 | 91 | 6 |
| Chrysene | 50 | 37.28 | 75 | 38.75 | 78 | 4 |
| Benzo(b)fluoranthene | 50 | 48.09 | 96 | 44.67 | 89 | 7 |
| Benzo(k)fluoranthene | 50 | 46.43 | 93 | 44.84 | 90 | 4 |
| Benzo(e)pyrene | 50 | 47.46 | 95 | 45.32 | 91 | 5 |
| Benzo(a)pyrene | 50 | 47.14 | 94 | 44.61 | 89 | 6 |
| Indeno(1,2,3-cd)pyrene | 50 | 48.18 | 96 | 45.77 | 92 | 5 |
| Dibenzo(a,h)anthracene | 50 | 48.08 | 96 | 45.39 | 91 | 6 |
| Benzo(g,h,i)perylene | 50 | 48.06 | 96 | 45.69 | 91 | 5 |

Table 2.1 (cont.) Results of BS/BSD Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support,
and Underwater Survey Activities

Sample ID: Lot Blank filter

| Compound | Spike Added µg | BS Rec. µg | % Rec. | BSD Rec. µg | % Rec. | RPD |
|-------------------------|----------------------|------------------|--------|-------------------|--------|-----|
| Naphthalene | 50 | 42.80 | 86 | 20.62 | 41 | 70 |
| 2-Methylnaphthalene | 50 | 46.56 | 93 | 33.82 | 68 | 32 |
| 1-Methylnaphthalene | 50 | 47.38 | 95 | 36.54 | 73 | 26 |
| Biphenyl | 50 | 48.02 | 96 | 39.62 | 79 | 19 |
| 2,6-Dimethylnaphthalene | 50 | 47.60 | 95 | 39.74 | 79 | 18 |
| Acenaphthylene | 50 | 48.36 | 97 | 41.68 | 83 | 15 |
| Acenaphthene | 50 | 49.52 | 99 | 43.70 | 87 | 12 |
| Dibenzofuran | 50 | 49.68 | 99 | 44.26 | 89 | 11 |
| Fluorene | 50 | 49.36 | 99 | 44.44 | 89 | 11 |
| Phenanthrene | 50 | 48.38 | 97 | 42.82 | 86 | 12 |
| Anthracene | 50 | 50.60 | 101 | 46.20 | 92 | 9 |
| Carbazole | 50 | 48.48 | 97 | 43.24 | 86 | 12 |
| Fluoranthene | 50 | 50.52 | 101 | 45.06 | 90 | 11 |
| Pyrene | 50 | 50.02 | 100 | 44.44 | 89 | 12 |
| Benzo(a)anthracene | 50 | 48.50 | 97 | 42.82 | 86 | 12 |
| Chrysene | 50 | 50.08 | 100 | 46.74 | 93 | 7 |
| Benzo(b)fluoranthene | 50 | 47.16 | 94 | 40.98 | 82 | 14 |
| Benzo(k)fluoranthene | 50 | 45.04 | 90 | 40.14 | 80 | 12 |
| Benzo(e)pyrene | 50 | 47.02 | 94 | 41.60 | 83 | 12 |
| Benzo(a)pyrene | 50 | 44.00 | 88 | 39.32 | 79 | 11 |
| Indeno(1,2,3-cd)pyrene | 50 | 44.26 | 89 | 39.74 | 79 | 11 |
| Dibenzo(a,h)anthracene | 50 | 43.44 | 87 | 40.00 | 80 | 8 |
| Benzo(g,h,i)perylene | 50 | 43.82 | 88 | 40.08 | 80 | 9 |

QA/QC for Metals in Air

Results of the BS/BSD Analysis for Metals in Air

A blank spike/blank spike duplicate analysis (BS/BSD) was run. The percent recoveries, listed in Table 2.2, ranged from 34 to 125. Fifty out of fifty-six values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.2, ranged from 0 (zero) to 7. QC limits are not available for this criterion.

Results of the Analysis of the Laboratory Control Sample for Metals in Air

Laboratory control samples were also analyzed. The percent recoveries ranged from 62 to 116 and are listed in Table 2.3. Twenty-five out of twenty-eight concentrations were within the acceptable QC limits.

Table 2.2 Results of the BS/BSD Analysis for Metals in Air
WA # D-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Metal | Sample Conc µg/filter | Original Conc | | Recovered Conc | | % Recovery | | RPD | Recommended QC Limit % Rec |
|------------|--------------------------|--------------------|------------------|--------------------|------------------|------------|-----|-----|-------------------------------|
| | | Spike µg/filter | Dup µg/filter | Spike µg/filter | Dup µg/filter | Spike | Dup | | |
| Aluminum | 2.3292 | 40.00 | 40.00 | 52.26 | 48.966 | 125 | 117 | 7 | 75-125 |
| Arsenic | U | 40.00 | 40.00 | 40.839 | 40.458 | 102 | 101 | 1 | 75-125 |
| Beryllium | U | 1.00 | 1.00 | 1.0592 | 1.046 | 106 | 105 | 1 | 75-125 |
| Cadmium | U | 1.00 | 1.00 | 1.0498 | 1.0388 | 105 | 104 | 1 | 75-125 |
| Calcium | 5.5364 | 1000 | 1000 | 1085.9 | 1074.6 | 108 | 107 | 1 | 75-125 |
| Chromium | 0.4428 | 4.00 | 4.00 | 4.9778 | 5.2212 | 113 | 120 | 6 | 75-125 |
| Cobalt | U | 10.00 | 10.00 | 10.184 | 10.04 | 102 | 100 | 1 | 75-125 |
| Copper | U | 5.00 | 5.00 | 5.47 | 5.41 | 109 | 108 | 1 | 75-125 |
| Iron | 0.4518 | 20.00 | 20.00 | 21.624 | 21.132 | 106 | 103 | 2 | 75-125 |
| Lead | U | 10.00 | 10.00 | 10.885 | 10.719 | 109 | 107 | 2 | 75-125 |
| Lithium | U | 40.00 | 40.00 | 46.163 | 46.319 | 115 | 116 | 0 | 75-125 |
| Magnesium | U | 1000 | 1000 | 1088.4 | 1078.6 | 109 | 108 | 1 | 75-125 |
| Manganese | U | 10.00 | 10.00 | 10.385 | 10.242 | 104 | 102 | 1 | 75-125 |
| Molybdenum | U | 40.00 | 40.00 | 42.838 | 42.729 | 107 | 107 | 0 | 75-125 |
| Nickel | U | 10.00 | 10.00 | 10.473 | 10.37 | 105 | 104 | 1 | 75-125 |
| Phosphorus | U | 40.00 | 40.00 | 20.931 | 20.372 | 52 | 51 | 3 | 75-125 |
| Platinum | U | 40.00 | 40.00 | 40.866 | 39.58 | 102 | 99 | 3 | 75-125 |
| Selenium | U | 40.00 | 40.00 | 40.464 | 39.955 | 101 | 100 | 1 | 75-125 |
| Silver | U | 1.00 | 1.00 | 1.0042 | 0.9952 | 100 | 100 | 1 | 75-125 |
| Sodium | 9.9974 | 1000 | 1000 | 1050.3 | 1039.6 | 104 | 103 | 1 | 75-125 |
| Tellurium | U | 40.00 | 40.00 | 39.822 | 38.988 | 100 | 97 | 2 | 75-125 |
| Thallium | U | 40.00 | 40.00 | 45.497 | 44.952 | 114 | 112 | 1 | 75-125 |
| Tin | U | 40.00 | 40.00 | 14.508 | 13.513 | 36 | 34 | 7 | 75-125 |
| Titanium | U | 40.00 | 40.00 | 41.055 | 40.774 | 103 | 102 | 1 | 75-125 |
| Vanadium | U | 10.00 | 10.00 | 10.586 | 10.46 | 106 | 105 | 1 | 75-125 |
| Yttrium | U | 40.00 | 40.00 | 42.336 | 41.878 | 106 | 105 | 1 | 75-125 |
| Zinc | 0.1208 | 10.00 | 10.00 | 10.784 | 10.535 | 107 | 104 | 2 | 75-125 |
| Zirconium | U | 40.00 | 40.00 | 14.671 | 14.075 | 37 | 35 | 4 | 75-125 |

Table 2.3 Results of the Analysis of the
Laboratory Control Sample for Metals in Air
WA # 0-110 Air Monitoring, Sampling, Analysis,
and Modeling Support, and Underwater Survey Activities

| Metal | Analyzed Value µg/L | Accepted Value µg/L | % Rec | QC Limits % Rec |
|------------|---------------------------|---------------------------|----------|--------------------|
| Aluminum | 4617.55 | 4000 | 115 | 80-120 |
| Arsenic | 4116.9 | 4000 | 103 | 80-120 |
| Beryllium | 107.5 | 100 | 107 | 80-120 |
| Cadmium | 103.18 | 100 | 103 | 80-120 |
| Calcium | 53193 | 50000 | 106 | 80-120 |
| Chromium | 439.2 | 400 | 110 | 80-120 |
| Cobalt | 1006.11 | 1000 | 101 | 80-120 |
| Copper | 541.74 | 500 | 108 | 80-120 |
| Iron | 2144.08 | 2000 | 107 | 80-120 |
| Lead | 1061.38 | 1000 | 106 | 80-120 |
| Lithium | 2319.62 | 2000 | 116 | 80-120 |
| Magnesium | 53566.16 | 50000 | 107 | 80-120 |
| Manganese | 1018.99 | 1000 | 102 | 80-120 |
| Molybdenum | 2127.76 | 2000 | 106 | 80-120 |
| Nickel | 1032.58 | 1000 | 103 | 80-120 |
| Phosphorus | 1451.11 | 2000 | 73 | 80-120 |
| Platinum | 2068.2 | 2000 | 103 | 80-120 |
| Selenium | 4062.26 | 4000 | 102 | 80-120 |
| Silver | 100.04 | 100 | 100 | 80-120 |
| Sodium | 50626.04 | 50000 | 101 | 80-120 |
| Tellurium | 1994.5 | 2000 | 100 | 80-120 |
| Thallium | 4513.77 | 4000 | 113 | 80-120 |
| Tin | 1250.53 | 2000 | 63 | 80-120 |
| Titanium | 2044.16 | 2000 | 102 | 80-120 |
| Vanadium | 1040.05 | 1000 | 104 | 80-120 |
| Yttrium | 2113.37 | 2000 | 106 | 80-120 |
| Zinc | 1028.63 | 1000 | 103 | 80-120 |
| Zirconium | 1244.98 | 2000 | 62 | 80-120 |

QA/QC for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

Results of the Internal Standard Recoveries for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

The results of the internal standard recoveries, listed in Table 2.4, ranged from 62 to 146. One hundred and two out of one hundred and eight values were within the acceptable QC limits.

Results of the BS/BSD Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

A blank was spiked in duplicate and analyzed. The percent recoveries ranged from 81 to 122 and are listed in Table 2.5. All thirty-four values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.5, ranged from 0 (zero) to 14. QC limits are not available for this analysis.

Table 2.4 Results of the Internal Standard Recoveries for Polychlorinated Dibenzodioxin
and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support
and Underwater Survey Activities

| Sample ID | Method Blank | 28080 | 28081 | 28082 | 28083 | 28084 | QC Limits |
|-------------------------|-----------------|----------|----------|----------|----------|----------|--------------|
| Location | | 0-1 | 0-2 | 0-3 | 0-4 | 0-5 | |
| Matrix Units | Air % | Air % | Air % | Air % | Air % | Air % | Percent |
| Internal Standard | | | | | | | |
| 13C-2,3,7,8-TCDD | 79 | 100 | 79 | 86 | 94 | 84 | 40-135 |
| 13C-1,2,3,6,7,8-HxCDD | 85 | 96 | 80 | 113 | 113 | 111 | 40-135 |
| 13C-2,3,7,8-TCDF | 79 | 103 | 80 | 108 | 109 | 109 | 40-135 |
| 13C-1,2,3,4,7,8-HxCDF | 81 | 91 | 77 | 119 | 112 | 110 | 40-135 |
| 13C-1,2,3,7,8-PeCDD | 123 | 130 | 126 | 100 | 95 | 93 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDD | 80 | 86 | 68 | 85 | 107 | 72 | 40-135 |
| 13C-1,2,3,7,8-PeCDF | 117 | 127 | 122 | 98 | 104 | 97 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDF | 86 | 99 | 83 | 115 | 107 | 106 | 40-135 |
| 13C-OCDD | 115 | 122 | 107 | 80 | 89 | 76 | 40-135 |

| Sample ID | 28085 | 28086 | 28088 | 28089 | Blank Spike | Blank Spike Duplicate | QC Limits |
|-------------------------|----------|----------|---------------|---------------|----------------|-----------------------------|--------------|
| Location | 10-UW1 | 10-UW2 | Trip Blank | Trip Blank | | | |
| Matrix Units | Air % | Air % | Air % | Air % | Air % | Air % | Percent |
| Internal Standard | | | | | | | |
| 13C-2,3,7,8-TCDD | 63 | 85 | 83 | 88 | 75 | 79 | 40-135 |
| 13C-1,2,3,6,7,8-HxCDD | 94 | 109 | 110 | 113 | 111 | 110 | 40-135 |
| 13C-2,3,7,8-TCDF | 114 | 146 | 138 | 143 | 142 | 128 | 40-135 |
| 13C-1,2,3,4,7,8-HxCDF | 98 | 119 | 128 | 129 | 124 | 124 | 40-135 |
| 13C-1,2,3,7,8-PeCDD | 81 | 85 | 77 | 83 | 75 | 70 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDD | 110 | 103 | 112 | 113 | 100 | 103 | 40-135 |
| 13C-1,2,3,7,8-PeCDF | 91 | 106 | 91 | 102 | 91 | 88 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDF | 119 | 126 | 139 | 141 | 132 | 127 | 40-135 |
| 13C-OCDD | 62 | 85 | 91 | 95 | 75 | 82 | 40-135 |

Table 2.5 Results of the BS/BSD Analysis
for Polychlorinated Dibenzodioxin and Polychlorinated Dibenzofurans in Air
WA # D-110 Air Monitoring, Sampling, Analysis, and Modeling Support
and Underwater Survey Activities

Sample ID Blank

| Parameter | Spike Added pg | Sample Conc pg | BS Conc pg | % Rec | BSD Conc pg | % Rec | RPD | QC Limits (% Rec) |
|---------------|----------------------|----------------------|------------------|----------|-------------------|----------|-----|-------------------------|
| 2378-TCDD | 200 | U | 239 | 120 | 244 | 122 | 2 | 60-140 |
| 12378-PeCDD | 200 | U | 214 | 107 | 230 | 115 | 7 | 60-140 |
| 123478-HxCDD | 500 | U | 550 | 110 | 507 | 101 | 8 | 60-140 |
| 123678-HxCDD | 500 | U | 473 | 95 | 493 | 99 | 4 | 60-140 |
| 123789-HxCDD | 500 | U | 429 | 86 | 418 | 84 | 3 | 60-140 |
| 1234678-HpCDD | 500 | U | 494 | 99 | 525 | 105 | 6 | 60-140 |
| OCDD | 1000 | 38.2 | 969 | 93 | 1040 | 100 | 7 | 60-140 |
| 2378-TCDF | 200 | U | 165 | 83 | 189 | 95 | 14 | 60-140 |
| 12378-PeCDF | 200 | U | 218 | 109 | 239 | 120 | 9 | 60-140 |
| 23478-PeCDF | 200 | U | 232 | 116 | 233 | 116 | 0 | 60-140 |
| 123478-HxCDF | 500 | U | 455 | 91 | 461 | 92 | 1 | 60-140 |
| 123678-HxCDF | 500 | U | 466 | 93 | 469 | 94 | 1 | 60-140 |
| 123789-HxCDF | 500 | U | 435 | 87 | 421 | 84 | 3 | 60-140 |
| 234678-HxCDF | 500 | U | 513 | 103 | 498 | 100 | 3 | 60-140 |
| 1234678-HpCDF | 500 | 12.4 | 418 | 81 | 461 | 90 | 10 | 60-140 |
| 1234789-HpCDF | 500 | U | 407 | 81 | 434 | 87 | 6 | 60-140 |
| OCDF | 1000 | 24.3 | 1090 | 107 | 1090 | 107 | 0 | 60-140 |

QA/QC for Inorganic Acids in Air

Results of the BS/BSD Analysis for Inorganic Acids in Air

A blank spike/blank spike duplicate analysis (BS/BSD) was run. The percent recoveries, listed in Table 2.6, ranged from 93 to 100. All twelve values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.6, ranged from 0 (zero) to 1. QC limits are not available for this criterion.

Table 2.6 Results of the BS/BSD Analysis for Inorganic Acids in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Analyte | Original Conc | | Recovered Conc | | % Recovery | | RPD | Recommended QC Limits % Rec |
|-------------------|---------------|-----------|----------------|-----------|------------|-----|-----|-----------------------------------|
| | Spike mg | Dup mg | Spike mg | Dup mg | Spike | Dup | | |
| Hydrobromic acid | 0.1053 | 0.1053 | 0.0983 | 0.0983 | 93 | 93 | 0 | 75-125 |
| Hydrochloric acid | 0.2054 | 0.2054 | 0.2042 | 0.2042 | 99 | 99 | 0 | 75-125 |
| Hydrofluoric acid | 0.4049 | 0.4049 | 0.3925 | 0.3937 | 97 | 97 | 0 | 75-125 |
| Nitric acid | 0.4067 | 0.4067 | 0.3901 | 0.3919 | 96 | 96 | 0 | 75-125 |
| Phosphoric acid | 0.5914 | 0.5914 | 0.5884 | 0.592 | 99 | 100 | 1 | 75-125 |
| Sulfuric acid | 0.4085 | 0.4085 | 0.3949 | 0.3937 | 97 | 96 | 0 | 75-125 |

LOCKHEED MARTIN

Southwest Research Institute
PO Box 28510, 6220 Culebra Road
San Antonio, TX 78228-0510

Attn: Jo Ann Boyd

19 November 1999

Project # RIA-00011 APG Burn Support

As per Lockheed Martin / REAC Purchase Order GA91969J73, please analyze samples according to the following parameters:

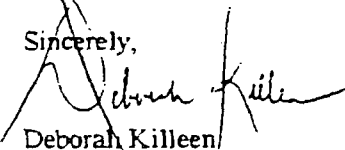
| Analysis/Method | Matrix | # of samples |
|---|--------|--------------|
| Dioxin/ Furans / Modified TO9 | Air | 20 |
| Inorganic Acids / NIOSH 7903 | Air | 20 |
| Metals/ NIOSH 7300 | Air | 20 |
| Data package: Package with Diskette Deliverable | | |

Samples are expected to arrive at your laboratory between November 23-December 31, 1999. All applicable QA/QC (BS/BSD) analysis as per method, will be performed on our sample matrix. Preliminary sample and QC result tables plus a signed copy of our Chain of Custody must be faxed to REAC 10 business days after receipt of the last samples. The complete data package is due 21 business days after receipt of the last samples. The complete data package must include all items on the deliverables checklist. **Expect all samples to be difficult matrix and all raw data must be included in final analytical report.**

All sample and QC results(ie: BS/BSD, LCS, Duplicates, and Blanks) must be summarized in a ExCel diskette deliverable.

Please submit all reports and technical questions concerning this project to John Johnson at (732) 321-4248 or fax to (732) 494-4020.

Sincerely,


Deborah Killeen
Data Validation and Report Writing Group Leader
Lockheed Martin / REAC Project

DK:jj Attachments

cc. R. Singhvi
D. Michunas
0011\non\mem\9911\sub\0011Con

D. Miller
Subcontracting File
D. Angwenyi

C. Lentini
A. DuBois
D. Killeen

732 (908) 321-4200

EPA Contract 68-C4-0022-4D

099-223

Project Name:

APG BURN SUPPORT

Project Number:

R1A00110

RFW Contact:

John Vikstrom

Phone:

732-321-4248

No:

03215

SHEET NO. 1 OF 1

Sample Identification

DIOXINS

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Volume (L) | Dioxins |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|------------|---------|
| | 28080 | 0-1 | A | 12/3/99 | 1 | AMIBGE JAR/KE | 0.1 | ✓ |
| | 28081 | 0-2 | | | 1 | | 0.87 L | ✓ |
| | 28082 | 0-3 | | | 1 | | 513.3 | ✓ |
| | 28083 | 0-4 | | | 1 | | 693 | ✓ |
| | 28084 | 0-5 | | | 1 | | 646.8 | ✓ |
| | 28085 | 0-UW1 | | | 1 | | 406 | ✓ |
| | 28086 | 0-UW2 | | | 1 | | 612 | ✓ |
| | 28089 | TRIP Blank | | | 1 | | 0 | ✓ |
| | 28088 | TRIP Blank | | | 1 | | 0 | ✓ |
| | - | BS/BSD | ↓ | ↓ | 2 | ↓ | 0 | ✓ |

Matrix:

SD - Sediment

PW - Potable Water

S - Soil

DS - Drum Solids

GW - Groundwater

W - Water

DL - Drum Liquids

SW - Surface Water

O - Oil

X - Other

SL - Sludge

A - Air

Special Instructions:

Puf plugs + Quartz
Filters certified
Clean by Southwest
Research Institute

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF CUSTODY #

Southwest Research Institute
16220 Culebra Road
San Antonio, TX 78228

Will use as FB-no flow through plug + filter

| Items/Reason | Relinquished By | Date | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|---------|-------------|---------|------|--------------|-----------------|------|-------------|------|------|
| 14/1/1999 | [Signature] | 12/4/99 | [Signature] | 12/4/99 | 0735 | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |

FORM #4

8/94

EPA Contract 68-~~C4-0022~~ (41)
C44-223

Project Name: APG - Ben Support

Project Number: 24A00110

RPW Contact: John Johnson

Phone: 732-321-4248

No: 03132

SHEET NO. / OF /

Sample Identification

Inorganic Acids

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Volume (L) | Inorganic Acids |
|--------|------------|-------------------|--------|----------------|--------------|--------------------------------|------------|-----------------|
| | 28060 | 0-1 | A | 12/3/99 | 1 | unw/peck / 10% ^{100%} | 59.2 | ✓ |
| | 28061 | 0-2 | | | 1 | | 41.2 | ✓ |
| | 28062 | 0-3 | | | 1 | | 57.8 | ✓ |
| | 28063 | 0-4 | | | 1 | | 52.0 | ✓ |
| | 28064 | 0-5 | | | 1 | | 58.0 | ✓ |
| | 28065 | 0-UW1 | | | 1 | | 53.0 | ✓ |
| | 28066 | 0-UW2 | | | 1 | | 43.2 | ✓ |
| | 28090 | Field Blank | | | 1 | | 0 | ✓ |
| | 28091 | Trip Blank | | | 1 | | 0 | ✓ |
| | 28092 | LOT Blank | | | 1 | | 0 | ✓ |
| | — | BS/BSD | | | 2 | | 0 | ✓ |

(100)

scale temp. 12.6°C

Matrix:

| | | |
|-------------------|--------------------|-----------|
| SD - Sediment | PW - Potable Water | S - Soil |
| DS - Drum Solids | GW - Groundwater | W - Water |
| DL - Drum Liquids | SW - Surface Water | O - Oil |
| X - Other | SL - Sludge | A - Air |

Special instructions:

① Inorganic Acids by NIOSH 7903
All Inorganic acids samples
used Lot # 1038 silica tubes

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF

Southwest Research Institute
6220 Culebra Rd
San Antonio, TX 78228

[illegible]Post-It[®] Fax Note 7671

7671

| | |
|-----------------------|----------------------|
| Date <u>1/19/2000</u> | No of pages <u>5</u> |
|-----------------------|----------------------|

| | |
|---------------|---|
| N of pages | 2 |
|---------------|---|

To John Johnson

From Joe Morin

Co/Dept. Lockheed Martin Aerospace

| | |
|-----|-------|
| CO. | SW RI |
|-----|-------|

Phone #

Phone 6

Fax: 771-494-4020

Fax # 210-522-6628

Time

732) REAC, Edison, NJ
 (800) 321-4200
 EPA Contract 68-C4-0022-AD
 09-223

CHAIN OF CUSTODY RECORD

Project Name: APG Burn Support
 Project Number: 21A00110
 RFW Contact: John Johnson Phone: 732-321-4248

No: 03133

SHEET NO. 1 OF 1

Sample Identification METALS

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Volume (L) | Notes |
|---|------------|-------------------|--------|----------------|--------------|------------------------|------------|-------|
| | 28050 | O-1 | H | 12/3/99 | 1 | W-Vickie/ice | 678 | ✓ |
| | 28051 | O-2 | | | | | 698.5 | ✓ |
| | 28052 | O-3 | | | | | 713 | ✓ |
| | 28053 | O-4 | | | | | 669.9 | ✓ |
| | 28054 | O-5 | | | | | 693 | ✓ |
| | 28055 | O-UW1 | | | | | 636 | ✓ |
| | 28056 | O-UW2 | | | | | 648 | ✓ |
| | 28057 | FIELD BLANK | | | | | 0 | ✓ |
| | 28058 | Trip Blank | | | | | 0 | ✓ |
| | 28059 | LOT BLANK | | | | | 0 | ✓ |
| | | BS/BSD | ✓ | ✓ | 2 | ✓ | 0 | ✓ |
| <div>000035</div> <div>RED</div> <div>Cooler Temp. 12°C</div> | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |

Matrix:
 SD - Sediment
 DS - Drum Solids
 DL - Drum Liquids
 X - Other
 PW - Potable Water
 GW - Groundwater
 SW - Surface Water
 SL - Sludge
 S - Soil
 W - Water
 O - Oil
 A - Air

Special Instructions:
 All metals (elements) samples
 collected using MCE Filter
 LOT # 88900

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF CUSTODY #

Southwest Research Institute
 6220 Culebra Rd
 San Antonio, TX 78228

| Item/Reason | Relinquished By | Date | Received By | Date | Time | Item/Reason | Relinquished By | Date | Received By | Date | Time |
|-------------|-----------------|---------|-------------|---------|-------|-------------|-----------------|------|-------------|------|------|
| 11/10/99 | [Signature] | 12/4/99 | [Signature] | 12/4/99 | 09:55 | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |

APPENDIX C
SBC COM Clearances for GB, GD, VX, and HD
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

From: Smith Sandra D SBCCOM <sandra.smith@SBCCOM.APGEA.ARMY.MIL>
To: Alfreda Dean <alfreda.dean@SBCCOM.APGEA.ARMY.MIL>, ...
Date: 12/8/99 4:57pm
Subject: EPA Clearances

| | | |
|--------------|----------------|--------------------------|
| POC: DuBois, | 732-494-4013 | O-FLD |
| Item# | GVH BKGD | taken 12/06/99 |
| UW-1 | 9912060122-M01 | Clear for GB, GD, VX, HD |
| UW-2 | 9912060123-M01 | Clear for GB, GD, VX, HD |
| O-1 | 9912060124-M01 | Clear for GB, GD, VX, HD |
| O-2 | 9912060125-M01 | Clear for GB, GD, VX, HD |
| O-3 | 9912060126-M01 | Clear for GB, GD, VX, HD |
| O-4 | 9912060127-M01 | Clear for GB, GD, VX, HD |
| O-5 | 9912060128-M01 | Clear for GB, GD, VX, HD |

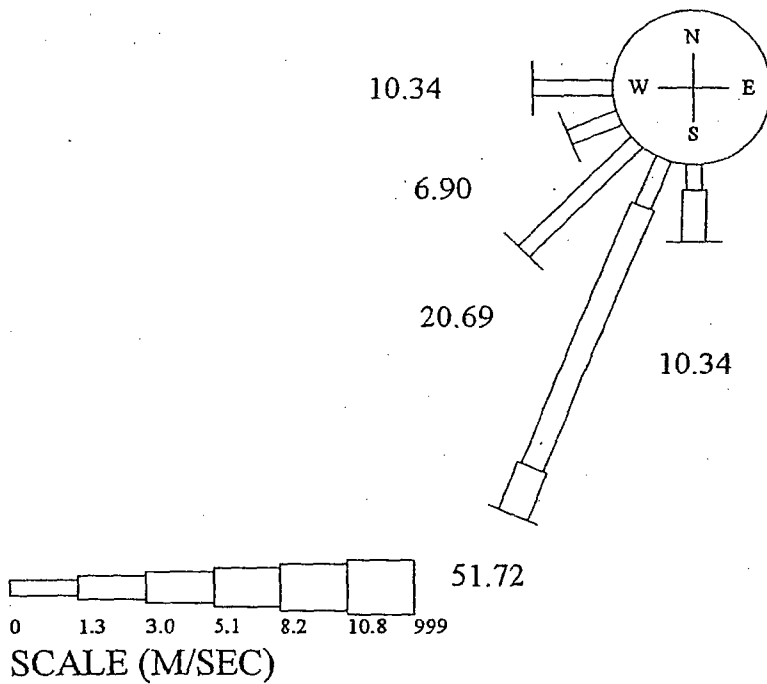
Sandra D. Smith (Sam)

APPENDIX D

Windroses

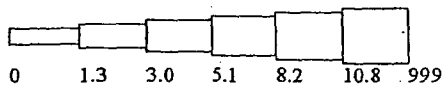
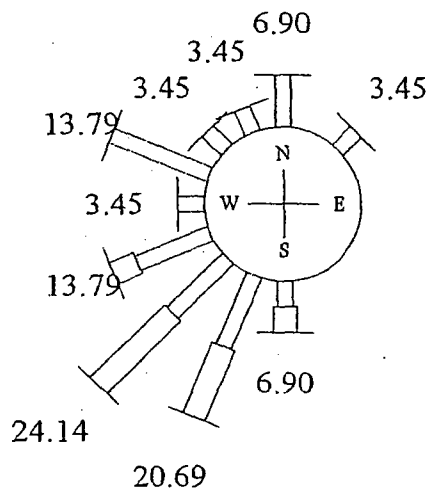
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

Aberdeen Proving Grounds Test Burn
Wind Rose Generated From H-Field Meteorological Data
12/3/99 14:00 - 21:00



| WIND SPEED (M/SEC) PERCENT OCCURRENCE | | | | | | | WIND SPEED (M/SEC) PERCENT OCCURRENCE | | | | | | |
|---------------------------------------|-------|---------|---------|---------|----------|-------|---------------------------------------|-------|---------|---------|---------|----------|-------|
| | 0-1.3 | 1.3-3.0 | 3.0-5.1 | 5.1-8.2 | 8.2-10.8 | >10.8 | | 0-1.3 | 1.3-3.0 | 3.0-5.1 | 5.1-8.2 | 8.2-10.8 | >10.8 |
| N | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | S | 3.45 | 6.90 | 0.00 | 0.00 | 0.00 | 0.00 |
| NNE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | SSW | 6.90 | 37.93 | 6.90 | 0.00 | 0.00 | 0.00 |
| NE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | SW | 20.69 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| ENE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | WSW | 6.90 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| E | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | W | 10.34 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| ESE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | WNW | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| SE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | NW | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| SSE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | NNW | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |

Aberdeen Proving Grounds Test Burn
Wind Rose Generated From Poverty Island Meteorological Data
12/3/99 14:00 - 21:00



SCALE (M/SEC)

| WIND SPEED (M/SEC) PERCENT OCCURRENCE | | | | | | | WIND SPEED (M/SEC) PERCENT OCCURRENCE | | | | | | |
|---------------------------------------|-------|---------|---------|---------|----------|-------|---------------------------------------|-------|---------|---------|---------|----------|-------|
| | 0-1.3 | 1.3-3.0 | 3.0-5.1 | 5.1-8.2 | 8.2-10.8 | >10.8 | | 0-1.3 | 1.3-3.0 | 3.0-5.1 | 5.1-8.2 | 8.2-10.8 | >10.8 |
| N | 6.90 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | S | 3.45 | 3.45 | 0.00 | 0.00 | 0.00 | 0.00 |
| NNE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | SSW | 10.34 | 10.34 | 0.00 | 0.00 | 0.00 | 0.00 |
| NE | 3.45 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | SW | 10.34 | 13.79 | 0.00 | 0.00 | 0.00 | 0.00 |
| ENE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | WSW | 10.34 | 3.45 | 0.00 | 0.00 | 0.00 | 0.00 |
| E | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | W | 3.45 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| ESE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | WNW | 13.79 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| SE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | NW | 3.45 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| SSE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | NNW | 3.45 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |

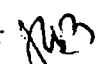
APPENDIX A-2

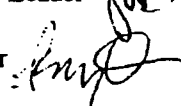
O-FIELD TRIP REPORT – BURN 2 (JULY 2000)

LOCKHEED MARTIN 

DATE: July 5, 2000

TO: David Mickunas, U.S. EPA/ERTC Work Assignment Manager

THROUGH: Jeff Bradstreet, REAC Air Group Leader 

FROM: Amy DuBois, REAC Task Leader 

SUBJECT: AIR MONITORING AND SAMPLING AT THE AIR MONITORING SAMPLING, ANALYSIS, AND MODELING SUPPORT, AND UNDERWATER SURVEY ACTIVITIES SITE, ABERDEEN PROVING GROUND, ABERDEEN, MD. WORK ASSIGNMENT #0-110 - TRIP REPORT - O-FIELD - BURN 2

BACKGROUND

The United States Environmental Protection Agency/Environmental Response Team Center (U.S. EPA/ERTC) issued Work Assignment Number 0-110 to Lockheed Martin under the Response, Engineering, and Analytical Contract (REAC) to provide air monitoring and air sampling during two controlled burns in the Edgewood Area of Aberdeen Proving Ground (APG). One burn was to be conducted at O-Field and one at J-Field. After problems igniting the marsh area during the O-Field burn, a second controlled burn was scheduled at O-Field.

Ordnance firing, ongoing test activities, and lightning strikes occasionally cause accidental fires in the test range areas at APG. Because of APG's long history of weapons testing and disposal practices, there is concern that contaminants have accumulated in the surface soils and vegetation at these locations and could be transported in the smoke plumes produced by such fires, posing a health risk to exposed individuals on and off the installation.

The scope of work for this work assignment included air sampling for dioxins, metals, polynuclear aromatic hydrocarbons (PAHs), inorganic acids, volatile organic compounds (VOCs) and chemical warfare agents (CWAs). Particulate monitoring was conducted utilizing an MIE DataRAM at each location.

OBSERVATIONS AND ACTIVITIES

REAC personnel mobilized to APG on December 17, 1999. Air sampling and monitoring were conducted at 5 downwind and 2 upwind locations (Figure 1).

VOC sampling and analysis was conducted following EPA Method TO-14A: *Determination of Volatile Organic Compounds in Ambient Air Using SUMMA Passivated Canister Sampling and Gas Chromatographic Mass Spectrometric (GC/MS) Analysis*. A sampling orifice was connected to each SUMMA canister to control the flow at 15 cubic centimeters per minute (cc/min). A solenoid valve was then connected to the SUMMA orifice. A battery operated timer was attached to each solenoid valve to trigger the solenoid at the anticipated start time for the burn.

PAH sampling and analysis was conducted following National Institute for Occupational Safety and Health (NIOSH) Method # 5515: *Polynuclear Aromatic Hydrocarbons*. Samples were collected utilizing a personal sampling pump

(SKC) to draw a measured volume of air (2 Liters per minute (L/min)) through a sampling train containing a teflon prefilter cassette and an XAD-2 sorbent tube. The pumps were programmed for a delayed start with a 4-hour sampling period.

Sampling and analysis for inorganic acids was conducted following NIOSH Method # 7903: *Acids, Inorganic*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (250 cc/min) through a sampling train containing a silica gel sorbent tube. The pumps were programmed for a delayed start with a 4-hour sampling period.

Sampling and analysis for dioxins was conducted following modified U.S. EPA Method TO9A, *Determination of Polychlorinated, Polybrominated and Brominated/Chlorinated Dibenzo-p-Dioxins and Dibenzofurans in Ambient Air*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (3 L/min) through a sampling train containing a polyurethane foam (PUF) plug and quartz filter. The pumps were programmed for a delayed start with a 4-hour sampling period. PUF glassware, plugs, and quartz filters were cleaned and certified by Southwest Research Institute in San Antonio, Texas prior to use.

Sampling and analysis for metals was conducted following modified NIOSH Method # 7300: *Elements (ICP)*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (3 L/min) through a sampling train containing a mixed cellulose ester filter cassette. The pumps were programmed for a delayed start with a 4-hour sampling period.

Samples were collected for CWAs utilizing a personal sampling pump (SKC) to draw a measured volume of air (100 cc/min) through a sampling train containing two Depot Area Air Monitoring System (DAAMS) sorbent tubes in a dual-sampling manifold. The pumps were programmed for a delayed start with a 4-hour sampling period. Tubes and analysis were provided by Soldiers Biological and Chemical Command (SBC COM).

Air monitoring for total particulates was performed utilizing an MIE DataRAM portable real-time aerosol monitor. Concentration data was logged every 10 seconds for the duration of the burn.

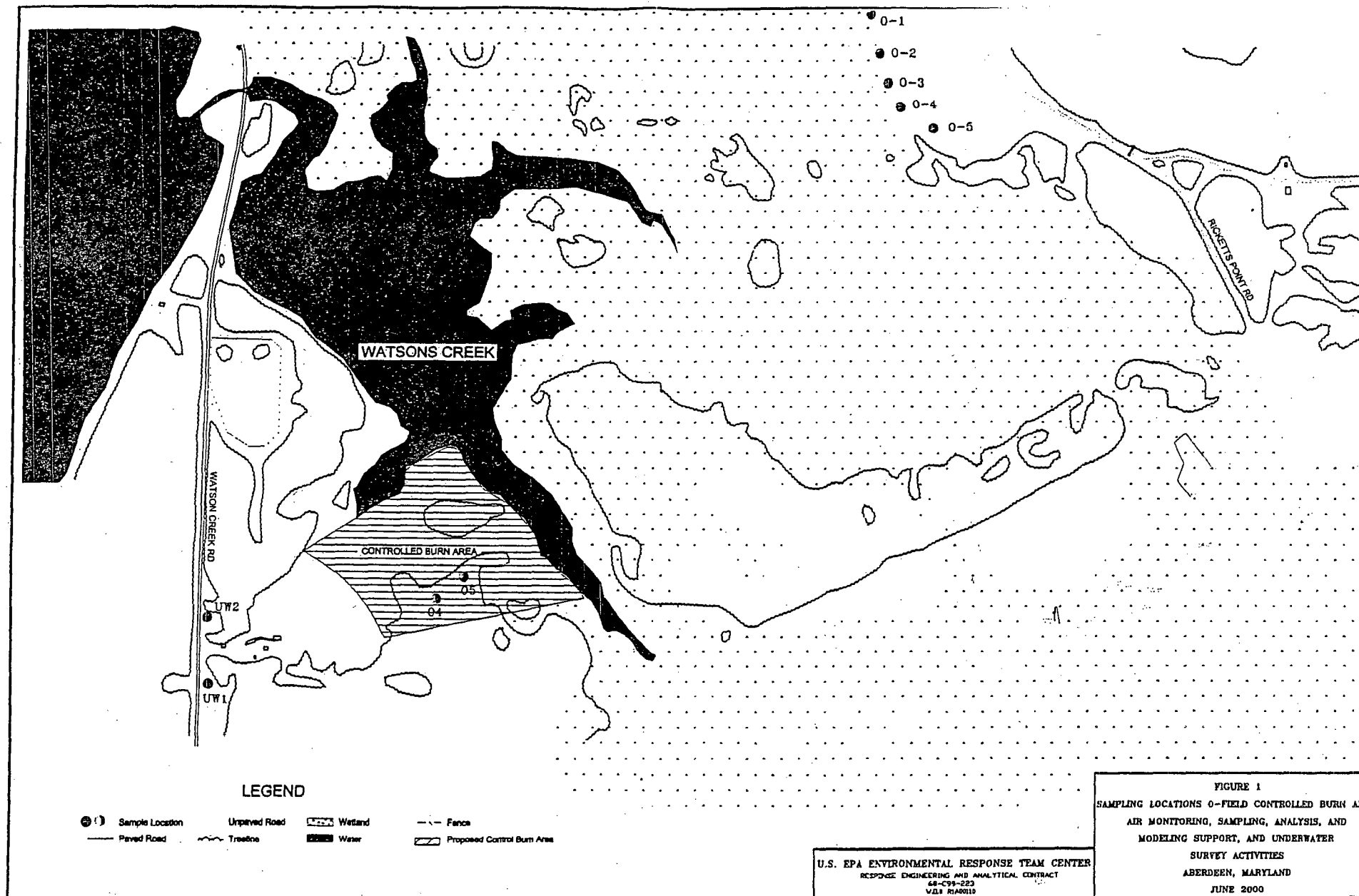
APG personnel positioned bridge sanctions at three downwind locations in Watsons Creek, prior to REAC's mobilization to the site. The two other downwind locations were positioned in trees along the edge of the marsh. REAC personnel set the samplers on the bridge sanctions and hoisted them into the trees with all timers set for a delayed start at 1345. When all personnel were out of the area, the APG Fire Department initiated the burn. In an attempt to propagate the burn through the marsh, approximately 8 to 10 gallons of kerosene were sprayed on the marsh vegetation. The fire still did not spread through the marsh and burned itself out after approximately 30 minutes.

RESULTS

Due to the short duration of the burn the decision was made between APG's Directorate of Safety, Health, and the Environment (DSHE) and the U.S. EPA/ERTC not to analyze the samples.

FUTURE ACTIVITIES

There are no future sampling activities planned for O-Field at this time.



APPENDIX A-3

J-FIELD TRIP REPORT
(JULY 2000)

LOCKHEED MARTIN

DATE: July 5, 2000

TO: David Mickunas, U.S. EPA/ERTC Work Assignment Manager

THROUGH: Jeff Bradstreet, REAC Air Group Leader

FROM: Amy DuBois, REAC Task Leader

SUBJECT: AIR MONITORING AND SAMPLING AT THE AIR MONITORING SAMPLING, ANALYSIS, AND MODELING SUPPORT, AND UNDERWATER SURVEY ACTIVITIES SITE, ABERDEEN PROVING GROUND, ABERDEEN, MD, WORK ASSIGNMENT #0-110 - TRIP REPORT - J-FIELD

BACKGROUND

The United States Environmental Protection Agency/Environmental Response Team Center (U.S. EPA/ERTC) issued Work Assignment Number 0-110 to Lockheed Martin under the Response, Engineering, and Analytical Contract (REAC) to provide air monitoring and air sampling during two controlled burns in the Edgewood Area of Aberdeen Proving Ground (APG). One burn was to be conducted at O-Field and one at J-Field.

Ordnance firing, ongoing test activities, and lightning strikes occasionally cause accidental fires in the test range areas at APG. Because of APG's long history of weapons testing and disposal practices, there is concern that contaminants have accumulated in the surface soils and vegetation at these locations and could be transported in the smoke plumes produced by such fires, posing a health risk to exposed individuals on and off the installation.

The scope of work for this work assignment included air sampling for dioxins, metals, polynuclear aromatic hydrocarbons (PAHs), inorganic acids, volatile organic compounds (VOCs) and chemical warfare agents (CWAs). Particulate monitoring was conducted utilizing an MIE DataRAM at five locations.

OBSERVATIONS AND ACTIVITIES

REAC personnel mobilized to APG on April 6, 2000. Air sampling and monitoring was conducted at 5 downwind and 2 upwind locations (see Figure 1).

VOC sampling and analysis was conducted following EPA Method TO-14A: *Determination of Volatile Organic Compounds in Ambient Air Using SUMMA Passivated Canister Sampling and Gas Chromatographic Mass Spectrometric (GC/MS) Analysis*. A sampling orifice was connected to each SUMMA canister to control the flow at 15 cubic centimeters per minute (cc/min). A solenoid valve was then connected to the SUMMA orifice. A battery operated timer was attached to each solenoid valve to trigger the solenoid at the anticipated start time for the burn.

PAH sampling and analysis was conducted following National Institute for Occupational Safety and Health (NIOSH) Method # 5515: *Polynuclear Aromatic Hydrocarbons*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (2 Liters per minute (L/min)) through a sampling train containing a teflon

prefilter cassette and an XAD-2 sorbent tube. The pumps were programmed for a delayed start with a 3-hour sampling period.

Sampling and analysis for inorganic acids was conducted following NIOSH Method # 7903: *Acids, Inorganic*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (250 cc/min) through a sampling train containing a silica gel sorbent tube. The pumps were programmed for a delayed start with a 3-hour sampling period.

Sampling and analysis for dioxins was conducted following modified U.S. EPA Method TO9A, *Determination of Polychlorinated, Polybrominated and Brominated/Chlorinated Dibenzo-p-Dioxins and Dibenzofurans in Ambient Air*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (3 L/min) through a sampling train containing a polyurethane foam (PUF) plug and quartz filter. The pumps were programmed for a delayed start with a 3-hour sampling period. PUF glassware, plugs, and quartz filters were cleaned and certified by Southwest Research Institute in San Antonio, Texas prior to use.

Sampling and analysis for metals was conducted following modified NIOSH Method # 7300: *Elements (ICP)*. Samples were collected utilizing a personal sampling pump (SKC) to draw a measured volume of air (3 L/min) through a sampling train containing a mixed cellulose ester filter cassette. The pumps were programmed for a delayed start with a 3-hour sampling period.

Samples were collected for CWAs utilizing a personal sampling pump (SKC) to draw a measured volume of air (100 cc/min) through a sampling train containing two Depot Area Air Monitoring System (DAAMS) sorbent tubes in a dual-sampling manifold. The CWAs analyzed for included: Sarin (GB), Soman (GD), Mustard (HD), and VX. The pumps were programmed for a delayed start with a 3-hour sampling period. Tubes and analysis were provided by Soldiers Biological and Chemical Command (SBC COM).

Air monitoring for total particulates was performed utilizing an MIE DataRAM portable real-time aerosol monitor. Concentration data was logged every 10 seconds for the duration of the burn. DataRAMs were positioned at locations DW1, DW2, DW3, DW4 and UW2.

The sampling devices were suspended 15 feet above the ground from trees and/or support poles, this positioned the samplers in the plume but out of the potential burn path of the fire. The collection of sampling devices was hoisted off the ground after setting the timers on the individual pumps and SUMMA canisters. The timers for the pumps controlled the start time and duration of the sampling period. The SUMMA timers only controlled the start of the sampling period. When all personnel were out of the area, the APG Fire Department initiated the burn.

RESULTS

VOCs: A summary of VOCs sampling results can be found in Table 1. Benzene and toluene were the only compounds detected above their quantitation limit in any of the samples. Benzene was detected at locations DW3 and DW4, and toluene was detected at locations DW3 and DW5. For complete analytical results for VOCs, see the Analytical Report in Appendix A.

PAHs: No PAHs were detected above the method detection limit in any of the samples.

Inorganic Acids: A summary of inorganic acids sampling results can be found in Table 2. Hydrochloric acid (HCl) was detected in samples DW1, DW3, DW5, and the Lot Blank. The detected HCl concentration ranged from 0.0176 to 0.1230 parts per million by volume (ppmv). The Lot Blank contained 0.0031 milligrams (mg) HCl. Hydrofluoric acid (HF) was detected in samples DW1 through DW5 at concentrations ranging from 0.0292 to 0.1030 ppmv. For complete analytical results for inorganic acids see the Analytical Report in Appendix B.

Dioxins/Furans: A summary of dioxins/furans results can be found in Table 3. Dioxins/furans were detected at six of the seven sampling locations. The OCDD results for samples DW3, DW5, UW1, and UW2 should be considered not detected because the concentration in the sample was less than five times that detected in the trip blank. The total dioxins/furans detected at each location after adjusting the OCDD results are as follows: DW1(not detected), DW2(1.920 picograms per cubic meter (pg/m³)), DW3(not detected), DW4(1.003 pg/m³), DW5(0.126 pg/m³), UW1(not detected), UW2(not detected), Trip Blank(0.0122 pg), Field Blank(not detected), and Lot Blank(0.070 pg). For complete analytical results for dioxins/furans, see the Analytical Report in Appendix B.

Metals: A summary of metals results are shown in Table 4. Aluminum was detected in samples DW1, DW2, DW3, DW5, UW1, and UW2 at concentrations ranging from 1.9 to 31.0 micrograms per cubic meter (µg/m³). Copper was detected at DW2 at 0.2 µg/m³. Lead was detected at DW2 at 0.3 µg/m³. Magnesium was detected at locations DW1, DW2, DW3, and UW2 at concentrations ranging from 2.0 to 30.0 µg/m³. Manganese was detected at locations DW2, DW3, and UW2 at concentrations ranging from 0.8 to 1.0 µg/m³. Phosphorous was detected at locations DW1, DW2, DW3, DW4, UW1, and UW2 at concentrations ranging from 0.8 to 2.2 µg/m³. Titanium was detected at locations UW1 and UW2 at 1.0 and 2.1 µg/m³, respectively. Sodium was detected in the method blank and should be regarded as not detected in all of the samples. Calcium, chromium, and zinc were detected in the lot blank and should be regarded as not detected in the rest of the samples because the concentrations were less than five times that detected in the lot blank. Iron was also detected in the lot blank. Locations DW3 and UW2 both had iron concentrations greater than 5 times the lot blank, iron should be regarded as not detected in the rest of the samples. Nickel was detected in the trip blank. Location UW1 had a nickel concentration greater than 5 times the trip blank, nickel should be regarded as not detected in the rest of the samples. For complete analytical results for metals, see the Analytical Report in Appendix B.

CWAs: No chemical warfare agents were detected in any of the samples. CWA results are provided by SBC COM, see Appendix C.

Particulates: Particulates results are shown in Figures 2 through 5. The DataRAM at location UW2 did not log data. The overall maximum concentration of 407,574.9 micrograms per cubic meter (µg/m³) was detected at location DW2 at 17:16 eastern standard time.

Meteorological data: Windroses representing local wind speed and wind direction during the burn period are provided in Appendix D. The data was collected at H-Field using a 10-meter tower, and at Poverty Island using a 5-meter tower. Winds were predominantly out of the west northwest. Times shown are in eastern standard time.

Analysis for VOCs and PAHs were provided by REAC, Edison, NJ. Analysis for dioxins/furans, inorganic acids, and metals were provided by Southwest Research Institute, San Antonio, TX. Analysis for CWAs was provided by SBC COM, APG, MD.

FUTURE ACTIVITIES

There are no future sampling activities planned at this time.

Table 1
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
Summary of VOCs Sampling Results - J-Field Controlled Burn - April 6, 2000

| Sample Number | 17747 | 17740 | 17741 | 17742 | 17743 | 17744 | 17744 Dup | 17745 | 17746 |
|-----------------|------------|-------|-------|-------|-------|-------|-----------|-------|-------|
| Sample Location | trip blank | DW3 | DW2 | DW1 | DW4 | DW5 | DW5 | UW1 | UW2 |
| concentration | ppbv | ppbv | ppbv | ppbv | ppbv | ppbv | ppbv | ppbv | ppbv |
| Chloromethane | U | 2 J | 9 J | 2 J | 3 J | 1 J | 1 J | U | 1 J |
| Benzene | U | 7 | U | 3 J | 5 | 3 J | 3 J | U | U |
| Toluene | U | 4 | U | 1 J | 3 J | 6 | 6 | U | U |
| Ethylbenzene | U | U | U | U | U | 1 J | 1 J | U | U |
| m & p-Xylenes | U | 1 J | U | U | U | 4 J | 4 J | U | U |
| o-Xylene | U | U | U | U | U | 1 J | 1 J | U | U |

VOCs - Volatile organic compounds
ppbv - parts per billion by volume
J - Below 1.00 nL Quantitation Limit
U - Not Detected

Table 2
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
Summary of Inorganic Acids Sampling Results - J-Field Controlled Burn - April 6, 2000

| Sample Number | 17734 | 17732 | 17733 | 17700 | 17701 | 17702 | 17703 | 17704 | 17705 | 17706 |
|----------------------------|-----------|-------------|------------|--------|--------|--------|--------|--------|--------|--------|
| Sample Location | Lot Blank | Field Blank | Trip Blank | DW3 | DW2 | DW1 | DW4 | DW5 | UW1 | UW2 |
| concentration | mg | mg | mg | ppmv | ppmv | ppmv | ppmv | ppmv | ppmv | ppmv |
| Hydrobromic Acid | U | U | U | U | U | U | U | U | U | U |
| Hydrochloric Acid | 0.0031 | U | U | 0.1230 | U | 0.0387 | U | 0.0188 | U | 0.0176 |
| Hydrofluoric Acid | U | U | U | 0.0724 | 0.1030 | 0.0292 | 0.0439 | 0.0389 | U | U |
| Nitric Acid | U | U | U | U | U | U | U | U | U | U |
| Phosphoric Acid | U | U | U | U | U | U | U | U | U | U |
| Sulfuric Acid ¹ | 0.0050 | U | 0.0013 | 0.0225 | 0.0262 | 0.0182 | 0.0161 | 0.0404 | 0.0175 | 0.0217 |

mg - total milligrams

ppmv - parts per million by volume

U - Not detected

¹ Due to the sulfuric acid concentration detected in the Trip Blank, the results for samples 17700 through 17706 are considered not detected.

Table 3
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
Summary of Dioxins/Furans Sampling Results - J-Field Controlled Burn - April 6, 2000

| Sample Number Sample Location | 17677 Trip Blank | 17678 Field Blank | 17679 Lot Blank | 17670 DW3 | 17671 DW2 | 17672 DW1 | 17673 DW4 | 17674 DW5 | 17675 UW1 | 17676 UW2 |
|-------------------------------------|---------------------|----------------------|--------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| Adjusted concentration ¹ | pg | pg | pg | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ | pg/m ³ |
| 2,3,7,8-TCDD | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,7,8-PeCDD | U | U | U | U | 1.88 | U | U | U | U | U |
| 1,2,3,4,7,8-HxCDD | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,6,7,8-HxCDD | U | U | U | U | U | U | 0.42 | U | U | U |
| 1,2,3,7,8,9-HxCDD | U | U | 0.066 | U | U | U | U | U | U | U |
| 1,2,3,4,6,7,8-HpCDD | U | U | U | U | U | U | 0.277 | 0.126 | U | U |
| OCDD ² | 0.0122 | U | U | 0.035 | U | U | 0.0738 | 0.094 | 0.033 | 0.03 |
| 2,3,7,8-TCDF | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,7,8-PeCDF | U | U | U | U | U | U | U | U | U | U |
| 2,3,4,7,8-PeCDF | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,4,7,8-HxCDF | U | U | U | U | U | U | 0.232 | U | U | U |
| 1,2,3,6,7,8-HxCDF | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,7,8,9-HxCDF | U | U | U | U | U | U | U | U | U | U |
| 2,3,4,6,7,8-HxCDF | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,4,6,7,8-HpCDF | U | U | U | U | U | U | U | U | U | U |
| 1,2,3,4,7,8,9-HpCDF | U | U | U | U | 0.0396 | U | U | U | U | U |
| OCDF ³ | U | U | 0.00442 | U | U | U | U | U | U | U |
| Total | 0.0122 | U | 0.07042 | 0.035 | 1.9196 | U | 1.0028 | 0.22 | 0.033 | 0.03 |

pg - picograms

pg/m³ - picograms per cubic meter

¹ Adjusted concentration - detected concentration multiplied by the toxicity equivalency factor (TEF) for each compound.

² The OCDD results for samples 17670, 17674, 17675, and 17676 are considered not detected because the concentration in the sample was less than five times that found in the trip blank.

³ The OCDF result for sample 17679 is considered estimated because the method blank contained 13 pg OCDF.

Table 4
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities
Summary of Metals Sampling Results - J-Field Controlled Burn - April 6, 2000

| Sample Number | 17687 | 17688 | 17689 | 17680 | 17681 | 17682 | 17683 | 17684 | 17685 | 17686 |
|-----------------------|-------------|------------|-----------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| Sample Location | Field Blank | Trip Blank | Lot Blank | DW3 | DW2 | DW1 | DW4 | DW5 | UW1 | UW2 |
| concentration | ug/filter | ug/filter | ug/filter | ug/m ³ | ug/m ³ | ug/m ³ | ug/m ³ | ug/m ³ | ug/m ³ | ug/m ³ |
| Aluminum | U | U | U | 3.8 | 4.0 | 1.9 | U | 1.9 | 4.0 | 31.0 |
| Calcium ² | 6.6 | 7.6 | 6.2 | 34.0 | 40.0 | 15.0 | 11.0 | 13.0 | 16.0 | 22.0 |
| Chromium ² | 0.6 | 0.9 | 0.5 | 1.0 | 1.1 | 1.0 | 1.1 | 1.0 | 1.1 | 1.2 |
| Copper | U | U | U | U | 0.2 | U | U | U | U | U |
| Iron ³ | 4.0 | 2.7 | 1.6 | 25.0 | 9.2 | 3.6 | 2.6 | 2.0 | 5.7 | 58.0 |
| Lead | U | U | U | U | 0.3 | U | U | U | U | U |
| Magnesium | U | U | U | 4.2 | 5.6 | 2.0 | U | U | U | 30.0 |
| Manganese | U | U | U | 1.0 | 1.0 | U | U | U | U | 0.8 |
| Nickel ⁴ | 0.4 | 0.2 | U | U | 0.8 | U | 0.4 | U | 6.9 | 0.6 |
| Phosphorous | U | U | U | 1.3 | 2.2 | 1.3 | 0.8 | U | 1.0 | 2.1 |
| Sodium ¹ | 9.8 | 9.0 | 7.3 | 17.0 | 18.0 | 17.0 | 16.0 | 11.0 | 14.0 | 18.0 |
| Titanium | U | U | U | U | U | U | U | U | 0.6 | 0.7 |
| Zinc ² | 2.1 | 2.4 | 1.1 | 3.4 | 2.0 | 0.9 | 1.1 | 0.8 | 1.7 | 1.2 |

ug/filter - micrograms per filter

ug/m³ - micrograms per cubic meter

U - not detected

¹ The method blank contained 11.81 ug/filter sodium, the sodium results for all samples should be considered not detected.

² The Calcium, Chromium and Zinc results for samples 17680 through 17688 are considered not detected because the concentration in the sample is less than 5 times that of the lot blank.

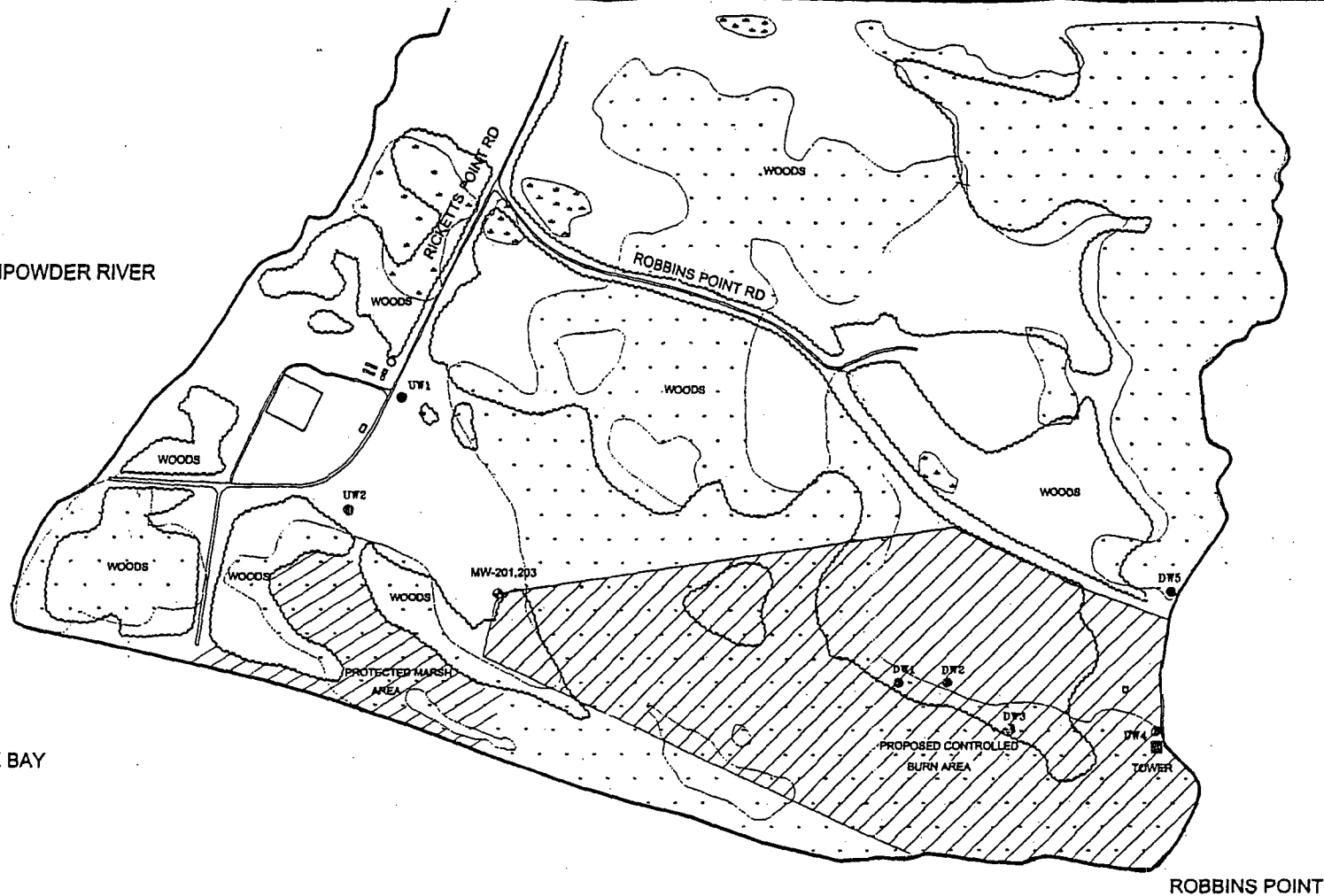
³ The Iron results for samples 17681 through 17685, 17687 and 17688 are considered not detected because the concentration in the sample is less than 5 times that of the lot blank.

⁴ The Nickel results for samples 17681, 17683, 17686, and 17687 are considered not detected because the concentration in the sample is less than 5 times that of the trip blank.



GUNPOWDER RIVER

CHESAPEAKE BAY



LEGEND

- | | | | |
|-----|------------------|-------|----------------------------|
| ● ● | Sample Locations | — | Unpaved Road |
| ⊙ | Monitoring Well | ▨ | Proposed Control Burn Area |
| ▧ | Water | ▤ | Wetland |
| | | - - - | Treeline |

U.S. EPA ENVIRONMENTAL RESPONSE TEAM CENTER
RESPONSE ENGINEERING AND ANALYTICAL CONTRACT
68-C99-223
V.D. 9/12/00

FIGURE 1
SAMPLING LOCATIONS J-FIELD CONTROLLED BURN AREA
AIR MONITORING, SAMPLING, ANALYSIS, AND
MODELING SUPPORT, AND UNDERWATER
SURVEY ACTIVITIES
ABERDEEN, MARYLAND
JUNE 2000

Figure 2
Aberdeen Proving Ground
J-Field Burn Data - Particulates
April 6, 2000
Location DW1 - Peak Burn Period

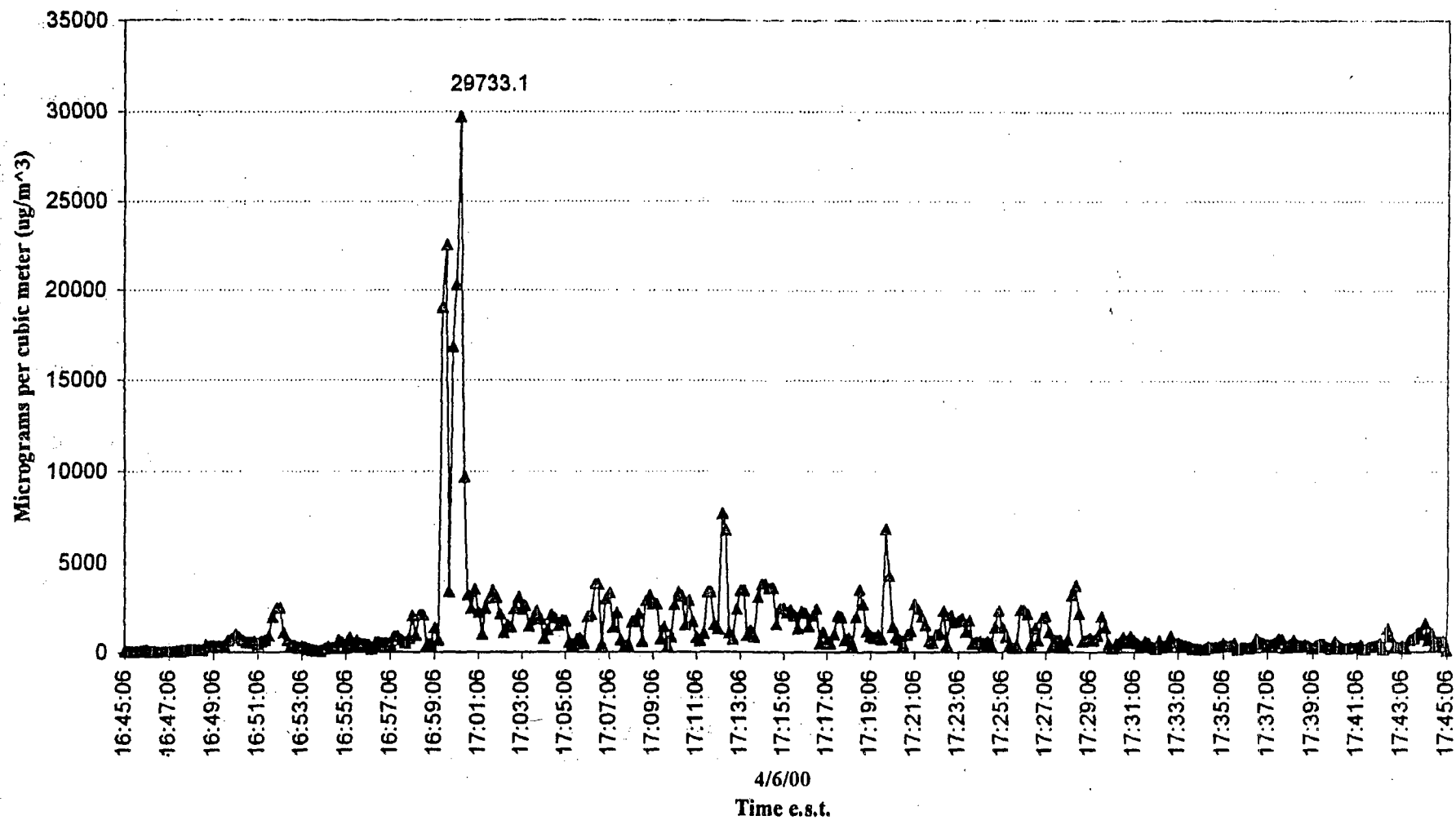


Figure 3
Aberdeen Proving Ground
J-Field Burn Data - Particulates
April 6, 2000
Location DW2 - Peak Burn Period

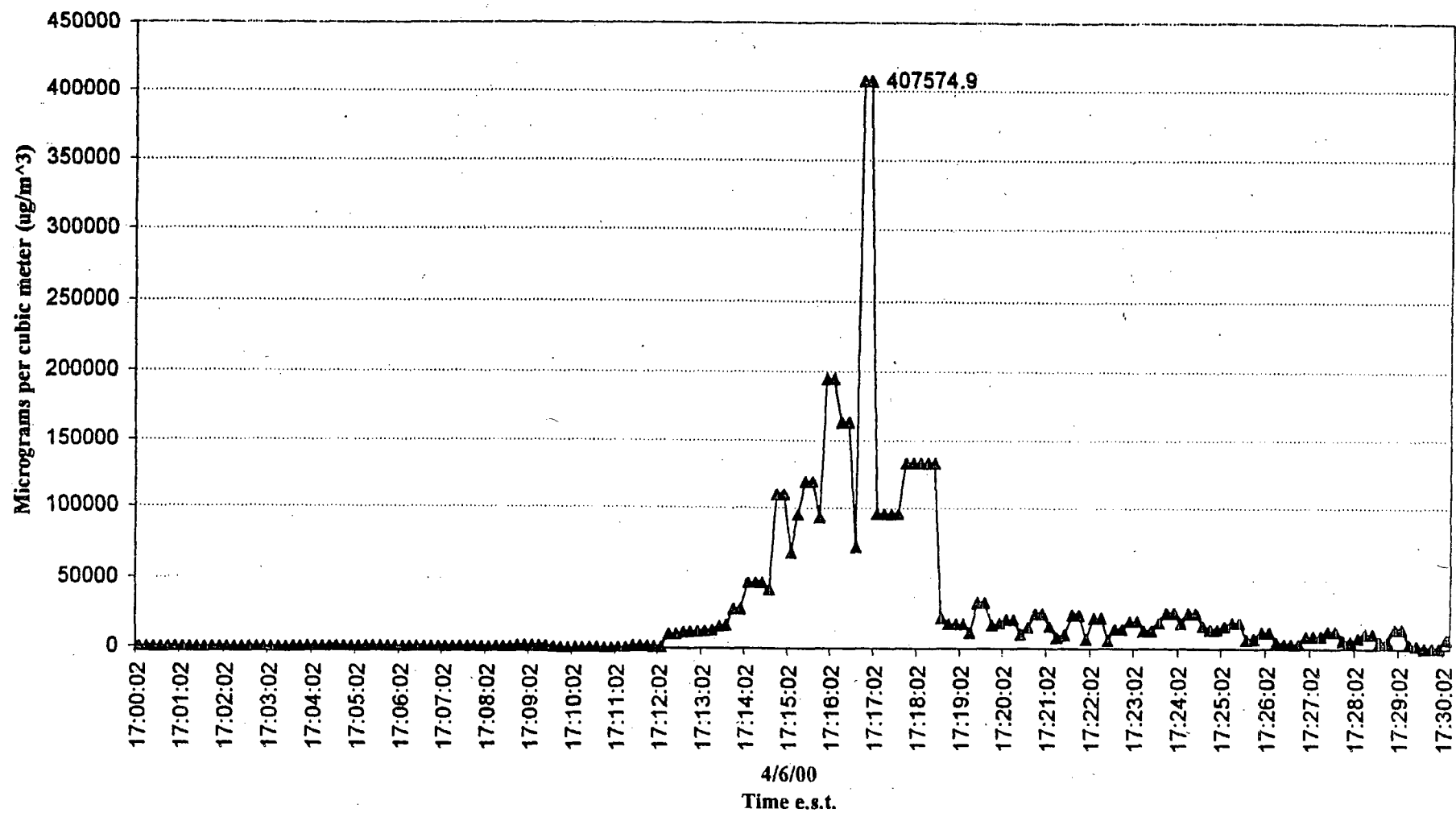


Figure 4
Aberdeen Proving Ground
J-Field Burn Data - Particulates
April 6, 2000
Location DW3 - Peak Burn Period

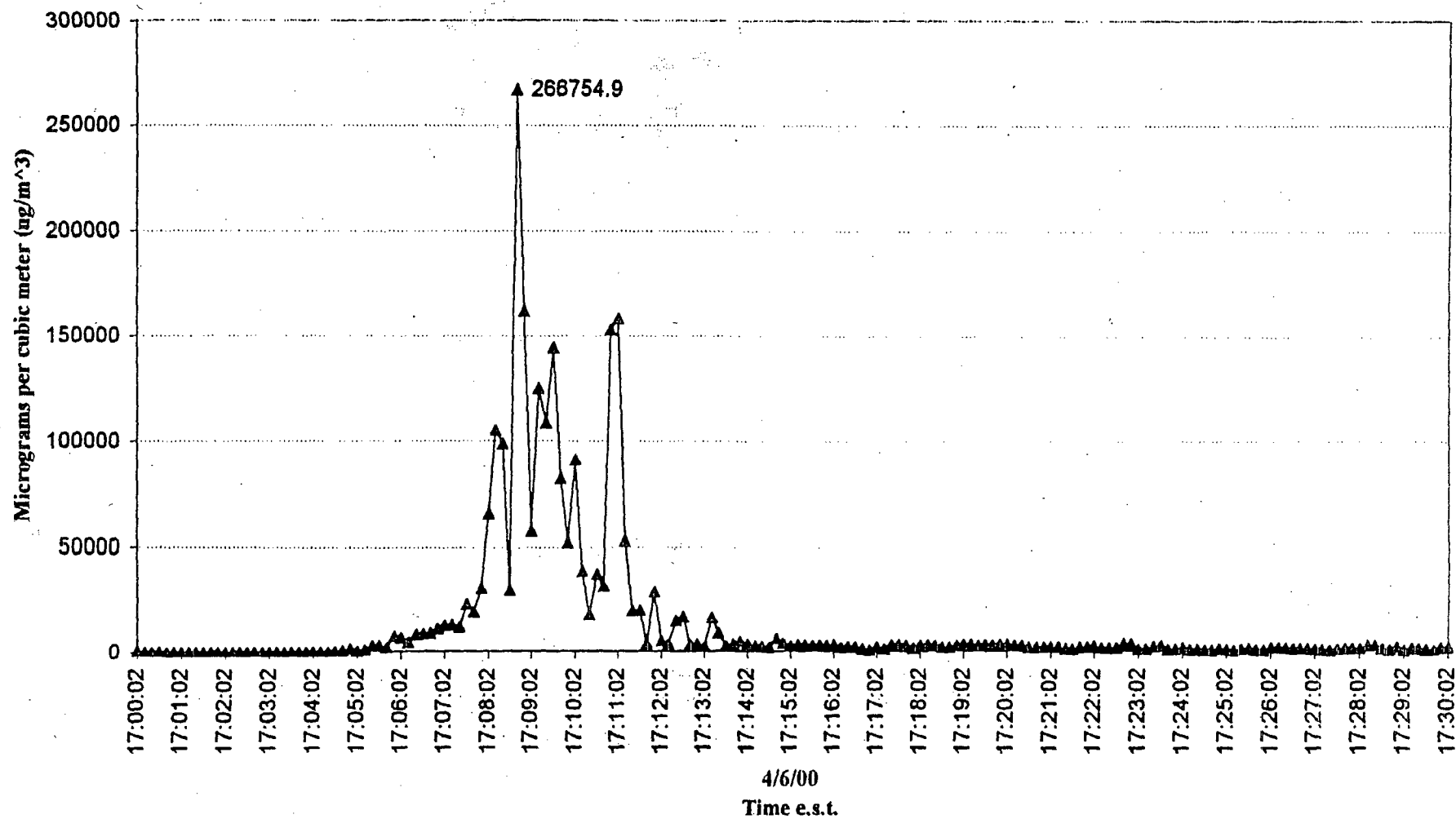
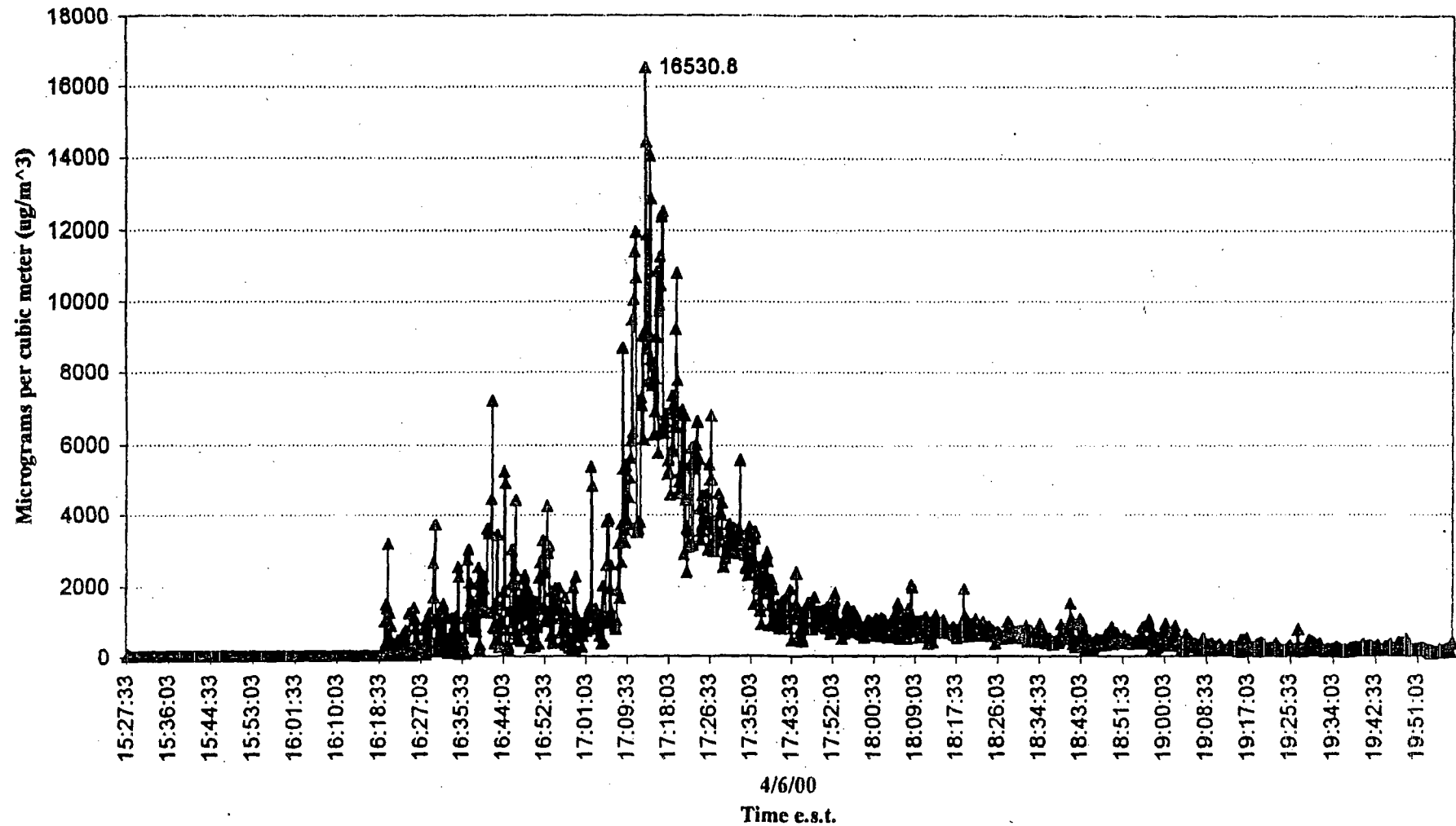


Figure 5
Aberdeen Proving Ground
J-Field Burn Data - Particulates
April 6, 2000
Location DW4



APPENDIX A
Analytical Report (VOCs)
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

ANALYTICAL REPORT

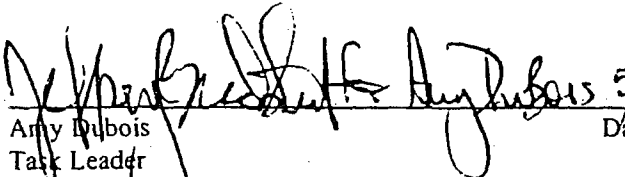
Prepared by
Lockheed Martin Technology Services Group

Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
Aberdeen Proving Ground, Aberdeen, MD


May 2000

EPA Work Assignment No. 0-110
Lockheed Martin Work Order No. R1A00110
EPA Contract No. 68-C99-223

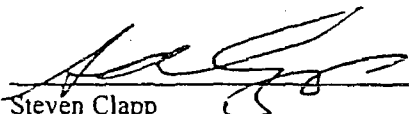
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1.0 INTRODUCTION

Summa canister samples were collected in support of the Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities work assignment at the Aberdeen Proving Ground, Aberdeen, MD on 06 April 2000. A total of seven (7) samples and a field blank were collected in 6-liter passivated Summa canisters. The samples were transported back to the Environmental Response Team Center (ERTC) facility in Edison, New Jersey. These samples were analyzed by the Response Engineering and Analytical Contract (REAC) using gas chromatography/mass spectrometry (GC/MS) on 10 and 11 April 2000.

2.0 GC/MS CANISTER PROCEDURES

2.1 Sample Pressurization

The Summa canisters used for sampling were cleaned by REAC using REAC Standard Operating Procedure (SOP) #1703 and were selected from clean batches certified by REAC. Before analysis, all canisters were pressurized. A pressurizing train was setup with a pressure gauge accurate to ± 0.1 pounds per square inch absolute (psia). The gauge and train were purged with nitrogen gas (Ultra High Pure grade) for 5 minutes. The train was then connected to the canister, an initial reading was taken. Nitrogen was added to all canister samples as followed:

| <u>Sample</u> | <u>Location</u> | <u>Initial Pressure (psia)</u> | <u>Final Pressure (psia)</u> |
|---------------|-----------------|------------------------------------|----------------------------------|
| 17740 | DW3 | 14.8 | 29.6 |
| 17741 | DW2 | 2.0 | 16.0 |
| 17742 | DW1 | 15.2 | 30.4 |
| 17743 | DW4 | 8.8 | 17.6 |
| 17744 | DW5 | 14.0 | 28.0 |
| 17745 | UW1 | 8.4 | 16.8 |
| 17746 | UW2 | 8.4 | 16.8 |
| 17747 | Trip/Field | 0.3 | 20.0 |

2.2 Summa Canister Analysis

Samples were analyzed by cryogenic trapping of aliquots from Summa canisters via a canister using a Hewlett-Packard 5890 gas chromatography (GC) and 5971A mass selective detector (MSD) running ChemStation software. Table 1 lists cryogenic trap and GC/MS conditions.

All canisters were attached to the Summa canister autosampler. Sample analysis began by cooling the first cryotrap, module -1 (M-1), to -160 degree Celsius ($^{\circ}\text{C}$). Once M-1 was cooled, a specified aliquot of sample or standard was cryotrapped. This aliquot was transferred to a Tenax trap, M-2, to eliminate most of the water, and then cryofocused at a third trap, M-3, before injection by direct heating.

2.3 Calibration and Sample Spiking

Standard mixture containing twenty-five (25) compounds was provided in compressed gas cylinder No ALM009519 by Scott Specialty Gases, Inc. These standard concentrations are .97 to 1.05 parts per million in volume (ppmv) and are listed in Table 2. The standards were diluted to a nominal concentration of 20 parts per billion (ppbv) in a Silco canister. An initial calibration range was obtained by varying the volume of the nominal 20 ppbv standard from 50 to 1250 milliliters (mL), equivalent to 1 nanoliter (nL) to 25 nL. Daily standards were obtained by analyzing the 20 ppbv standard at 500 mL (equivalent to 10 nL).

Bromochloromethane (BCM) and p-bromofluorobenzene (BFB) were added to both samples and standards. Both standards were provided in compressed gas cylinder No. ALM046281 by Scott Specialty Gases. These standard concentrations were 1.06 ppmv. BCM was used as an internal standard and BFB was used as a surrogate standard. This standard was diluted from a nominal concentration of 1 ppmv to 100 ppbv in a Silco canister. An aliquot of 100 mL (equivalent to 10 nL) was added to all standards and samples. To validate the mass spectrometer tuning, an aliquot of 70 mL (equivalent to 50 nanograms of BFB) was analyzed alone. Standard cylinder I.D. numbers, concentrations, and their quantitation ions are listed in Table 2.

2.4 Compound Identification/Quantitation

Target Compounds in samples were identified and quantitated using ChemStation software. This software was used to tentatively identify and quantitate target compounds using reconstructed and extracted ion chromatogram which were matched with retention time windows. The report format includes the identified compound mass spectra (both raw and background subtracted), quantitation, and qualifier ion chromatogram.

Target compound results are originally reported in nL. The limit of quantitation (LOQ) for all the target compounds is estimated to be 1 nL, being the lowest volume of standard on the calibration curve. Any target compound detected at 4 times lower than the LOQ is not reported. The target compound results are calculated in ppbv using the following equation:

$$\text{Concentration (ppbv)} = \frac{\text{Quant Result (nL)} \times 1000}{\text{Undiluted Sample Volume (mL)}}$$

Non-target compounds were identified by a library search of all peaks in a chromatogram. The library search report prints out the sample spectrum along with the ten best library matches and the three best library match spectra. These matches were used along with mass spectral interpretation techniques to tentatively identify the unknowns. Concentrations were calculated based on the total ion response of bromochloromethane in the daily standard. All compounds appearing in the method blank as well as other background compounds commonly found in Summa canister GC/MS analyses (siloxanes, carbon dioxide, etc.) were deleted from the sample results to provide a true listing of the compounds in the samples.

2.5 QA/QC

The following QA/QC procedures were performed for this analysis:

- ▶ The HP 5971A was tuned daily for perfluorotributylamine (PFTBA) to meet abundance criteria for p-bromofluorobenzene as listed in EPA Method 624. Tuning results are included in the QA/QC data section (Appendix B). The tune was adjusted when necessary.
- ▶ An initial calibration by automated injection from a Silco canister standard at 20 ppbv was performed on 24 March 2000. All compounds met the acceptance criteria of having relative standard deviations (RSD) of less than 25%.
- ▶ Continuing calibrations were performed on 10 and 11 April 2000 to satisfy the 12 hour requirement. All compounds met the acceptance criteria of having relative percent difference (RPD) less than 25%, except chloroethane (43.8%) on 11 April 2000. This compound was not detected in the associated samples; the data are not affected.
- ▶ A surrogate standard of BFB was added to all standards and samples. Percent recoveries were calculated against the daily standards, and are listed in Table 3. Recoveries should be within 70% to 130% for BFB.

- ▶ Method blanks were analyzed after each continuing calibration to ensure that the system was clean.
- ▶ A replicate was analyzed on sample 17744 (DW5).
- ▶ A set of matrix spike and matrix spike duplicates (MS/MSD) was analyzed on sample 17746 (UW2) by spiking the samples with 500 mL of the 20 ppbv standard. There is no specific recovery range established according to SOP # 1705.

3.0 RESULTS

Summa canister target and non-target results are listed in Tables 3 and 4, respectively. The recoveries for the MS/MSD are presented in Table 5. All results are reported in ppbv for Summa canister samples and blanks. The chain-of-custody is in Appendix A. The Summa canister data are in Appendix B.

In Appendix B, the Analysis Log is followed by the calibration package for each day of analysis. The calibration package includes the daily analysis log, canister pressurization log, BFB tune, and initial or continuing calibration quant report. The quant report lists the retention time, quantitation ion, peak area, and concentration in nL. Concentrations listed on the quant reports are generated by using the average response factors of the initial calibration and the response factors of the continuing calibrations.

The following is a list of the QA/QC flags used in qualifying the results:

- A - Assumed volume for method blank.
- B - Concentration less than 3 times method blank value.
- C - Compound calibration relative standard deviation (RSD) >25% (concentrations calculated by average response factor only).
- E - Exceeds calibration range.
- J - Below 1.0 nL quantitation limit.
- U - Not detected.

4.0 DATA ASSESSMENT

A total of 7 samples and a field blank were collected on 4/6/00 on chain of custody numbers 03310 from the Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities in Aberdeen Proving Ground, Aberdeen, MD under U.S.EPA WA# 0-110. The samples were received on 4/10/00.

The data contained in this report has been validated to two significant figures. Any other interpretation of the data is the responsibility of the user.

The samples were treated with procedures consistent with those described in SOP # 1008.

The reported year on the raw data for the acquisition time is incorrectly reported as "100" and the reported year on the raw data for the quantitation time is incorrectly reported as "19100". This is due to a software problem related to the year 2000.

In the continuing calibration on 2/11/00 the percent difference for chloroethane (44%) exceeded the QC limits. This compound was not detected in the associated samples; the data are not affected.

TABLE 1 - GC/MS Instrument Conditions

A. Preconcentrator Conditions:

| | |
|---------------------------------|---------------|
| M-1 Cryotrap Temperature | : -160°C |
| Internal Standard Trap Time | : 1.0 minute |
| Sample flow | : 150 mL/min |
| M-1 Cryotrap Desorb Temperature | : 20°C |
| M-2 Cryotrap Temperature | : -10°C |
| Transfer (M-1 to M-2) Time | : 4.5 minutes |
| M-2 Cryotrap Desorb Temperature | : 240°C |
| M-3 Cryotrap Temperature | : -160°C |
| Transfer (M-2 to M-3) Time | : 3.5 minutes |
| Injection Time | : 2.0 minutes |

B. GC/MS Conditions, Sample Analysis:

| | |
|---------------------|-----------------|
| Initial Temperature | : 40.0°C |
| Initial Time | : 6.0 minutes |
| Ramp Rate | : 8.0°C/min |
| Final Temperature | : 220.0°C |
| Final Time | : 9.5 minutes |
| Run Time | : 35.03 minutes |
| Mass Scan Range: | : 35 to 250 AMU |

Column: 0.25 mm x 30 meter Restek RTx-VOA, 3.0 µm film thickness (Restek Corporation)

TABLE 2 - Air Toxic Standards (Concentrations and Quantitation Ions)

| <u>Compound</u> | <u>Cylinder</u> | <u>Conc. (ppmv)</u> | <u>Quant. Ion</u> |
|----------------------------|-----------------|---------------------|-------------------|
| chloromethane | ALM009519 | 0.98 | 50 |
| vinyl chloride | ALM009519 | 0.97 | 62 |
| chloroethane | ALM009519 | 1.00 | 64 |
| trichlorofluoromethane | ALM009519 | 1.04 | 101 |
| 1,1-dichloroethene | ALM009519 | 1.02 | 61 |
| dichloromethane | ALM009519 | 1.00 | 49 |
| trans-1,2-dichloroethene | ALM009519 | 1.00 | 61 |
| 1,1-dichloroethane | ALM009519 | 1.02 | 63 |
| trichloromethane | ALM009519 | 1.02 | 83 |
| 1,1,1-trichloroethane | ALM009519 | 1.01 | 97 |
| 1,2-dichloroethane | ALM009519 | 1.02 | 62 |
| benzene | ALM009519 | 1.00 | 78 |
| carbon tetrachloride | ALM009519 | 0.98 | 117 |
| trichloroethene | ALM009519 | 1.00 | 130 |
| dibromomethane | ALM009519 | 0.98 | 174 |
| bromodichloromethane | ALM009519 | 1.01 | 83 |
| toluene | ALM009519 | 1.01 | 91 |
| 1,1,2-trichloroethane | ALM009519 | 0.98 | 97 |
| tetrachloroethene | ALM009519 | 1.00 | 166 |
| ethylbenzene | ALM009519 | 1.01 | 91 |
| meta-xylene | ALM009519 | 1.02 | 91 |
| styrene | ALM009519 | 1.04 | 104 |
| ortho-xylene | ALM009519 | 1.04 | 91 |
| 1,1,2,2-tetrachloroethane | ALM009519 | 1.00 | 83 |
| 1,3,5-trimethylbenzene | ALM009519 | 1.05 | 120 |
| <u>Surrogate Standards</u> | | | |
| bromochloromethane | ALM046281 | 1.06 | 49 |
| p-bromofluorobenzene | ALM046281 | 1.06 | 95 |

Table 3 - Air Toxic Target Compound Results for Summa Canister Samples
WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
(concentrations in ppbv)

Page 1 of 3

| Sample Number | Method | 17747 | 17740 | 17741 | 17742 |
|--------------------------------|----------|------------|----------|----------|----------|
| Sample Location | Blank | Trip/Field | DW3 | DW2 | DW1 |
| Date Sampled | N/A | 04/06/00 | 04/06/00 | 04/06/00 | 04/06/00 |
| Date Analyzed | 04/10/00 | 04/10/00 | 04/10/00 | 04/10/00 | 04/10/00 |
| Data File | AGS003 | AGS004 | AGS005 | AGS006 | AGS007 |
| Chloromethane | 4 U | 4 U | 2 J | 9 J | 2 J |
| Vinyl Chloride | 4 U | 4 U | 4 U | 16 U | 4 U |
| Chloroethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Trichlorofluoromethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| 1,1-Dichloroethene | 4 U | 4 U | 4 U | 16 U | 4 U |
| Methylene Chloride | 4 U | 4 U | 4 U | 16 U | 4 U |
| trans-1,2-Dichloroethylene | 4 U | 4 U | 4 U | 16 U | 4 U |
| 1,1-Dichloroethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Trichloromethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| 1,1,1-Trichloroethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Carbon Tetrachloride | 4 U | 4 U | 4 U | 16 U | 4 U |
| 1,2-Dichloroethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Benzene | 4 U | 4 U | 7 | 16 U | 3 J |
| Trichloroethylene | 4 U | 4 U | 4 U | 16 U | 4 U |
| Bromodichloromethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Dibromomethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Toluene | 4 U | 4 U | 4 | 16 U | 1 J |
| 1,1,2-Trichloroethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| Tetrachloroethylene | 4 U | 4 U | 4 U | 16 U | 4 U |
| Ethylbenzene | 4 U | 4 U | 4 U | 16 U | 4 U |
| m & p-Xylenes | 4 U | 4 U | 1 J | 16 U | 4 U |
| o-Xylene | 4 U | 4 U | 4 U | 16 U | 4 U |
| Styrene | 4 U | 4 U | 1 J | 16 U | 4 U |
| 1,1,2,2-Tetrachloroethane | 4 U | 4 U | 4 U | 16 U | 4 U |
| 1,3,5-Trimethylbenzene | 4 U | 4 U | 4 U | 16 U | 4 U |
| p-Bromofluorobenzene (% Rec) | 103 | 100 | 107 | 101 | 109 |
| Pressurized Sample Volume (mL) | 250 | 250 | 500 | 500 | 500 |
| Initial Pressure (psia) | N/A | N/A | 14.8 | 2.0 | 15.2 |
| Final Pressure (psia) | N/A | N/A | 29.6 | 16.0 | 30.4 |
| Quantitation Limit (ppbv) | 4 | 4 | 4 | 16 | 4 |

A - Assumed volume for Blanks
B - <3 times Method Blank value
C - Compound Calibration >25% RSD
D - Compound Calibration Check >25% RPD
E - Concentration exceeded calibration limit (25nL)
J - Below 1.00 nL Quantitation Limit
U - Not Detected

0006

Table 3 - Air Toxic Target Compound Results for Summa Canister Samples
WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
(concentrations in ppbv)

Page 2 of 3

| Sample Number | 17743 | 17744 | 17744 Rep | 17745 | 17746 |
|--------------------------------|----------|----------|-----------|----------|----------|
| Sample Location | DW4 | DW5 | DW5 | UW1 | UW2 |
| Date Sampled | 04/06/00 | 04/06/00 | 04/06/00 | 04/06/00 | 04/06/00 |
| Date Analyzed | 04/10/00 | 04/10/00 | 04/11/00 | 04/10/00 | 04/10/00 |
| Data File | AGS008 | AGS009 | AGS016 | AGS010 | AGS012 |
| Chloromethane | 3 J | 1 J | 1 J | 4 U | 1 J |
| Vinyl Chloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| Chloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Trichlorofluoromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1-Dichloroethene | 4 U | 4 U | 4 U | 4 U | 4 U |
| Methylene Chloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| trans-1,2-Dichloroethylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1-Dichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Trichloromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1,1-Trichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Carbon Tetrachloride | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,2-Dichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Benzene | 5 | 3 J | 3 J | 4 U | 4 U |
| Trichloroethylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| Bromodichloromethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Dibromomethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Toluene | 3 J | 6 | 6 | 4 U | 4 U |
| 1,1,2-Trichloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| Tetrachloroethylene | 4 U | 4 U | 4 U | 4 U | 4 U |
| Ethylbenzene | 4 U | 1 J | 1 J | 4 U | 4 U |
| m & p-Xylenes | 4 U | 4 J | 4 J | 4 U | 4 U |
| o-Xylene | 4 U | 1 J | 1 J | 4 U | 4 U |
| Styrene | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,1,2,2-Tetrachloroethane | 4 U | 4 U | 4 U | 4 U | 4 U |
| 1,3,5-Trimethylbenzene | 4 U | 4 U | 4 U | 4 U | 4 U |
| p-Bromofluorobenzene (% Rec) | 110 | 112 | 112 | 106 | 105 |
| Pressurized Sample Volume (mL) | 500 | 500 | 500 | 500 | 500 |
| Initial Pressure (psia) | 8.8 | 14.0 | 14.0 | 8.4 | 8.4 |
| Final Pressure (psia) | 17.6 | 28.0 | 28.0 | 16.8 | 16.8 |
| Quantitation Limit (ppbv) | 4 | 4 | 4 | 4 | 4 |

A - Assumed volume for Blanks
 B - <3 times Method Blank value
 C - Compound Calibration >25% RSD
 D - Compound Calibration Check >25% RPD
 E - Concentration exceeded calibration limit (25nL)
 J - Below 1.00 nL Quantitation Limit
 U - Not Detected
 N/A - Not Applicable

Table 3 - Air Toxic Target Compound Results for Summa Canister Samples
WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
(concentrations in ppbv)

Page 3 of 3

Sample Number Method
Sample Location Blank
Date Sampled N/A
Date Analyzed 04/11/00
Data File AGS015

| | |
|--------------------------------|-----|
| Chloromethane | 4 U |
| Vinyl Chloride | 4 U |
| Chloroethane | 4 U |
| Trichlorofluoromethane | 4 U |
| 1,1-Dichloroethene | 4 U |
| Methylene Chloride | 4 U |
| trans-1,2-Dichloroethylene | 4 U |
| 1,1-Dichloroethane | 4 U |
| Trichloromethane | 4 U |
| 1,1,1-Trichloroethane | 4 U |
| Carbon Tetrachloride | 4 U |
| 1,2-Dichloroethane | 4 U |
| Benzene | 4 U |
| Trichloroethylene | 4 U |
| Bromodichloromethane | 4 U |
| Dibromomethane | 4 U |
| Toluene | 4 U |
| 1,1,2-Trichloroethane | 4 U |
| Tetrachloroethylene | 4 U |
| Ethylbenzene | 4 U |
| m & p-Xylenes | 4 U |
| o-Xylene | 4 U |
| Styrene | 4 U |
| 1,1,2,2-Tetrachloroethane | 4 U |
| 1,3,5-Trimethylbenzene | 4 U |
| | |
| p-Bromofluorobenzene (% Rec) | 100 |
| | |
| Pressurized Sample Volume (mL) | 250 |
| Initial Pressure (psia) | N/A |
| Final Pressure (psia) | N/A |
| Quantitation Limit (ppbv) | 4 |

A - Assumed volume for Blanks
B - <3 times Method Blank value
C - Compound Calibration >25% RSD
D - Compound Calibration Check >25% RPD
E - Concentration exceeded calibration limit (25nL)
J - Below 1.00 nL Quantitation Limit
U - Not Detected
N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 1 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | |
|---------------------|----------|--|
| Sample Number: | Method | Reference Standard: Bromochloromethane |
| Sample Location: | Blank | Reference Std Conc. (ppbv): 21.2 |
| Sample Volume (mL): | 250 | Reference Std Volume (mL): 500 |
| Date Sampled: | N/A | Reference Std Area: 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): N/A |
| Data File: | AGS003 | Final Pressure (psig): N/A |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|----------------------------|----------------|------|----------------------|
| No non-targets were found. | | | |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples

Page 2 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|------------|-----------------------------|--------------------|
| Sample Number: | 17747 | Reference Standard: | Bromochloromethane |
| Sample Location: | Trip/Field | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 250 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | N/A |
| Data File: | AGS004 | Final Pressure (psig): | N/A |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|----------------------------|----------------|------|----------------------|
| No non-targets were found. | | | |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 3 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17740 | Reference Standard: | Bromochloromethane |
| Sample Location: | DW3 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 14.8 |
| Data File: | AGS005 | Final Pressure (psig): | 29.6 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|-------------------------|----------------|---------|----------------------|
| cycloalkane/alkene | 2.792 | 2442275 | 8 |
| cycloalkane/alkene | 3.611 | 962606 | 3 * |
| acetaldehyde | 3.914 | 916963 | 3 * |
| furan + unknown | 6.675 | 2271909 | 7 |
| acetone | 6.937 | 1719096 | 6 |
| 2methyl-furan + unknown | 10.765 | 1956463 | 6 |
| aldehyde | 19.271 | 2398466 | 8 |

* - Below 4 ppbv Limit of Quantitation

N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 4 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17741 | Reference Standard: | Bromochloromethane |
| Sample Location: | DW2 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 2.0 |
| Data File: | AGS006 | Final Pressure (psig): | 16.0 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|---------------|----------------|---------|----------------------|
| acetone | 6.999 | 1058843 | 13 * |

* - Below 16 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples

Page 5 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17742 | Reference Standard: | Bromochloromethane |
| Sample Location: | DW1 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 15.2 |
| Data File: | AGS007 | Final Pressure (psig): | 30.4 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|---------------|----------------|--------|----------------------|
| unknown | 8.192 | 987161 | 3 * |
| aldehyde | 19.288 | 940362 | 3 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 6 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17743 | Reference Standard: | Bromochloromethane |
| Sample Location: | DW4 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 8.8 |
| Data File: | AGS008 | Final Pressure (psig): | 17.6 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|---------|----------------------|
| alkene | 2.800 | 1499003 | 5 |
| acetone | 6.968 | 1622183 | 5 |
| unknown | 8.193 | 1208501 | 4 * |
| alkane | 9.626 | 942198 | 3 * |
| 2-methyl-furan + unknown | 10.772 | 924300 | 3 * |
| aldehyde | 19.283 | 936177 | 3 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 7 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17744 | Reference Standard: | Bromochloromethane |
| Sample Location: | DW5 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 14.0 |
| Data File: | AGS009 | Final Pressure (psig): | 28.0 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|---------|----------------------|
| cycloalkane/alkene | 2.800 | 834872 | 3 * |
| alkane + alkane | 3.238 | 1004830 | 3 * |
| n-butane + alkane | 3.628 | 1397502 | 4 |
| n-hexane | 9.635 | 1440931 | 5 |
| toluene + siloxane | 16.720 | 2327997 | 7 |
| trimethyl-benzene isomer | 22.913 | 939916 | 3 * |

* - Below 4 ppbv Limit of Quantitation

N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 8 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17745 | Reference Standard: | Bromochloromethane |
| Sample Location: | UW1 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 8.4 |
| Data File: | AGS010 | Final Pressure (psig): | 16.8 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|----------------------------|----------------|------|----------------------|
| No non-targets were found. | | | |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples**

Page 9 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|-----------------------------|--------------------|
| Sample Number: | 17746 | Reference Standard: | Bromochloromethane |
| Sample Location: | UW2 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 13322670 |
| Date Analyzed: | 04/10/00 | Initial Pressure (psig): | 8.4 |
| Data File: | AGS012 | Final Pressure (psig): | 16.8 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|----------------------------|----------------|------|----------------------|
| No non-targets were found. | | | |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples

Page 10 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|----------|--|---------|
| Sample Number: | Method | Reference Standard: Bromochloromethane | |
| Sample Location: | Blank | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 250 | Reference Std Volume (mL): | 500 |
| Date Sampled: | N/A | Reference Std Area: | 8543457 |
| Date Analyzed: | 04/11/00 | Initial Pressure (psig): | N/A |
| Data File: | AGS015 | Final Pressure (psig): | N/A |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|----------------------------|----------------|------|----------------------|
| No non-targets were found. | | | |

* - Below 4 ppbv Limit of Quantitation

N/A - Not Applicable

Table 4 - Air Toxic Non-target Compound Results
Summa Canister Samples

Page 11 of 11

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| | | | |
|---------------------|-----------|-----------------------------|--------------------|
| Sample Number: | 17744 Rep | Reference Standard: | Bromochloromethane |
| Sample Location: | DW5 | Reference Std Conc. (ppbv): | 21.2 |
| Sample Volume (mL): | 500 | Reference Std Volume (mL): | 500 |
| Date Sampled: | 04/06/00 | Reference Std Area: | 8543457 |
| Date Analyzed: | 04/11/00 | Initial Pressure (psig): | 14.0 |
| Data File: | AGS016 | Final Pressure (psig): | 28.0 |

| Compound Name | Retention Time | Area | Concentration (ppbv) |
|--------------------------|----------------|---------|----------------------|
| unknown + alkane | 3.198 | 696075 | 4 * |
| alkane | 3.596 | 898837 | 5 |
| alkene | 4.956 | 613106 | 3 * |
| acetone | 6.890 | 892721 | 4 |
| toluene + siloxane | 16.677 | 1684447 | 8 |
| trimethyl-benzene isomer | 22.853 | 716418 | 4 * |

* - Below 4 ppbv Limit of Quantitation
N/A - Not Applicable

**Table 5 - Air Toxic MS/MSD Recovery Summary for Summa Canister Samples
APG Burn Site, Edgewood, MD WA # 0-110**

| Sample Number | | 17746 | 17746 MS | | 17746 MSD | |
|-------------------------------|-----------------|----------|----------|----------|-----------|----------|
| Sample Location | | UW2 | UW2 | | UW2 | |
| Date Sampled | | 04/06/00 | 04/06/00 | | 04/06/00 | |
| Date Analyzed | | 04/10/00 | 04/10/00 | % | 04/10/00 | % |
| Data File | Spike Amount | AGS012 | AGS013 | Recovery | AGS014 | Recovery |
| Chloromethane | 9.8 | 0.19 | 10.95 | 110 | 10.99 | 110 |
| Vinyl Chloride | 9.7 | 0.00 | 11.12 | 115 | 11.50 | 119 |
| Chloroethane | 10.0 | 0.00 | 11.91 | 119 | 11.99 | 120 |
| Trichlorofluoromethane | 10.4 | 0.00 | 12.87 | 124 | 13.33 | 128 |
| 1,1-Dichloroethene | 10.2 | 0.00 | 10.65 | 104 | 10.79 | 106 |
| Methylene Chloride | 10.0 | 0.00 | 10.14 | 101 | 10.23 | 102 |
| trans-1,2-Dichloroethene | 10.0 | 0.00 | 10.34 | 103 | 10.58 | 106 |
| 1,1-Dichloroethane | 10.2 | 0.00 | 10.56 | 104 | 10.79 | 106 |
| Trichloromethane | 10.2 | 0.00 | 10.22 | 100 | 10.54 | 103 |
| 1,1,1-Trichloroethane | 10.1 | 0.00 | 10.17 | 101 | 10.64 | 105 |
| Carbon Tetrachloride | 9.8 | 0.00 | 10.40 | 106 | 10.68 | 109 |
| 1,2-Dichloroethane | 10.2 | 0.00 | 10.61 | 104 | 10.74 | 105 |
| Benzene | 10.0 | 0.00 | 10.06 | 101 | 10.19 | 102 |
| Trichloroethylene | 10.0 | 0.00 | 9.95 | 99 | 10.12 | 101 |
| Bromodichloromethane | 10.1 | 0.00 | 10.16 | 101 | 10.43 | 103 |
| Dibromomethane | 9.8 | 0.00 | 10.10 | 103 | 10.23 | 104 |
| Toluene | 10.1 | 0.00 | 10.07 | 100 | 10.19 | 101 |
| 1,1,2-Trichloroethane | 9.8 | 0.00 | 10.12 | 103 | 10.45 | 107 |
| Tetrachloroethylene | 10.0 | 0.00 | 10.35 | 103 | 10.49 | 105 |
| Ethylbenzene | 10.1 | 0.00 | 10.13 | 100 | 10.17 | 101 |
| meta & para-Xylenes | 10.2 | 0.00 | 10.20 | 100 | 10.25 | 100 |
| ortho-Xylene | 10.4 | 0.00 | 10.31 | 99 | 10.42 | 100 |
| Styrene | 10.4 | 0.00 | 10.51 | 101 | 10.52 | 101 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 0.00 | 10.70 | 107 | 10.96 | 110 |
| 1,3,5-trimethylbenzene | 10.5 | 0.00 | 10.67 | 102 | 10.82 | 103 |
| 1-Bromofluorobenzene (% Rec.) | N/A | 105 | 105 | N/A | 103 | N/A |

N/A - Not Applicable

APPENDIX A
CHAIN-OF-CUSTODY

Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sampled on 06 April 2000

WA#: R1A00110

00110/del/ar/0005/APGBurnar

132-~~1008~~ 321-4200 *RB*

EPA Contract ~~68-C4-0022~~

68099-223

CHA OF CUSTODY RECORD

Project Name: APG Burn

Project Number: 14A00/10

UN-RFW Contact: Amy DeBelle

Phone: 494 4013

No: 03310

SHEET NO. 1 OF 1

041000

Sample Identification

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | VOCs | Volume |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|------|---------|
| 426 | 17740 | DW3 | A | 4/16/00 | 1 | none | ✓ | 6(L) |
| 427 | 17741 | DW2 | | | | | ✓ | 8(L) |
| 428 | 17742 | DW1 | | | | | ✓ | 10(L) |
| 429 | 17743 | DW4 | | | | | ✓ | 4.5(L) |
| 430 | 17744 | DW5 | | | | | ✓ | 6(L) |
| 431 | 17745 | WW1 | | | | | ✓ | 3.96(L) |
| 432 | 17746 | WW2 | | | | | ✓ | 4.8(L) |
| 433 | 17747 | Trip Field | ↓ | ↓ | ↓ | ↓ | ✓ | 8(L) |
| (AED) | | | | | | | | |
| (AED) | | | | | | | | |
| 0022 | | | | | | | | |

Matrix:

SD - Sediment
DS - Drum Solids
DL - Drum Liquids
X - Other

PW - Potable Water
GW - Groundwater
SW - Surface Water
SL - Sludge

| | |
|-----|-------|
| S - | Soil |
| W - | Water |
| O - | Oil |
| A - | Air |

Special Instructions:

VOCs - Volatile Organic
Compounds Analysis
by Modified TO-14

(L) - Liters

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF
CUSTODY #

| Items/Reason | Relinquished By | Date | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|---------|-------------|---------|------|--------------|-----------------|---------|-------------|---------|------|
| Standby | James J. | 4/10/00 | Dan H. Weny | 4/10/00 | 1410 | ALL/Analysis | Dan H. Weny | 4/10/00 | James J. | 4/10/00 | 1510 |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |

APPENDIX B
Analytical Report (PAH, Inorganic Acids, Metals, and Dioxins/Furans)
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

ANALYTICAL REPORT

Prepared by
Lockheed Martin

Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities
Aberdeen Proving Ground, Aberdeen, MD

May 2000

EPA Work Assignment No. 0-110
LOCKHEED MARTIN Work Order No. R1A00110
EPA Contract No. 68-C99-223

Submitted to
D. Mickunas
EPA-ERT

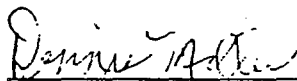


A. Dubois
Task Leader

5/31/00

Date

Analysis by:
REAC
SWRI

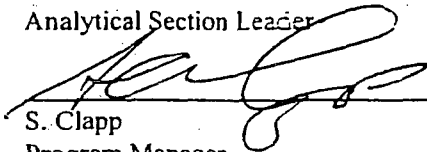


D. Miller
Analytical Section Leader

6/1/2000

Date

Prepared by:
M. Bernick



S. Clapp
Program Manager

6/1/00

Date

Reviewed by:
D. Killeen

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Appendices will be furnished on request.

Introduction

REAC in response to WA # 0-110, provided analytical support for environmental samples collected from Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities, located in Aberdeen Proving Ground, Aberdeen, MD as described in the following table. The support also included QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

The samples were treated with procedures consistent with those described in SOP # 1008 and are summarized in the following table:

| COC # | Number of Samples | Sampling Date | Date Received | Matrix | Analysis | Laboratory |
|-------|-------------------|---------------|---------------|--------|-----------------|------------|
| 06966 | 10 | 4/6/00 | 4/11/00 | Air | Dioxin | SWRI* |
| 05254 | 10 | 4/6/00 | 4/11/00 | | Metals | |
| 06965 | 10 | 4/6/00 | 4/11/00 | | Inorganic Acids | |
| 05654 | 10 | 4/6/00 | 4/10/00 | | PAH | REAC |

* SWRI denotes Southwest Research Institute

Case Narrative

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

PAH in Air Package J142

The data were examined and were found to be satisfactory.

Inorganic Acids in Air Package J159

All sample results were lot blank subtracted.

Sample 17733, the trip blank, contained 0.0013 mg of sulfuric acid. The sulfuric acid results for samples 17700 through 17706 are considered not detected.

Metals in Air Package J158

The method blank contained 11.81 µg/filter sodium (Na). The Na results for samples 17680 through 17689 are considered not detected.

Sample 17689, the lot blank, contained 6.2 µg/filter calcium (Ca), 0.51 µg/filter chromium (Cr), 1.6 µg/filter iron (Fe) and 1.1 µg/filter zinc (Zn). The Ca, Cr and Zn results for samples 17680 through 17688, and the Fe results for samples 17681 through 17685, 17687 and 17688 are considered not detected because the concentration in the

00110\DEL\AR\0005\APGBurnar

sample is less than five times the concentration in the lot blank.

Sample 17688, the trip blank, contained 0.23 µg/filter nickel (Ni). The Ni results for samples 17681, 17683, 17686 and 17687 are considered not detected because the concentration in the sample is less than five times the concentration in the trip blank.

The LCS percent recovery exceeded the QC limits for tellurium (Te) (6.7%). The BS/BSD percent recovery exceeded the QC limits for Te (BS 8.2%, BSD 8.6%), phosphorus (P) (BS 126%, BSD 127%), tin (Sn) (BS 131%, BSD 132%), and zirconium (Zr) (BS 134%, BSD 131%). The Te results for samples 17680 through 17689 are considered unusable.

Dioxins in Air Package J 160

The method blank contained 13 pg OCDF. The OCDF result for sample 17679 is considered estimated.

Sample 17679, the lot blank, contained 0.660 pg 123789-HxCDD. This compound was not detected in the associated samples; the data are not affected.

Sample 17677, the trip blank, contained 12.2 pg OCDD. The OCDD results for samples 17670, 17675, 17674 and 17676 are considered not detected because the sample concentrations were less than five times that found in the trip blank.

Lock mass ion 342 (penta dioxins and furans) exhibited a loss of sensitivity during the calibration verification on 4/17/00 (6:41 am) on instrument H. None of the associated samples exhibited a sensitivity loss for this ion during analysis; the data are not affected.

In the ending calibration verification standard of 4/17/00 (6:41 am), the acceptable percent difference QC limits were exceeded for ¹³C-12378-PeCDD (45%) and ¹³C-OCDD (46%). As required by the method criteria, the subcontracted laboratory used the two continuing calibrations bracketing the samples to calculate average relative response factors for quantitation. Samples 17673, 17674, 17675, 17676, 17677 and 17678 were quantitated using these average response factors. The percent relative standard deviation of these average response factors exceeded the QC limits for OCDF (21), ¹³C-12378-PeCDD (43) and ¹³C-OCDD (33). The OCDD results for samples 17673, 17674, 17675, 17676 and 17677 are considered estimated.

The acceptable QC limits were exceeded for the percent recovery for internal standard ¹³C-12378-PeCDD (138%) for sample 17676. 12378-PeCDD was not detected in this sample; the data are not affected.

The acceptable QC limits were exceeded for the percent recovery for internal standard ¹³C-12378-PeCDF (137%) for sample 17676. Pentafurans were not detected in this sample; the data are not affected.

Summary of Abbreviations

| | |
|----------------|--|
| AA | Atomic Absorption |
| B | The analyte was found in the blank |
| BFB | Bromofluorobenzene |
| C | Centigrade |
| D | (Surrogate Table) this value is from a diluted sample and was not calculated (Result Table) this result was obtained from a diluted sample |
| Dioxin | denotes Polychlorinated Dibenzo-p-dioxins and Polychlorinated Dibenzofurans and/or PCDD and PCDF |
| CLP | Contract Laboratory Protocol |
| COC | Chain of Custody |
| CONC | Concentration |
| CRDL | Contract Required Detection Limit |
| CRQL | Contract Required Quantitation Limit |
| DFTPP | Decafluorotriphenylphosphine |
| DL | Detection Limit |
| E | The value is greater than the highest linear standard and is estimated |
| EMPC | Estimated maximum possible concentration |
| ICAP | Inductively Coupled Argon Plasma |
| ISTD | Internal Standard |
| J | The value is below the method detection limit and is estimated |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| MDL | Method Detection Limit |
| MI | Matrix Interference |
| MS | Matrix Spike |
| MSD | Matrix Spike Duplicate |
| MW | Molecular Weight |
| NA | either Not Applicable or Not Available |
| NC | Not Calculated |
| NR | Not Requested |
| NS | Not Spiked |
| % D | Percent Difference |
| % REC | Percent Recovery |
| PPB | Parts per billion |
| PPBV | Parts per billion by volume |
| PPMV | Parts per million by volume |
| PQL | Practical Quantitation Limit |
| QA/QC | Quality Assurance/Quality Control |
| QL | Quantitation Limit |
| RPD | Relative Percent Difference |
| RSD | Relative Standard Deviation |
| SIM | Selected Ion Monitoring |
| TCLP | Toxic Characteristics Leaching Procedure |
| U | Denotes not detected |
| W | Weathered analyte; the results should be regarded as estimated |
| m ³ | cubic meter kg kilogram µg microgram |
| L | liter g gram pg picogram |
| mL | milliliter mg milligram ng nanogram |
| µL | microliter |
| * | denotes a value that exceeds the acceptable QC limit Abbreviations that are specific to a particular table are explained in footnotes on that table |

Revision 1/5/00

Analytical Procedure for PAH in Air (XAD-2 Tubes)

XAD-2 Tube Preparation

The XAD-2 tubes were analyzed for polycyclic aromatic hydrocarbons (PAH) using modified NIOSH Method 5515. The front, back and filter portions of the tubes were analyzed separately by extracting them with 2.0 mL methylene chloride. A preweighed filter was also collected with these tubes and this filter was extracted with 4.0 mL methylene chloride. One mL of extract was spiked with 20 µL of a 2000 ppm XAD internal standards solution consisting of naphthalene-d₈, acenaphthene-d₁₀, phenanthrene-d₁₀, chrysene-d₁₂, and perylene-d₁₂, resulting in a 40.0 ppm concentration and analyzed.

GC/MS Analysis

An HP 6890 MSD, equipped with a 6890 autosampler and controlled by a personal computer equipped with HP-Enviroquant software was used to analyze the samples.

The instrument conditions were:

| | |
|--|---|
| Column | Restek Rtx-5 (cross bonded SE-54) 30 meter x 0.25 mm ID, 0.50 µm film thickness. |
| Flow Rate | 1 mL/min, EPC enabled |
| Injection Temperature | 280° C |
| Transfer Temperature | 280° C |
| Source Temperature & Analyzer Temperature | Controlled by thermal transfer of heat from Transfer Line 280° C |
| Temperature Program | 70° C for 0.5 min 30° C/min to 295° C; hold for 8 minutes 30° C/min to 315° C; hold for 7 min |
| Pulsed Splitless Injection | Pressure Pulse = 16 psi for 1.0 min, then normal flow 8:1 Split Ratio |
| Injection Volume | 1µL |

The GC/MS was calibrated using 6 PAH standards at 10, 25, 50, 75, 100 and 150 ppm. Before analysis each day the system was tuned with 50-ng decafluorotriphenylphosphine (DFTPP) and passed a continuing calibration check by analyzing a 50µg/mL daily standard. The QC limit for the initial calibration is %RSD less than 30 and %D less than 25 for the daily check. Sample quantification is based on the average response factor of the calibration curve or the response factor of the daily calibration check.

The XAD-2 tube PAH results are listed in Table 1.1. Tentatively identified compounds are listed in Table 1.2. The following equations were used to calculate the analyte - total µg/sample:

$$\mu\text{g/sample} = C_u \times V \times DE = \frac{A_u \times C_{is} \times V \times DE}{A_{is} \times RRF}$$

where

- C_u = Concentration of the analyte (µg/mL)
- V = Extraction Volume (mL)
- DE = Desorption Efficiency = 100/(% Recovery)
- A_u = Area of the analyte
- C_{is} = Concentration of the internal standard (µg/mL)
- A_{is} = Area of the internal standard

The Relative Response Factor, RRF, is calculated from the calibration standard mixture using

$$RRF = \frac{A_u \times C_{is}}{A_{is} \times C_u}$$

where

- RRF = Relative Response Factor (unitless)
- A_u = Area of Analyte in the standard mixture
- C_{is} = Concentration of Internal Standard in the standard mixture (µg/mL)
- A_{is} = Area of Internal Standard in the standard mixture
- C_u = Concentration of Analyte in the standard mixture (µg/mL)

The concentration of the analyte in mg/m³ and ppbv (parts per billion by volume) is calculated using the following:

$$\text{mg/m}^3 = \frac{(\text{Total } \mu\text{gfront} + \text{Total } \mu\text{gback})}{\text{Liters Sampled}}$$

$$\text{ppbv} = \frac{\text{mg/m}^3 \times 24.45 \times 1000}{MW}$$

where MW is the molecular weight of the analyte

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Analytical Procedure for Inorganic Acids in Air

The subcontract laboratory determined the concentration of inorganic acids in the samples by analyzing them according to NIOSH Method 7903. The results of the analysis for the air samples are listed in Table 1.3.

Analytical Procedure for Metals in Air

The subcontract laboratory determined the concentration of Metals in the samples by analyzing them according to NIOSH Method 7300. The results of the analysis for the air samples are listed in Table 1.4.

Analytical Procedure for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

The subcontract laboratory determined the concentration of polychlorinated dibenzodioxins and polychlorinated dibenzofurans in the samples by analyzing them according to USEPA SW-846 Method 8290. The results of the analysis are listed in Table 1.5.

Table 1.1 Results of the Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| Sample No. Sampling Location Volume (L) | 17699 Lot Blank 0 | | 17690 DW3 360 | | 17691 DW2 351 | | 17692 DW1 360 | | 17693 DW4 369 | |
|---|-------------------------|-----------|---------------------|-------------|---------------------|-------------|---------------------|-------------|---------------------|-------------|
| Compound Name | Conc. µg | MDL µg | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv |
| Naphthalene | U | 11 | U | 5.6 | 1.7 J | 5.7 | U | 5.6 | U | 5.4 |
| 2-Methylnaphthalene | U | 11 | U | 5.3 | U | 5.4 | U | 5.3 | U | 5.2 |
| 1-Methylnaphthalene | U | 11 | U | 5.2 | U | 5.3 | U | 5.2 | U | 5.0 |
| Biphenyl | U | 11 | U | 5.0 | U | 5.1 | U | 5.0 | U | 4.8 |
| 2,6-Dimethylnaphthalene | U | 11 | U | 4.9 | U | 5.0 | U | 4.9 | U | 4.7 |
| Acenaphthylene | U | 12 | U | 5.2 | U | 5.3 | U | 5.2 | U | 5.1 |
| Acenaphthene | U | 11 | U | 4.7 | U | 4.9 | U | 4.7 | U | 4.6 |
| Dibenzofuran | U | 11 | U | 4.6 | U | 4.7 | U | 4.6 | U | 4.5 |
| Fluorene | U | 11 | U | 4.7 | U | 4.8 | U | 4.7 | U | 4.6 |
| Phenanthrene | U | 11 | U | 4.1 | U | 4.2 | U | 4.1 | U | 4.0 |
| Anthracene | U | 10 | U | 3.9 | U | 4.0 | U | 3.9 | U | 3.8 |
| Carbazole | U | 12 | U | 4.8 | U | 4.9 | U | 4.8 | U | 4.6 |
| Fluoranthene | U | 12 | U | 4.0 | U | 4.1 | U | 4.0 | U | 3.9 |
| Pyrene | U | 12 | U | 4.0 | U | 4.1 | U | 4.0 | U | 3.9 |
| Benzo(a)anthracene | U | 12 | U | 3.5 | U | 3.6 | U | 3.5 | U | 3.4 |
| Chrysene | U | 14 | U | 4.3 | U | 4.4 | U | 4.3 | U | 4.2 |
| Benzo(b)fluoranthene | U | 11 | U | 2.9 | U | 3.0 | U | 2.9 | U | 2.8 |
| Benzo(k)fluoranthene | U | 11 | U | 3.1 | U | 3.2 | U | 3.1 | U | 3.0 |
| Benzo(e)pyrene | U | 12 | U | 3.2 | U | 3.3 | U | 3.2 | U | 3.1 |
| Benzo(a)pyrene | U | 12 | U | 3.3 | U | 3.4 | U | 3.3 | U | 3.2 |
| Indeno(1,2,3-cd)pyrene | U | 13 | U | 3.1 | U | 3.2 | U | 3.1 | U | 3.0 |
| Dibenzo(a,h)anthracene | U | 12 | U | 3.0 | U | 3.1 | U | 3.0 | U | 2.9 |
| Benzo(g,h,i)perylene | U | 12 | U | 3.1 | U | 3.1 | U | 3.1 | U | 3.0 |

Table 1.1 (cont.) Results of the Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

| Sample No. | 17694 | | 17695 | | 17696 | | 17697 | | 17698 | |
|-------------------------|---------------|-------------|---------------|-------------|---------------|-------------|-------------|-----------|-------------|-----------|
| Sampling Location | DW5 | | UW1 | | UW2 | | Field Blank | | Trip Blank | |
| Volume (L) | 360 | | 332 | | 360 | | 0 | | 0 | |
| Compound Name | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. ppbv | MDL ppbv | Conc. µg | MDL µg | Conc. µg | MDL µg |
| Naphthalene | U | 5.6 | U | 6.1 | U | 5.6 | U | 11 | U | 11 |
| 2-Methylnaphthalene | U | 5.3 | U | 5.8 | U | 5.3 | U | 11 | U | 11 |
| 1-Methylnaphthalene | U | 5.2 | U | 5.6 | U | 5.2 | U | 11 | U | 11 |
| Biphenyl | U | 5.0 | U | 5.4 | U | 5.0 | U | 11 | U | 11 |
| 2,6-Dimethylnaphthalene | U | 4.9 | U | 5.3 | U | 4.9 | U | 11 | U | 11 |
| Acenaphthylene | U | 5.2 | U | 5.6 | U | 5.2 | U | 12 | U | 12 |
| Acenaphthene | U | 4.7 | U | 5.1 | U | 4.7 | U | 11 | U | 11 |
| Dibenzofuran | U | 4.6 | U | 5.0 | U | 4.6 | U | 11 | U | 11 |
| Fluorene | U | 4.7 | U | 5.1 | U | 4.7 | U | 11 | U | 11 |
| Phenanthrene | U | 4.1 | U | 4.4 | U | 4.1 | U | 11 | U | 11 |
| Anthracene | U | 3.9 | U | 4.2 | U | 3.9 | U | 10 | U | 10 |
| Carbazole | U | 4.8 | U | 5.2 | U | 4.8 | U | 12 | U | 12 |
| Fluoranthene | U | 4.0 | U | 4.3 | U | 4.0 | U | 12 | U | 12 |
| Pyrene | U | 4.0 | U | 4.3 | U | 4.0 | U | 12 | U | 12 |
| Benzo(a)anthracene | U | 3.5 | U | 3.8 | U | 3.5 | U | 12 | U | 12 |
| Chrysene | U | 4.3 | U | 4.6 | U | 4.3 | U | 14 | U | 14 |
| Benzo(b)fluoranthene | U | 2.9 | U | 3.1 | U | 2.9 | U | 11 | U | 11 |
| Benzo(k)fluoranthene | U | 3.1 | U | 3.4 | U | 3.1 | U | 11 | U | 11 |
| Benzo(e)pyrene | U | 3.2 | U | 3.5 | U | 3.2 | U | 12 | U | 12 |
| Benzo(a)pyrene | U | 3.3 | U | 3.6 | U | 3.3 | U | 12 | U | 12 |
| Indeno(1,2,3-cd)pyrene | U | 3.1 | U | 3.4 | U | 3.1 | U | 13 | U | 13 |
| Dibenzo(a,h)anthracene | U | 3.0 | U | 3.2 | U | 3.0 | U | 12 | U | 12 |
| Benzo(g,h,i)perylene | U | 3.1 | U | 3.3 | U | 3.1 | U | 12 | U | 12 |

Table 1.2 Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17699 Lot Blank

LabFile # APG059

Con. Factor

2.0

| | CAS# | Compound | Q | RT | Conc ** Total µg |
|----|---------|------------------------|----|-------|---------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 95 | 9.73 | 28 |
| 2 | | Unknown | | 15.14 | 13 |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
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| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17690

LabFile # APG062

Con. Factor

5.6

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|-------|------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 91 | 9.73 | 97 |
| 2 | | Unknown | | 15.15 | 48 |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
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| 12 | | | | | |
| 13 | | | | | |
| 14 | | | | | |
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| 17 | | | | | |
| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17691

LabFile # APG065

Con. Factor

5.7

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|-------|------------------|
| 1 | | Unknown phenol | | 4.96 | 28 |
| 2 | 85-68-7 | Butyl benzyl phthalate | 91 | 7.93 | 100 |
| 3 | | Unknown | | 15.15 | 45 |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
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| 11 | | | | | |
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| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17692

LabFile # APG068

Con. Factor

5.6

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|------|------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 70 | 9.73 | 93 |
| 2 | | | | | |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
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| 13 | | | | | |
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| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17693

LabFile # APG071

Con. Factor

5.4

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|------|------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 95 | 9.73 | 92 |
| 2 | | | | | |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
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| 10 | | | | | |
| 11 | | | | | |
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| 13 | | | | | |
| 14 | | | | | |
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| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17694

LabFile # APG074

Con. Factor

5.6

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|------|------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 91 | 9.73 | 91 |
| 2 | | | | | |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
| 9 | | | | | |
| 10 | | | | | |
| 11 | | | | | |
| 12 | | | | | |
| 13 | | | | | |
| 14 | | | | | |
| 15 | | | | | |
| 16 | | | | | |
| 17 | | | | | |
| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17695

LabFile # APG077

Con. Factor

6.0

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|------|------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 94 | 9.73 | 95 |
| 2 | | | | | |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
| 9 | | | | | |
| 10 | | | | | |
| 11 | | | | | |
| 12 | | | | | |
| 13 | | | | | |
| 14 | | | | | |
| 15 | | | | | |
| 16 | | | | | |
| 17 | | | | | |
| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17696

LabFile # APG080

Con. Factor

5.6

| | CAS# | Compound | Q | RT | Conc ** µg/m3 |
|----|---------|------------------------|----|------|------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 70 | 9.73 | 85 |
| 2 | | | | | |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
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| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17697

LabFile # APG083

Con. Factor

2.0

| | CAS# | Compound | Q | RT | Conc ** Total µg |
|----|---------|------------------------|----|-------|---------------------|
| 1 | 85-68-7 | Butyl benzyl phthalate | 94 | 9.73 | 30 |
| 2 | | Unknown | | 15.13 | 12 |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
| 9 | | | | | |
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| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |

**Estimated Concentration (Response Factor = 1.0)

Table 1.2 (cont.) Results of the TIC for PAH in Air

WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample # 17698

LabFile # APG086

Con. Factor

2.0

| | CAS# | Compound | Q | RT | Conc ** |
|----|---------|------------------------|----|-------|----------|
| | | | | | Total µg |
| 1 | 85-68-7 | Butyl benzyl phthalate | 94 | 9.73 | 33 |
| 2 | | Unknown | | 15.14 | 14 |
| 3 | | | | | |
| 4 | | | | | |
| 5 | | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
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**Estimated Concentration (Response Factor = 1.0)

Table 1.3 Results of the Analysis for Inorganic Acids in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| | | | | | | | | | | |
|-------------------|------------|-----------|-------------|-----------|------------|-----------|---------------|--------------|--------------|-------------|
| Sample ID: | 17734 | | 17732 | | 17733 | | 17700 | | | |
| Location: | Lot Blank | | Field Blank | | Trip blank | | DW-3 | | | |
| Air Volume (L): | 0 | | 0 | | 0 | | 45.9 | | | |
| | | | | | | | | | | |
| Analyte | Conc mg | MDL mg | Conc mg | MDL mg | Conc mg | MDL mg | Conc mg/m³ | MDL mg/m³ | Conc ppmv | MDL ppmv |
| Hydrobromic acid | U | 0.0010 | U | 0.0010 | U | 0.0010 | U | 0.0220 | U | 0.0067 |
| Hydrochloric acid | 0.0031 | 0.0010 | U | 0.0010 | U | 0.0010 | 0.184 | 0.0224 | 0.123 | 0.0150 |
| Hydrofluoric acid | U | 0.0011 | U | 0.0011 | U | 0.0011 | 0.0592 | 0.0229 | 0.0724 | 0.0280 |
| Nitric acid | U | 0.0045 | U | 0.0045 | U | 0.0045 | U | 0.0980 | U | 0.0380 |
| Phosphoric acid | U | 0.0032 | U | 0.0032 | U | 0.0032 | U | 0.0689 | U | 0.0172 |
| Sulfuric acid | 0.0050 | 0.0010 | U | 0.0010 | 0.0013 | 0.0010 | 0.0901 | 0.0222 | 0.0225 | 0.0055 |

| | | | | | | | | | | | | |
|-------------------|---------------------------|--------------------------|--------------|-------------|---------------------------|--------------------------|--------------|-------------|---------------------------|--------------------------|--------------|-------------|
| Sample ID | 17701 | | | | 17702 | | | | 17703 | | | |
| Location | DW2 | | | | DW1 | | | | DW4 | | | |
| Air Volume (L): | 45.0 | | | | 45.9 | | | | 45.5 | | | |
| | <hr/> | | | | | | | | | | | |
| Analyte | Conc mg/m ³ | MDL mg/m ³ | Conc ppmv | MDL ppmv | Conc mg/m ³ | MDL mg/m ³ | Conc ppmv | MDL ppmv | Conc mg/m ³ | MDL mg/m ³ | Conc ppmv | MDL ppmv |
| | <hr/> | | | | | | | | | | | |
| Hydrobromic acid | U | 0.0225 | U | 0.0068 | U | 0.0220 | U | 0.0067 | U | 0.0222 | U | 0.0067 |
| Hydrochloric acid | U | 0.0223 | U | 0.0153 | 0.0578 | 0.0224 | 0.0387 | 0.0150 | U | 0.0226 | U | 0.0152 |
| Hydrofluoric acid | 0.0842 | 0.0234 | 0.103 | 0.0286 | 0.0239 | 0.0229 | 0.0292 | 0.0280 | 0.0359 | 0.0231 | 0.0439 | 0.0283 |
| Nitric acid | U | 0.1000 | U | 0.0388 | U | 0.0980 | U | 0.0380 | U | 0.0989 | U | 0.0384 |
| Phosphoric acid | U | 0.0702 | U | 0.0175 | U | 0.0689 | U | 0.0172 | U | 0.0695 | U | 0.0173 |
| Sulfuric acid | 0.105 | 0.0227 | 0.0262 | 0.0057 | 0.0732 | 0.0222 | 0.0182 | 0.0055 | 0.0644 | 0.0224 | 0.0161 | 0.0056 |

| Sample ID | 17704 | | | | 17705 | | | | 17706 | | | |
|-------------------|---------------|--------------|--------------|-------------|---------------|--------------|--------------|-------------|---------------|--------------|--------------|-------------|
| Location | DW5 | | | | UW1 | | | | UW2 | | | |
| Air Volume (L): | 45.0 | | | | 41.8 | | | | 45.5 | | | |
| Analyte | Conc mg/m³ | MDL mg/m³ | Conc ppmv | MDL ppmv | Conc mg/m³ | MDL mg/m³ | Conc ppmv | MDL ppmv | Conc mg/m³ | MDL mg/m³ | Conc ppmv | MDL ppmv |
| Hydrobromic acid | U | 0.0225 | U | 0.0068 | U | 0.0242 | U | 0.0073 | U | 0.0222 | U | 0.0067 |
| Hydrochloric acid | 0.0281 | 0.0223 | 0.0188 | 0.0153 | U | 0.0246 | U | 0.0165 | 0.0262 | 0.0226 | 0.0176 | 0.0152 |
| Hydrofluoric acid | 0.0318 | 0.0234 | 0.0389 | 0.0286 | U | 0.0252 | U | 0.0308 | U | 0.0231 | U | 0.0283 |
| Nitric acid | U | 0.1000 | U | 0.0388 | U | 0.1077 | U | 0.0418 | U | 0.0989 | U | 0.0384 |
| Phosphoric acid | U | 0.0702 | U | 0.0175 | U | 0.0756 | U | 0.0189 | U | 0.0695 | U | 0.0173 |
| Sulfuric acid | 0.162 | 0.0227 | 0.0404 | 0.0057 | 0.0703 | 0.0244 | 0.0175 | 0.0061 | 0.087 | 0.0224 | 0.0217 | 0.0056 |

All sample results are lot blank subtracted.

Table 1.4 Results of the Analysis for Metals in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Client ID | Method Blank | | 17680 | | 17681 | | 17682 | | 17683 | | 17684 | |
|----------------|--------------|-----------|---------------|--------------|---------------|--------------|---------------|--------------|---------------|--------------|---------------|--------------|
| Location | - | | DW3 | | DW2 | | DW1 | | DW4 | | DW5 | |
| Air Volume (L) | 0 | | 540 | | 540 | | 540 | | 540 | | 540 | |
| Parameter | Conc µg | MDL µg | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ |
| Aluminum | U | 1.0 | 3.8 | 1.9 | 4.0 | 1.9 | 1.9 | 1.9 | U | 1.9 | 1.9 | 1.9 |
| Arsenic | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Beryllium | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Cadmium | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Calcium | U | 2.0 | 34 | 3.7 | 40 | 3.7 | 15 | 3.7 | 11 | 3.7 | 13 | 3.7 |
| Chromium | U | 0.10 | 1.0 | 0.19 | 1.1 | 0.19 | 1.0 | 0.19 | 1.1 | 0.19 | 0.98 | 0.19 |
| Cobalt | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Copper | U | 0.10 | U | 0.19 | 0.22 | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Iron | U | 1.0 | 25 | 1.9 | 9.2 | 1.9 | 3.6 | 1.9 | 2.6 | 1.9 | 2.0 | 1.9 |
| Lead | U | 0.10 | U | 0.19 | 0.33 | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Lithium | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Magnesium | U | 1.0 | 4.2 | 1.9 | 5.6 | 1.9 | 2.0 | 1.9 | U | 1.9 | U | 1.9 |
| Manganese | U | 0.10 | 1.0 | 0.19 | 1.0 | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Molybdenum | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Nickel | U | 0.20 | U | 0.37 | 0.81 | 0.37 | U | 0.37 | 0.38 | 0.37 | U | 0.37 |
| Phosphorus | U | 0.40 | 1.3 | 0.74 | 2.2 | 0.74 | 1.3 | 0.74 | 0.81 | 0.74 | U | 0.74 |
| Platinum | U | 1.0 | U | 1.9 | U | 1.9 | U | 1.9 | U | 1.9 | U | 1.9 |
| Selenium | U | 0.20 | U | 0.37 | U | 0.37 | U | 0.37 | U | 0.37 | U | 0.37 |
| Silver | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Sodium | 11.8 | 6.0 | 17 | 11 | 18 | 11 | 17 | 11 | 16 | 11 | 11 | 11 |
| Tellurium | U | 1.0 | U | 1.9 | U | 1.9 | U | 1.9 | U | 1.9 | U | 1.9 |
| Thallium | U | 0.40 | U | 0.74 | U | 0.74 | U | 0.74 | U | 0.74 | U | 0.74 |
| Tin | U | 0.40 | U | 0.74 | U | 0.74 | U | 0.74 | U | 0.74 | U | 0.74 |
| Titanium | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Vanadium | U | 0.10 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 | U | 0.19 |
| Yttrium | U | 0.20 | U | 0.37 | U | 0.37 | U | 0.37 | U | 0.37 | U | 0.37 |
| Zinc | U | 0.30 | 3.4 | 0.56 | 2.0 | 0.56 | 0.93 | 0.56 | 1.1 | 0.56 | 0.81 | 0.56 |
| Zirconium | U | 0.20 | U | 0.37 | U | 0.37 | U | 0.37 | U | 0.37 | U | 0.37 |
| Tungsten | U | 0.40 | U | 0.74 | U | 0.74 | U | 0.74 | U | 0.74 | U | 0.74 |

Table 1.4 (cont.) Results of the Analysis for Metals in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Client ID | 17685 | | 17686 | | 17687 | | 17688 | | 17689 | |
|----------------|---------------|--------------|---------------|--------------|-------------------|------------------|-------------------|------------------|-------------------|------------------|
| Location | UW1 | | UW2 | | Field Blank | | Trip Blank | | Lot Blank | |
| Air Volume (L) | 498 | | 540 | | 0 | | 0 | | 0 | |
| Parameter | Conc µg/m³ | MDL µg/m³ | Conc µg/m³ | MDL µg/m³ | Conc µg/filter | MDL µg/filter | Conc µg/filter | MDL µg/filter | Conc µg/filter | MDL µg/filter |
| Aluminum | 4.0 | 2.0 | 31 | 1.9 | U | 1.0 | U | 1.0 | U | 1.0 |
| Arsenic | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Beryllium | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Cadmium | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Calcium | 16 | 4.0 | 22 | 3.7 | 6.6 | 2.0 | 7.6 | 2.0 | 6.2 | 2.0 |
| Chromium | 1.1 | 0.20 | 1.2 | 0.19 | 0.62 | 0.10 | 0.88 | 0.10 | 0.51 | 0.10 |
| Cobalt | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Copper | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Iron | 5.7 | 2.0 | 58 | 1.9 | 4.0 | 1.0 | 2.7 | 1.0 | 1.6 | 1.0 |
| Lead | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Lithium | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Magnesium | U | 2.0 | 30 | 1.9 | U | 1.0 | U | 1.0 | U | 1.0 |
| Manganese | U | 0.20 | 0.77 | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Molybdenum | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Nickel | 6.9 | 0.40 | 0.55 | 0.37 | 0.44 | 0.20 | 0.23 | 0.20 | U | 0.20 |
| Phosphorus | 1.0 | 0.80 | 2.1 | 0.74 | U | 0.40 | U | 0.40 | U | 0.40 |
| Platinum | U | 2.0 | U | 1.9 | U | 1.0 | U | 1.0 | U | 1.0 |
| Selenium | U | 0.40 | U | 0.37 | U | 0.20 | U | 0.20 | U | 0.20 |
| Silver | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Sodium | 14 | 12 | 18 | 11 | 9.8 | 6.0 | 9.0 | 6.0 | 7.3 | 6.0 |
| Tellurium | U | 2.0 | U | 1.9 | U | 1.0 | U | 1.0 | U | 1.0 |
| Thallium | U | 0.80 | U | 0.74 | U | 0.40 | U | 0.40 | U | 0.40 |
| Tin | U | 0.80 | U | 0.74 | U | 0.40 | U | 0.40 | U | 0.40 |
| Titanium | 0.62 | 0.20 | 0.71 | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Vanadium | U | 0.20 | U | 0.19 | U | 0.10 | U | 0.10 | U | 0.10 |
| Yttrium | U | 0.40 | U | 0.37 | U | 0.20 | U | 0.20 | U | 0.20 |
| Zinc | 1.7 | 0.60 | 1.2 | 0.56 | 2.1 | 0.30 | 2.4 | 0.30 | 1.1 | 0.30 |
| Zirconium | U | 0.40 | U | 0.37 | U | 0.20 | U | 0.20 | U | 0.20 |
| Tungsten | U | 0.80 | U | 0.74 | U | 0.40 | U | 0.40 | U | 0.40 |

Table 1.5 Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID Location Volume of Air (L) | Blank 04/12/00 0 | | | | 17670 DW3 540 | | | | |
|--|------------------------|------------|-----------|-----------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-------|
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | TEF |
| 2,3,7,8-TCDD | | 0.72 | 10.0 | U | | 3.89 | 18.5 | U | 1 |
| 1,2,3,7,8-PeCDD | | 1.40 | 10.0 | U | | 2.92 | 18.5 | U | 0.5 |
| 1,2,3,4,7,8-HxCDD | U | | 25.0 | U | | 4.85 | 46.3 | U | 0.1 |
| 1,2,3,6,7,8-HxCDD | | 3.16 | 25.0 | U | | 1.22 | 46.3 | U | 0.1 |
| 1,2,3,7,8,9-HxCDD | | 2.20 | 25.0 | U | | 3.37 | 46.3 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | | 4.66 | 25.0 | U | | 2.63 | 46.3 | U | 0.01 |
| OCDD | | 16.7 | 50.0 | U | 35.7 J | | 92.5 | 0.035 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | U | | | | U | | | | |
| Total Hexa-Dioxins | U | | | | U | | | | |
| Total Hepta-Dioxins | U | | | | U | | | | |
| 2,3,7,8-TCDF | U | | 10.0 | U | | 2.04 | 18.5 | U | 0.1 |
| 1,2,3,7,8-PeCDF | U | | 10.0 | U | U | | 18.5 | U | 0.05 |
| 2,3,4,7,8-PeCDF | U | | 10.0 | U | | 1.44 | 18.5 | U | 0.5 |
| 1,2,3,4,7,8-HxCDF | | 2.72 | 25.0 | U | | 0.259 | 46.3 | U | 0.1 |
| 1,2,3,6,7,8-HxCDF | | 1.38 | 25.0 | U | | 0.592 | 46.3 | U | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | | 25.0 | U | | 1.41 | 46.3 | U | 0.1 |
| 2,3,4,6,7,8-HxCDF | U | | 25.0 | U | | 1.41 | 46.3 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | | 12.5 | 25.0 | U | | 7.33 | 46.3 | U | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | | 25.0 | U | U | | 46.3 | U | 0.01 |
| OCDF | 13.0 J | | 50.0 | 0.0130 | | 5.11 | 92.5 | U | 0.001 |
| Total Tetra-Furans | U | | | | U | | 46.3 | | |
| Total Penta-Furans | U | | | | U | | 92.5 | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | U | | | | U | | | | |
| Total | | | | 0.0130 | | | | 0.035 | |

Table 1.5 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 17673 | | | | 17674 | | | | |
|---------------------|--------------|------------|-----------|-----------------------|-----------------|---------------|--------------|--------------------------|-------|
| Location | DW4 | | | | DW5 | | | | |
| Volume of Air (L) | 0 | | | | 540 | | | | |
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | Result pg/m³ | EMPC pg/m³ | MDL pg/m³ | Adjusted Conc (pg/m³) | TEF |
| 2,3,7,8-TCDD | | 0.400 | 10.0 | U | U | | 18.5 | U | 1 |
| 1,2,3,7,8-PeCDD | | 0.440 | 10.0 | U | | 0.888 | 18.5 | U | 0.5 |
| 1,2,3,4,7,8-HxCDD | | 3.72 | 25.0 | U | | 1.96 | 46.3 | U | 0.1 |
| 1,2,3,6,7,8-HxCDD | 4.20 J | | 25.0 | 0.420 | | 6.66 | 46.3 | U | 0.1 |
| 1,2,3,7,8,9-HxCDD | | 3.12 | 25.0 | U | | 2.48 | 46.3 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | 27.7 | | 25.0 | 0.277 | 12.6 J | | 46.3 | 0.126 | 0.01 |
| OCDD | 73.8 | | 50.0 | 0.0738 | 94.6 | | 92.5 | 0.094 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | U | | | | U | | | | |
| Total Hexa-Dioxins | 4.20 | | | | U | | | | |
| Total Hepta-Dioxins | 27.7 | | | | 12.6 | | | | |
| 2,3,7,8-TCDF | | 2.10 | 10.0 | U | U | | 18.5 | U | 0.1 |
| 1,2,3,7,8-PeCDF | | 0.980 | 10.0 | U | | 0.629 | 18.5 | U | 0.05 |
| 2,3,4,7,8-PeCDF | | 0.700 | 10.0 | U | | 0.851 | 18.5 | U | 0.5 |
| 1,2,3,4,7,8-HxCDF | 2.32 J | | 25.0 | 0.232 | | 0.207 | 46.3 | U | 0.1 |
| 1,2,3,6,7,8-HxCDF | | 1.76 | 25.0 | U | U | | 46.3 | U | 0.1 |
| 1,2,3,7,8,9-HxCDF | | 1.56 | 25.0 | U | | 0.666 | 46.3 | U | 0.1 |
| 2,3,4,6,7,8-HxCDF | | 2.58 | 25.0 | U | U | 2.04 | 46.3 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | | 9.64 | 25.0 | U | | 14.5 | 46.3 | U | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | | 6.36 | 25.0 | U | | 6.11 | 46.3 | U | 0.01 |
| OCDF | | 21.3 | 50.0 | U | | 14.0 | 92.5 | U | 0.001 |
| Total Tetra-Furans | U | | | | U | | | | |
| Total Penta-Furans | U | | | | 3.18 | | | | |
| Total Hexa-Furans | 2.32 | | | | U | | | | |
| Total Hepta-Furans | U | | | | U | | | | |
| Total | | | | 1.00 | | | | 0.221 | |

Table 1.5 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 17675 | | | | 17676 | | | | |
|---------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-----------------------------|---------------------------|--------------------------|---------------------------------------|-------|
| Location | UW1 | | | | UW2 | | | | |
| Volume of Air (L) | 498 | | | | 540 | | | | |
| Analyte | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | Result pg/m ³ | EMPC pg/m ³ | MDL pg/m ³ | Adjusted Conc (pg/m ³) | TEF |
| 2,3,7,8-TCDD | | 1.73 | 20.1 | U | U | | 18.5 | U | 1 |
| 1,2,3,7,8-PeCDD | U | | 20.1 | U | | 2.04 | 18.5 | U | 0.5 |
| 1,2,3,4,7,8-HxCDD | | 1.53 | 50.2 | U | U | | 46.3 | U | 0.1 |
| 1,2,3,6,7,8-HxCDD | | 3.01 | 50.2 | U | U | | 46.3 | U | 0.1 |
| 1,2,3,7,8,9-HxCDD | | 3.14 | 50.2 | U | | 1.48 | 46.3 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | | 10.0 | 50.2 | U | | 4.18 | 46.3 | U | 0.01 |
| OCDD | 33.0 J | | 100 | 0.033 | 30.9 J | | 92.5 | 0.030 | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | | |
| Total Penta-Dioxins | U | | | | U | | | | |
| Total Hexa-Dioxins | U | | | | U | | | | |
| Total Hepta-Dioxins | U | | | | U | | | | |
| 2,3,7,8-TCDF | | 1.33 | 20.1 | U | | 1.11 | 18.5 | U | 0.1 |
| 1,2,3,7,8-PeCDF | | 1.65 | 20.1 | U | | 0.481 | 18.5 | U | 0.05 |
| 2,3,4,7,8-PeCDF | | 0.804 | 20.1 | U | | 0.888 | 18.5 | U | 0.5 |
| 1,2,3,4,7,8-HxCDF | | 1.81 | 50.2 | U | | 1.37 | 46.3 | U | 0.1 |
| 1,2,3,6,7,8-HxCDF | | 1.85 | 50.2 | U | | 1.44 | 46.3 | U | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | | 50.2 | U | | 0.925 | 46.3 | U | 0.1 |
| 2,3,4,6,7,8-HxCDF | | 2.17 | 50.2 | U | U | | 46.3 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | | 28.9 | 50.2 | U | | 7.84 | 46.3 | U | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | | 50.2 | U | | 0.703 | 46.3 | U | 0.01 |
| OCDF | | 12.0 | 100 | U | | 9.47 | 92.5 | U | 0.001 |
| Total Tetra-Furans | U | | | | U | | | | |
| Total Penta-Furans | U | | | | U | | | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | U | | | | U | | | | |
| Total | | | | 0.033 | | | | 0.030 | |

Table 1.5 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID Location Volume of Air (L) | | 17677 Trip Blank 0 | | | 17678 Field Blank 0 | | | | |
|--|--------------|--------------------------|-----------|-----------------------|---------------------------|------------|-----------|-----------------------|-------|
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | TEF |
| 2,3,7,8-TCDD | | 0.720 | 10.0 | U | | 0.840 | 10.0 | U | 1 |
| 1,2,3,7,8-PeCDD | U | | 10.0 | U | | 0.880 | 10.0 | U | 0.5 |
| 1,2,3,4,7,8-HxCDD | | 0.280 | 25.0 | U | U | | 25.0 | U | 0.1 |
| 1,2,3,6,7,8-HxCDD | | 0.920 | 25.0 | U | | 1.80 | 25.0 | U | 0.1 |
| 1,2,3,7,8,9-HxCDD | U | | 25.0 | U | | 0.840 | 25.0 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | | 1.84 | 25.0 | U | | 2.16 | 25.0 | U | 0.01 |
| OCDD | 12.2 J | | 50.0 | 0.0122 | | 21.3 | 50.0 | U | 0.001 |
| Total Tetra-Dioxins | U | | | | U | | | U | |
| Total Penta-Dioxins | U | | | | U | | | U | |
| Total Hexa-Dioxins | U | | | | U | | | U | |
| Total Hepta-Dioxins | U | | | | U | | | U | |
| 2,3,7,8-TCDF | | 0.780 | 10.0 | U | U | | 10.0 | U | 0.1 |
| 1,2,3,7,8-PeCDF | | 0.680 | 10.0 | U | | 1.00 | 10.0 | U | 0.05 |
| 2,3,4,7,8-PeCDF | | 0.520 | 10.0 | U | | 0.240 | 10.0 | U | 0.5 |
| 1,2,3,4,7,8-HxCDF | U | | 25.0 | U | U | | 25.0 | U | 0.1 |
| 1,2,3,6,7,8-HxCDF | U | | 25.0 | U | U | | 25.0 | U | 0.1 |
| 1,2,3,7,8,9-HxCDF | | 0.600 | 25.0 | U | | 0.260 | 25.0 | U | 0.1 |
| 2,3,4,6,7,8-HxCDF | | 0.520 | 25.0 | U | | 1.04 | 25.0 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | | 3.38 | 25.0 | U | | 2.64 | 25.0 | U | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | | 25.0 | U | | 0.720 | 25.0 | U | 0.01 |
| OCDF | | 2.82 | 50.0 | U | | 6.70 | 50.0 | U | 0.01 |
| Total Tetra-Furans | U | | | | U | | | | |
| Total Penta-Furans | U | | | | 1.96 | | | | |
| Total Hexa-Furans | U | | | | U | | | | |
| Total Hepta-Furans | U | | | | U | | | | |
| Total | | | | 0.0122 | | | | U | |

Table 1.5 (cont.) Results of the Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Sample ID | 17679 | | | | |
|---------------------|--------------|------------|-----------|-----------------------|-------|
| Location | Lot Blank | | | | |
| Volume of Air (L) | 0 | | | | |
| Analyte | Result pg | EMPC pg | MDL pg | Adjusted Conc (pg) | TEF |
| 2,3,7,8-TCDD | | 1.70 | 10.0 | U | 1 |
| 1,2,3,7,8-PeCDD | | 1.02 | 10.0 | U | 0.5 |
| 1,2,3,4,7,8-HxCDD | | 1.06 | 25.0 | U | 0.1 |
| 1,2,3,6,7,8-HxCDD | | 0.500 | 25.0 | U | 0.1 |
| 1,2,3,7,8,9-HxCDD | 0.660 J | | 25.0 | 0.0660 | 0.1 |
| 1,2,3,4,6,7,8-HpCDD | | 0.330 | 25.0 | U | 0.01 |
| OCDD | | 20.9 | 50.0 | U | 0.001 |
| Total Tetra-Dioxins | U | | | | |
| Total Penta-Dioxins | U | | | | |
| Total Hexa-Dioxins | 0.660 | | | | |
| Total Hepta-Dioxins | U | | | | |
| 2,3,7,8-TCDF | U | | 10.0 | U | 0.1 |
| 1,2,3,7,8-PeCDF | U | | 10.0 | U | 0.05 |
| 2,3,4,7,8-PeCDF | U | | 10.0 | U | 0.5 |
| 1,2,3,4,7,8-HxCDF | | 1.20 | 25.0 | U | 0.1 |
| 1,2,3,6,7,8-HxCDF | | 0.440 | 25.0 | U | 0.1 |
| 1,2,3,7,8,9-HxCDF | U | | 25.0 | U | 0.1 |
| 2,3,4,6,7,8-HxCDF | U | | 25.0 | U | 0.1 |
| 1,2,3,4,6,7,8-HpCDF | | 5.64 | 25.0 | U | 0.01 |
| 1,2,3,4,7,8,9-HpCDF | U | | 25.0 | U | 0.01 |
| OCDF | 4.22 J | | 50.0 | 0.00442 | 0.001 |
| Total Tetra-Furans | U | | | | |
| Total Penta-Furans | U | | | | |
| Total Hexa-Furans | U | | | | |
| Total Hepta-Furans | U | | | | |
| Total | | | | 0.0702 | |

QA/QC for PAH in Air

Results of the BS/BSD Analysis for PAH in Air

An XAD lot blank and a lot blank filter were chosen for the blank spike/blank spike duplicate (BS/BSD) analyses. The percent recoveries, for the XAD lot blank, ranging from 95 to 132, are listed in Table 2.1. The relative percent differences, also listed in Table 2.1, ranged from zero (0) to 6. The percent recoveries, for the lot blank filter, ranging from 82 to 96, are also listed in Table 2.1. The relative percent differences, also listed in Table 2.1, ranged from zero (0) to 4. QC limits are not available for either the percent recoveries or the relative percent differences for this analysis.

Table 2.1 Results of BS/BSD Analysis for PAH in Air
WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample ID: XAD Spike

| Compound | Spike Added µg | BS Rec. µg | % Rec. | BSD Rec. µg | % Rec. | RPD |
|-------------------------|----------------------|------------------|--------|-------------------|--------|-----|
| Naphthalene | 50 | 50.02 | 100 | 50.73 | 101 | 1 |
| 2-Methylnaphthalene | 50 | 50.11 | 100 | 50.87 | 102 | 2 |
| 1-Methylnaphthalene | 50 | 52.95 | 106 | 53.81 | 108 | 2 |
| Biphenyl | 50 | 50.63 | 101 | 51.75 | 103 | 2 |
| 2,6-Dimethylnaphthalene | 50 | 49.41 | 99 | 50.51 | 101 | 2 |
| Acenaphthylene | 50 | 49.07 | 98 | 49.70 | 99 | 1 |
| Acenaphthene | 50 | 49.56 | 99 | 49.02 | 98 | 1 |
| Dibenzofuran | 50 | 50.37 | 101 | 50.47 | 101 | 0 |
| Fluorene | 50 | 50.12 | 100 | 49.76 | 100 | 1 |
| Phenanthrene | 50 | 47.69 | 95 | 47.58 | 95 | 0 |
| Anthracene | 50 | 47.56 | 95 | 48.93 | 98 | 3 |
| Carbazole | 50 | 49.10 | 98 | 48.48 | 97 | 1 |
| Fluoranthene | 50 | 50.81 | 102 | 51.52 | 103 | 1 |
| Pyrene | 50 | 50.59 | 101 | 50.28 | 101 | 1 |
| Benzo(a)anthracene | 50 | 51.55 | 103 | 51.53 | 103 | 0 |
| Chrysene | 50 | 65.94 | 132 | 66.08 | 132 | 0 |
| Benzo(b)fluoranthene | 50 | 51.85 | 104 | 51.52 | 103 | 1 |
| Benzo(k)fluoranthene | 50 | 51.71 | 103 | 54.93 | 110 | 6 |
| Benzo(e)pyrene | 50 | 52.94 | 106 | 53.19 | 106 | 1 |
| Benzo(a)pyrene | 50 | 55.93 | 112 | 52.47 | 105 | 6 |
| Indeno(1,2,3-cd)pyrene | 50 | 53.51 | 107 | 53.66 | 107 | 0 |
| Dibenzo(a,h)anthracene | 50 | 54.32 | 109 | 53.84 | 108 | 1 |
| Benzo(g,h,i)perylene | 50 | 53.69 | 107 | 53.45 | 107 | 1 |

Table 2.1 (cont.) Results of BS/BSD Analysis for PAH in Air
 WA # 0-110 Air Monitoring, Sampling, and Modeling Support, and Underwater Survey Activities

Sample ID: Filter Spike

| Compound | Spike Added μg | BS Rec. μg | % Rec. | BSD Rec. μg | % Rec. | RPD |
|-------------------------|----------------------|------------------|--------|-------------------|--------|-----|
| Naphthalene | 50 | 46.32 | 93 | 45.88 | 92 | 1 |
| 2-Methylnaphthalene | 50 | 45.10 | 90 | 44.80 | 90 | 0.7 |
| 1-Methylnaphthalene | 50 | 48.10 | 96 | 48.24 | 96 | 0 |
| Biphenyl | 50 | 45.90 | 92 | 45.20 | 90 | 2 |
| 2,6-Dimethylnaphthalene | 50 | 44.46 | 89 | 44.94 | 90 | 1 |
| Acenaphthylene | 50 | 44.10 | 88 | 44.40 | 89 | 1 |
| Acenaphthene | 50 | 46.78 | 94 | 44.98 | 90 | 4 |
| Dibenzofuran | 50 | 45.64 | 91 | 44.70 | 89 | 2 |
| Fluorene | 50 | 45.00 | 90 | 44.88 | 90 | 0 |
| Phenanthrene | 50 | 45.10 | 90 | 45.50 | 91 | 1 |
| Anthracene | 50 | 47.32 | 95 | 47.26 | 95 | 0 |
| Carbazole | 50 | 45.02 | 90 | 44.36 | 89 | 1 |
| Fluoranthene | 50 | 44.94 | 90 | 44.38 | 89 | 1 |
| Pyrene | 50 | 45.80 | 92 | 44.82 | 90 | 2 |
| Benzo(a)anthracene | 50 | 44.36 | 89 | 43.46 | 87 | 2 |
| Chrysene | 50 | 46.96 | 94 | 46.74 | 93 | 1 |
| Benzo(b)fluoranthene | 50 | 42.18 | 84 | 40.78 | 82 | 3 |
| Benzo(k)fluoranthene | 50 | 43.64 | 87 | 45.02 | 90 | 3 |
| Benzo(e)pyrene | 50 | 43.28 | 87 | 42.50 | 85 | 2 |
| Benzo(a)pyrene | 50 | 44.12 | 88 | 42.68 | 85 | 3 |
| Indeno(1,2,3-cd)pyrene | 50 | 43.02 | 86 | 41.78 | 84 | 3 |
| Dibenzo(a,h)anthracene | 50 | 43.04 | 86 | 41.74 | 83 | 3 |
| Benzo(g,h,i)perylene | 50 | 42.92 | 86 | 42.88 | 86 | 0 |

QA/QC for Inorganic Acids in Air

Results of the BS/BSD Analysis for Inorganic Acids in Air

A blank spike/blank spike duplicate analysis (BS/BSD) was run. The percent recoveries, listed in Table 2.2, ranged from 83 to 105. All twelve values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.2, ranged from 0 (zero) to 9. QC limits are not available for the RPD.

Table 2.2 Results of the BS/BSD Analysis for Inorganic Acids in Air
 WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Analyte | Original Conc | | Recovered Conc | | % Recovery | | RPD | Recommended QC Limits % Rec |
|-------------------|---------------|-----------|----------------|-----------|------------|-----|-----|-----------------------------------|
| | Spike mg | Dup mg | Spike mg | Dup mg | Spike | Dup | | |
| Hydrobromic acid | 0.0808 | 0.0808 | 0.0843 | 0.0845 | 104 | 105 | 0 | 75-125 |
| Hydrochloric acid | 0.0413 | 0.0413 | 0.0423 | 0.0414 | 102 | 100 | 2 | 75-125 |
| Hydrofluoric acid | 0.0211 | 0.0211 | 0.0191 | 0.0175 | 91 | 83 | 9 | 75-125 |
| Nitric acid | 0.0812 | 0.0812 | 0.0808 | 0.0808 | 100 | 100 | 0 | 75-125 |
| Phosphoric acid | 0.118 | 0.118 | 0.110 | 0.115 | 93 | 97 | 4 | 75-125 |
| Sulfuric acid | 0.0817 | 0.0817 | 0.0846 | 0.0852 | 104 | 104 | 1 | 75-125 |

QA/QC for Metals in Air

Results of the BS/BSD Analysis for Metals in Air

A blank spike/blank spike duplicate analysis (BS/BSD) was run. The percent recoveries, listed in Table 2.3, ranged from 8 to 134. Fifty out of fifty-eight values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.3, ranged from 0 (zero) to 20. QC limits are not available for the RPD.

Results of the Analysis of the Laboratory Control Sample for Metals in Air

A laboratory control sample was also analyzed. The percent recoveries ranged from 7 to 116 and are listed in Table 2.4. Twenty-eight out of twenty-nine concentrations were within the acceptable QC limits.

Table 2.3 Results of the BS/BSD Analysis for Metals in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities

| Metal | Lot | Original Conc | | Recovered Conc | | % Recovery | | RPD | Recommended QC Limit % Rec |
|------------|------|----------------------------|---------------------------|---------------------------|------------------|------------|-----|-----|----------------------------------|
| | | Blank Conc µg/filter | Spike Dup µg/filter | Spike Dup µg/filter | Dup µg/filter | Spike | Dup | | |
| Aluminum | U | 40.00 | 40.00 | 44.49 | 44.56 | 111 | 111 | 0 | 75-125 |
| Arsenic | U | 40.00 | 40.00 | 43.98 | 44.17 | 110 | 110 | 0 | 75-125 |
| Beryllium | U | 1.00 | 1.00 | 1.04 | 1.05 | 104 | 105 | 0 | 75-125 |
| Cadmium | U | 1.00 | 1.00 | 1.09 | 1.09 | 109 | 109 | 0 | 75-125 |
| Calcium | 6.2 | 1000 | 1000 | 1058 | 1062 | 105 | 106 | 0 | 75-125 |
| Chromium | 0.51 | 4.00 | 4.00 | 5.07 | 5.29 | 114 | 120 | 5 | 75-125 |
| Cobalt | U | 10.00 | 10.00 | 10.29 | 10.37 | 103 | 104 | 1 | 75-125 |
| Copper | U | 5.00 | 5.00 | 5.40 | 5.43 | 108 | 109 | 1 | 75-125 |
| Iron | 1.6 | 20.00 | 20.00 | 21.29 | 21.99 | 98 | 102 | 3 | 75-125 |
| Lead | U | 10.00 | 10.00 | 10.69 | 10.74 | 107 | 107 | 0 | 75-125 |
| Lithium | U | 40.00 | 40.00 | 43.52 | 44.06 | 109 | 110 | 1 | 75-125 |
| Magnesium | U | 1000 | 1000 | 1069 | 1072 | 107 | 107 | 0 | 75-125 |
| Manganese | U | 10.00 | 10.00 | 10.65 | 10.72 | 107 | 107 | 1 | 75-125 |
| Molybdenum | U | 40.00 | 40.00 | 44.04 | 44.30 | 110 | 111 | 1 | 75-125 |
| Nickel | U | 10.00 | 10.00 | 10.52 | 10.50 | 105 | 105 | 0 | 75-125 |
| Phosphorus | U | 40.00 | 40.00 | 50.42 | 50.81 | 126 | 127 | 1 | 75-125 |
| Platinum | U | 40.00 | 40.00 | 42.03 | 41.60 | 105 | 104 | 1 | 75-125 |
| Selenium | U | 40.00 | 40.00 | 42.56 | 42.44 | 106 | 106 | 0 | 75-125 |
| Silver | U | 1.00 | 1.00 | 1.10 | 1.11 | 110 | 111 | 0 | 75-125 |
| Sodium | 7.3 | 1000 | 1000 | 974 | 983 | 97 | 98 | 1 | 75-125 |
| Tellurium | U | 40.00 | 40.00 | 3.29 | 3.44 | 8 | 9 | 4 | 75-125 |
| Thallium | U | 40.00 | 40.00 | 44.27 | 44.02 | 111 | 110 | 1 | 75-125 |
| Tin | U | 40.00 | 40.00 | 52.36 | 52.70 | 131 | 132 | 1 | 75-125 |
| Titanium | U | 40.00 | 40.00 | 42.70 | 42.51 | 107 | 106 | 0 | 75-125 |
| Vanadium | U | 10.00 | 10.00 | 10.50 | 10.56 | 105 | 106 | 1 | 75-125 |
| Yttrium | U | 40.00 | 40.00 | 43.51 | 43.31 | 109 | 108 | 0 | 75-125 |
| Zinc | 1.1 | 10.00 | 10.00 | 11.43 | 11.41 | 103 | 103 | 0 | 75-125 |
| Zirconium | U | 40.00 | 40.00 | 53.59 | 52.21 | 134 | 131 | 3 | 75-125 |
| Tungsten | U | 40.00 | 40.00 | 42.05 | 34.50 | 105 | 86 | 20 | 75-125 |

**Table 2.4 Results of the Analysis of the
Laboratory Control Sample for Metals in Air
WA # 0-110 Air Monitoring, Sampling, Analysis,
and Modeling Support, and Underwater Survey Activities**

| Metal | Analyzed Value µg/L | Accepted Value µg/L | % Rec | QC Limits % Rec |
|------------|---------------------------|---------------------------|----------|--------------------|
| Aluminum | 4301.16 | 4000 | 108 | 80-120 |
| Arsenic | 4299.25 | 4000 | 108 | 80-120 |
| Beryllium | 103.75 | 100 | 104 | 80-120 |
| Cadmium | 107.18 | 100 | 107 | 80-120 |
| Calcium | 52777.01 | 50000 | 106 | 80-120 |
| Chromium | 430.77 | 400 | 108 | 80-120 |
| Cobalt | 1025.19 | 1000 | 103 | 80-120 |
| Copper | 528.69 | 500 | 106 | 80-120 |
| Iron | 2038.85 | 2000 | 102 | 80-120 |
| Lead | 1058.99 | 1000 | 106 | 80-120 |
| Lithium | 2162.79 | 2000 | 108 | 80-120 |
| Magnesium | 53539.27 | 50000 | 107 | 80-120 |
| Manganese | 1055.69 | 1000 | 106 | 80-120 |
| Molybdenum | 2202.15 | 2000 | 110 | 80-120 |
| Nickel | 1024.3 | 1000 | 102 | 80-120 |
| Phosphorus | 2197.74 | 2000 | 110 | 80-120 |
| Platinum | 2123 | 2000 | 106 | 80-120 |
| Selenium | 4167.54 | 4000 | 104 | 80-120 |
| Silver | 107.28 | 100 | 107 | 80-120 |
| Sodium | 48612.46 | 50000 | 97 | 80-120 |
| Tellurium | 133.3 | 2000 | 7 | 80-120 |
| Thallium | 4446.75 | 4000 | 111 | 80-120 |
| Tin | 2324.71 | 2000 | 116 | 80-120 |
| Titanium | 2122.45 | 2000 | 106 | 80-120 |
| Vanadium | 1047.41 | 1000 | 105 | 80-120 |
| Yttrium | 2178.35 | 2000 | 109 | 80-120 |
| Zinc | 1061.42 | 1000 | 106 | 80-120 |
| Zirconium | 2269.74 | 2000 | 113 | 80-120 |
| Tungsten | 2035.00 | 2000 | 102 | 80-120 |

QA/QC for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

Results of the Internal Standard Recoveries for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

The internal standard percent recoveries, listed in Table 2.5, ranged from 70 to 138. One hundred and fifteen out of one hundred and seventeen values were within the acceptable QC limits.

Results of the BS/BSD Analysis for Polychlorinated Dibenzodioxins and Polychlorinated Dibenzofurans in Air

A blank was spiked in duplicate and analyzed. The percent recoveries, listed in Table 2.6, ranged from 76 to 114. All thirty-four values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.6, ranged from 0 (zero) to 16. All 17 RPDs were within the acceptable QC limits.

**Table 2.5 Results of the Internal Standard Recoveries for Polychlorinated Dibenzodioxin
and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support
and Underwater Survey Activities**

| Sample ID | Method Blank | 17670 | 17671 | 17672 | 17673 | 17674 | QC Limits |
|----------------------------|-----------------|-------|-------|-------|-------|-------|--------------|
| Units Internal Standard | % | % | % | % | % | % | |
| 13C-2,3,7,8-TCDD | 90 | 94 | 102 | 98 | 92 | 98 | 40-135 |
| 13C-1,2,3,6,7,8-HxCDD | 107 | 109 | 116 | 117 | 116 | 117 | 40-135 |
| 13C-2,3,7,8-TCDF | 91 | 103 | 103 | 108 | 108 | 112 | 40-135 |
| 13C-1,2,3,4,7,8-HxCDF | 111 | 114 | 104 | 100 | 121 | 122 | 40-135 |
| 13C-1,2,3,7,8-PeCDD | 98 | 121 | 118 | 118 | 135 | 133 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDD | 94 | 107 | 98 | 102 | 101 | 90 | 40-135 |
| 13C-1,2,3,7,8-PeCDF | 104 | 106 | 109 | 112 | 127 | 137 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDF | 84 | 105 | 92 | 90 | 108 | 108 | 40-135 |
| 13C-OCDD | 99 | 103 | 111 | 88 | 128 | 128 | 40-135 |

| Sample ID | 17675 | 17676 | 17677 | 17678 | 17679 | Blank Spike | Blank Spike Duplicate | QC Limits |
|----------------------------|-------|-------|-------|-------|-------|----------------|--------------------------|--------------|
| Units Internal Standard | % | % | % | % | % | % | % | |
| 13C-2,3,7,8-TCDD | 97 | 94 | 85 | 83 | 88 | 86 | 98 | 40-135 |
| 13C-1,2,3,6,7,8-HxCDD | 113 | 114 | 114 | 115 | 122 | 126 | 119 | 40-135 |
| 13C-2,3,7,8-TCDF | 112 | 110 | 100 | 96 | 91 | 95 | 109 | 40-135 |
| 13C-1,2,3,4,7,8-HxCDF | 117 | 127 | 118 | 118 | 109 | 105 | 101 | 40-135 |
| 13C-1,2,3,7,8-PeCDD | 126 | 138 | 114 | 113 | 110 | 111 | 115 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDD | 87 | 90 | 70 | 107 | 116 | 104 | 114 | 40-135 |
| 13C-1,2,3,7,8-PeCDF | 132 | 132 | 110 | 111 | 102 | 102 | 105 | 40-135 |
| 13C-1,2,3,4,6,7,8-HpCDF | 110 | 119 | 104 | 104 | 108 | 109 | 113 | 40-135 |
| 13C-OCDD | 127 | 121 | 130 | 110 | 116 | 128 | 123 | 40-135 |

Table 2.6 Results of the BS/BSD Analysis
for Polychlorinated Dibenzodioxin and Polychlorinated Dibenzofurans in Air
WA # 0-110 Air Monitoring, Sampling, Analysis, and Modeling Support
and Underwater Survey Activities

| Sample ID Blank Parameter Units | Spike pg | Blank Conc pg | BS Conc pg | Rec % | BSD Conc pg | Rec % | RPD | QC Limits % Rec | RPD |
|---------------------------------------|-------------|---------------------|------------------|----------|-------------------|----------|-----|--------------------------|-----|
| 2378-TCDD | 200 | U | 221 | 111 | 227 | 114 | 3 | 60-140 | 50 |
| 12378-PeCDD | 200 | U | 204 | 102 | 203 | 102 | 0 | 60-140 | 50 |
| 123478-HxCDD | 500 | U | 542 | 108 | 553 | 111 | 2 | 60-140 | 50 |
| 123678-HxCDD | 500 | U | 447 | 89 | 452 | 90 | 1 | 60-140 | 50 |
| 123789-HxCDD | 500 | 0.66 | 385 | 77 | 385 | 77 | 0 | 60-140 | 50 |
| 1234678-HpCDD | 500 | U | 555 | 111 | 513 | 103 | 8 | 60-140 | 50 |
| OCDD | 1000 | U | 940 | 94 | 873 | 87 | 7 | 60-140 | 50 |
| 2378-TCDF | 200 | U | 173 | 87 | 166 | 83 | 4 | 60-140 | 50 |
| 12378-PeCDF | 200 | U | 190 | 95 | 179 | 90 | 6 | 60-140 | 50 |
| 23478-PeCDF | 200 | U | 198 | 99 | 204 | 102 | 3 | 60-140 | 50 |
| 123478-HxCDF | 500 | U | 499 | 100 | 521 | 104 | 4 | 60-140 | 50 |
| 123678-HxCDF | 500 | U | 444 | 89 | 450 | 90 | 1 | 60-140 | 50 |
| 123789-HxCDF | 500 | U | 427 | 85 | 465 | 93 | 9 | 60-140 | 50 |
| 234678-HxCDF | 500 | U | 486 | 97 | 512 | 102 | 5 | 60-140 | 50 |
| 1234678-HpCDF | 500 | U | 468 | 94 | 459 | 92 | 2 | 60-140 | 50 |
| 1234789-HpCDF | 500 | U | 520 | 104 | 445 | 89 | 16 | 60-140 | 50 |
| OCDF | 1000 | 4.22 | 767 | 76 | 815 | 81 | 6 | 60-140 | 50 |

LOCKHEED MARTIN

Southwest Research Institute
PO Box 28510, 6220 Culebra Road
San Antonio, TX 78228-0510

Attn: Jo Ann Boyd

19 November 1999

Project # RIA-00011 APG Burn Support

As per Lockheed Martin / REAC Purchase Order GA91969J73, please analyze samples according to the following parameters:

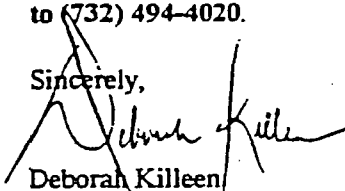
| Analysis Method | Matrix | # of samples |
|---|--------|--------------|
| Dioxin/ Furans / Modified TO9 | Air | 20 |
| Inorganic Acids / NIOSH 7903 | Air | 20 |
| Metals/ NIOSH 7300 | Air | 20 |
| Data package: Package with Diskette Deliverable | | |

Samples are expected to arrive at your laboratory between November 23-December 31, 1999. All applicable QA/QC (BS/BSD) analysis as per method, will be performed on our sample matrix. Preliminary sample and QC results tables plus a signed copy of our Chain of Custody must be faxed to REAC 10 business days after receipt of the last samples. The complete data package is due 21 business days after receipt of the last samples. The complete data package must include all items on the deliverables checklist. **Expect all samples to be difficult matrix and all raw data must be included in final analytical report.**

All sample and QC results (ie: BS/BSD, LCS, Duplicates, and Blanks) must be summarized in a Excel diskette deliverable.

Please submit all reports and technical questions concerning this project to John Johnson at (732) 321-4248 or fax to (732) 494-4020.

Sincerely,


Deborah Killeen
Data Validation and Report Writing Group Leader
Lockheed Martin / REAC Project

DK:jj Attachments

cc. R. Singhvi
D. Michunas
0011\mon\mem\9911\sub\0011Con

D. Miller
Subcontracting File
D. Angwenyi

C. Lentini
A. DuBois
D. Killeen

000010

APG Burn Compound List

NIOSH 7300 Metals

Aluminum
Arsenic
Beryllium
Cadmium
Calcium
Chromium
Cobalt
Copper
Iron
Lead
Lithium
Magnesium
Manganese
Molybdenum
Nickel
Phosphorus
Platinum
Selenium
Silver
Sodium
Tellurium
Tungsten
Thallium
Titanium
Vanadium
Yttrium
Zinc
Zirconium

Inorganic Acids
Hydrofluoric Acids
Hydrochloric Acids
Phosphoric acid
Hydrobromic acid
Nitric acid
Sulfuric acid

00040

~~(908)~~ 321-4200

68-099-223

Project Name: HPG Burn Support

Project Number: 24A00110

LM-RPW Contact: Amy D. B. S.

Phone: 732-494-4013

No: 06966

SHEET NO. 1 OF 1

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Volume(L) | Dioxins |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-----------|---------|
| ✓ | 17670 | DW3 | A | 4/6/02 | 1 | AMBER JAR/KE | 540 | ✓ |
| ✓ | 17671 | DW2 | | | | | 540 | ✓ |
| ✓ | 17672 | DW1 | | | | | 540 | ✓ |
| ✓ | 17673 | DW4 | | | | | Ø | ✓ |
| ✓ | 17674 | DW5 | | | | | 540 | ✓ |
| ✓ | 17675 | UW1 | | | | | 498 | ✓ |
| ✓ | 17676 | UW2 | | | | | 540 | ✓ |
| ✓ | 17677 | TRIP BLANK | | | | | Ø | ✓ |
| ✓ | 17678 | FIELD BLANK | | | | | Ø | ✓ |
| ✓ | 17679 | LOT BLANK | | | | | Ø | ✓ |
| — | — | BBS/BSD | ↓ | ↓ | 2 | ↓ | Ø | ✓ |

| | | | |
|------|--------------|------|---------------|
| SD - | Sediment | PW - | Potable Water |
| DS - | Drum Solids | GW - | Groundwater |
| DL - | Drum Liquids | SW - | Surface Water |
| X - | Other | SL - | Sludge |

Southwest Research Institute

S. Soil
W. Water
O. Oil
A. Air

Analysis for Dioxins/Furans

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF
CUSTODY #

| Items/Reason | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|---------|-------------|------|------|
| V/Anderson | [Signature] | 4/21/00 | | | |
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

| Items/Reason | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|------|-------------|------|------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

00017

68-099-223

4/11 RFW Contact: Amey Dubois

Phone: 732-494-4013

SHEET NO. 1 OF 1

Analyses Requested

FROM CHAIN OF
CUSTODY #

[illegible]

73-24908) 321-4200

EPA Contract ~~68-G4-0022~~ (12)

68-099-223

Project Name: APG Burn

Project Number: 21A00110

RFV Contact: Amey DuBois

Phone: 732-494-4013

No: 05654

SHEET NO. / OF /

041000

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | PAH | Volume(L) |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-----|-----------|
| 415 | 17690 | DW3 | A | 4/6/00 | 1 | WHIRLPAK/ICE/DARK | ✓ | 360 |
| 416 | 17691 | DW2 | ↓ | ↓ | ↓ | ↓ | ✓ | 351 |
| 417 | 17692 | DW1 | ↓ | ↓ | ↓ | ↓ | ✓ | 360 |
| 418 | 17693 | DW4 | ↓ | ↓ | ↓ | ↓ | ✓ | 369 |
| 419 | 17694 | DW5 | ↓ | ↓ | ↓ | ↓ | ✓ | 360 |
| 420 | 17695 | UW1 | ↓ | ↓ | ↓ | ↓ | ✓ | 332 |
| 421 | 17696 | UW2 | ↓ | ↓ | ↓ | ↓ | ✓ | 360 |
| 422 | 17697 | FIELD BLANK | ↓ | ↓ | ↓ | ↓ | ✓ | 0 |
| 423 | 17698 | TRIP BLANK | ↓ | ↓ | ↓ | ↓ | ✓ | 0 |
| 424 | 17699 | LOT BLANK | ↓ | ↓ | ↓ | ↓ | ✓ | 0 |
| | — | BS/BS D | ↓ | ↓ | 2 | ↓ | ✓ | 0 |

SD - Sediment
DS - Drum Solids
DL - Drum Liquids
X - Other

PW - Potable Water
GW - Groundwater
SW - Surface Water
SL - Sludge

S - Soil
W - Water
O - Oil
~~A - Air~~

Analysis for PAHs by NIOSH
5515

ALL tubes 226-30-06 LOT 1041

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF
CUSTODY #

DW - DOWNWIND
 LW - UPWIND

L - Liters

| Items/Reason | Relinquished By | Date | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|---------|--------------|---------|------|--------------|-----------------|---------|-------------|---------|------|
| All analysis | J McCall | 4/10/00 | David Angewy | 4/10/00 | 0950 | All Analysis | David Angewy | 4/10/00 | Chris Green | 4/10/00 | 1015 |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |

32 (908) 321-4200

EPA Contract 68-C4-0022-~~101~~
68-C99-223

CHAIN OF CUSTODY RECORD

Project Name: APG BURN Support

Project Number: R1A00110

LM ~~REV~~ Contact: Amy DuBois

Phone: 732-494-4023

No: 05254

SHEET NO. 1 OF 1

Sample Identification

METALS

Analyses Requested

[illegible]

Matrix:

SD - Sediment
DS - Drum Solids
DL - Drum Liquids
X - Other

PW - Potable Water
GW - Groundwater
SW - Surface Water
SL - Sludge

S - Soil
W - Water
O - Oil
A - Air

Special Instructions:

DW - DOWN WIND
 UW - UP WIND
 L - LIFE'S

FOR SUBCONTRACTING USE ONLY

FROM CHAIN OF
CUSTODY #

Analysis for Metals by NIOSH 7300
LOT # 517

[illegible]

APPENDIX C
SBC COM Clearances for GB, GD, VX, and HD
Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

Dubois, Amy E

From: Snyder Juanita A SBCCOM [juanita.snyder@SBCCOM.APGEA.ARMY.MIL]
Wednesday, April 12, 2000 12:55 PM
To: DUBOIS.AMY@EPAMAIL.EPA.GOV; salford@genphysics.com
Cc: axdean@CBDCOM-EMH1.APGEA.ARMY.MIL; dghall@CBDCOM-EMH1.APGEA.ARMY.MIL; fglattin@CBDCOM-EMH1.APGEA.ARMY.MIL; jasnyder@CBDCOM-EMH1.APGEA.ARMY.MIL; jefranch@CBDCOM-EMH1.APGEA.ARMY.MIL; rdmoore@CBDCOM-EMH1.APGEA.ARMY.MIL; sds smith@CBDCOM-EMH1.APGEA.ARMY.MIL; tablades@CBDCOM-EMH1.APGEA.ARMY.MIL; thomas.rosso@SBCCOM.APGEA.ARMY.MIL
Subject: J-field Clearances

DOC: Dubois, x (732)494-4013 J-Field
Item GVH background monitoring
04/06/00
W3 0004070050-M01 Clear for GB GD VX & HD
W2 0004070051-M01 Clear for GB GD VX & HD
W1 0004070052-M01 Clear for GB GD VX & HD
W4 0004070053-M01 Clear for GB GD VX & HD
W5 0004070054-M01 Clear for GB GD VX & HD
W3 0004070055-M01 Clear for GB GD VX & HD
W4 0004070056-M01 Clear for GB GD VX & HD
PA 0004070057-M01 Clear for GB GD VX & HD
P1 0004070058-M01 Clear for GB GD VX & HD

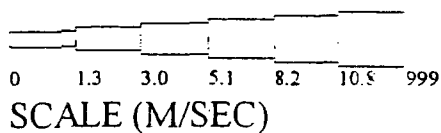
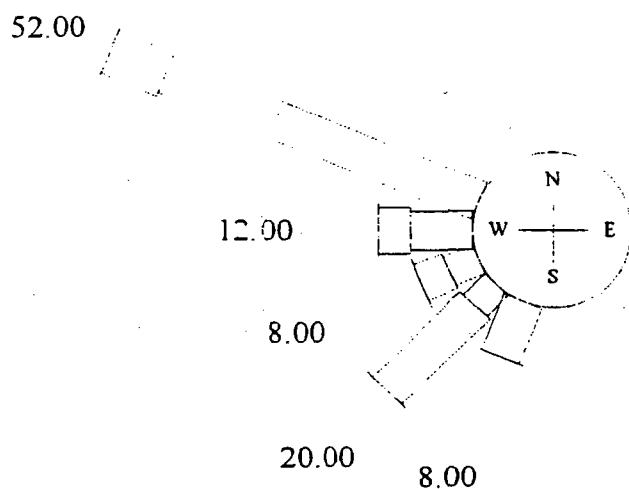
ita Snyder
ample Team

APPENDIX D

Windroses

Air Monitoring, Sampling, Analysis, and Modeling Support, and Underwater Survey Activities Site
July 2000

Aberdeen Proving Grounds Test Burn
Wind Rose Generated From H-Field Meteorological Data
4/6/00 14:30-20:30



| WIND SPEED (M/SEC) PERCENT OCCURRENCE | | | | | | | WIND SPEED (M/SEC) PERCENT OCCURRENCE | | | | | | |
|---------------------------------------|-------|---------|---------|---------|----------|-------|---------------------------------------|-------|---------|---------|---------|----------|-------|
| | 0-1.3 | 1.3-3.0 | 3.0-5.1 | 5.1-8.2 | 8.2-10.8 | >10.8 | | 0-1.3 | 1.3-3.0 | 3.0-5.1 | 5.1-8.2 | 8.2-10.8 | >10.8 |
| N | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | S | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| NNE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | SSW | 0.00 | 0.00 | 0.00 | 8.00 | 0.00 | 0.00 |
| NE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | SW | 0.00 | 0.00 | 4.00 | 16.00 | 0.00 | 0.00 |
| ENE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | WSW | 0.00 | 0.00 | 4.00 | 4.00 | 0.00 | 0.00 |
| E | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | W | 0.00 | 0.00 | 0.00 | 8.00 | 4.00 | 0.00 |
| ESE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | WNW | 0.00 | 0.00 | 0.00 | 44.00 | 8.00 | 0.00 |
| SE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | NW | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| SSE | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | NNW | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |

APPENDIX B

AIR SAMPLING EQUIPMENT PHOTOGRAPHS

Arrangement of Air Sampling Equipment



APG Controlled Burn Project – Analytes of Interest:

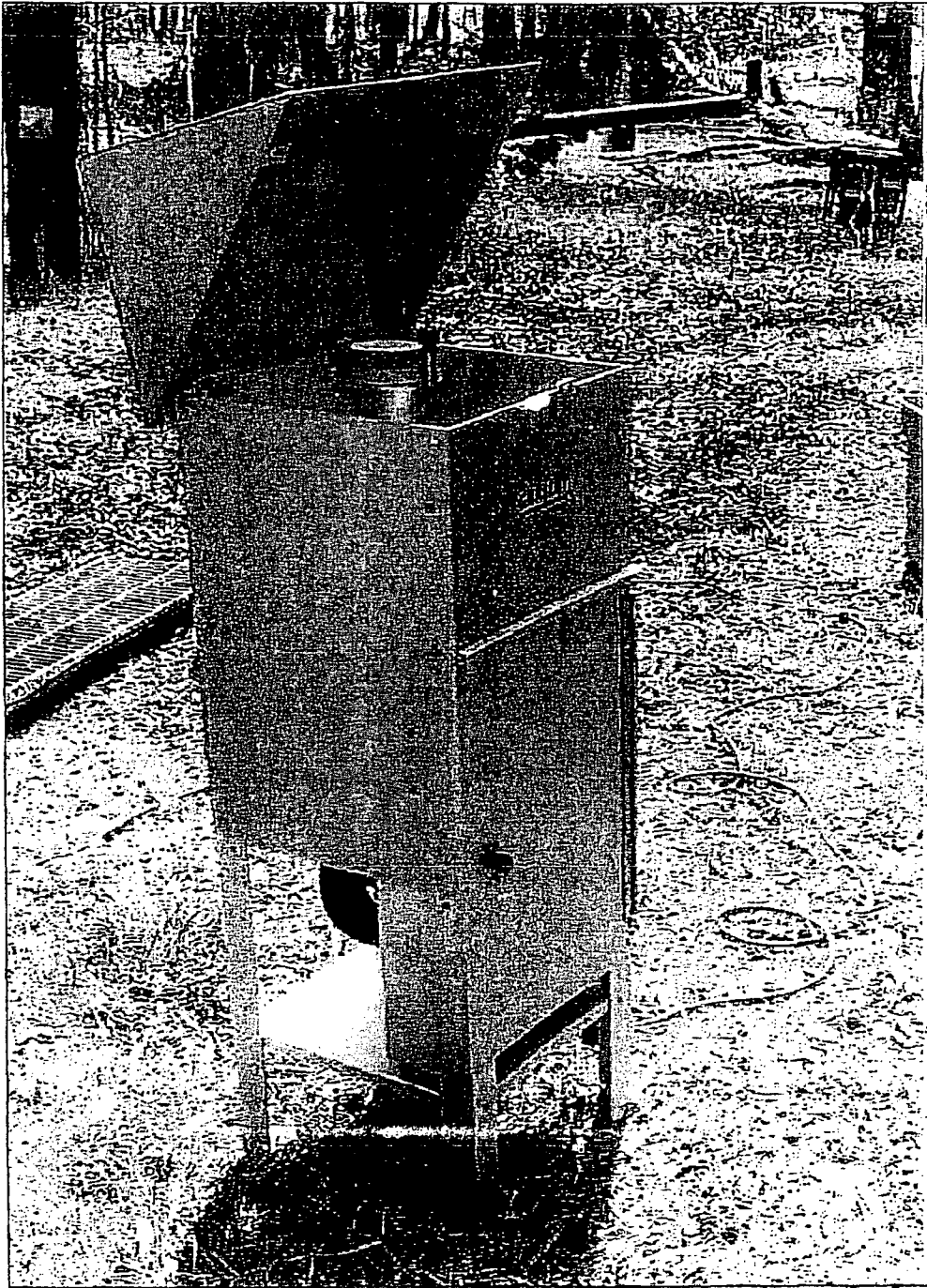
- Volatile Organic Compounds (VOCs)
- Pesticides and Polychlorinated Biphenyls (PCBs)
- Explosives
- Inorganics
- Radiologicals
- Chemical Agents

Summa Canister (left) and DAAMS Tube (right)



Summa canister for collection of air samples for volatile organic analysis. DAAMS tubes used for collection of samples for chemical agent analysis.

High Volume PUF Sampler

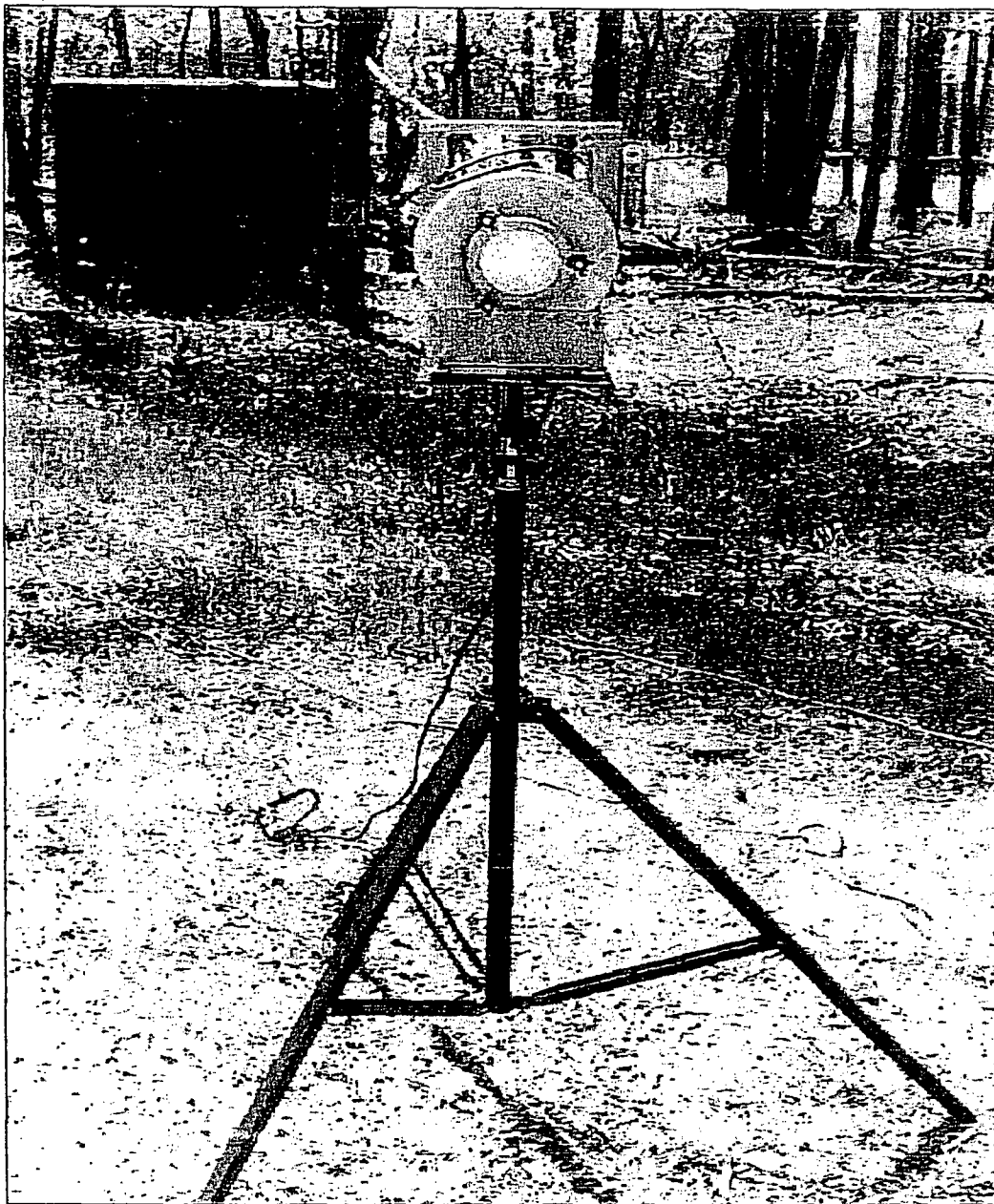


Sampler for collection of air samples for Pesticides, PCBs, and Explosives.

PUF – Polyurethane foam

PCB – Polychlorinated biphenyls

High Volume Sampler (Handi-Vol)



Sampler for collection of air samples for Inorganics and Radionuclides analyses

APPENDIX C

LIST OF ANALYTES FOR THE ABERDEEN PROVING GROUND CONTROLLED BURN PROJECT

Table C-1. TARGET ANALYTES FOR EACH ANALYSIS PERFORMED ON AIR SAMPLES FROM THE MAIN FRONT, NEW O-FIELD, AND J-FIELD CONTROLLED BURNS

| Volatiles ¹ | Pesticides/PCBs ² | Explosives ² | Metals (Total) | Radiologicals | Gamma Spectroscopy | Chemical Agents |
|---------------------------|--------------------------------------|----------------------------|----------------------|---------------|--------------------------|-----------------|
| Freon 12 | 4,4'-DDD | 1,3,5-Trinitrobenzene | Aluminum | Gross Alpha | Actinium-228 | Mustard (HD) |
| Chloromethane | 4,4'-DDE | 1,3-Dinitrobenzene | Antimony | Gross Beta | Bismuth-212 | Sarin (GB) |
| Freon 114 | 4,4'-DDT | 2,4,6-Trinitrotoluene | Arsenic | | Bismuth-214 | Soman (GD) |
| Chloroethene | Aldrin | 2,4-Dinitrotoluene | Barium | | Cesium-137 | VX |
| Bromomethane | Dieldrin | 2,6-Dinitrotoluene | Beryllium | | Cobalt-60 | |
| Chloroethane | Endosulfan I | 2-Amino-4,6-dinitrotoluene | Cadmium | | Lead-210 | |
| Freon 11 | Endosulfan II | 2-Nitrotoluene | Calcium | | Lead-212 | |
| 1,1-Dichloroethene | Endosulfan Sulfate | 3-Nitrotoluene | Chromium | | Lead-214 | |
| Methylene chloride | Endrin | 4-Amino-2,6-dinitrotoluene | Cobalt | | Potassium-40 | |
| Freon 113 | Endrin Aldehyde | 4-Nitrotoluene | Copper | | Protactinium-231 | |
| 1,1-Dichloroethane | Endrin Ketone | HMX | Iron | | Protactinium-234 | |
| cis-1,2-Dichloroethylene | Heptachlor | Nitrobenzene | Lead | | Radium-223 | |
| Chloroform | Heptachlor Epoxide | RDX | Magnesium | | Radium-224 | |
| 1,2-Dichloroethane | Lindane (gamma-BHC) | Tetryl | Manganese | | Radium-226 | |
| 1,1,1-Trichloroethane | Methoxychlor | | Mercury | | Uranium-235 | |
| Benzene | Toxaphene | | Nickel | | Uranium-233/234 | |
| Carbon tetrachloride | alpha-BHC | | Potassium | | Uranium-235/236 | |
| 1,2-Dichloropropane | alpha-Chlordane | | Selenium | | Uranium-238 ⁴ | |
| Trichloroethene | beta-BHC | | Silver | | | |
| cis-1,3-Dichloropropene | delta-BHC | | Sodium | | | |
| trans-1,3-Dichloropropene | gamma-Chlordane | | Thallium | | | |
| 1,1,2-Trichloroethane | 2-Chlorobiphenyl | | Uranium ³ | | | |
| Toluene | 2,3-Dichlorobiphenyl | | Vanadium | | | |
| 1,2-Dibromoethane | 2,2',5'-Trichlorobiphenyl | | Zinc | | | |
| Tetrachloroethene | 2,4',5'-Trichlorobiphenyl | | | | | |
| Chlorobenzene | 2,2',5,5'-Tetrachlorobiphenyl | | | | | |
| Ethylbenzene | 2,2',3,5'-Tetrachlorobiphenyl | | | | | |
| m-/p- Zylenes | 2,3',4,4'-Tetrachlorobiphenyl | | | | | |
| Styrene | 2,2',4,5,5'-Pentachlorobiphenyl | | | | | |
| o-Xylene | 2,2',3,4,5'-Pentachlorobiphenyl | | | | | |
| 1,1,2,2-Tetrachloroethane | 2,3,3',4',6-Pentachlorobiphenyl | | | | | |
| 1,3,5-Trimethylbenzene | 2,2',4,4',5,5',6-Hexachlorobiphenyl | | | | | |
| 1,2,4-Trimethylbenzene | 2,2',4,4',5,5'-Hexachlorobiphenyl | | | | | |
| 1,3-Dichlorobenzene | 2,2',3,4,5,5'-Hexachlorobiphenyl | | | | | |
| 1,4-Dichlorobenzene | 2,2',3,4,4'-Hexachlorobiphenyl | | | | | |
| 1,2-Dichlorobenzene | 2,2',3,4',5,5',6-Heptachlorobiphenyl | | | | | |
| 1,2,4-Trichlorobenzene | 2,2',3,4,4',5',6-Heptachlorobiphenyl | | | | | |
| Hexachlorobutadiene | 2,2',3,4,4',5,5'-Heptachlorobiphenyl | | | | | |
| | 2,2',3,3',4,4',5-Heptachlorobiphenyl | | | | | |
| | 2,2',3,3',4,4',5,6-Nonchlorobiphenyl | | | | | |

¹ Volatiles analysis on air samples collected during New O-Field and Main Front controlled burns was performed for the purposes of identifying only non-target peaks (a.k.a., Tentatively Identified Compounds (TICs)) and not the full range of TO-14 compounds.

² Analysis for these analytes was performed on PUF and filter samples from the Main Front controlled burn; analysis for these analytes on the air samples from the New O-Field and J-Field controlled burns was performed only on filter samples.

³ Analysis for Total Uranium was performed only on the air samples from the Main Front and New O-Field controlled burns.

⁴ Analysis for these radionuclides was performed only on the air samples from the J-Field controlled burn.

Table C-2. Volatile Organic Compound Detection Limits

| Compound | Method Detection Limit (MDL)* ppb | Compound | Method Detection Limit (MDL) * ppb |
|--------------------------|--------------------------------------|---------------------------|---------------------------------------|
| Freon 12 | ~ 0.2 | Cis-1,3-Dichloropropene | ~ 0.2 |
| Chloromethane | ~ 0.2 | Trans-1,3-Dichloropropene | ~ 0.2 |
| Freon 114 | ~ 0.2 | 1,1,2-Trichloroethane | ~ 0.2 |
| Chloroethene | ~ 0.2 | Toluene | ~ 0.2 |
| Bromomethane | ~ 0.2 | 1,2-Dibromoethane | ~ 0.2 |
| Chloroethane | ~ 0.2 | Tetrachloroethene | ~ 0.2 |
| Freon 11 | ~ 0.2 | Chlorobenzene | ~ 0.2 |
| 1,1-Dichloroethene | ~ 0.2 | Ethylbenzene | ~ 0.2 |
| Methylene chloride | ~ 0.2 | m-/p-Xylenes | ~ 0.2 |
| Freon 113 | ~ 0.2 | Styrene | ~ 0.2 |
| 1,1-Dichloroethane | ~ 0.2 | o-Xylene | ~ 0.2 |
| Cis-1,2-Dichloroethylene | ~ 0.2 | 1,1,2,2-Tetrachloroethane | ~ 0.2 |
| Chloroform | ~ 0.2 | 1,3,5-Trimethylbenzene | ~ 0.2 |
| 1,2-Dichloroethane | ~ 0.2 | 1,2,4-Trimethylbenzene | ~ 0.2 |
| 1,1,1-Trichloroethane | ~ 0.2 | 1,3-Dichlorobenzene | ~ 0.2 |
| Benzene | ~ 0.2 | 1,4-Dichlorobenzene | ~ 0.2 |
| Carbon tetrachloride | ~ 0.2 | 1,2-Dichlorobenzene | ~ 0.2 |
| 1,2-Dichloropropane | ~ 0.2 | 1,2,4-Trichlorobenzene | ~ 0.2 |
| Trichloroethene | ~ 0.2 | Hexachlorobutadiene | ~ 0.2 |

- Laboratory reports MDL of approximately 0.2 ppb for all compounds as adjusted for flow and sample volume.

Table C-3. Explosive Detection Limits

| Compound | Method Detection Limit (MDL) ug/L * | Compound | Method Detection Limit (MDL) ug/L * |
|-----------------------|--|----------------------------|--|
| 1,3,5-Trinitrobenzene | 0.11 | 3-Nitrotoluene | 0.16 |
| 2-Nitrotoluene | 0.09 | HMX | 0.23 |
| 4-Nitrotoluene | 0.21 | Tetryl | 0.22 |
| RDX | 0.23 | 2,4,6-Trinitrotoluene | 0.10 |
| 2,4-Dinitrotoluene | 0.10 | 2-Amino-4,6-Dinitrotoluene | 0.09 |
| 1,3-Dinitrobenzene | 0.09 | 4-Amino-2,6-Dinitrotoluene | 0.11 |
| 2,6-Dinitrotoluene | 0.13 | Nitrobenzene | 0.08 |

* The sample being analyzed is in a liquid matrix; therefore, the analytical MDL is expressed as ug/L.

Table C-4. Pesticides/Polychlorinated Biphenyls Detection Limits

| Compound | Method Detection Limit (MDL) ug/L * | Compound | Method Detection Limit (MDL) ug/L * |
|---------------------|--|--|--|
| 4,4'-DDD | 0.016 | Methoxychlor | 0.0096 |
| 4,4'-DDE | 0.0076 | Toxaphene | 0.39 |
| 4,4'-DDT | 0.0069 | 2,3-Dichlorobiphenyl | 0.021 |
| Aldrin | 0.010 | 2,2',5-Trichlorobiphenyl | 0.015 |
| alpha-BCH | 0.0085 | 2,4',5-Trichlorobiphenyl | 0.0080 |
| alpha-Chlordane | 0.0084 | 2,2',5,5'-Tetrachlorobiphenyl | 0.0084 |
| beta-BHC | 0.012 | 2,2',3,5'-Tetrachlorobiphenyl | 0.012 |
| delta-BHC | 0.0073 | 2,3',4,4'-Tetrachlorobiphenyl | 0.0070 |
| Dieldrin | 0.012 | 2,2',4,5,5'-Pentachlorobiphenyl | 0.0030 |
| Endosulfan I | 0.038 | 2,2',3,4,5'-Pentachlorobiphenyl | 0.0036 |
| Endosulfan II | 0.0097 | 2,3,3',4',6-Pentachlorobiphenyl | 0.0046 |
| Endosulfan Sulfate | 0.012 | 2,2',3,5,5',6-Hexachlorobiphenyl | |
| Endrin | 0.0068 | 2,2',4,4',5,5'-Hexachlorobiphenyl | 0.0045 |
| Endrin Aldehyde | 0.024 | 2,2',3,4,5,5'-Hexachlorobiphenyl | 0.0067 |
| Endrin Ketone | 0.0089 | 2,2',3,4,4',5'-Hexachlorobiphenyl | 0.0081 |
| gamma-BHC (Lindane) | 0.0065 | 2,2',3,4',5,5',6-Heptachlorobiphenyl | 0.014 |
| gamma-Chlordane | 0.020 | 2,2',3,4,4',5',6-Heptachlorobiphenyl | 0.012 |
| Heptachlor | 0.017 | 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 0.0033 |
| Heptachlor Epoxide | 0.0065 | 2,2',3,3',4,4',5-Heptachlorobiphenyl | 0.9992 |
| | | 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 0.0031 |

* The sample being analyzed is in a liquid matrix; therefore, the analytical MDL is expressed as ug/L.

Table C-5. Metals Detection Limits

| Compound | Method Detection Limit (MDL) ug/L * | Compound | Method Detection Limit (MDL) ug/L * |
|-----------|--|-----------|--|
| Aluminum | 176.0 | Magnesium | 82.4 |
| Antimony | 3.7 | Manganese | 3.5 |
| Arsenic | 1.5 | Mercury | 0.2 |
| Barium | 17.9 | Nickel | 6.4 |
| Beryllium | 0.7 | Potassium | 68.3 |
| Cadmium | 0.5 | Selenium | 2.0 |
| Calcium | 86.4 | Silver | 1.4 |
| Chromium | 11.9 | Sodium | 281.0 |
| Cobalt | 4.4 | Thallium | 3.0 |
| Copper | 1.9 | Vanadium | 4.5 |
| Iron | 114.0 | Zinc | 3.0 |
| Lead | 1.1 | | |

* The sample being analyzed is in a liquid matrix; therefore, the analytical MDL is expressed as ug/L.

Table C-6. Chemical Agents Detection Limits

| Compound | Method Detection Limit (MDL) (mg/m ³)* |
|--|---|
| Sarin (GB) | 0.0003 |
| Soman (GD) | 0.0003 |
| O-ethyl-S-(2-diisopropylaminoethyl)- methylphosphonothiolate (VX) | 0.0003 |
| Mustard (HD) | 0.003 |

* MDL/sensitivity is not uniformly defined or reported. The above MDL/sensitivity is based on a 2 – 3 hour sampling time and represents the information currently available. (Reference: Site Monitoring Concept Plan, U.S. Army Chemical Materiel Destruction Agency, 15 September 1993).

APPENDIX D

BURN EVENT PHOTOGRAPHS

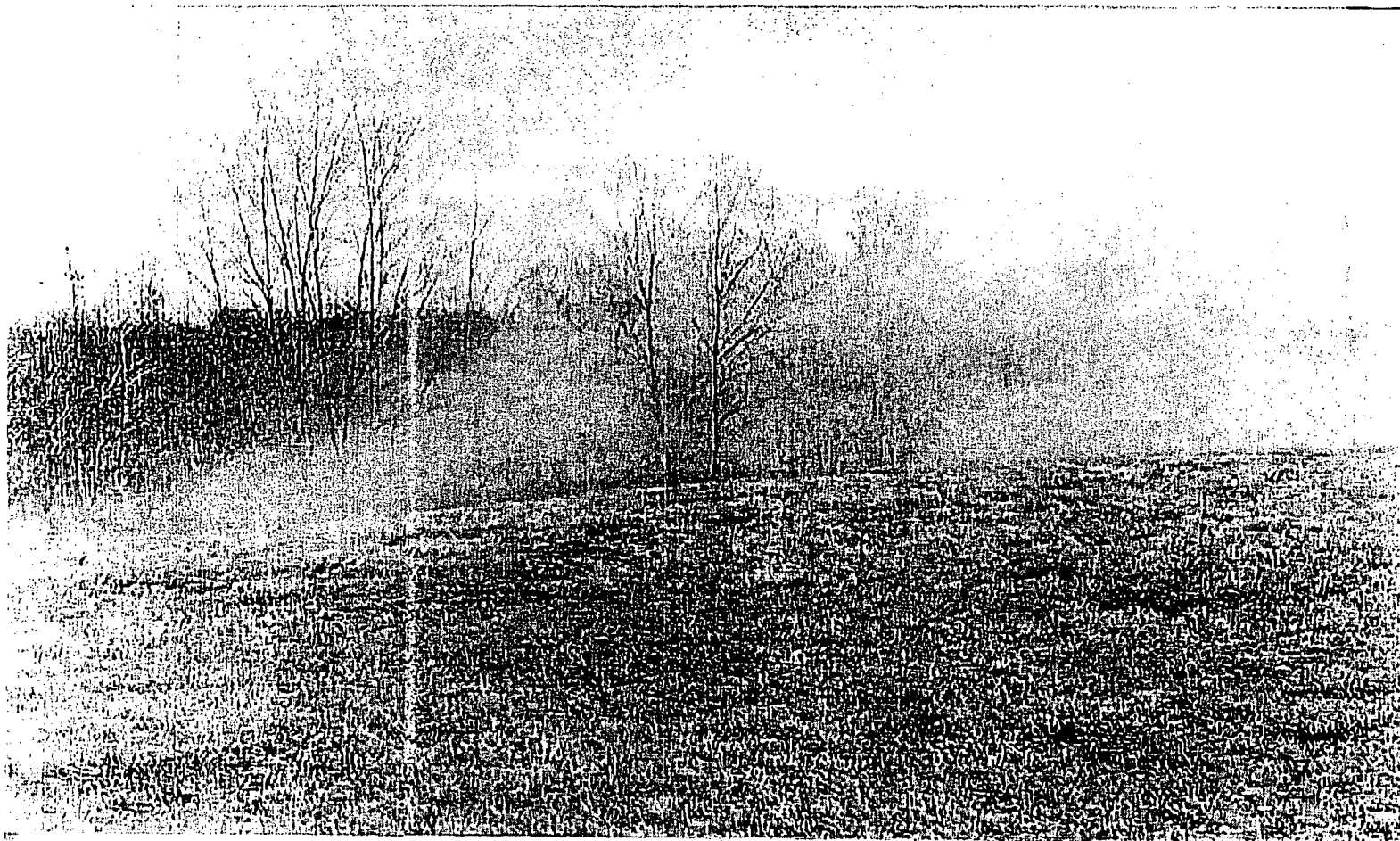
APPENDIX D-1

MAIN FRONT – 1999

Smoke Plume at the Beginning of the Main Front Range Controlled Burn – April 1999



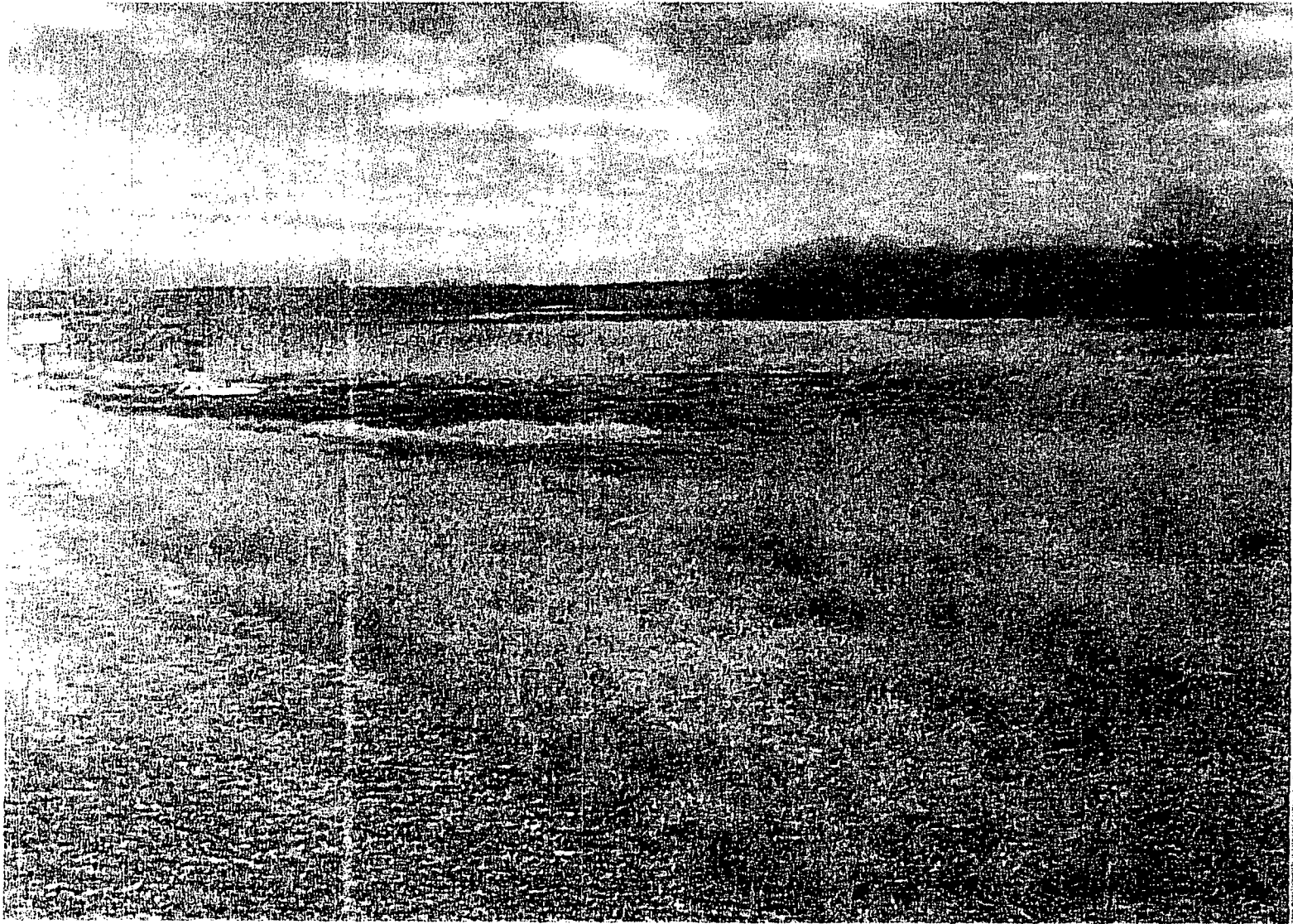
Smoke Plume at the Beginning of the Main Front Range Controlled Burn – April 1999



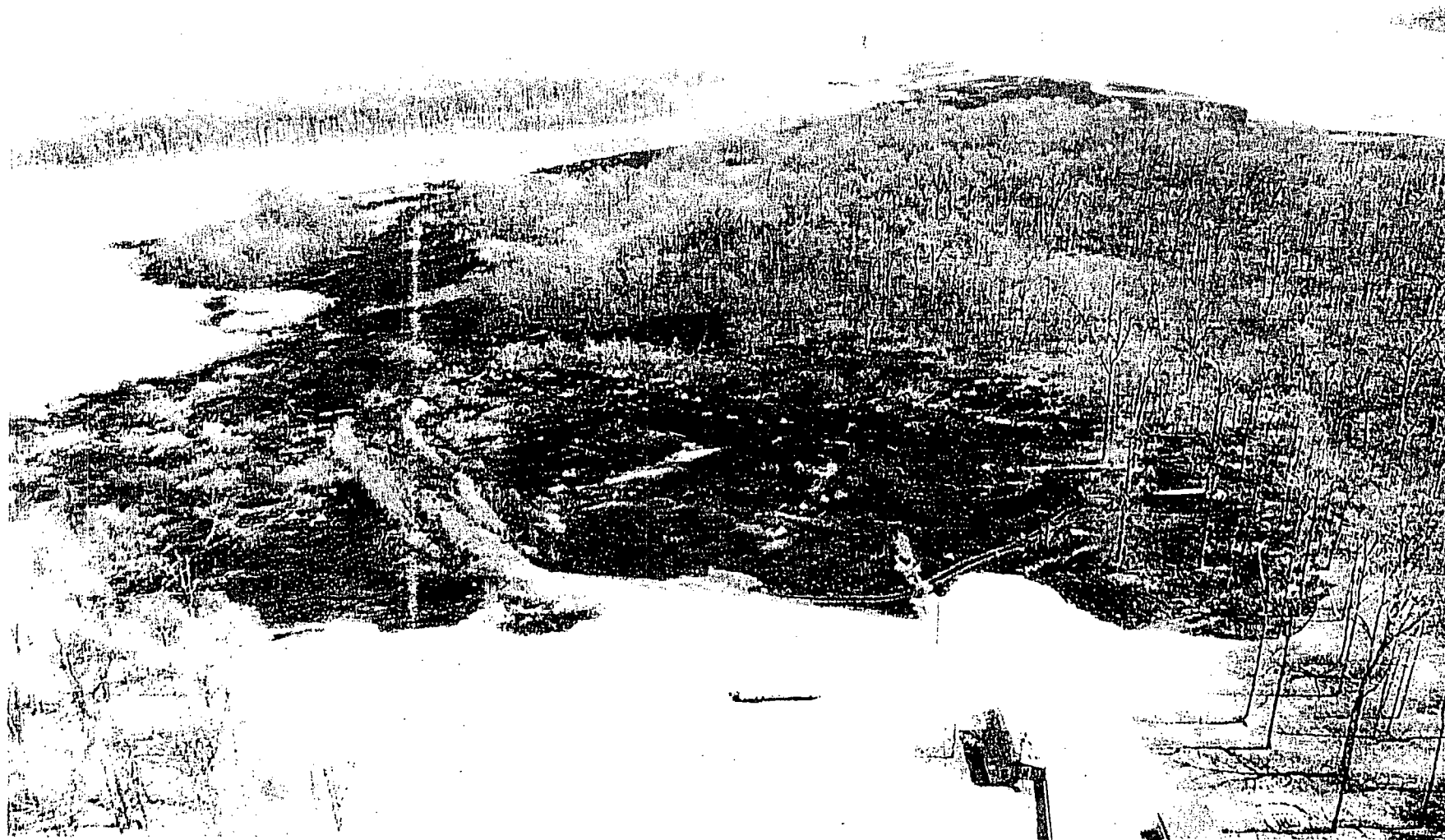
APPENDIX D-2

NEW O-FIELD – 1999

New O-Field Burn Area – Facing Watson Creek (Northeast)



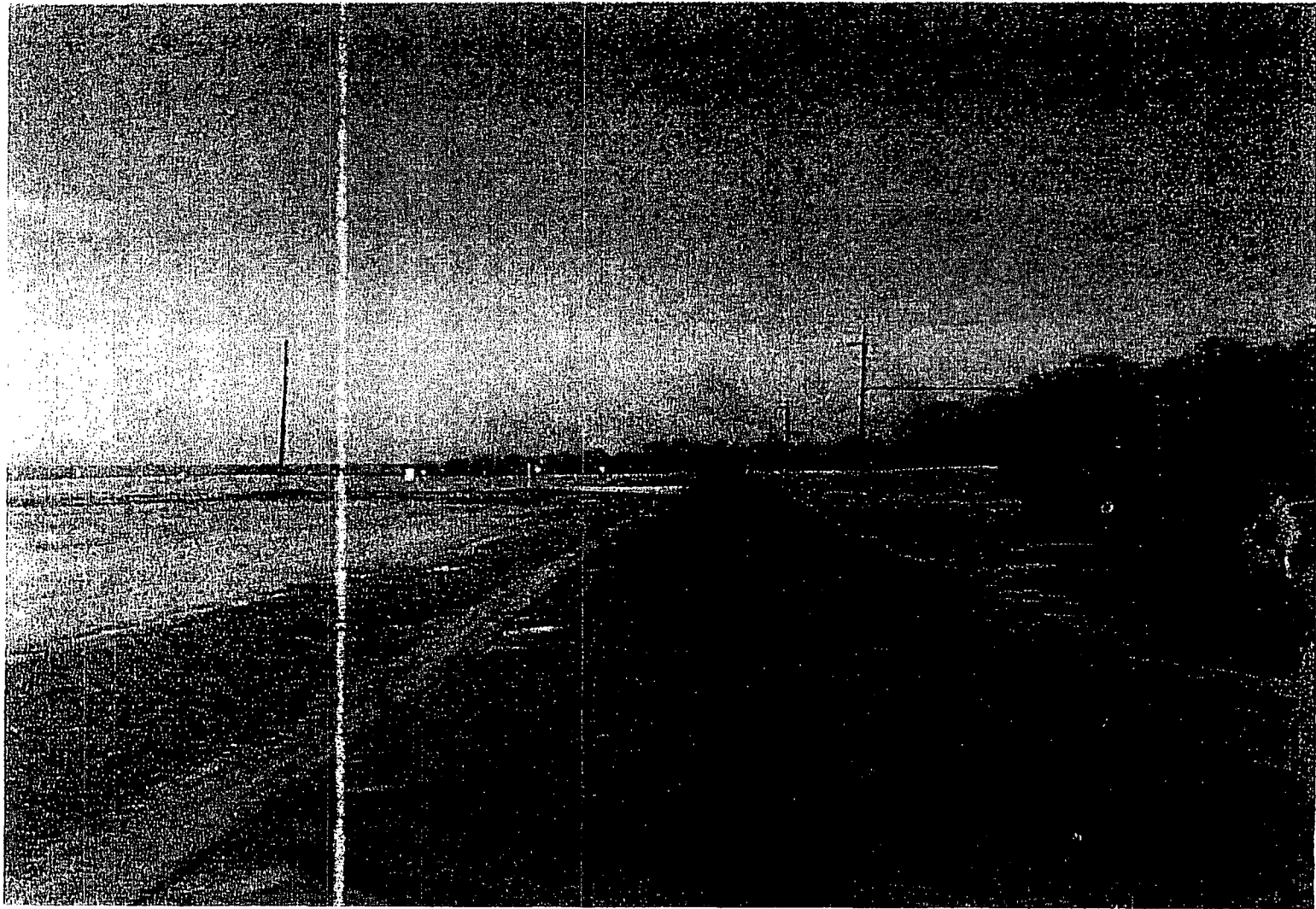
Aerial View – New O-Field Controlled Burn Area – December 1999



APPENDIX D-3

J-FIELD – 2000

Initial Smoke Plume Produced by the J-Field Controlled Burn – April 2000



Smoke Plume Produced by the J-Field Controlled Burn – April 2000



Aerial View of Area Burned During J-Field Controlled Burn – April 2000



APPENDIX E

CONTROLLED BURN DATA TABLES

APPENDIX E-1

DATA TABLES FOR THE MAIN FRONT CONTROLLED BURN – APRIL 1999

Table E-1. Main Front Controlled Burn Air Samples - April 1999

Volatile Organic Compound Analysis Results from Summa Canisters

| Analytes | Toxic Air Pollutant (Hour Screening Level) Concentration (ppm) | EPA Region III Ambient Air RBC Concentration (µg/m ³) | OSHA PELs Concentration (µg/m ³) | SP1 | | SP2 | | SP3 | | SP4 | |
|----------------------------|--|---|---|------------------------|---------------------------------------|------------------------|---------------------------------------|------------------------|---------------------------------------|------------------------|---------------------------------------|
| | | | | Concentration (ppb) | Concentration (µg/m ³) | Concentration (ppb) | Concentration (µg/m ³) | Concentration (ppb) | Concentration (µg/m ³) | Concentration (ppb) | Concentration (µg/m ³) |
| Acetone | 17,820 | 37 | 2,400,000 | 3.04 | 7.22 | 3.01 | 7.15 | 5.16 | 12.26 | 3.08 | 7.32 |
| Methylene chloride | NA | 3.8 | 86.843 | ND | ND | ND | ND | 1.69 | 5.87 | 1.85 | 6.43 |
| Toluene | NA | 42 | 753,703 | 0.91 | 3.42 | 2.31 | 8.71 | 42.83 | 161.4 | 8.56 | 32.26 |
| Octane | 17500 | NA | 2,350,000 | ND | ND | ND | ND | 2.94 | 13.7 | ND | ND |
| Nonane | NA | NA | NA | ND | ND | 4.1 | 21.5 | 20.2 | 106 | 4.05 | 21.25 |
| Decane | NA | NA | NA | ND | ND | 4.22 | 24.56 | 4.08 | 23.7 | 2.75 | 16 |
| m-/p-xylene | 6,510 | 730 | 435,000 | 1.77 | - | 7.12 | - | 32.13 | - | 30.8 | - |
| o-xylene | 6,510 | 730 | 435,000 | ND | ND | ND | ND | 3.89 | 16.9 | 3.83 | 16.63 |
| Unknown hydrocarbon | - | - | - | ND | ND | 24.59 | - | 18.05 | - | 4.37 | - |
| Unknown hydrocarbon | - | - | - | ND | ND | 2.71 | - | 3.48 | - | 18.75 | - |
| Unknown hydrocarbon | - | - | - | ND | ND | 2.11 | - | 3.76 | - | 2.23 | - |
| Unknown hydrocarbon | - | - | - | ND | ND | 1.61 | - | 3.13 | - | 6.13 | - |
| Unknown hydrocarbon | - | - | - | ND | ND | 1.54 | - | 2.56 | - | 8.08 | - |
| Unknown hydrocarbon | - | - | - | ND | ND | ND | ND | 3.25 | - | ND | ND |
| Benzaldehyde | NA | 73 | NA | 1.01 | 4.38 | ND | ND | ND | ND | ND | ND |
| Ethylhexanol | NA | NA | NA | 4.86 | - | ND | ND | ND | ND | ND | ND |
| Methylheptanone | NA | NA | NA | 0.87 | 4.56 | ND | ND | ND | ND | ND | ND |
| Methylbutane | NA | NA | NA | ND | ND | ND | ND | ND | ND | 19.92 | 58.78 |
| Hexamethylcyclotrisiloxane | NA | NA | NA | ND | ND | ND | ND | ND | ND | 7.49 | 68.01 |
| Trimethylbenzene | NA | 0.62 | NA | ND | ND | ND | ND | ND | ND | 2.55 | 12.54 |
| Dichlorobenzene | 60.12 | 0.28 | 300,000 | ND | ND | ND | ND | ND | ND | 1.56 | 9.40 |
| Total VOC | | | | 12.46 | | 53.32 | | 147.15 | | 126 | |

* The detected analytes were reported as Tentatively Identified Compounds (TICs)

NA - Screening criteria not available or does not apply

ND - nondetected

Shadowed cells indicate detected concentrations above screening criteria

Table E-2. Main Front Controlled Burn Air Samples - April 1999

Chemical Agent Analysis Results

| Chemical Agent | SP1 | | SP2 | | SP3 | | SP4 | |
|----------------|---------------------|--|---------------------|--|---------------------|--|---------------------|--|
| | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) |
| Sarin (GB) | ND | ND | ND | ND | ND | ND | ND | ND |
| Soman (GD) | ND | ND | ND | ND | ND | ND | ND | ND |
| VX | ND | ND | ND | ND | ND | ND | ND | ND |
| Mustard (HD) | ND | ND | ND | ND | ND | ND | ND | ND |

ND - nondetected

Analysis provided by Edgewood Chemical Biological Center.

Table E-3. Main Front Controlled Burn Air Samples - April 1999

Radiological Analysis Results

| Analytes | Toxic Air Pollutant 1-Hour Screening Level Concentration (ug/m ³) | EPA Region III Ambient Air RBCs Concentration (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|---------------------------------------|---|--|---|-------------------------|---|-------------------------|---|-------------------------|---|-------------------------|---|-------------------------|
| | | | | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) |
| Gross Alpha | NA | NA | NA | 6.8 | 0.0231 | 6.2 | 0.0210 | 1 | 0.0122 | 1.4 | 0.0133 | 5.6 |
| Gross Beta | NA | NA | NA | 37 | 0.1257 | 42 | 0.1424 | 0.6 | 0.0073 | 1.6 | 0.0152 | 32 |
| Actinium-228 | NA | NA | NA | -2.1 | -0.0071 | 1.6 | 0.0054 | 15 | 0.1833 | 2.1 | 0.0199 | 13 |
| Bismuth-212 | NA | NA | NA | 43 | 0.1461 | 42 | 0.1424 | -72 | -0.8797 | 13 | 0.1233 | 21 |
| Bismuth-214 | NA | NA | NA | -8.5 | -0.0289 | 3.7 | 0.0125 | 0.64 | 0.0078 | -7.4 | -0.0702 | -6.1 |
| Cesium-137 | NA | NA | NA | -0.72 | -0.0024 | -3.8 | -0.0129 | 3.6 | 0.0440 | 1.3 | 0.0123 | 3.7 |
| Cobalt-60 | NA | NA | NA | 0.88 | 0.0030 | 1.2 | 0.0041 | -0.41 | -0.0050 | 1.3 | 0.0123 | 0.71 |
| Lead-210 | NA | NA | NA | -46 | -0.1563 | -14 | -0.0475 | -70 | -0.8552 | -95 | -0.9012 | 10 |
| Lead-212 | NA | NA | NA | 0.37 | 0.0013 | 16 | 0.0543 | -2.5 | -0.0305 | -1.4 | -0.0133 | 5.6 |
| Lead-214 | NA | NA | NA | -3.6 | -0.0122 | 7.4 | 0.0251 | -5.3 | -0.0648 | -0.18 | -0.0017 | -3.6 |
| Potassium-40 | NA | NA | NA | 52 | 0.1767 | 4.1 | 0.0139 | -47 | -0.5742 | 22 | 0.2087 | 11 |
| Protactinium-231 | NA | NA | NA | -7.8 | -0.0265 | 20 | 0.0678 | 120 | 1.4661 | 97 | 0.9202 | 4.6 |
| Protactinium-234 | NA | NA | NA | 3.5 | 0.0119 | -23 | -0.0780 | 6.3 | 0.0770 | 3.4 | -0.0323 | -4.9 |
| Radium-223 | NA | NA | NA | -5.4 | -0.0184 | 1.4 | 0.0047 | -6.9 | -0.0843 | -1.4 | -0.0133 | 4.6 |
| Radium-224 | NA | NA | NA | 9 | 0.0306 | 180 | 0.6104 | 120 | 1.4661 | -53 | -0.5028 | 160 |
| Radium-226 | NA | NA | NA | -8.3 | -0.0282 | 3.6 | 0.0122 | 0.62 | 0.0076 | -7.2 | -0.0683 | -5.9 |
| Uranium-235 | NA | NA | NA | -5.6 | -0.0190 | -3.3 | -0.0112 | -14 | -0.1710 | 4.4 | 0.0417 | 12 |
| | | | | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) |
| Total Uranium by Mass Spectrometry | 6 | 11 | 50 | 2 | 0.0068 | 2 | 0.0068 | ND | ND | ND | ND | 2.2 |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|----------------------|-------------|-------------------------------------|
| SP1 | TSP3 | 294.26 |
| SP2 | TSP6 | 294.87 |
| SP3 | Handi Vol 6 | 81.85 |
| SP4 (Upwind) | Handi Vol 7 | 105.41 |

NA - Screening criteria not available or does not apply

ND - Nondetected

Table E-4. Main Front Controlled Burn Air Samples - April 1999

Inorganics Analysis Results

| Analytes | Toxic Air Pollutant 1-Hour Screening Level Concentration (ug/m ³) | EPA Region III Ambient Air RBCs Concentration (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|-----------|---|--|---|------------------------|--|------------------------|--|------------------------|--|------------------------|--|------------------------|
| | | | | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) |
| Mercury | 0.3 | 0.031 | 100 (acceptable ceiling) | 0.01 | 0.00003 | 0.01 | 0.00003 | BQL | - | BQL | - | BQL |
| Silver | NA | 1.8 | 10 | 2 | 0.0068 | 1.7 | 0.0058 | BQL | - | BQL | - | 3.1 |
| Aluminum | NA | 0.37 | 5,000 | 15000 | 51.1910 | 11800 | 40.0489 | 43.7 | 0.5955 | 53.5 | 0.4598 | 21700 |
| Arsenic | NA | 0.00041 | 500 | 4.3 | 0.0147 | 3.6 | 0.0122 | BQL | - | BQL | - | 6.5 |
| Barium | NA | 0.051 | 500 | 36900 | 125.9300 | 29000 | 98.4252 | 2.5 | 0.0341 | 2.7 | 0.0232 | 48900 |
| Beryllium | 0.1 | 0.00075 | 2 | 0.14 | 0.0005 | 0.11 | 0.0004 | BQL | - | BQL | - | 0.17 |
| Calcium | NA | NA | NA | 10400 | 35.4925 | 8340 | 28.3057 | 67.5 | 0.9199 | 71.5 | 0.6145 | 13800 |
| Cadmium | NA | 0.00099 | 5 | BQL | - | BQL | - | 0.04 | 0.0005 | BQL | - | BQL |
| Cobalt | NA | 22 | 100 | BQL | - | BQL | - | BQL | - | 0.06 | 0.0005 | BQL |
| Chromium | NA | 0.00015 | 500 | 11.4 | 0.0389 | 9.1 | 0.0309 | 0.59 | 0.0080 | 0.52 | 0.0045 | 18.3 |
| Copper | NA | 15 | 100 | 16.2 | 0.0553 | 11.9 | 0.0404 | 22.8 | 0.3107 | 4.9 | 0.0421 | BQL |
| Iron | NA | 110 | NA | 274 | 0.9351 | 230 | 0.7806 | 38.3 | 0.5219 | 42.1 | 0.3618 | 315 |
| Potassium | NA | NA | NA | 22200 | 75.7627 | 17100 | 58.0369 | 23.8 | 0.3243 | 30.8 | 0.2647 | 28800 |
| Magnesium | NA | NA | NA | 991 | 3.3820 | 799 | 2.7118 | 15.2 | 0.2071 | 19.4 | 0.1667 | 1300 |
| Manganese | NA | 0.0052 | 5000 | 10.3 | 0.0352 | 9.2 | 0.0312 | 1.4 | 0.0191 | 1.7 | 0.0146 | 11.3 |
| Sodium | NA | NA | NA | 66600 | 227.2882 | 48500 | 164.6077 | 186 | 2.5348 | 166 | 1.4267 | 81400 |
| Nickel | NA | 7.3 | 1,000 | 1.7 | 0.0058 | 1.4 | 0.0048 | 0.45 | 0.0061 | 0.41 | 0.0035 | 1.7 |
| Lead | NA | NA | 50 | 5 | 0.0171 | 3.9 | 0.0132 | 0.45 | 0.0061 | 0.44 | 0.0038 | 4.9 |
| Antimony | NA | 0.15 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Selenium | NA | 1.8 | 200 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Thallium | NA | 0.026 | 100 | 1.5000 | 0.0051 | 1.3 | 0.0044 | BQL | - | BQL | - | 2.5 |
| Vanadium | NA | 2.6 | 500 | 0.6900 | 0.0024 | 0.68 | 0.0023 | 0.29 | 0.0040 | 0.33 | 0.0028 | 0.58 |
| Zinc | 100 | 110 | NA | 29300.0000 | 99.9932 | 22500 | 76.3644 | 3.5 | 0.0477 | 4.1 | 0.0352 | 38100 |

| Sampling Location | Sampler ID | Total Air Flow (m ³ /min) | |
|-------------------|-------------|--------------------------------------|---------|
| SP1 | TSP1 | 293.02 | Mercury |
| SP1 | TSP2 | 297.77 | |
| SP2 | TSP4 | 294.64 | Mercury |
| SP2 | TSP5 | 294.64 | |
| SP3 | Handi Vol 1 | 73.38 | Mercury |
| SP3 | Handi Vol 4 | 88.92 | |
| SP4 (Upwind) | Handi Vol 2 | 116.35 | Mercury |
| SP4 (Upwind) | Handi Vol 5 | 114.11 | |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shadowed cells indicate detected concentrations above screening criteria.

Table E-5. Main Front Controlled Burn Air Samples - April 1999
PCBs Analysis Results from Filters

| Analytes | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|------------------------------------|------------------------|--|------------------------|--|------------------------|--|------------------------|--|------------------------|
| | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) |
| 2,2',3,5'-Tetrachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',5,5'-Tetrachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',5-Trichlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,4'-Hexachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,5'-Pentachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,5,5'-Hexachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,5,5',6'-Hexachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',4,5,5',6'-Pentachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3',4,4'-Tetrachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3',3',4',6'-Pentachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3-Dichlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4',5-Trichlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Chlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22'33'44'5'-Heptachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22'33'44'55'6'-Nonachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22'34'55'6'-Heptachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22'344'5'6'-Heptachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22'344'55'-Heptachlorobiphenyl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|-------------------------------------|
| SP1 | Filter7 | 53.39 |
| SP2 | Filter6 | 46.95 |
| SP3 | Filter1 | 37.78 |
| SP4 (Upwind) | Filter3 | 61.1 |

BQL - Below Quantitation Limit

Table E-6. Main Front Controlled Burn Air Samples - April 1999

PCBs Analysis Results from PUF Samplers

| Analytes | Toxic Air Pollutant 1-Hour Screening Level (ug/m ³) | EPA Region III Ambient Air RBCs (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|---|---|--|---|---------------------|--|---------------------|--|---------------------|--|---------------------|--|---------------------|
| | | | | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) |
| 2,2',3,5'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',5,5'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',5-Trichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,4',5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,5'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | 0.03 | 0.0008 | BQL | - | BQL |
| 2,2',3,4,5,5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,5,5',6'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',4,5,5'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3',4,4'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3,3',4',6'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3-Dichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4',5-Trichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | 0.43 | 0.0110 | 0.073 | 0.0012 | BQL |
| 2-Chlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,3',4,4',5'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,3',4,4',5,5',6'-Nonachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4',5,5',6'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,4',5',6'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,4',5,5',6'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|----------------------------------|
| SP1 | PUF7 | 53.39 |
| SP2 | PUF6 | 46.95 |
| SP3 | PUF1 | 37.78 |
| SP4 (Upwind) | PUF3 | 61.1 |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shadowed cells indicated detected concentrations above screening criteria

Table E-7. Main Front Controlled Burn Air Samples - April 1999

Pesticides Analysis Results from Filters

| Analytes | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|---------------------|------------------------|--|------------------------|--|------------------------|--|------------------------|--|------------------------|
| | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) |
| alpha-BHC | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| beta-BHC | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| delta-BHC | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Lindane (gamma-BHC) | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Heptachlor | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Aldrin | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Heptachlor epoxide | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan I | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Dieldrin | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDE | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan II | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDD | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan sulfate | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDT | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Methoxychlor | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin ketone | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin aldehyde | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| alpha-Chlordane | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| gamma-Chlordane | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Toxaphene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|----------------------------------|
| SP1 | Filter9 | 60.17 |
| SP2 | Filter8 | 41.09 |
| SP3 | Filter2 | 33.16 |
| SP4 (Upwind) | Filter4 | 45.89 |

BQL - Below Quantitation Limit

Table E-8. Main Front Controlled Burn Air Samples - April 1999

Pesticides Analysis Results from PUF Samplers

| Analytes | Toxic Air Pollutant | EPA Region III | OSHA PEL's Concentration | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|---------------------|--|--|--------------------------|---------------------|---------------------------------------|---------------------|---------------------------------------|---------------------|---------------------------------------|---------------------|---------------------------------------|---------------------|
| | 1-Hour Screening Level (ug/m ³) | Ambient Air RBCs (ug/m ³) | (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) |
| alpha-BHC | NA | NA | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| beta-BHC | NA | NA | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| delta-BHC | NA | NA | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Lindane (gamma-BHC) | NA | NA | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Heptachlor | NA | 0.0014 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Aldrin | NA | 0.00037 | 250 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Heptachlor epoxide | NA | 0.00069 | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan I | NA | 2.2 | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Dieldrin | NA | 0.00039 | 250 | 0.032 | 0.0005 | 0.02 | 0.0005 | 0.1 | 0.0030 | BQL | - | BQL |
| 4,4'-DDE | NA | 0.018 | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin | NA | 0.11 | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan II | NA | 2.2 | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDD | NA | 0.026 | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan sulfate | NA | NA | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDT | NA | 0.018 | 1,000 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Methoxychlor | NA | 1.8 | 1,500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin ketone | NA | NA | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin aldehyde | NA | NA | NA | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| alpha-Chlordane | NA | NA | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| gamma-Chlordane | NA | NA | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Toxaphene | NA | 0.0057 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|----------------------------------|
| SP1 | PUF9 | 60.17 |
| SP2 | PUF8 | 41.09 |
| SP3 | PUF2 | 33.16 |
| SP4 (Upwind) | PUF4 | 45.89 |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shadowed cells indicate detected concentrations above screening criteria

| Analytes | SP1 | | | | SP2 | | | | SP3 | | | |
|--------------|----------------------|-----------|------------------|----------------------|----------------------|-----------|------------------|----------------------|----------------------|-----------|------------------|----------------------|
| | Results (pCi/filter) | Error +/- | MDA (pCi/filter) | Validation Qualifier | Results (pCi/filter) | Error +/- | MDA (pCi/filter) | Validation Qualifier | Results (pCi/filter) | Error +/- | MDA (pCi/filter) | Validation Qualifier |
| Gross Alpha | 6.8 | 2.4 | 1.8 | U3, J6 | 6.2 | 2.3 | 2 | U3, J6 | 1 | 0.8 | 1 | U1, U3, J6 |
| Gross Beta | 37 | 3.2 | 1.9 | D | 42 | 3.4 | 2 | D | 0.6 | 0.9 | 1.5 | U1, U2, U3 |
| Bismuth-212 | 43 | 43 | 85 | U1, U2 * | 42 | 75 | 130 | U1, U2 * | -72 | 73 | 110 | U1, U2 * |
| Lead-212 | 0.37 | 7.5 | 7.3 | U1, U2 D | 16 | 16 | 14 | U1, U2 D | -2.5 | 9.2 | 15 | U1, U2 * |
| Potassium-40 | 52 | 48 | 51 | U1, J6 D | 4.1 | 83 | 180 | U1, U2 * | -47 | 75 | 160 | U1, U2 * |
| Radium-223 | -5.4 | 6.5 | 11 | U1, U2 * | 1.4 | 10 | 17 | U1, U2 D | -6.9 | 11 | 17 | U1, U2 * |
| Uranium-235 | -5.6 | 14 | 23 | U1, J6 * | -3.3 | 30 | 50 | U1, U2 * | -14 | 30 | 49 | U1, U2 * |

| Analytes | SP4 (Upwind) | | | | Blank | | | |
|--------------|----------------------|-----------|------------------|----------------------|----------------------|-----------|------------------|----------------------|
| | Results (pCi/filter) | Error +/- | MDA (pCi/filter) | Validation Qualifier | Results (pCi/filter) | Error +/- | MDA (pCi/filter) | Validation Qualifier |
| Gross Alpha | 1.4 | 0.8 | 1 | U3, J6 | 5.6 | 2.1 | 1.7 | J6 |
| Gross Beta | 1.6 | 1 | 1.5 | U3, J6 | 32 | 2.9 | 1.6 | D |
| Bismuth-212 | 13 | 42 | 79 | U1, U2 * | 21 | 69 | 120 | U1, U2 |
| Lead-212 | -1.4 | 4.9 | 9 | U1, U2 * | 5.6 | 9 | 15 | U1, U2 * |
| Potassium-40 | 22 | 57 | 65 | U1, U2 D | 11 | 120 | 80 | U1, U2 |
| Radium-223 | -1.4 | 6.4 | 11 | U1, U2 * | 4.6 | 12 | 18 | U1, U2 * |
| Uranium-235 | 4.4 | 4.2 | 24 | U1, U2 D | 12 | 29 | 50 | U1, U2 * |

* reported nondetected in 8 results

MDA - minimum detectable amount

D reported detected in 8 results

Qualifier

U1 - results less than MDA

U2 - results less than error

U3 - results less than blank

J6 - error greater than 20%

Table E-9. Main Front Controlled Burn Air Samples - April 1999
Explosives Analysis Results from PUF Samplers

| Analytes | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|----------------------------|---------------------|---------------------------------------|---------------------|---------------------------------------|---------------------|---------------------------------------|---------------------|---------------------------------------|---------------------|
| | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) | Concentration (ug/m ³) | Results (ug/PUF) |
| 1,3,5-Trinitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 1,3-Dinitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4,6-Trinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4-Dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,6-Dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Amino-4,6-dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 3-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4-Amino-2,6-dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| HMX | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Nitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| RDX | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Tetryl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

BQL - Below Quantitation Limit

Table E-10 Main Front Controlled Burn Air Samples - April 1999
Explosives Analysis Results from Filters

| Analytes | SP1 | | SP2 | | SP3 | | SP4 (Upwind) | | Blank |
|----------------------------|---------------------|------------------------------------|---------------------|------------------------------------|---------------------|------------------------------------|---------------------|------------------------------------|---------------------|
| | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) | Concentration (ug/m ³) | Results (ug/filter) |
| 1,3,5-Trinitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 1,3-Dinitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4,6-Trinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4-Dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,6-Dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Amino-4,6-dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 3-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4-Amino-2,6-dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| HMX | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Nitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| RDX | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Tetryl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

BQL - Below Quantitation Limit

APPENDIX E-2

DATA TABLES FOR THE NEW O-FIELD CONTROLLED BURN – DECEMBER 1999

Table E-11. New O-Field Controlled Burn Air Samples - December 1999
Volatile Organic Compound Analysis Results from Summa Canisters

| Analytes | Toxic Air Pollutants 12-Hour Screening Level Concentration (ug/m ³) | EPA Region II Ambient Air RBCs Concentration (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP2 | | SP3 | | SP4-A (Background) | | SP4-B (Background) | |
|---------------------------------|--|--|---|---------------------|------------------------------------|---------------------|------------------------------------|---------------------|------------------------------------|---------------------|------------------------------------|---------------------|------------------------------------|
| | | | | Concentration (ppb) | Concentration (ug/m ³) | Concentration (ppb) | Concentration (ug/m ³) | Concentration (ppb) | Concentration (ug/m ³) | Concentration (ppb) | Concentration (ug/m ³) | Concentration (ppb) | Concentration (ug/m ³) |
| Acetone | 17,820 | 37 | 2,400,000 | 1.68 | 3.99 | 27.6 | 65.56 | 1.11 | 2.64 | ND | - | 1.51 | 3.59 |
| Benzene | 80 | 0.22 | 3,195 | 0.806 | 2.58 | 6.23 | 19.9 | 0.727 | 2.32 | 1.42 | 4.54 | 1.54 | 4.92 |
| Benzonitrile | NA | NA | NA | 0.609 | 2.57 | ND | ND | ND | ND | ND | ND | ND | ND |
| Carbon Dioxide | 2,088,000 | NA | 9,000,000 | ND | ND | ND | ND | ND | ND | ND | ND | ND | ND |
| Carbon Disulfide | NA | 73 | 62,275 | ND | ND | 5.94 | 18.5 | ND | ND | ND | ND | ND | ND |
| Dodecene | NA | NA | NA | 2.9 | 19.96 | ND | ND | ND | ND | ND | ND | ND | ND |
| Hexane | NA | 21 | 1,800,000 | ND | ND | 11.3 | 39.83 | ND | ND | ND | ND | ND | ND |
| Methylene Chloride | NA | 3.8 | 86,843 | ND | ND | 7.27 | 25.25 | ND | ND | ND | ND | ND | ND |
| Pinene Isomer | NA | NA | NA | ND | ND | ND | ND | ND | ND | 1.49 | - | ND | ND |
| Toluene | NA | 42 | 753,703 | 1.48 | 5.58 | 23.6 | 88.94 | 1.6 | 6.03 | 3.58 | 13.49 | 1.21 | 4.56 |
| Xylene Isomer | 6,510 | 730 | NA | 0.796 | - | 81.94 | - | 0.797 | - | 1.6 | - | 0.831 | - |
| Unknown C11 Hydrocarbon | - | - | - | ND | ND | ND | ND | ND | ND | 1.33 | - | ND | ND |
| Unknown C12 Hydrocarbon | - | - | - | ND | ND | ND | ND | ND | ND | 1.97 | - | ND | ND |
| Unknown Chlorofluorohydrocarbon | - | - | - | ND | ND | ND | ND | ND | ND | 1.85 | - | ND | ND |
| Unknown | - | - | - | ND | ND | ND | ND | ND | ND | 3.26 | - | ND | ND |
| Unknown | - | - | - | ND | ND | ND | ND | ND | ND | 1.83 | - | ND | ND |
| Unknown | - | - | - | ND | ND | ND | ND | ND | ND | 1.89 | - | ND | ND |
| Total VOC | | | | 8.271 | | 81.940 | | 4.2340 | | 20.220 | | 5.091 | |

* The detected analytes were reported as Tentatively Identified Compounds (TICs)

NA - Screening criteria not available or does not apply

ND - nondetected

Shadowed cells indicated detected concentrations above criteria

Table E-12. New O-Field Controlled Burn Air Samples - December 1999
Pesticide Analysis Results from PUF Samplers

| Analytes | SP1 | | SP2 | | SP3 | | SP4 (Background) | | Blank |
|---------------------|------------------|-------------------------------------|------------------|-------------------------------------|------------------|-------------------------------------|------------------|-------------------------------------|------------------|
| | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) |
| alpha-BHC | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| beta-BHC | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| delta-BHC | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Lindane (gamma-BHC) | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Heptachlor | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Aldrin | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Heptachlor epoxide | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan I | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Dieldrin | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDE | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan II | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDD | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endosulfan sulfate | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4,4'-DDT | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Methoxychlor | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin ketone | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Endrin aldehyde | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| alpha-Chlordane | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| gamma-Chlordane | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Toxaphene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|----------------------------------|
| SP1 | PUF9 | 36.59 |
| SP2 | PUF8 | 42.36 |
| SP3 | PUF2 | 34.4 |
| SP4 (Background) | PUF4 | 46.37 |

BQL - Below Quantitation Limit

Table E-13 New O-Field Controlled Burn Air Samples - December 1999
PCBs Analysis Results from PUF Samplers

| Analyses | Toxic Air Pollutant 1-Hour Screening Level (ug/m ³) | EPA Region III Ambient Air RBCs (ug/m ³) | OSHA/PELs Concentration (ug/m ³) | SP1 Results (ug/PUF) | SP1 Concentrations (ug/m ³) | SP2 Results (ug/PUF) | SP2 Concentrations (ug/m ³) | SP3 Results (ug/PUF) | SP3 Concentrations (ug/m ³) | SP4 (Background) Results (ug/PUF) | SP4 (Background) Concentrations (ug/m ³) | Blank Results (ug/PUF) |
|--------------------------------------|---|--|---|----------------------------|---|----------------------------|---|----------------------------|---|---|--|------------------------------|
| 2,2',3,5'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',5,5'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | 0.017 |
| 2,2',5-Trichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,4'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,4,5'-Pentachlorobiphenyl | NA | 0.0031 | 500 | 0.042 | 0.0010 | 0.0160 | 0.0004 | 0.0130 | 0.0004 | BQL | - | BQL |
| 2,2',3,4,5,5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',3,5,5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | 0.0050 | 0.0001 | 0.0200 | 0.0006 | 0.009 | 0.0002 | BQL |
| 2,2',4,4',5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,2',4,5,5'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3',4,4'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | 0.091 | - | 0.004 |
| 2,3,3',4'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,3-Dichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | 0.11 |
| 2,4',5-Trichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Chlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22',33',44',5-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22',33',44',55',6-Nonachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22',34',55',6-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22',34',4',5',6-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 22',34',4',55',6-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|-------------------------------------|
| SP1 | PUF9 | 36.59 |
| SP2 | PUF6 | 42.36 |
| SP3 | PUF2 | 34.4 |
| SP4 (Background) | PUF4 | 46.37 |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Table E-14 New O-Field Controlled Burn Air Samples - December 1999
Explosives Analysis Results from PUF Samplers

| Analytes | SP1 | | SP2 | | SP3 | | SP4 (Background) | | Blank |
|----------------------------|------------------|-------------------------------------|------------------|-------------------------------------|------------------|-------------------------------------|------------------|-------------------------------------|------------------|
| | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) |
| 1,3,5-Trinitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 1,3-Dinitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4,6-Trinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,4-Dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2,6-Dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Amino-4,6-dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 2-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 3-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4-Amino-2,6-dinitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| 4-Nitrotoluene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| HMX | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Nitrobenzene | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| RDX | BQL | - | BQL | - | BQL | - | BQL | - | BQL |
| Tetryl | BQL | - | BQL | - | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|----------------------------------|
| SP1 | PUF7 | 36.15 |
| SP2 | PUF6 | 42.95 |
| SP3 | PUF1 | 38.14 |
| SP4 (Background) | PUF3 | 54.08 |

BQL - Below Quantitation Limit

Table E-15 New O-Field Controlled Burn Air Samples - December 1999
Inorganics Analysis Results

| Analytes | Toxic Air Pollutant 1-Hour Screening Level (ug/m ³) | EPA Region III Ambient Air RBCs (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP2 | | SP3 | | SP4 (Background) | | Blank | Blank |
|-----------|---|--|---|------------------------|--|------------------------|--|------------------------|--|------------------------|--|-------------------------------|-------------------------------------|
| | | | | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) TSP | Results (ug/filter) HANDI VOL |
| Mercury | 0.3 | 0.03 | 100 (acceptable ceiling) | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Silver | NA | 1.8 | 10 | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Aluminum | NA | 0.37 | 5,000 | 110 | 0.5200 | 55.7 | 0.2000 | 44.9 | 0.3800 | 39.9 | 0.3100 | 24.6 | BQL |
| Arsenic | NA | 0.00041 | 500 | 3.2 | 0.0100 | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Barium | NA | 0.051 | 500 | 24.7 | 0.1100 | 18.4 | 0.0700 | 2.3 | 0.0200 | 2.2 | 0.0200 | 20.1 | 1.4 |
| Beryllium | 0.1 | 0.00075 | 2 | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Calcium | NA | NA | NA | 595 | 2.5900 | 362.0 | 1.3300 | 89.2 | 0.7200 | 77.2 | 0.6000 | 373 | BQL |
| Cadmium | NA | 0.00090 | 5 | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Cobalt | NA | 22 | 100 | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Chromium | NA | 0.00015 | 500 | 2.1 | 0.0100 | 1.4 | 0.0010 | BQL | - | BQL | - | 1.3 | BQL |
| Copper | NA | 15 | 100 | 87.7 | 0.3800 | 31.4 | 0.1200 | 6.4 | 0.0500 | 10.3 | 0.0800 | 0.88 | BQL |
| Iron | NA | 110 | NA | 188 | 0.8100 | 88.1 | 0.2500 | 49.6 | 0.4000 | 55 | 0.4300 | 13.2 | BQL |
| Potassium | NA | NA | NA | 197 | 0.8600 | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Magnesium | NA | NA | NA | 80 | 0.3500 | 44.5 | 0.1800 | BQL | - | BQL | - | 35.3 | BQL |
| Manganese | NA | 0.0052 | 5000 | 4.5 | 0.0200 | 1.6 | 0.0100 | 1.0 | 0.0080 | 1 | 0.0080 | BQL | BQL |
| Sodium | NA | NA | NA | 1270 | 5.5300 | 993.0 | 3.6400 | 230.0 | 1.8500 | 225 | 1.7800 | 1150 | 213 |
| Nickel | NA | 7.3 | 1,000 | 3 | 0.0100 | 1.0 | 0.0040 | BQL | - | BQL | - | BQL | BQL |
| Lead | NA | NA | 50 | 17.1 | 0.0700 | 6.0 | 0.0220 | 5.5 | 0.0400 | 12.5 | 0.1000 | BQL | BQL |
| Antimony | NA | 0.15 | 500 | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Selenium | NA | 1.8 | 200 | 1.80 | 0.0100 | 0.5 | 0.0020 | 0.6 | 0.0050 | 0.52 | 0.0040 | BQL | BQL |
| Thallium | NA | 0.026 | 100 | BQL | - | BQL | - | BQL | - | BQL | - | BQL | BQL |
| Vanadium | NA | 2.6 | 500 | 6.9 | 0.0300 | 1.4 | 0.0050 | 1.4 | 0.0110 | 1.6 | 0.0100 | BQL | BQL |
| Zinc | 100 | 110 | NA | 33.9 | 0.1500 | 12.2 | 0.0450 | 6.5 | 0.0500 | 6.2 | 0.0500 | 2.1 | BQL |

| Sampling Location | Sample ID | Total Air Flow (m ³) | |
|-------------------|-------------|----------------------------------|---------|
| SP1 | TSP3 | 222.12 | Mercury |
| SP1 | TSP2 | 229.68 | |
| SP2 | TSP6 | 271.08 | Mercury |
| SP2 | TSP5 | 272.62 | |
| SP3 | Handl Vol 6 | 138.96 | Mercury |
| SP3 | Handl Vol 4 | 124.06 | |
| SP4 (Background) | Handl Vol 5 | 116.58 | Mercury |
| SP4 (Background) | Handl Vol 7 | 127.04 | |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shaded cells indicate detected concentrations above screening criteria

Table E-16. New O-Field Controlled Burn Air Samples - December 1999
Chemical Agent Analysis Results

| Chemical Agent | SP1 | | SP2 | | SP3 | | SP4 | |
|----------------|---------------------|--|---------------------|--|---------------------|--|---------------------|--|
| | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) | Concentration (ppb) | Concentration ($\mu\text{g}/\text{m}^3$) |
| Sarin (GB) | ND | ND | ND | ND | ND | ND | ND | ND |
| Soman (GD) | ND | ND | ND | ND | ND | ND | ND | ND |
| VX | ND | ND | ND | ND | ND | ND | ND | ND |
| Mustard (HD) | ND | ND | ND | ND | ND | ND | ND | ND |

ND - nondetected

Analysis provided by Edgewood Chemical Biological Center

Table E-17. New O-Field Controlled Burn Air Samples - December 1999
Radiological Analysis Results

| Analytes | EPA Region II Ambient Air RBCs (ug/m ³) | Toxic Air Pollutant/ 1-Hour Screening Level (ug/m ³) | OSRA PELs Concentration (ug/m ³) | SP1 | | SP2 | | SP3 | | SP4 (Background) | |
|---------------------------------------|---|--|---|-------------------------|---|-------------------------|---|-------------------------|---|-------------------------|---|
| | | | | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) |
| Gross Alpha | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Gross Beta | NA | NA | NA | 23 | 0.1013 | ND | ND | 3.4 | 0.0255 | 4.3 | 0.0363 |
| Actinium-228 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Bismuth-212 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Bismuth-214 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Cesium-137 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Cobalt-60 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Lead-210 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Lead-212 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Lead-214 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Potassium-40 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Protactinium-231 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Protactinium-234 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Radium-223 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Radium-224 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Radium-226 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Uranium-235 | NA | NA | NA | ND | ND | ND | ND | ND | ND | ND | ND |
| Total Uranium by Mass Spectrometry | 6 | 11 | 50 | ND | ND | ND | ND | ND | ND | ND | ND |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|----------------------|-------------|-------------------------------------|
| SP1 | TSP1 | 228.98 |
| SP2 | TSP4 | 267.89 |
| SP3 | Handi Vol 1 | 133.45 |
| SP4 (Background) | Handi Vol 2 | 118.58 |

NA - Screening criteria not available or does not apply
ND - nondetected

APPENDIX E-3

DATA TABLES FOR THE J-FIELD CONTROLLED BURN – APRIL 2000

Table E-18. J-Field Controlled Burn Air Samples - April 2000
Volatile Organic Compound Analysis Results from Summa Canisters

| Analytes | Toxic Air Pollutant Hour Screening Level Concentration (ug/m ³) | EPA Regional Ambient Air RBCs Concentration (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SPA (Background) | |
|---------------------------|---|--|---|------------------------|---------------------------------------|---------------------|---------------------------------------|
| | | | | Concentration (ppb) | Concentration (ug/m ³) | Concentration (ppb) | Concentration (ug/m ³) |
| Acetaldehyde * | 450 | 0.81 | 360,000 | 2.21 | 3.98 | ND | ND |
| Acetic Acid * | 370 | NA | 25,000 | ND | ND | 1.07 | 2.63 |
| Acetone * | 17,820 | 37 | 2,400,000 | 6.05 | 14.37 | 3.29 | 7.82 |
| Acetonitrile * | 1,010 | 62 | 70,000 | 1.73 | 2.9 | ND | ND |
| Benzene | 80 | 0.22 | 3,195 | 6.44 | 20.57 | 0.746 J | 2.38 |
| Chloroethene | NA | 0.21 | 2,556 | 0.332 J | 0.849 | ND | ND |
| Chloromethane | 525 | 1.80 | NA | 1.65 | 3.41 | 0.757 | 1.56 |
| Ethylbenzene | 5,430 | 110 | 435,000 | 5.91 | 25.66 | 1.62 | 7.03 |
| Ethylhexanol * | NA | NA | NA | 1.88 | 10.01 | ND | ND |
| Freon 12 | NA | NA | NA | ND | ND | 0.385 J | 1.90 |
| Furan * | NA | 0.37 | NA | 3.08 | 8.58 | ND | ND |
| Furfural * | NA | 3.70 | 20,000 | 6.56 | 25.78 | ND | ND |
| m/p-Xylenes | 6,510 | 730 | 435,000 | 3.43 | 14.89 | 0.967 J | 4.2 |
| Methylester Acetic Acid * | 7,570 | NA | NA | 1.21 | 3.67 | ND | ND |
| Methylfuran * | NA | NA | NA | 2.49 | 8.36 | ND | ND |
| Methylpropene * | NA | NA | NA | 1.89 | 4.34 | ND | ND |
| o-Xylene | 6,510 | 730 | 435,000 | 0.335 J | 1.45 | ND | ND |
| Styrene | 1,700 | 100 | 42,598 | 9.01 | 38.38 | 2.54 | 10.82 |
| Toluene | NA | 42 | 753,703 | 5.93 | 22.35 | 1.42 | 5.35 |
| Unknown C8 Hydrocarbon * | - | - | - | 1.92 | - | 0.973 | - |
| Unknown C4 Alkene * | - | - | - | 0.922 | - | ND | ND |
| Unknown * | - | - | - | 0.855 | - | ND | ND |
| Unknown * | - | - | - | 1.52 | - | ND | ND |
| Total VOC | | | | 62.477 | | 11 | |

* Analyte Identified as a Tentatively Identified Compound (TIC)

NA - Screening criteria not available or does not apply

ND - nondetected

J - Estimate concentration. Target detected at greater than the detection limit, but less than the quatitation limit (i.e., detection limit x5)

Shadowed cells indicate detected concentrations above screening criteria

Table E-19. J-Field Controlled Burn Air Samples - April 2000
Pesticide Analysis Results from PUF Samplers

| Analytes | Toxic Air Pollutant 1-Hour Screening Level (ug/m ³) | EPA Region III Ambient Air RBCs (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP4 (Background) | | Blank |
|---------------------|---|--|---|---------------------|--|---------------------|--|---------------------|
| | | | | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) |
| alpha-BHC | NA | NA | NA | BQL | - | BQL | - | BQL |
| beta-BHC | NA | NA | NA | BQL | - | BQL | - | BQL |
| delta-BHC | NA | NA | NA | BQL | - | BQL | - | BQL |
| Lindane (gamma-BHC) | NA | NA | 500 | BQL | - | BQL | - | BQL |
| Heptachlor | NA | 0.0014 | 500 | BQL | - | 0.078 | 0.0020 | BQL |
| Aldrin | NA | 0.00037 | 250 | BQL | - | BQL | - | BQL |
| Heptachlor epoxide | NA | 0.00069 | NA | BQL | - | BQL | - | BQL |
| Endosulfan I | NA | 2.2 | NA | BQL | - | BQL | - | BQL |
| Dieldrin | NA | 0.00039 | 250 | BQL | - | BQL | - | BQL |
| 4,4'-DDE | NA | 0.018 | NA | BQL | - | BQL | - | BQL |
| Endrin | NA | 0.11 | NA | BQL | - | BQL | - | BQL |
| Endosulfan II | NA | 2.2 | NA | BQL | - | BQL | - | BQL |
| 4,4'-DDD | NA | 0.026 | NA | BQL | - | BQL | - | BQL |
| Endosulfan sulfate | NA | NA | NA | BQL | - | BQL | - | BQL |
| 4,4'-DDT | NA | 0.018 | 1,000 | BQL | - | BQL | - | BQL |
| Methoxychlor | NA | 1.8 | 1,500 | BQL | - | BQL | - | BQL |
| Endrin ketone | NA | NA | NA | BQL | - | BQL | - | BQL |
| Endrin aldehyde | NA | NA | NA | BQL | - | BQL | - | BQL |
| alpha-Chlordane | NA | NA | 500 | BQL | - | BQL | - | BQL |
| gamma-Chlordane | NA | NA | 500 | BQL | - | BQL | - | BQL |
| Toxaphene | NA | 0.0057 | 500 | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|----------------------------------|
| SP1 | PUF9 | 47.59 |
| SP4 (Background) | PUF4 | 39.52 |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shadowed cells indicate detected concentrations above screening criteria

Table E-20. J-Field Controlled Burn Air Samples - April 2000
PCBs Analysis Results from PUF Samplers

| Analytes | Toxic Air Pollutant | EPA Region III | OSHA PELs Concentration (ug/m3) | SP1 | | SP4 (Background) | | Blank |
|-----------------------------------|-----------------------------------|-----------------------------|------------------------------------|---------------------|---------------------------|---------------------|---------------------------|---------------------|
| | 1-Hour Screening Level (ug/m³) | Ambient Air RBCs (ug/m³) | | Results (ug/PUF) | Concentrations (ug/m³) | Results (ug/PUF) | Concentrations (ug/m³) | Results (ug/PUF) |
| 2,2',3,5'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2',5,5'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2',5-Trichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2'3,4,4'5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2'3,4,5'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2'3,4,5,5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2'3,5,5'6'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2'4,4'5,5'-Hexachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,2'4,5,5'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,3',4,4'-Tetrachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,3,3'4'6'-Pentachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,3-Dichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2,4',5-Trichlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 2-Chlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 22'33'44'5'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 22'33'44'55'6'-Nonachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 22'34'55'6'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 22'344'5'6'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |
| 22'344'55'-Heptachlorobiphenyl | NA | 0.0031 | 500 | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m³) |
|-------------------|------------|---------------------|
| SP1 | PUF9 | 47.86 |
| SP4 (Background) | PUF4 | 39.52 |

NA - Screening criteria not available or does not apply
BQL - Below Quantitation Limit

Table E-21. J-Field Controlled Burn Air Samples - April 2000
Explosives Analysis Results from PUF Samplers

| Analytes | Toxic Air Pollutant | EPA Region III | OSHA PELs Concentration (ug/m ³) | SP1 | | SP4 (Background) | | Blank |
|----------------------------|--|--|---|---------------------|--|---------------------|--|---------------------|
| | 1-Hour Screening Level (ug/m ³) | Ambient Air RBCs (ug/m ³) | | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) | Concentrations (ug/m ³) | Results (ug/PUF) |
| 1,3,5-Trinitrobenzene | NA | 11 | NA | BQL | - | BQL | - | BQL |
| 1,3-Dinitrobenzene | NA | 0.037 | 1,000 | BQL | - | BQL | - | BQL |
| 2,4,6-Trinitrotoluene | NA | 0.21 | 1,500 | BQL | - | BQL | - | BQL |
| 2,4-Dinitrotoluene | 50 | 0.73 | 1,500 | BQL | - | BQL | - | BQL |
| 2,6-Dinitrotoluene | NA | 0.37 | 1,500 | BQL | - | BQL | - | BQL |
| 2-Amino-4,6-dinitrotoluene | NA | NA | NA | 21.3 | 0.4570 | BQL | - | BQL |
| 2-Nitrotoluene | NA | NA | 30,000 | BQL | - | BQL | - | BQL |
| 3-Nitrotoluene | NA | NA | 30,000 | BQL | - | BQL | - | BQL |
| 4-Amino-2,6-dinitrotoluene | NA | NA | NA | 5.9 | 0.1266 | BQL | - | BQL |
| 4-Nitrotoluene | NA | NA | 30,000 | BQL | - | BQL | - | BQL |
| HMX | NA | 18 | NA | BQL | - | BQL | - | BQL |
| Nitrobenzene | NA | 0.22 | 5,000 | BQL | - | BQL | - | BQL |
| RDX | NA | 0.0057 | NA | BQL | - | BQL | - | BQL |
| Tetryl | NA | 3.7 | 1,500 | BQL | - | BQL | - | BQL |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|------------|-------------------------------------|
| SP1 | PUF7 | 46.61 |
| SP4 (Background) | PUF3 | 47.27 |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shadowed cells indicated detected concentrations above screening criteria

Table E-22. J-Field Controlled Burn Air Samples - April 2000

Inorganics Analysis Results

| Analytes | Toxic Air Pollutant 1-Hour Screening Level (ug/m ³) | EPA Region III Ambient Air RBCs (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP4 (Background) | | Blank | Blank |
|-----------|---|--|---|------------------------|--|------------------------|--|-------------------------------|------------------------------------|
| | | | | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) | Concentrations (ug/m ³) | Results (ug/filter) TSP | Results (ug/filter) HANDIVOL |
| Mercury | 0.3 | 0.03 | 100 (acceptable ceiling) | 0.1 | 0.00043 | BQL | - | BQL | BQL |
| Silver | NA | 1.8 | 10 | 0.1 | 0.0004 | 0.08 | 0.0009 | BQL | BQL |
| Aluminum | NA | 0.37 | 5,000 | 159 | 0.6802 | 80.8 | 0.8831 | 6.8 | 2.3 |
| Arsenic | NA | 0.00041 | 500 | 0.59 | 0.0025 | BQL | - | BQL | BQL |
| Barium | NA | 0.051 | 500 | 18.3 | 0.0783 | 2.6 | 0.0284 | 3.6 | 1.2 |
| Beryllium | 0.1 | 0.00075 | 2 | BQL | - | BQL | - | BQL | BQL |
| Calcium | NA | NA | NA | 1910 | 8.1711 | 187 | 2.0437 | 161 | 43.4 |
| Cadmium | NA | 0.00099 | 5 | 0.84 | 0.0036 | 0.05 | 0.0005 | BQL | BQL |
| Cobalt | NA | 22 | 100 | 0.21 | 0.0009 | BQL | - | BQL | BQL |
| Chromium | NA | 0.00015 | 500 | 0.82 | 0.0035 | 0.51 | 0.0056 | 0.52 | 0.13 |
| Copper | NA | 15 | 100 | 16.2 | 0.0693 | 11 | 0.1202 | 0.22 | 0.23 |
| Iron | NA | 110 | NA | 148 | 0.6332 | 63.7 | 0.6962 | 5.5 | 3.2 |
| Potassium | NA | NA | NA | 740 | 3.1658 | 26.1 | 0.2852 | 14.4 | 8.5 |
| Magnesium | NA | NA | NA | 321 | 1.3733 | 35.3 | 0.3858 | 17.1 | 6.2 |
| Manganese | NA | 0.0052 | 5000 | 128 | 0.5476 | 2.4 | 0.0262 | 0.19 | 0.1 |
| Sodium | NA | NA | NA | 618 | 2.6439 | 158 | 1.7268 | 413 | 131 |
| Nickel | NA | 7.3 | 1,000 | 0.95 | 0.0041 | 0.35 | 0.0038 | BQL | BQL |
| Lead | NA | NA | 50 | 7.6 | 0.0325 | 0.7 | 0.0077 | BQL | BQL |
| Antimony | NA | 0.15 | 500 | BQL | - | BQL | - | BQL | BQL |
| Selenium | NA | 1.8 | 200 | 0.52 | 0.0022 | BQL | - | BQL | BQL |
| Thallium | NA | 0.026 | 100 | BQL | - | BQL | - | BQL | BQL |
| Vanadium | NA | 2.6 | 500 | 0.6 | 0.0026 | 0.29 | 0.0032 | BQL | BQL |
| Zinc | 100 | 110 | NA | 30.9 | 0.1322 | 4 | 0.0437 | 0.48 | 0.61 |

| Sampling Location | Sampler ID | Total Air Flow (m ³) | |
|-------------------|-------------|----------------------------------|---------|
| SP1 | TSP2 | 233.75 | Mercury |
| SP1 | TSP3 | 230.2 | |
| SP4 (Background) | Handi Vol 5 | 91.5 | Mercury |
| SP4 (Background) | Handi Vol 7 | 107.26 | |

NA - Screening criteria not available or does not apply

BQL - Below Quantitation Limit

Shadowed cells indicate detected concentrations above screening criteria

Table E-23. J-Field Controlled Burn Air Samples - April 2000
Chemical Agent Analysis Results

| Chemical Agent | SP1 | | SP4 | |
|----------------|---------------------|------------------------------------|---------------------|------------------------------------|
| | Concentration (ppb) | Concentration (ug/m ³) | Concentration (ppb) | Concentration (ug/m ³) |
| Sarin (GB) | ND | ND | ND | ND |
| Soman (GD) | ND | ND | ND | ND |
| VX | ND | ND | ND | ND |
| Mustard (HD) | ND | ND | ND | ND |

ND - nondetected

Analysis provided by Edgewood Chemical Biological Center

Table E-24. J-Field Controlled Burn Air Samples - April 2000
Radiological Analysis Results

| Pollutants | Toxic Air Pollutant 1-Hour Screening Level (ug/m ³) | EPA Region III Ambient Air RBCs (ug/m ³) | OSHA PELs Concentration (ug/m ³) | SP1 | | SP4 (Background) | |
|------------------|---|--|---|-------------------------|---|-------------------------|---|
| | | | | Results (pCi/filter) | Concentrations (pCi/m ³) | Results (pCi/filter) | Concentrations (pCi/m ³) |
| Gross Alpha | NA | NA | NA | ND | ND | ND | ND |
| Gross Beta | NA | NA | NA | ND | ND | ND | ND |
| Actinium-228 | NA | NA | NA | ND | ND | ND | ND |
| Bismuth-212 | NA | NA | NA | ND | ND | ND | ND |
| Bismuth-214 | NA | NA | NA | ND | ND | ND | ND |
| Cesium-137 | NA | NA | NA | ND | ND | ND | ND |
| Cobalt-60 | NA | NA | NA | ND | ND | ND | ND |
| Lead-210 | NA | NA | NA | ND | ND | ND | ND |
| Lead-212 | NA | NA | NA | ND | ND | ND | ND |
| Lead-214 | NA | NA | NA | ND | ND | ND | ND |
| Potassium-40 | NA | NA | NA | ND | ND | ND | ND |
| Protactinium-231 | NA | NA | NA | ND | ND | ND | ND |
| Protactinium-234 | NA | NA | NA | ND | ND | ND | ND |
| Radium-223 | NA | NA | NA | ND | ND | ND | ND |
| Radium-224 | NA | NA | NA | ND | ND | ND | ND |
| Radium-226 | NA | NA | NA | ND | ND | ND | ND |
| Uranium-235 | NA | NA | NA | ND | ND | ND | ND |
| U-233/234 | NA | NA | NA | ND | ND | ND | ND |
| U-235/236 | NA | NA | NA | 0.12 | 0.0005 | ND | ND |
| U-238 | NA | NA | NA | ND | ND | ND | ND |

| Sampling Location | Sampler ID | Total Air Flow (m ³) |
|-------------------|-------------|----------------------------------|
| SP1 | TSP1 | 230.95 |
| SP4 (Background) | Handi Vol 2 | 83.84 |

NA - Screening criteria not available or does not apply
ND - nondetected