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Decision Document Study Area CS-8/FS-21

> Final October 2000

Prepared for: AFCEE/MMR Installation Restoration Program 322 E. Inner Road Otis ANGB, MA 02542 DSN: 557-4670 Comm: 508-968-4670

> Prepared by: HARDING ESE

Submitted by: Advanced Infrastructure Management Technologies Oak Ridge, Tennessee 37831-7606 Managed by: LOCKHEED MARTIN ENERGY SYSTEMS for the U.S. DEPARTMENT OF ENERGY under contract DE-AC05-84OR21400



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December 6, 2000

Mr. Robert M. Gill Remediation Program Manager HQ AFCEE/MMR 322 East Inner Road Otis ANG Base, MA 02542-5028

Subject: CS-8/FS-21 Decision Document Massachusetts Military Reservation

Dear Mr. Gill:

For the Air Force Center for Environmental Excellence, and as directed by Advanced Infrastructure Management Technologies, Harding ESE is hereby submitting copies of the Final Decision Document for Study Area CS-8/FS-21. This final deliverable includes the executed signature pages provided under HQ AFCEE/MMR transmittal letter dated November 28, 2000.

Please contact me at (207) 775-5401 if you have any questions.

Sincerely,

Harding ESE, Inc.

In INV.

Jøhn W. Peterson MMR Project Manager

Enclosures (16 bound, 1 unbound, 1 digital)

cc: P. Marchessault, USEPA (2 bound)
J. Murphy, USEPA (letter only)
L. Pinaud, MADEP (2 bound)
E. Grillo, MADEP (letter only)
M. Maly, SFIM-AEC-IRP (1 bound)
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File

INSTALLATION RESTORATION PROGRAM

DECISION DOCUMENT STUDY AREAS CS-8/FS-21 OPERATIONAL MOTOR POOL, OMS-22/CURRENT PRODUCT TANK NO. 90

FINAL OCTOBER 2000

Prepared for: Air Force Center for Environmental Excellence and Air National Guard

> Prepared by: HARDING ESE Portland, Maine Project No. 49240

Advanced Infrastructure Management Technologies Oak Ridge, Tennessee 37831-7606

Managed by: LOCKHEED MARTIN ENERGY SYSTEMS for the U.S. DEPARTMENT OF ENERGY under contract DE-AC05-84OR21400

DECISION DOCUMENT STUDY AREAS CS-8/FS-21

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EXECUTIVE SUMMARY

As part of the preliminary assessment conducted in 1986 for the Installation Restoration Program at the Massachusetts Military Reservation, Study Area CS-8 (Operational Motor Pool, OSM-22), and adjoining Study Area FS-21 (Current Product Tank No. 90), were identified as potential sites of past uncontrolled disposal of hazardous substances. Because of this finding, contamination in these areas was investigated and characterized during two site inspection programs (1988, 1989), a drainage structure removal program (1996), underground storage tank removals (1996, 1999), and additional groundwater sampling (1999). This Decision Document provides a summary of these activities, an assessment of the contaminants detected, and a preliminary risk evaluation.

Potential contaminant sources for the CS-8/FS-21 Study Areas consisted of:

- Vehicle maintenance wastes prior to 1970.
- Battery electrolyte disposed of in an on-site cesspool from 1970 to 1985.
- A former 12,500-gallon underground storage tank used to store diesel fuel.
- The former location of Current Product Tank No. 90, a 5,000-gallon underground storage tank used to store motor vehicle gasoline.

The deposition of these potential contaminant sources is as follows:

- Wastes generated at CS-8 are now transported to the Army's 3500 Area at the MMR for temporary storage prior to shipment off-site to a licensed Treatment, Storage, and Disposal Facility.
- The cesspool was removed in 1996 during the Drainage Structure Removal Program. Clean closure was achieved.
- 12,500-gallon underground storage tank used to store diesel fuel was removed in 1988 and replaced with a new double wall tank. Pressure testing of the original tank in 1985 indicated it to be tight (i.e., no leaks). This new tank was removed in 1998 and not replaced. Soil contamination during underground storage tank excavations was not detected.
- A leak test conducted on Current Product Tank No. 90 (5,000-gallon tank) in 1985 was inconclusive due to a bent fill pipe. The tank was removed in 1988 and replaced with a double wall tank of the same capacity, west of the vehicle repair shop. This new tank was later removed in 1996. Soil contamination during underground storage tank excavations was not detected.

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Results from the human health and ecological preliminary risk evaluation suggest that unacceptable levels of risk are not anticipated. Therefore, the Air Force Center for Environmental Excellence recommends no further action for Study Areas CS-8 and FS-21 via this Decision Document.

1.0 INTRODUCTION

The objectives of this decision document are to (1) describe the histories of the Operational Motor Pool, OMS-22 (Study Area CS-8), and the Current Product Tank No. 90 (Study Area FS-21) at the Massachusetts Military Reservation (MMR); (2) present results of site characterization investigations at Study Areas Chemical Spill (CS)-8/Fuel Spill (FS) FS-21; (3) present results of human health and ecological preliminary risk evaluations (PREs); and (4) explain why no further action is recommended for these study areas. Study Areas CS-8/FS-21 initially were identified in the Task 6 records search, a preliminary assessment (PA) of MMR conducted as part of the U.S. Department of Defense Installation Restoration Program (IRP) (E. C. Jordan Co., 1986).

The IRP is a program at military facilities to identify, evaluate, and remediate wastedisposal and spill sites that were contaminated through past practices. Study Areas CS-8/FS-21 were investigated further in the Tasks 2-3A and 2-3C site investigations (SIs) (E. C. Jordan Co., 1989a and 1990b).

In 1988, Current Product Tank No. 90 (5,000-gallon motor vehicle gasoline [MOGAS] underground storage tank [UST]), and a 12,500-gallon Diesel-fuel UST were replaced at the study areas with two double-walled USTs of the same size, and storing the same petroleum product (5,000-gallon MOGAS, 12,500-gallon Diesel fuel). The 5,000-gallon UST and the 12,500-gallon UST were removed in 1996 and 1999, respectively. In 1996, a cesspool was also removed from the study area as part of the Drainage Structure Removal Program.

In 1999, confirmational sampling was conducted to confirm the finding that no additional remedial actions were required.

2.0 STUDY AREAS DESCRIPTION AND HISTORY

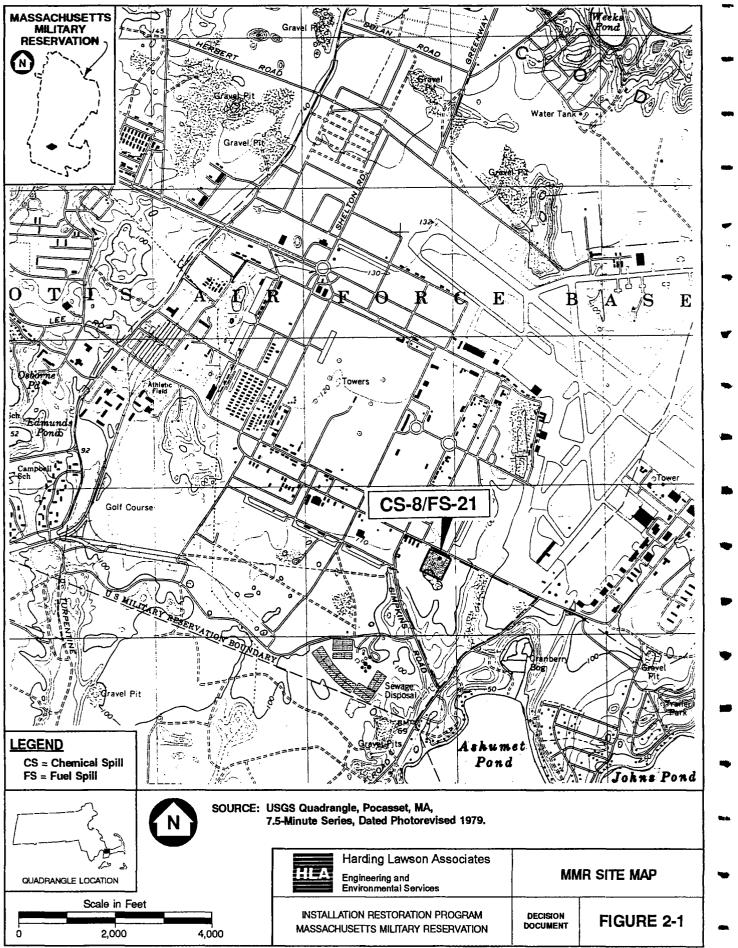
Study Areas CS-8/FS-21 are located next to each other along the south-central boundary (Figure 2-1). Because of this geographical proximity, Study Areas CS-8/FS-21 were combined for purposes of SI field activities.

2.1 Study Area CS-8

Study Area CS-8 includes an active and an abandoned concrete wash pad, a cesspool, and a 12,500-gallon diesel-fuel UST and pump island located west of the Vehicle Repair Shop. The study area has been operating as the current Army National Guard Vehicle Repair Shop at MMR since 1950. The maintenance shop has been responsible for the maintenance of from 20 vehicles (1950) to 300 vehicles (1986). In November 1998, maintenance operations at the Study Area were shifted to the Unit Training Equipment Storage (UTES) site off Greenway Road. Minimal maintenance activities still occur at the Study Area by reservists during weekend activities.

Summary of Waste Management Practices: Since 1970, all vehicle maintenance wastes, except used battery electrolyte, have been disposed of off-site through a hazardous-waste contractor. Until 1985, battery electrolyte had been discharged to the cesspool. No records exist regarding pre- 1970 waste-disposal practices. It is likely that before 1970, wastes either were taken to the base landfill or dumped onto the ground at the study area. Reportedly up to 250 gallons per year (gal/yr) of waste oil (mixed with solvents), 20 gal/yr of paint thinner, and 50 gal/yr of waste solvents were generated by vehicle-maintenance activities before 1970 (E. C. Jordan Co., 1989a). Degreasing agents likely included halogenated solvents such as TCE and DCE, and petroleum distillates. The quantities actually disposed of on-site are unknown. Until November 1998, wastes generated at Study Area CS-8 were drummed, sent to Building 4600 for temporary storage, prior to shipment off-site to a licensed Treatment, Storage, and Disposal Facility (TSDF). Currently, wastes generated at Study Area CS-8 are transported to the Army's 3500 Area at the MMR for temporary storage, prior to shipment off-site to a TSDF.

Cesspool: Details of the original volume and design of the cesspool are not available, nor is it known if it was connected to a leachfield. Upon its removal in 1996, the cesspool was observed to have been constructed of concrete barrel block walls, with an open dirt bottom



ten feet below grade surface, and a 4-inch-diameter ceramic inlet pipe with no outlet pipe. The walls had previously collapsed (Jacobs, 1996).

12,500-Gallon Diesel Fuel UST: The 12,500-gallon diesel-fuel UST at Study Area CS-8 was pressure-tested in May 1985 and found to be tight. In 1988, this tank was removed and replaced with a double-walled tank of the same capacity at the same location (Pesce and Berube, 1990). During removal of the tank in 1988, no contaminated soil was encountered (Berube, 1990). In 1998, this new 12,500-gallon diesel-fuel tank was removed and not replaced. See Appendix C for details on the 1998 UST removal. Although documentation of the 1988 removal is based on verbal conversation, and not written documentation, the location of the 1998 and 1988 removals were in the same location, and no contamination was detected within the 1998 UST excavation.

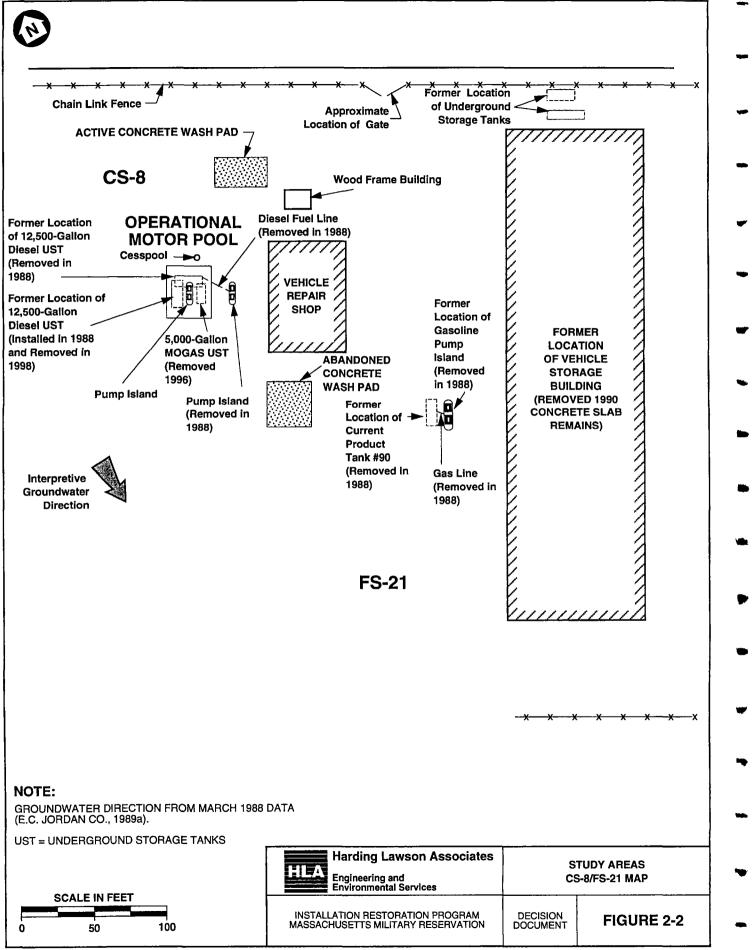
2.2 Study Area FS-21

Study Area FS-21, southwest of the Vehicle Repair Shop, is the former location of Current Product Tank No. 90, a 5,000-gallon MOGAS UST installed in 1954 (Figure 2-2). Study Area FS-21 also includes the former location of two USTs at the northern end of the Vehicle Storage Building. Details concerning the size, construction, and former use of the two USTs are not available.

Current Product Tank No. 90: This tank was pressure-tested in May 1985 to assess potential leakage, but the test was inconclusive because of a bent fill pipe. In 1988, the tank was removed and replaced with a double-walled MOGAS tank of the same capacity, installed west of the Vehicle Repair Shop, and adjacent to the 12,500 gallon diesel-fuel UST (Pesce and Berube, 1990) (see Figure 2-2). At the time of the removal of Current Product Tank No. 90 in 1988, the associated piping and gasoline pump island were removed and the area was re-graded. No contamination was noted in the surrounding soil, which is consistent with the interpretation of investigation results (Pesce and Berube, 1990). The new 5,000-gallon UST was subsequently removed in 1996. See Appendix C for details on the 1996 UST removal.

2.3 Task 6 Records Search

Potential contaminants identified in the Task 6 Records Search were waste solvents and oils, battery electrolyte, and MOGAS. The primary constituents of MOGAS are hydrocarbons. Study Areas CS-8 and FS-21 were assigned Hazard Assessment Rating Methodology (HARM) ratings of 66.6 and 61.9, respectively. The HARM system is



designed to prioritize sites at an installation to indicate the relative need for a SI. HARM considers factors such as site characteristics, waste characteristics, potential for contaminant migration, and waste management practices. Based on findings of the Task 6 Records Search, Study Areas CS-8 and FS-21 were recommended for SI studies.

Historically, a chain-link fence enclosed Study Areas CS-8/FS-21. In the late 1980s, the fence was temporarily opened along the southern boundary for site improvements, including grading of the vehicle parking area. The fence has not been reinstalled along the southern boundary.

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3.0 SUMMARY OF STUDY AREAS INVESTIGATIONS

The SI field program for Study Areas CS-8/FS-21 was designed to evaluate whether past maintenance activities, waste-disposal methods, and a potentially leaking gasoline UST (i.e., Current Product Tank No. 90) affected soil and groundwater quality.

Task 2-3A SI: This 1987 investigation consisted of a metal-detector survey, a soil-gas survey, excavation of 10 test pits, completion of six soil borings with installation of monitoring wells in four of the borings, field gas chromatograph (GC) screening of soil samples, and laboratory analysis of soil and groundwater samples. Selected samples were submitted to a U.S. Environmental Protection Agency (USEPA) Contract Laboratory Program (CLP) laboratory and analyzed for selected Target Compound List (TCL) and Target Analyte List (TAL) chemicals. Table 3-1 lists the specific analyses for each sample.

Task 2-3C SI: This 1989 investigation was intended to resample groundwater from the existing study-area monitoring wells and sample contents of the on-site cesspool. However, only one of the four on-site wells was sampled (CS8/MW2). Wells CS8/MW1 and FS21/MW2 had been destroyed, and well FS21/MWI had been damaged (PVC riser was bent slightly) during previous site construction activities designed to improve surface-water drainage. The exact location of the cesspool could not be determined at the time of the investigation, because of newly graded gravel.

Laboratory analytical data from Tasks 2-3A and 2-3C are USEPA CLP data. Although the data were not validated according to USEPA functional guidelines or USEPA Region I guidelines, they were evaluated for holding times and the presence of blank contamination during preparation of this report and for suitability for use in risk assessments and to support engineering studies. Results of this data evaluation are in Appendix A and as discussed in this evaluation, the data were found suitable. Lack of full validation likely adds conservatism to the characterization of site conditions, because full validation could result in rejection of a maximum concentration that otherwise would have been used in the contamination assessment and PRE. Additional details concerning data-quality evaluation are in the Tasks 2-3A and 2-3C reports.

Phase I Sump Removal Program: In 1992, four TerraProbe borings were completed in the cesspool area to further evaluate the presence of contamination as part of the Phase I

TABLE 3-1 CLP ANALYTICAL SAMPLING PROGRAM

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| | | | SOIL S | SAMPLES | | | WATER SAMPLES | | | | | | | |
|-------------------|----------------|---------------------|---------|----------|-----------------|------------|----------------|----------|----------|-----------------|------------|-------------|-----|--|
| EXPLORATION ID | SAMPLE DATE | DEPTH (feet bgs) | TCL VOC | TCL SVOC | TCL Pest/PCB | TAL Inorg. | SAMPLE DATE | TCL VOC | TCL SVOC | TCL Pest/PCB | TAL inorg. | EDB MTBE | ТРН | |
| FS-21/TB-1 | 9/1/87 | 13-15 | Y | N | N | N | | | | | | | | |
| | | 18-20 | Y | N | N | N | | | | | 1 | | | |
| | | 28-30 | Y | (Y | Y | Y | | 1 | | | [[| | | |
| | | 33-35 | Y | N | N | N | | | | i | | | | |
| | | 63-65 | Y | Y | Y | Y | | | | | | | | |
| | | 63-65 (dup) | Y | Y | Y | Y | | ļ | 1 | | | | | |
| FS-21/MW-1 | 9/1/87 | 52-62 | | | - | | 10/28/87 | Y | Y | N | Y | N | N | |
| | | (screen) | | | | | 3/10/99 | Y | N | N | Y | Y | N | |
| FS-21/TB-2 | 9/9/87 | 0-2 | Y | Y | Y | Y | | | | | | | | |
| | | 10-12 | Y | Y | Y | Y | | | | | 1 1 | | | |
| | | 39-41 | Y | N | N | N | | | | | | | | |
| | | 59-61 | Y | Y | Y | Y | | | | | | | | |
| FS-21/MW-2 | 9/9/87 | 51-66 (screen) | - | - | | - | 1/7/88 | N | Y | N | Y | N | N | |
| CS-8/TP-1 | 8/29/87 | 1-2 | Y | Y | Y | Y | | | | | | | | |
| CS-8/TP-6 | 8/31/87 | 5-6 | Y | Y | Y | Y | | | | | | | | |
| CS-8/TB-1 | 8/30/87 | 23-25 | Y | Y | Y | Y | | | | | | | | |
| | | 43-45 | Y | Y | Y | Y | | | | | 1 1 | | | |
| | | 43-45 (dup) | Y | Y | Y | Y | | | | | | | | |
| | | 58-60 | Y | Y | Y | Y | | | | | | | | |
| CS-8/TB-2 | 8/31/87 | 19-21 | Y | Y | Y | Y | | <u> </u> | | | | | | |
| | | 54-56 | Y | Y | Y | Y | | | | | 1 1 | | | |

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continued

TABLE 3-1 CLP ANALYTICAL SAMPLING PROGRAM

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| | | | SOIL S | AMPLES | | | WATER SAMPLES | | | | | | | |
|-------------------|----------------|---|-----------------------|-------------|------------------|------------------|--|---------|-----------------------|-----------------|-----------------------|-----------------------|-----------------------|--|
| EXPLORATION ID | SAMPLE DATE | DEPTH (feet bgs) | TCL VOC | TCL SVOC | TCL Pest/PCB | TAL inorg. | SAMPLE DATE | TCL VOC | TCL SVOC | TCL Pest/PCB | TAL Inorg. | EDB MTBE | ТРН | |
| CS-8/TB-2 | 8/31/87 | 54-56 (dup) | Y | Y | Y | Y | | | 1 | | | | | |
| CS-8/MW-2 | 8/31/87 | 52-62 (screen) | - | - | | _ | 1/7/88 1/7/88 (dup) 6/28/89 6/28/89(dup) 3/10/99 3/10/99(dup) | Y | Y Y Y N N | 2 Z Z Z Z Z | Y Y Y Y Y | N N N Y Y | N Y Y N N | |
| CS-8/TB-3 | 8/31/87 | 8-10 18-20 58-60 | Y Y Y | N Y Y | N Y Y | N Y Y | | | | | | | | |
| CS-8/TB-4 | 9/1/87 | 0-2 5-7 8-10 10-12 34-36 54-56 | Y Y Y Y Y | N N N Y Y | N N Y Y | N N Y Y | | | | | | | | |
| CS-8/MW-4 | 9/1/87 | 52-62 (screen) | ~ | - | - | - | 10/28/87 10/28/87 (dup) | Y Y | Y Y | N N | Y Y | N N | N N | |
| 04BH0001 | 12/14/99 | 35-40 | _ | - | | _ | 12/14/99 | Y | Y | N | Y | Ŷ | N | |

Υ

N

dup

screen

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Notes:

TCL=Target Compound ListTAL=Target Analyte ListVOCs=volatile organic compoundsSVOCs=semivolatile organic compounds

PCBs = polychlorinated biphenyls

TB = test boring TP = test pit

Ξ

=

CLP

MW

TPH = total petroleum hydrocarbons

monitoring well

- sample analyzed for parameter
- sample not analyzed for parameter
- well-screen interval
- duplicate sample

Page 2 of 2

Contract Laboratory Program

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sump removal program with field GC screening of the soil samples collected (ABB Environmental Services, Inc. [ABB-ES], 1992a).

Confirmational Groundwater Sampling: In March 1999, monitoring wells CS8/MW2 and FS21/MW1 were resampled using the low flow purge technique. Monitoring well FS21/MW1, which couldn't be sampled during the Task 2-3C Site Investigation due to a bent riser, was accessed with a submersible pump. Groundwater samples were analyzed off-site for inorganics, and VOCs including methyl-tert-butyl ether (MTBE) and ethylene dibromide (EDB). In December 1999, a groundwater sample (Sample I.D. 04BH0001) was collected using the USEPA Region One's low flow purge and sample technique, at the top of the groundwater table from a borehole advanced at the former location of FS-21/MW-2. The groundwater sample was analyzed for VOCs, SVOCs, EDB, and inorganics.

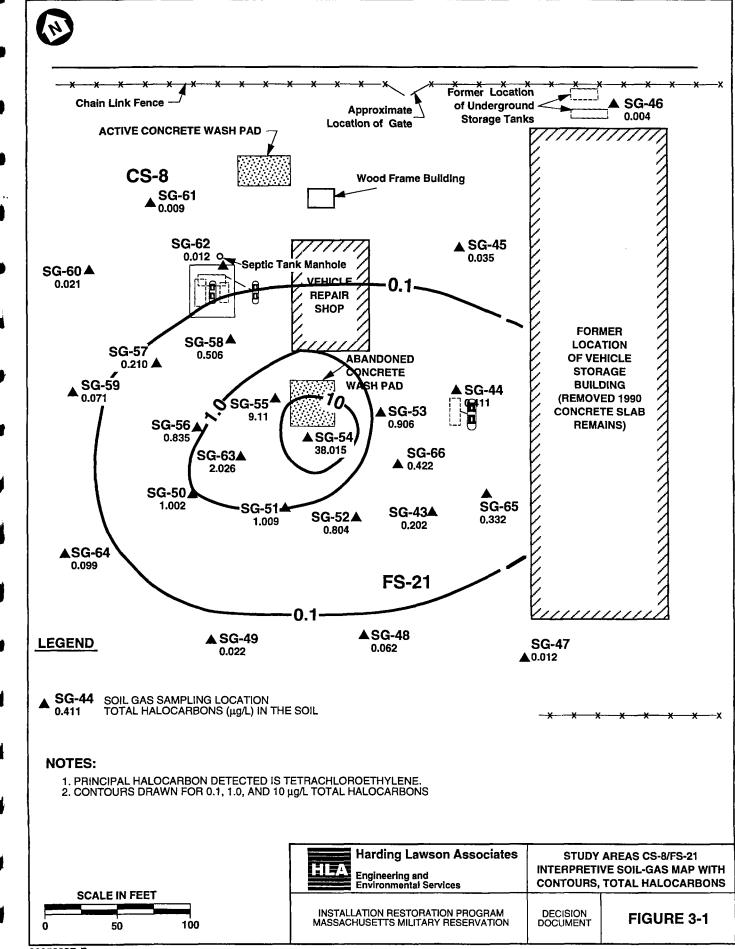
3.1 METAL-DETECTOR SURVEY

A metal-detector survey was conducted at Study Areas CS-8/FS-21 on August 26, 1987, to locate USTs and underground fuel and utility lines so that invasive work would be done in areas likely to show contamination, while avoiding underground structures. Field procedures consisted of traverses spaced at intervals of 10 feet or less. The survey located the manhole cover of the cesspool and the diesel-fuel UST (see Figure 2-2). The supply line from the tank to the diesel pump island could not be located.

3.2 SOIL-GAS SURVEY

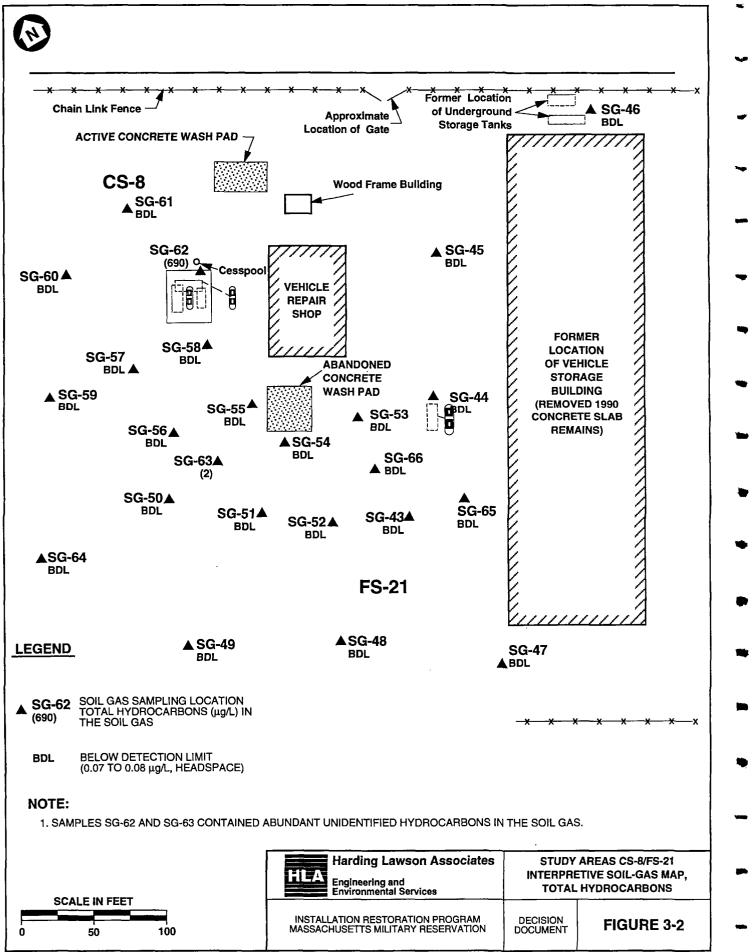
To assist in locating test pits and borings, a soil-gas survey was conducted on August 27 and 28, 1987, in areas considered likely to have received fuel or solvent spills and leakage. These areas included the abandoned wash pad, abandoned and active UST sites, and cesspool. Probes were completed in 24 locations to approximately 4 feet below ground surface (bgs) (Figure 3-1). Target analytes included trichloroethene (TCE), tetrachloroethene (PCE), trichloroethane, chloroform benzene, toluene, xylenes, and total non-methane hydrocarbons. All probes detected at least trace amounts of halocarbons, with the highest concentrations detected along the southern edge of the abandoned concrete wash pad, south of the Vehicle Repair Shop. The suspected source of these detections is minor spills of runoff from mechanical parts that had been cleaned with solvents. Halocarbon concentrations generally decreased with distance from the abandoned wash pad. Only two soil-gas probes detected the presence of hydrocarbons: The highest concentration was adjacent to the cesspool north of the diesel pump island (690 micrograms per liter [µg/L] headspace); the second highest concentration was 2 µg/L headspace, approximately I 00 feet south of the diesel pump island (Figure 3-2).

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3.3 TEST-PIT PROGRAM

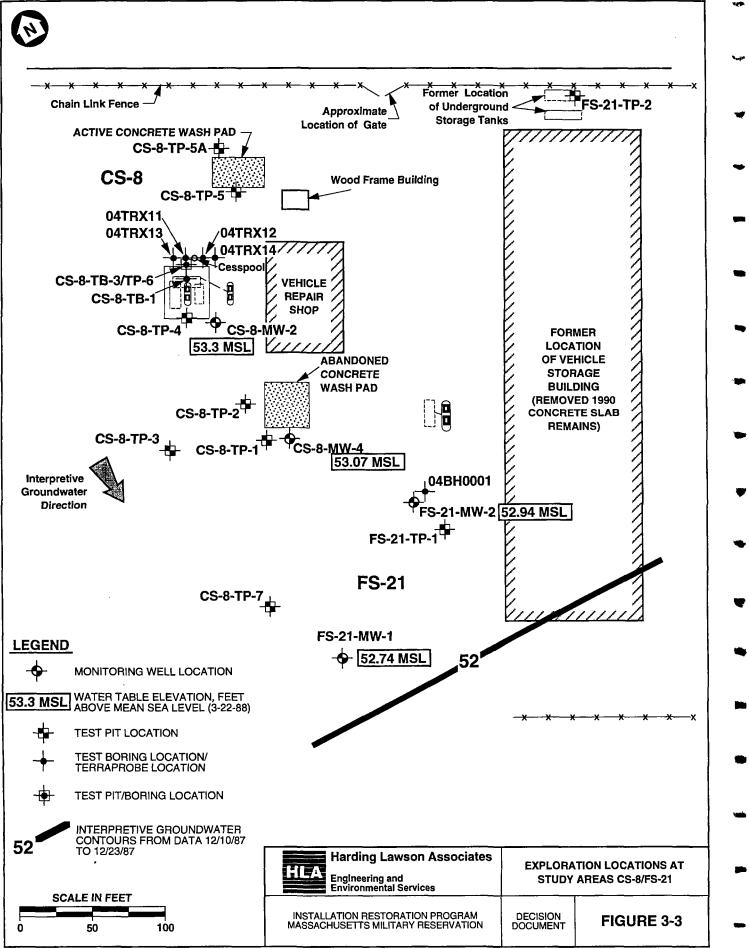
Ten test pits were excavated between August 29 and 31, 1987, to investigate shallow subsurface soil at suspected contaminant source locations. Logs for these test pits are in the Task 2-3A SI appendices (E. C. Jordan Co., 1989a).

Study Area FS-21: Two test pits were completed at Study Area FS-21 (Figure 3-3). Test pit FS-21/TP-1, located in the area of a geophysical anomaly approximately 50 feet south of the MOGAS pump island, was completed to a depth of approximately 7.5 feet bgs. During excavation, an iron pipe and a clay pipe were encountered; it was determined that neither pipe was in use. Soil surrounding these pipes at 3-5 feet bgs had photoionization (PI) readings at background levels and a soil sample just below the pipes was screened using a field GC; target VOCs were not detected. A 1981 utilities map for MMR does not specifically show pipes in the vicinity of FS-21 TP1. However, the map legend indicates that generally all building sanitary sewer connections were 6-inch diameter vituperative clay pipe, and all building water connections were 4-inch diameter cast iron, or 2.5-inch diameter wrought iron pipe. Given these diameters and piping materials, along with SI observations, these pipes were likely associated with water and/or sewer feeds to the Vehicle Storage Building.

The second test pit (FS-21/TP-2) was located at the former location of the USTs north of the Vehicle Storage Building to characterize shallow soil that may have been contaminated by spills or leaks from those tanks. During completion of both test pits, soil staining was not observed and PI meter readings were below background levels. Two soil samples (i.e., 5 and 7 feet bgs) from each test pit were screened with the field GC; target volatile organic compounds (VOCs) were not detected in any of the samples.

Study Area CS-8: Eight test pits (designated CS-8/TP-1 through CS-8/TP-7 and CS-8/TP-5A) were completed at Study Area CS-8 (see Figure 3-3); only one (i.e., CS8/TP-1) had PI meter readings above background levels. This test pit was located adjacent to the abandoned concrete wash pad and had PI meter readings of 13 parts per million (ppm). Eleven soil samples from the eight test pits were screened using a field GC; two VOC constituents were detected at trace concentrations. PCE was detected in CS-8/TP-1 (0.8 foot bgs) at a concentration of 1.1 micrograms per kilogram (μ g/kg); benzene was detected in CS-8/TP-6 (5 feet bgs) at a concentration of 0.3 μ g/kg. Two soil samples (CS8/TP-1, CS8/TP-6) were collected on the basis of field observations and submitted for off-site CLP analyses (see Table 3-1).

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3.4 SOIL BORINGS AND MONITORING WELLS

Six soil borings were completed and four monitoring wells were installed between August 30 and September 1, 1987. Four borings and two wells were positioned to investigate Study Area CS-8, and two borings and two wells were located to assess conditions at Study Area FS-21 (see Figure 3-3). Logs for these borings and wells are in the Task 2-3A SI appendices (E. C. Jordan, Co., 1989a). "CS-8/MW-2" and "CS-8/TB-2" are the designations for a monitoring well and test boring, respectively, completed in one location. If a test boring was completed but no monitoring well installed at a specific location, then a "TB" does not have a corresponding "MW".

A 2-foot-long split-spoon sample was collected from each boring at intervals from near the ground surface to the bottom of the boring. These samples were screened in the field by a PI meter and a subset of the samples was screened with the field GC. On the basis of PI meter readings and field GC results, the apparently most contaminated soil samples (a total of 23 different sample intervals) were submitted for off-site CLP analyses from the test borings completed at CS-8 and FS-21. The explorations IDs are listed in Table 3-1.

Study Area CS-8: At Study Area CS-8, two monitoring wells (CS-8/MW-2 and CS-8/MW-4) were installed to evaluate soil and groundwater quality. (Note that soil and groundwater samples are identified as TB and MW samples, respectively.) CS-8/MW-2 was positioned adjacent to the diesel pump island west of the Vehicle Repair Shop and downgradient of the cesspool. CS-8/MW-4 was positioned downgradient of the abandoned concrete wash pad. PI meter readings for the split spoon soil samples did not exceed background levels; however, several negative PI meter readings suggest that humidity may have reduced instrument sensitivity. No soil staining was observed during monitoring-well installation.

Field GC screening detected target VOCs in the soil boring samples. In CS-8/MW-2, trace concentrations (i.e., less than 1.5 μ g/kg) of benzene and toluene were detected in subsurface soil samples over 29 feet bgs. In CS-8/MW-4, field GC screening detected up to 9.5 μ g/kg of PCE at the ground surface, where PI meter readings of 6 to 8 ppm were recorded. The suspected source of these detections is minor spills of runoff from mechanical parts that had been cleaned with solvents. Trace concentrations of TCE and toluene also were detected in the surface soil. Both soil and groundwater samples were collected at these two locations and submitted for CLP analyses (see Table 3-1). CS-8/MW-2 was re-sampled in 1989 and 1999 for confirmational analyses (see Table 3-1).

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Field GC screening also detected target VOCs in soil samples from the two borings in Study Area CS-8 in which wells were not installed (i.e., CS-8/TB-1 and CS-8/TB-3). In CS-8/TB-1, located adjacent to the 12,500-gallon diesel-fuel UST, field GC screening of most soil samples from 8 to 63 feet bgs detected trace concentrations of dichloroethene (DCE), TCE, PCE, benzene, and toluene. In CS-8/TB-3, located just south of the cesspool, benzene (up to 1.5 μ g/kg), TCE (up to 0.6 μ g/kg), and PCE (1.8 to 3.2 μ g/kg) were detected from 53 to 58 feet bgs. Toluene was detected throughout the soil column in CS-8/TB-3, with the highest concentration (4.1 μ g/kg) at 8 feet bgs, where soil staining also was observed.

Study Area FS-21: Two monitoring wells were installed at Study Area FS-21 to evaluate potential soil and groundwater contamination. FS-21/MW-1 was positioned downgradient of the study area along the southern boundary; FS-21/MW-2 was installed adjacent to the former location of Current Product Tank No. 90.

During installation of FS-21/MW-1, elevated PI meter readings of 7 ppm were observed at 28 to 30 feet bgs; readings at other depths were at background levels. Ten soil samples were screened using the field GC; traces of benzene, TCE, and toluene were detected, primarily within the unsaturated soil (13-53' bgs). Soil staining was not observed.

During installation of FS-21/MW-2, PI meter readings of 6 to 17 ppm and a two-inch layer of black stained soil was observed in the surface sample at this location. Soil staining was encountered in surface soil. Fifteen soil samples screened with the field GC showed the presence of low concentrations of VOCs. Low concentrations of VOCs also were detected at the water table: TCE ($11 \mu g/kg$) and toluene (2.3 $\mu g/kg$).

Groundwater samples from these two monitoring wells were submitted for CLP analysis (see Table 3-1). In 1999, FS-21/MW-1 was re-sampled for confirmational analyses (see Table 3-1). In December 1999, a groundwater sample (Sample I.D. 04BH0001) was collected using the USEPA Region One's low flow purge and sample technique, at the top of the groundwater table from a borehole advanced at the former location of FS-21/MW-2. The groundwater sample was analyzed for VOCs, SVOCs, EDB, and inorganics.

3.5 TERRAPROBE BORINGS

Four TerraProbe borings (i.e., 04TRXI1, 04TRX12, 04TRX13, and 04TRX14) were completed in February 1992 as part of the Phase I Sump Removal Action in the cesspool area along a line approximately perpendicular to the western wall of the Vehicle Repair Shop. Approximate locations of the borings are shown in Figure 3-3. Soil samples were

collected for field laboratory analysis at 4, 8, 12, 18, and 24 feet bgs from each boring. Each sample was analyzed for VOCs, semivolatile organic compounds (SVOCs), pesticides, polychlorinated biphenyls (PCBs), inorganics, and total petroleum hydrocarbons (TPH). A summary of the results and their corresponding Tier I and Tier II hazard equivalent concentrations (HECs) are presented on Table 4-2b, and a summary of the onsite analytical results are included in Appendix B.

Review of the TerraProbe data shows that with one exception, detection of target VOCs, SVOCs, pesticides, and inorganics was limited to samples from boring 04TRX12 (4-12 feet bgs). Possible fuel-related VOCs, including toluene (130 μ g/kg), ethylbenzene (190 μ g/kg), m/p-xylene (6.1 and 250 μ g/kg), and o-xylene (67 μ g/kg), were reported in one sample from this boring, but not its duplicate. The inorganics copper (270 milligrams per kilogram [mg/kg]) and zinc (380 mg/kg) also were reported in the same sample, but not its duplicate. Two pesticides, delta-benzene hexachloride (D-BHC) (150 to 530 μ g/kg) and heptachlor epoxide (140 to 1,800 μ g/kg), were reported in samples from this boring. The VOC m/p-xylene (6.1 μ g/kg) was reported in one sample from boring 04TRX14. TPH was detected in all samples at 40 to 980 mg/kg. TPH concentrations were apparently unrelated to both target analyte concentrations (except in samples from boring 04TRX12) and PI meter readings. SVOCs were not detected in any samples.

3.6 GEOLOGY

The geology of the study areas was interpreted from the test pits and test borings. The borings drilled at this site extended to depths of 65 to 66 feet bgs that is just below the water table.

Subsurface soil encountered was typical of outwash deposits found across the southern part of MMR. The soil was primarily poor- to well-graded, fine to medium sand, with trace amounts of coarse sand and fine gravel. In some locations, the top 1.4 feet of soil was fill over a thin, discontinuous ash layer. The ash was observed to be grey to black, loose to medium dense, and dry. The ash measured from 0.1 foot thick in FS21/TP1 to 0.7 feet thick in CS8/TP5. The typical study-area soil sample consisted of about 95 percent fine to medium sand, 5 percent fin gravel and coarse sand, and trace amounts of silt.

3.7 Hydrogeology

As interpreted from MMR-wide groundwater-elevation data, groundwater flows in a southerly direction in the vicinity of Study Areas CS-8/FS-21, toward the western portion of Study Areas SD-1 and Ashumet Pond (E. C. Jordan Co., 1989b). Groundwater elevations at Study Areas CS-8/FS-21 on March 22, 1988, ranged from 52.7 to 53.3 feet above mean sea level (MSL), or approximately 58 feet bgs.

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4.0 CONTAMINATION ASSESSMENT

In general, the contamination assessment of a site is based on CLP data and, to a lesser extent, field analytical data, depending on the methodologies and analyses performed. At Study Areas CS-8/FS-21, the field GC screening analytical data, in which concentrations were reported in micrograms per liter (μ g/L) headspace, and the PI meter readings are considered qualitative only. The CLP analytical data, which quantify concentrations of TCL contaminants in soil and groundwater, are most useful for contamination assessment

4.1 SURFACE SOIL

Three surface-soil samples (zero to 2 feet bgs) were analyzed for CLP analytes (Table 4-1). TCL VOCs were not detected. Two TCL SVOCs, both phthalates, were detected at concentrations of less dm half their Contract Required Quantitation Limits (CRQLs) in one of two surface-soil samples. Phthalates would not be expected at Study Areas CS-8/FS-21 as a result of manufacturing- or process-related activities because such activities did not occur on-site. Because of their ubiquitous presence in plastics, phthalates are recognized by USEPA as common laboratory and sampling artifacts (USEPA, 1991). Given the low concentrations at which they were detected and their nature as common artifacts, the two phthalates are not considered site contaminants. The PCB Aroclor-1260 was detected at 660 μ g/kg in one of two samples in which it was a target analyte. This single detection was in the sample collected adjacent to the abandoned wash pad. It is possible that very localized PCB contamination exists in this area and is related to some former activity at the abandoned wash pad. However, the measured concentration is below the clean-closure limit of 1,000 μ g/kg set by USEPA (USEPA, 1989c), and PCB-containing equipment is not known to be present in the study areas.

In the two samples analyzed for TAL inorganics, only barium, copper, thallium, and zinc were detected at concentrations above MMR maximum background concentrations for surface soil (ABB-ES, 1992b). All four inorganics were detected at concentrations less than twice MMR maximum background concentrations.

TABLE 4-1

SUMMARY OF ANALYTES DETECTED IN SURFACE SOIL - OFF-SITE ANALYSIS

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| "Exploration ID | | FS-21/TB-2 | CS-8/TP-1 | CS-8/TB-4 |
|--|------------------------|----------------------------|----------------------------|------------|
| "Sampled | | 10-SEPT-87 | 02-SEPT-87 | 04-SEPT-87 |
| ANALYTES "Depth | :Concentration (mg/kg) | 0 | <u> </u> | 1 |
| VOLATILE ORGANIC COMPOUNDS (µg/kg) | | _ | _ | _ |
| | | | | - |
| SEMIVOLATILE ORGANIC COMPOUNDS (µg/kg) | | | | |
| Bis(2-ethylhexyl)phthalate | NA | 64J | - | NR |
| Butylbenzylphthalate | NA | 140J | - | NR |
| | | | | |
| Pesticides/PCBs (µg/kg) | | | | |
| 4,4'-DDT | NA | - | - | NR |
| Aroclor 1260 | NA | - | 660 | NR |
| | | | | |
| INORGANIC COMPOUNDS (mg/kg) | | s autorization constrainte | 7.74 - Constitution 3, 107 | |
| Aluminum | 8,930 | 3220 🔤 | 1300 🖓 | |
| Arsenic | 3.6 | 2.2 | 0.69J | NR |
| Barium | 10.4 | 9.6J | 11J | NR |
| Beryllium | 0.65 | | 0.37J | NR |
| Calcium | 969 | 218J | 331J | NR |
| Chromium | 6.8 | 5.3 | 2:1J | NR |
| Cobalt | 4.1 | | 1.5J | NR |
| Copper | 5.2 | | 9.2J | NR |
| Iron | 12,400 | 4010J | 2490 | NR |
| Lead | 12.05 | 5.6 | 6.9 | NR |
| Magnesium | 794.5 | 178 J | 536J | NR |
| Manganese | 108 | 31J | 59J | NR |
| Selenium | 0.33 | - | | NR |
| Thallium | 0.25 | | 0.40J | NR |
| Vanadium | 15.2 | 6.4J | 3.7 | NR |
| Zinc | 16 | 30J | 17J | NR |

NOTES:

- = not detected
 bgs = below ground surface
 4,4'-DDT = 4,4'- dichlorodiphenyltrichloroethane
 j = estimated concentration
- mg/kg = milligrams per kilogram NR = not requested
- PCBs = polychlorinated biphenyls
- µg/kg = micrograms per kilogram

Depth indicates top of sample interval.

Shaded results are below MMR background concentrations.

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4.2 SUBSURFACE SOIL

Review of CLP analytical data indicates that 25 subsurface-soil samples collected from 5 to 65 feet bgs were analyzed for TCL VOCs; none were detected (Table 4-2a). Of the 18 samples analyzed for TCL SVOCs, one phthalate was detected in one sample only, at a concentration of 340 μ g/kg. This concentration is barely above the CRQL of 330 μ g/kg; as discussed previously, phthalates would not be expected at this site and are typical sampling and laboratory artifacts. Therefore, the one phthalate detection is not considered to represent site-related contamination.

Eighteen samples were analyzed for TCL pesticides and PCBs. The pesticide 4,4'dichlorodiphenyltrichloroethane (DDT) was detected in three samples only, at concentrations of 18, 24, and 45 μ g/kg, compared to the CRQL of 16 μ g/kg. Pesticide detections were in deep soil, occurring at depths from 19 to 61 feet bgs; they likely reflect residual concentrations from widespread pesticide applications for insect control on MMR in the 1960s and 1970s. At these depths, pesticides affect no human or ecological receptors; therefore, the low 4,4'-DDT concentrations are not considered to represent contamination of concern.

The TAL inorganics beryllium, chromium, and copper each were detected once at concentrations above the MMR maximum background concentrations for subsurface soil. These three detections occurred in three different samples; in two of these, a duplicate sample failed to replicate the background exceedance. Additionally, the chromium and copper detections were at concentrations lower than the MMR maximum background concentration for surface soil. For this reason, and because the detections were not reproducible in duplicate samples, site-related contamination of subsurface soil by inorganic analytes is not indicated.

Interpretation of field laboratory analytical results for VOCs, SVOCs, pesticides, PCBs, and inorganics in subsurface-soil samples from the TerraProbe borings in the cesspool area support the conclusions derived from the CLP data that contaminants are not widespread around the former cesspool (Table 4-2b). The one sample that showed the presence of fuel-related compounds was a duplicate sample (i.e., the 8-foot bgs sample from 04TRX12) and no fuel-related compounds were detected in the original sample. At this same sample location, the duplicate sample contained detectable concentrations of copper and zinc while neither of these analytes was detected in the original sample.

TABLE 4-22 SUMMARY OF ANALYTES DETECTED IN SUBSURFACE SOIL - OFF-SITE ANALYSIS

STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| "Exploration ID: | Maximum | FS-21/TB-1 | FS-21/TB-1 | FS-21/TB-1 | FS-21/TB-1 | FS-21/TB-1 | FS-21/TB-1* | FS-21/TB-2 | FS-21/TB-2 | FS-21/TB-2 | CS-8/TP-8 | CS-8/TB-1 | CS-8/TB-1 | CS-8/TB-1* | CS-8/TB-1 | CS-&/TB-2 |
|-----------------------------------|-----------------------------|------------------|------------------|-----------------------------|------------------|------------------|------------------|------------------|------------|------------|--------------------------------|-------------------|------------------|------------------|------------------|------------------|
| "Sampled: ANALYTES "Depth: | Background Conc. (mg/kg) | 03-SEPT-87 13 | 03-SEPT-87 18 | 03-SEPT-87 28 | 03-SEPT-87 33 | 03-SEPT-87 63 | 03-SEPT-87 63 | 10-SEPT-87 10 | 10-SEPT-87 | 10-SEPT-87 | 02-SEPT-87 | 01-SEPT-87 23 | 01-SEPT-87 43 | 01-SEPT-87 43 | 01-SEPT-87 58 | 01-SEPT-87 19 |
| ANALTIES Depti: | CONC. (MUKg) | | 10 | | | | 03 | | | | | | | | | 1.0 |
| VOLATILE ORGANIC COMPOUNDS (ug/kg | NA | - | - | - | - | - | - | - | - |] - | - | - | - | - | - | - |
| SEMIVOLATILE ORGANIC COMPOUNDS | | | | | | | | | | | | | | | | |
| Butylbenzylphthalate | NA | NR | NR | - | NR | - | - | - | NR | - | 340 | - | - | - | - | - |
| PESTICIDES/PCBs (µg/kg) | | | | | | | | | | } | | | | | | |
| 4.4'-DDT | NA | NR | NR | - | NR | - | - | - | NR | 45 | - 1 | - | - 1 | - | - | 18 |
| Aroclor 1260 | NA | NR | NR | - | NR | - | - | - | NR | - | - | - | - | - | - | - |
| INORGANIC COMPOUNDS (mg/kg) | | | | | | | | | 1 | | | | | U. | | |
| Aluminum | 1,980 | NR | NR | 877 | NR | 352 | 316 | 914 | NR | 432 | 647 | 259 | 401 | 925 | 483 | .420 |
| Arsenic | 2.3 | NR | NR | 0.66J | NR | $d = d_{2}$ | 0.65 | | NR | | | | in the second | | | |
| Barium | 14.7 | NR | NR | 3.2J | NR | 1.2J | | 7 4.0J | NR | 3.4J | 1.7J | 1.5J | 2.6J | 4.1J | 2.8J | 1.5J |
| Beryllium | 0.69 | NR | NR | 19 A 4 | NR | | 1.3 | | NR | | 「「「「「」」 | | | .0.40J | | 141.00 |
| Calcium | 933 | NR | NR | 85.1 | NR | 68J | 75J | 76J | NR | 136J | 27J | 27J | 36J | 56J | 42J | /i 27J |
| Chromium | 3.9 | NR | NR | | NR | | | 4.3 | NR | | and the property of the second | | | | | Second Strate |
| Cobalt | 2.6 | NR | NR | | NR | | | 70 25 25 | NR | - | | | | 1.30J | 1:5J" | |
| Copper | 4.3 | NR | NR | 17482-3 1948-3 1949-3 | NR | | | 0.94J | NR | | 4.1J | 3.1J | 3.2J | . 3.9J | 4.7J | 3.6J |
| Iron | 2,600 | NR | NR | 1300 | NR | 825 | 573 | 2020J | NR | 865J | 961 | 418 | 718 | 2170 | 6 979 ° | 690 . |
| Lead | 3.70 | NR | NR | 11 | NR | 0.72J | 0,79J | 1.8 | NR | ા નનું િ | 1.0J | 0.84J | 0.82J | 1.2 | 0.72J | 0.78J |
| Magnesium | 742.0 | NR | NR | 251J | NR | | 4 | 354J | NR | 95J | 129J | Shares and start | , 95J | 557J | . 196J | |
| Manganese | 587 | NR | NR | 41J | NR | 8.3J | 5.20J | 🧆 53J 🎋 | NR | 56J | 19J | 10J 🗇 | 15J | 23J | 53J | 22J |
| Selenium | 0.62 | NR | NR | 0.35J | NR | 0.49J | | | NR | 0,39J | | | | | | |
| Thallium | 1.00 | NR | NR | يرين الروادين | NR | | E. | 建品合物的 | NR | | . 0.41J 🖞 | 0.45J | 0.45J | 0.44J | 0.55J | 0.41J |
| Vanadium | 5.7 | NR | NR | 1.9J | NR | | | 2.0J | NR | 2,0J | | المراجعة المعادمة | | 2.9J | | |
| Zinc | 16 | NR | NR | 6.0 | NR | 6.6 | 4.1J | 8.7J | NR | 8.7J | 6.6J | 6.7J | 8.1J | 9.2J | 12J | ૣ ૼ ,6Ĵ |
| | | | | | | | | | 1 | | | | | | | |
| | | | | | | | | | | | 1 | | | | | |

NOTES:

= not detected

= duplicate sample

4,4'-DDT = 4,4'-dichlorodiphenyltrichloroethane = estimated concentration

= milligrams per kilogram

= not analyzed

not analyzed
 not requested
 polychlorinated biphenyls
 micrograms per kilogram
 Depth indicates top of sample interval.
 Shaded results are below MMR background concentrations.

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TABLE 4-2a SUMMARY OF ANALYTES DETECTED IN SUBSURFACE SOIL - OFF-SITE ANALYSIS

STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| Background Conc. (mg/kg) NA NA | 01-SEPT-87 54 - | 01-SEPT-87 54 - | 01-SEPT-87 8 - | 01-SEPT-87 18 | 01-SEPT-87 58 | 04-SEP-87 5 | 04-SEP-87 8 | 04-SEP-87 10 | 04-SEP-87 34 | 04-SEP-87 54 |
|---|--|---|--|---|--|--|--|---|--|--|
| NA | - | | - | | | | | | | |
| | - | - | - | | | | | | | |
| NA | | | r – | - | • | - | - | - | - | - |
| NA | | | | | | | | | | |
| | - | - | NR | - | - | NR | NR | - | - | - |
| | | | | | | | | | | Ļ |
| NA | - | 24 | NR | - | - | NR | NR | - | - | - 1 |
| NA | - | - | NR | - | - | NR | NR | - | - | - 1 |
| | | | | | | | | | | 1 |
| 1,980 | 409 | 469 | NR | 429 | 225 | NR | NR | 551 | 373 | .612.1 |
| 2.3 | | | NR | | | NR | NR | 0.55J - | 0.55J | 1 ³ 0.63J |
| 14.7 | 1.4J | 1.3J | NR | 1.7J | 1.4 | NR | NR | | -**21J | 3.6J |
| 0.69 | | | NR | | | NR | NR | See See See | The state of the | (1)保全。5 |
| 933 | 26J | 29J | NR | 23J | 23.J | NR | NR | : 81J | 63J | j, 93J) } |
| 3.9 | | | NR | | | NR | NR | 1.1 | A Start | 13-1-144 |
| 2.6 | | | NR | مەربىيە يەربىيە يەربىيە يەربىيە يەربىيە | | NR | NR | | | · 1642 小小小 |
| 4.3 | 4,3J | 4.3J | NR | 3.5J | | NR | NR | e | | |
| 2,600 | 822 | 1060 | NR | 821J | 418 | NR | NR | 893J | 583 | is it is reapy to |
| 3.70 | 0,9J: | , 0.81J | NR | 2.1J | 0.92J | NR | NR | 0.98J | 0.52J | 0.67J |
| 742.0 | 103J | 147J | NR | 83J | and the second secon | NR | NR | 171J | 82J 1 | ູ່ 148 ປີະ |
| 587 | 9.5J | 24J | NR | 27J | 5.4J | NR | NR | | 7 16J | 3.414J |
| 0.62 | | | NR | 1 N - 199 | | NR | NR | | ₩ 0,33J | A dish to see |
| 1.00 | 0.47J | 0.52J | NR | 0,52J | 0.54J | NR | NR | | | |
| 5.7 | | | NR | | 5.5 | NR | NR | 1.6J.+ | 1 13 13 1 | 1.1 |
| 16 | 7.1J | 8.8J | NR | 7.1J | 5.3 | NR | NR | 5.0 | 5.0 | 6.3 |
| | NA 1,980 2.3 14.7 0.69 933 3.9 2.6 4.3 2,600 3.70 742.0 587 0.62 1.00 5.7 | NA - 1,980 409 2.3 14.7 1.4J 0.69 933 26J 3.9 2.6 4.3 4,3J 2,600 822 3.70 0,9J 742.0 103J 587 9.5J 0.62 - 1.00 0.47J 5.7 | NA - - 1,980 409 469 2.3 14.7 1.4J 14.7 1.4J 1.3J 0.69 933 26J 29J 3.9 - - - 2.6 4.3 4.3J 4.3J 2.600 822 1060 3.70 3.70 0.9J 0.8IJ 147J 587 9.5J 24J 0.62 1.00 0.47J 0.52J 5.7 | NA - - NR 1,980 409 469 NR 2.3 NR NR 14.7 1.4J 1.3J NR 14.7 1.4J 1.3J NR 933 26J 29J NR 3.9 - NR NR 2.6 NR NR NR 4.3 4.3J NR NR 2.600 822 1060 NR 3.70 0.9J 0.81J NR 742.0 103J 147J NR 587 9.5J 24J NR 0.62 NR NR NR 1.00 0.47J 0.52J NR 5.7 - NR NR | NA - - NR - 1,980 409 469 NR 429 2.3 NR - NR - 14.7 1.4J 1.3J NR 1.7J 0.69 NR - NR - 933 26J 29J NR 23J 3.9 NR - NR - 4.3 4.3J 4.3J NR 3.5J 2,600 822 1060 NR 821J 3.70 0.9J 0.81J NR 2.1J 742.0 103J 147J NR 83J 587 9.5J 24J NR 27J 0.62 - NR 0.52J NR 1.00 0.47J 0.52J NR 0.52J | NA - - NR - - 1,980 409 469 NR 429 225 2.3 NR NR 1.41 1.31 NR 1.71 1.41 14.7 1.4J 1.31 NR 1.7J 1.41 1.41 0.69 NR NR 2.3J 2.3J 2.3J 3.9 1.41 2.6 NR NR - NR - - - 2.60 822 1060 NR 821.1 4.16; - - 2.60 822 1060 NR 821.1 4.16; - 2.60 822 1060 NR 821.1 4.16; - 3.70 0.9J 0.31.J NR 2.1J 0.92.J - 742.0 103.3 147.J NR 83.J - - NR 1.00 0.47.J 0.52.J NR 0.52.J 0.54.J | NA - NR - NR - NR 1,980 409 469 NR 429 225 NR 2.3 NR NR 1.429 225 NR 14.7 1.4J 1.3J NR 1.7J 1.4J NR 933 26J 29J NR 23J 23J NR 933 26J 29J NR 23J 23J NR 3.9 - NR - NR - NR 2.6 - NR - NR - NR 4.3 4.3J 4.3J NR 3.5J 3.6J NR 2.600 822 1060 NR 821J 416 NR 3.70 0.9J 0.81J NR 2.1J 0.92J NR 587 9.5J 24J NR 27J 5.4J NR 1.00 0.47J 0.52J NR <td>NA - NR - NR NR NR 1,980 409 469 NR 429 225 NR NR 2.3 NR NR NR 1.42 NR NR NR 14.7 1.4J 1.3J NR 1.7J 1.4L NR NR 933 26J 29J NR 23J 23J NR NR 3.9 NR NR - NR NR NR NR 2.6 4.3J NR - NR NR NR NR 2.60 822 1060 NR 8213 3.6J NR NR 2.600 822 1060 NR 8213 3.6J NR NR 2.600 822 1060 NR 8213 3.6J NR NR 3.70 0.9J 0.31J NR 2.1J 0.92J NR NR 742.0<td>NA - NR - NR NR NR . 1,980 409 469 NR 429 225 NR NR 551 2.3 NR NR 1.7J 1.4U NR NR 0.55J 14.7 1.4J 1.3J NR 1.7J 1.4U NR NR 2.1J 933 26J 29J NR 23J 23J NR NR 81J 3.9 - NR - NR NR NR 81J 2.6 NR - NR NR NR NR 81J 2.60 822 1060 NR 821J 3.6J NR NR 83J 2.600 822 1060 NR 821J 9.9J NR NR 9.98J 742.0 103J 147J NR 83J - NR NR 171J 587 9.5J 24J<td>NA - NR - NR NR NR NR - - NR NR NR - - - NR NR - - - - NR NR -</td></td></td> | NA - NR - NR NR NR 1,980 409 469 NR 429 225 NR NR 2.3 NR NR NR 1.42 NR NR NR 14.7 1.4J 1.3J NR 1.7J 1.4L NR NR 933 26J 29J NR 23J 23J NR NR 3.9 NR NR - NR NR NR NR 2.6 4.3J NR - NR NR NR NR 2.60 822 1060 NR 8213 3.6J NR NR 2.600 822 1060 NR 8213 3.6J NR NR 2.600 822 1060 NR 8213 3.6J NR NR 3.70 0.9J 0.31J NR 2.1J 0.92J NR NR 742.0 <td>NA - NR - NR NR NR . 1,980 409 469 NR 429 225 NR NR 551 2.3 NR NR 1.7J 1.4U NR NR 0.55J 14.7 1.4J 1.3J NR 1.7J 1.4U NR NR 2.1J 933 26J 29J NR 23J 23J NR NR 81J 3.9 - NR - NR NR NR 81J 2.6 NR - NR NR NR NR 81J 2.60 822 1060 NR 821J 3.6J NR NR 83J 2.600 822 1060 NR 821J 9.9J NR NR 9.98J 742.0 103J 147J NR 83J - NR NR 171J 587 9.5J 24J<td>NA - NR - NR NR NR NR - - NR NR NR - - - NR NR - - - - NR NR -</td></td> | NA - NR - NR NR NR . 1,980 409 469 NR 429 225 NR NR 551 2.3 NR NR 1.7J 1.4U NR NR 0.55J 14.7 1.4J 1.3J NR 1.7J 1.4U NR NR 2.1J 933 26J 29J NR 23J 23J NR NR 81J 3.9 - NR - NR NR NR 81J 2.6 NR - NR NR NR NR 81J 2.60 822 1060 NR 821J 3.6J NR NR 83J 2.600 822 1060 NR 821J 9.9J NR NR 9.98J 742.0 103J 147J NR 83J - NR NR 171J 587 9.5J 24J <td>NA - NR - NR NR NR NR - - NR NR NR - - - NR NR - - - - NR NR -</td> | NA - NR - NR NR NR NR - - NR NR NR - - - NR NR - - - - NR NR - |

NOTES:

= not detected

= duplicate sample

4,4'-DDT = 4,4'-dichlorodiphenyltrichloroethane

= estimated concentration

= milligrams per kilogram

= not analyzed

= not requested

= polychlorinated biphenyls

= micrograms per kilogram Depth indicates top of sample interval. Shaded results are below MMR background concentrations

TABLE 4-2b SUMMARY OF ANALYTES DETECTED IN SOIL - ON-SITE ANALYSIS

| EXPLORATION ID: SAMPLED: DEPTH: | Maximum Background Conc. (mg/kg) | 04TRX11 28-FEB-92 4 | 04TRX11 28-FEB-92 8 | 04TRX11 28-FEB-02 12 | 04TRX11 28-FEB-92 18 | 04TRX11 28-FEB-92 24 | 04TRX12 28-FEB-92 4 | 04TRX12 28-FEB-92 8 | 04TRX12 28-FEB-92 8 | 04TRX12 28-FEB-92 12 | 04TRX12 28-FEB-92 18 | 04TRX12 28-FEB-92 24 |
|---------------------------------------|--|---------------------------|---------------------------|----------------------------|----------------------------|----------------------------|---------------------------|---------------------------|---------------------------|----------------------------|----------------------------|----------------------------|
| ANALYTES | | | | | | | | | | | | |
| VOCs (ug/kg) | | | | | 1 | | | | | | | |
| Toluene | NA | - | - | - | - | - | - | 130 | - | - | | - |
| Ethylbenzene | NA | - 1 | - | - | - | - | - 1 | 190 | - | - | - | - (|
| o-Xylene | NA | | - | - | - | - | - 1 | 67 | - | - | - | - |
| m/p-Xylene | NA | - | - | - | - | - | - | 250 | - | - | - | - |
| Pesticides/PCBs (µg/kg) | | | | | | | } | | | | | |
| Delta-BHC | NA | - ' | - | - | - | - I | - | 330 | 530 | 150 | - | - 1 |
| Heptachlor epoxide | NA | - | - | - | - | - | 140 | 1700 | 1800 | 1300 | - | - |
| inorganics (mg/kg) | | | i . | | | i | 1 | | | | | |
| Lead | 3.7 | - | - | - | - | - | - | - 1 | - | - | - | - |
| Copper | 4.3 | - | - 1 | - | - | - | - | 270 | - | - | - | - |
| Zinc | 16.0 | - | - | • | - | - ' | - 1 | 380 | - | - | - 1 | - |
| Arsenic | 2.3 | - | - | - | - | - | l - | - | - | - | - | - |
| Chromium | 3.9 | - | - | - | - | - | - ' | - | - | - | - | - |
| TPH (mg/kg) | NA | 130 | 83 | 84 | 120 | 120 | 130 | 980 | 820 | 83 | 120 | 40 |

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

NOTES:

N/A = not detected N/A = not analyzed µg/kg = micrograms per kilogram mg/kg = milligrams per kilogram

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TABLE 4-2b SUMMARY OF ANALYTES DETECTED IN SOIL - ON-SITE ANALYSIS

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| EXPLORATION ID: SAMPLED: DEPTH: | | 04TRX13 04-MAR-92 4 | 04TRX13 04-MAR-92 8 | 04TRX13 04-MAR-92 8 | 98TRX13 04-MAR-92 12 | 98TRX13 04-MAR-92 18 | 98TRX13 04-MAR-92 24 | 98TRX14 04-MAR-92 4 | 98TRX14 04-MAR-92 8 | 98TRX14 04-MAR-92 12 | 98TRX14 04-MAR-92 18 | 98TRX14 04-MAR-92 24 |
|---------------------------------------|------|---------------------------|---------------------------|---------------------------|----------------------------|----------------------------|----------------------------|---------------------------|---------------------------|----------------------------|----------------------------|----------------------------|
| ANALYTES | | | | | | | | | | | | |
| VOCs (ug/kg) | | | | | | | | | | | | |
| Toluene | NA | - | - | - | - | - | - | - | - | - | - | - |
| Ethylbenzene | NA | - 1 | - | - | - | - | - | - | - | - | - | - |
| o-Xylene | NA | - | - | - | - | - | - | - | - | - | - | - |
| m/p-Xylene | NA | - | - | - | - | - | - | 6.1 | • | - | - | - |
| Pesticides/PCBs (µg/kg) | | | | | | | | | | | | |
| Delta-BHC | NA | - | - | - | - | - | | - | - | - | - | - |
| Heptachlor epoxide | NA | - | - | - | - | - | - | - | - | - | - | - |
| Inorganics (mg/kg) | | | | | | | | | | | | |
| Lead | 3.7 | - | - | - | - | - | - | - | - | - | - | - |
| Copper | 4.3 | - | - | - | - | - | - | - | - | - | - | - |
| Zinc | 16.0 | - | - | - | - | - | - 1 | - | - | - | - | - |
| Arsenic | 2.3 | - | - | - | - | - | - | - | - | - | - | - |
| Chromium | 3.9 | - | - | - | - | - | - | - | - | - | - | - |
| TPH (mg/kg) | NA | 170 | 330 | 320 | 290 | 410 | 410 | 180 | 160 | 120 | 210 | 410 |

NOTES:

- = not detected

N/A = not analyzed µg/kg = micrograms per kilogram mg/kg = milligrams per kilogram

4.3 GROUNDWATER

A total of twelve groundwater samples (including four duplicates) were collected from four monitoring wells and one borehole located at Study Area CS-8/FS-21; however, not all samples were analyzed for all parameters. The analytical sampling program is presented in Table 3-1.

Organic Analytes: Site-related TCL VOCs were detected above laboratory detection limits in the groundwater sample collected from borehole 04BH0001; ethylbenzene and xylenes were detected at concentrations of 1.14 and 3.04 μ g/l, respectively. Although 2-butanone was detected in groundwater samples collected from monitoring well CS8/MW2, the lack of 2-butanone detected in groundwater from adjacent wells and from groundwater collected from CS8/MW2 in 1999, indicates that its presence in 1989 was likely due to sample contamination introduced during collection and/or laboratory analysis (Haase, et al., 1988 and E.C. Jordan Co., 1991). As a result of these findings, 2-butanone was not considered a site-related contaminant of potential concern and was not evaluated in the PRE/PRA.

Site-related TCL SVOCs were detected above laboratory detection limits in the groundwater sample collected from borehole 04BH0001; Acenapthene and Fluorene were detected at concentrations of 0.16 and 0.18 μ g/l, respectively. Naphthalene was detected at an estimated concentration of 0.08 μ g/l (below its quantitation limit) from this same sample.

04BH0001 had 1999 Borehole been advanced in primarily confirm to benzo(b/k)fluoranthene and bis(2-ethylhexyl)phthalate concentrations detected below CROLs in a 1988 groundwater sample collected from monitoring well FS-21/MW-2, which had since been destroyed. These contaminants were not detected in the 1999 borehole sample that was advanced in the approximate same location as FS-21/MW-2. The presence of benzo(b/k)fluoranthene in the historical groundwater sample collected from well FS-21/MW-2 may be due to soil derived suspended solids (i.e., turbidity) that were introduced into the groundwater during sampling using non-low-flow techniques. The groundwater Field Sample Data Record for this 1988 sample stated that the groundwater sample appearance was "light brown with moderate suspended solids" (E.C. Jordan, 1990a). As a result of this data, benzo(b/k)fluoranthene was not considered a site-related contaminant of potential concern and was not evaluated in the PRE/PRA.

Inorganic Analytes: In a comprehensive study of inorganic concentrations in groundwater at the MMR conducted in 1996-1997, where 158 groundwater monitoring wells were sampled and analyzed for inorganic analytes (*MMR Inorganics Investigation Technical Memorandum, Jacobs Engineering, 1998*), it was determined that samples collected using the USEPA Region I Low Flow (Minimum Stress) Purging and Sampling Procedure (USEPA, 1996a) consistently have lower concentrations of suspended solids, and therefore, lower total metals concentrations than samples collected using more traditional (high flow) sampling techniques (Jacobs, 1998). The study also found that metals data representing MMR "background" conditions were probably biased on the high side by the method used to collect the groundwater samples, and that these data are suspect.

Because groundwater samples collected prior to 1999 at the site did not use the Low Flow (Minimum Stress) Purging and Sampling Procedure, this pre-1999 laboratory data was not considered reliable. Only data obtained during the March and December 1999 sampling rounds, which used the Low Flow (Minimum Stress) Purging and Sampling Procedure was used to evaluate the site. The March 1999 sampling round sampled wells CS-8/MW-2 and FS-21/MW-1. Manganese (708 μ g/l) and Beryllium (0.94J μ g/l) were detected at concentrations above groundwater screening criteria, although the Beryllium detected was below sample quantitation limits. The December 1999 sampling round collected groundwater from borehole 04BH001 and detected arsenic (7.4J μ g/l) above groundwater screening criteria, although it was also below sample quantitation limits.

TABLE 4-3 SUMMARY OF ANALYTES DETECTED IN GROUNDWATER - OFF-SITE ANALYSIS

STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| "Exploration ID: | | CS-8/MW-2* | | CS-8/MW-2* | | | | CS-8/MW-4* | | | | 04BH0001 |
|---------------------------------------|-----------|------------|------------|------------|---------------|-----------|-----------|------------|-----------|-----------|-----------|-----------|
| "Sampled: | 07-JAN-88 | 07-JAN-88 | 28-JUNE-89 | 28-JUNE-89 | 10-Mar-99 | 10-Mar-99 | 28-OCT-87 | 28-OCT-87 | 28-OCT-87 | 10-Mar-99 | 07-JAN-88 | 14-Dec-99 |
| VOLATILE ORGANIC COMPOUNDS (µg/L) | | | | | | | | | | | | |
| 2-Butanone | | _ | 6400 | 3800 | - | _ | | | _ | _ | NR | R |
| 2-Hexanone | | - | - | - | - | - | - | | - | - | NR | Ŕ |
| Acetone | - | - | í - ' | - | - | - 1 | - 1 | - | - | - | NR | R |
| Ethylbenzene | - | - | - | - | - | - | - | - | - | - | NR | 1.14 |
| Xylenes | - | - | - | - | - | - | - | - | - | - | NR | 3.04 |
| SEMIVOLATILE ORGANIC COMPOUNDS (µg/L) | | | | | | |] |) | | | | |
| Benzo(a)pyrene | - | - | - | - | - | - | - 1 | - | - | - | - 1 | R |
| Benzo(b)/(k)fluoranthene | - | - 1 | - 1 | - 1 | NR | NR | - | | - | NR | 3J | - |
| Chyrsene | - | - | - 1 | - | NR | NR | - | - 1 | - | NR | 2J | - |
| Naphthalene | - | _ | 2J | 2.1 | NR | NR | - 1 | | - | NR | - | |
| Fluoranthene | - | _ | | | NR | NR | - | - 1 | - | NR | 3J | - 1 |
| Pyrene | - | _ | - | - | NR | NR | - | _ | - | NR | 3.1 | - 1 |
| Bis (2-ethylhexyl)phthalate | 2J | 4J | - | - | NR | NR | - | - | - | NR | 13J | - |
| INORGANIC COMPOUNDS (µg/L) | | | | | | | | | | | | |
| Aluminum | R | R | R | R | - | - | R | R | R | 1180 | R | |
| Arsenic | R | R | R | R | _ | _ | R | R | R | | R | 7.4J |
| Barium | R | R | R | R | 17.2 J | 17.1 J | R | R | R | 70.4 | R | 7.40 |
| Beryllium | R | R | R | R | 11.20 | 17.10 | R | R | R | 0.94 J | R | |
| Cadmium | | R | R | R | 0.35 J | 0.62 J | R | | R | 0.4 J | | - |
| Calcium | | R | R | R | 2590 | 2620 | R | R | R | 6030 | R | - 1 |
| | <u> </u> | | R | | 2590 287 J | | R | R | R | | | - |
| Magnésium | ĸ | R | | R | | 280 J | | | | 1200 | R | - |
| Manganese | к | R | R | R | 34.1 | 35.2 | R | R | R | 708 | R | - |
| Potassium | R | R | R | R | 1120 J | 1180 J | R | R | R | 1300 J | R | - |
| Sodium | R | R | R | R | 18800 | 19000 | R | R | R | 21700 | R | - |
| Zinc | R | R | R | R | - | 8.4 J | R | R | R | 19.3 J | R | 29.6 |
| TOTAL PETROLEUM HYDROCARBONS (mg/L) | NR | NR | - | - | NR | NR | NR | NR | NR | NR | NR | NR |

NOTES:

not detected

* a duplicate sample

J = estimated concentration

mg/L. × milligrams per liter

MW ≈ montoring well

NR ≈ not requested

R ≈ rejected (unusable results)

µg/L ≈ micrograms per liter

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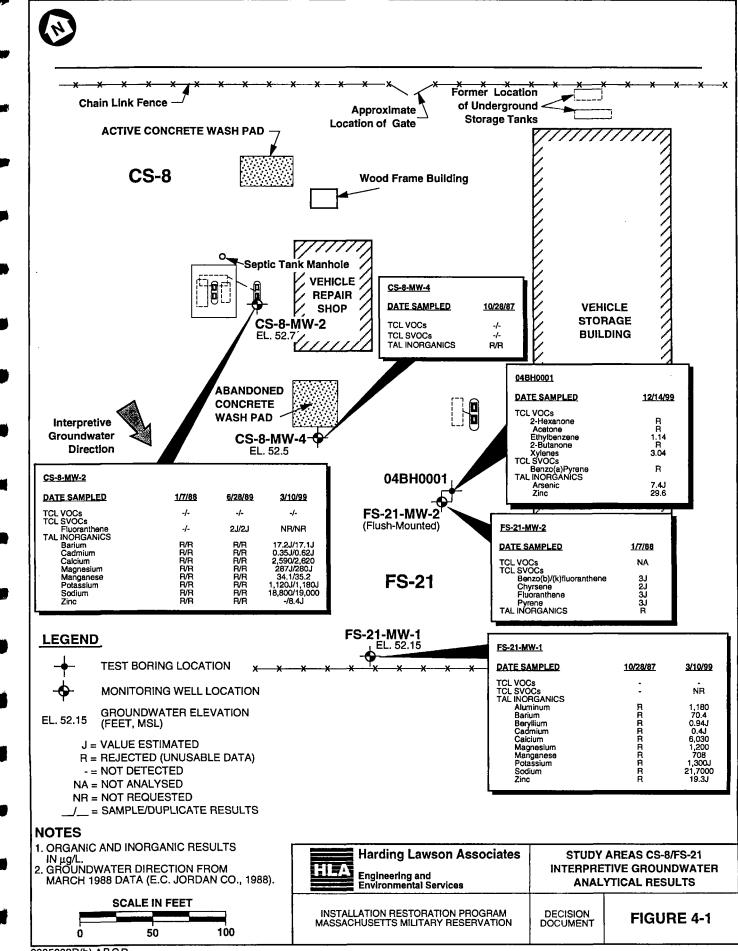
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5.0 PRELIMINARY RISK EVALUATION

This section describes how the human-health and ecological PREs were conducted and presents the results of each.

5.1 INTRODUCTION

A PRE was conducted at Study Areas CS-8/FS-21 to provide a screening-level assessment of potential risks to human and ecological receptors associated with exposure to environmental contaminants identified during site characterization efforts.

5.1.1 Approach

A PRE is a worst-case analysis that addresses the most sensitive receptors and all potential current and future pathways. The human health PRE was performed using current USEPA Region I risk assessment guidance (USEPA, 1989b; 1994; 1995; 1996; 1999a, 2000c). The ecological PRE followed the approach outlined in the MMR Risk Assessment Handbook (RAH) (Automated Sciences Group, Inc. [ASG], 1994).

Data were evaluated for surface soil, subsurface soil, and groundwater collected at Study Areas CS-8/FS-21 during the SI and subsequent sampling events. The data were determined to be sufficient to complete human-health and ecological PREs at these study areas. The ecological PRE considered exposures to surface soil only; the human-health PRE included an evaluation of exposures to surface soil, subsurface soil, and groundwater. No surface water or sediment exists at this site.

The human health PRE was performed as a two-step process. First, analytes that could pose a health risk of concern were identified by comparing the maximum detected concentrations to conservative, health-based screening values. Analytes with maximum detected concentrations that were less than the screening values were considered to not pose a health risk of concern (i.e., they posed a *de minimus* risk) and therefore, were not further evaluated. Analytes with maximum detected concentrations that exceeded the screening values were retained for further evaluation as chemicals of potential concern (CPCs). In the second step, analytes selected as CPCs were evaluated to assess whether potential exposure to the CPCs would be associated with acceptable cumulative cancer and noncancer risks. This evaluation was performed by identifying an exposure point

mmr/docs/cs8dd/text/cs8-fs21 Final DD2. doc

concentration (EPC) for each CPC, and then quantitatively evaluating exposure doses and risks using default residential exposure and toxicity assumptions recommended by USEPA. The cumulative risks were compared to USEPA and MADEP cumulative receptor risk limits to determine if the health risks meet risk management criteria.

The ecological PRE also followed a tiered approach. In the Tier I analysis, maximum reported analyte concentrations were compared to Hazard Equivalent Concentration (HECs) based on the most sensitive receptor species likely to inhabit the study areas. If maximum site concentrations exceed HECs for Tier I, a remedial investigation (RI) can be recommended without Tier II evaluation. Otherwise, a Tier II evaluation is completed to compare maximum concentrations to HECs for indicator species likely to inhabit the study areas based on future land-use plans.

5.1.2 Study Areas CS-8/FS-21 Objectives

The purposes of this PRE were as follows:

- select CPCs
- identify human and ecological exposure pathways
- identify human and ecological receptors
- evaluate possible human health risks by calculating quantitative risk estimates for potential exposures to CPCs and by comparing CPCs at Study Areas CS-8/FS-21 to applicable or relevant and appropriate requirements
- evaluate possible ecological risks by comparing CPCs at Study Areas CS-8/FS-21 to HECs

The PRE was conducted to support the recommendation of one of the following alternatives: (1) decision document (no further action), (2) RI/feasibility study, or (3) removal action.

5.2 HUMAN-HEALTH PRELIMINARY RISK EVALUATION

The purpose of the PRE for Study Areas CS-8/FS-21 was to identify potential human-health risks associated with exposures to study-area-related CPCs.

5.2.1 Data Evaluation

Analytical data available for Study Areas CS-8/FS-21 were evaluated to identify studyarea- and medium-specific CPCs. Frequencies of detection, ranges of CRQLs and CRDLs, and a range of minimum to maximum detected concentrations of analytes detected in surface soil, subsurface soil, and groundwater are listed in Tables 5-1, 5-2, and 5-3, respectively.

Surface soil data consisting of soil samples collected 0-2 ft bgs are identified in Table 4-1 and are summarized in Table 5-1. Surface soil data were collected in accordance with the SI workplan for CS-8/FS-21.

In accordance with USEPA Region I guidance (USEPA, 1995), subsurface soils considered for evaluation in health risk assessment should be represented by the depth intervals that could realistically be contacted by individuals under the current and foreseeable future land uses. USEPA Region I has defined this depth as up to 10 ft bgs. Subsurface soil data for soil samples collected 2-10 ft bgs identified in Table 4-2a were used to develop the data summary presented in Table 5-2.

As discussed in Section 4.3, 2-butanone was detected in groundwater samples collected in 1989 from monitoring well CS8/MW2. The lack of 2-butanone detected in groundwater from adjacent wells and from groundwater collected from CS8/MW2 in 1999 indicates that its presence in 1989 was likely due to sample contamination introduced during collection and/or laboratory analysis (Haase, et al., 1988 and E.C. Jordan Co., 1991). As a result of these findings, 2-butanone was not considered a site-related contaminant of potential concern and was not evaluated in the PRE/PRA.

Soil samples collected from the TerraProbe borings installed around the cesspool in 1992 were analyzed in an on-site laboratory to characterize the extent of soil contamination in support of a removal action. Samples were not submitted for off-site analysis. Therefore, the field screening soil data collected from the TerraProbe boring locations were qualitatively evaluated by a comparison of detected concentrations to USEPA Region IX residential soil preliminary remediation goals (PRGs). Detected concentrations of all VOCs, copper, and zinc are substantially lower than their respective residential soil PRGs. The greatest TPH concentrations (820 mg/kg and 980 mg/kg in the 8-foot bgs sample from 04TRX12) are slightly greater than the MADEP S-1 Method 1 standard for TPH (800 mg/kg) provided in the Massachusetts Contingency Plan, 310 CMR 40.0985(6)

(MADEP, 1999). The concentrations of delta-BHC and heptachlor epoxide exceed the USEPA Region IX residential soil PRGS. In 1996, the cesspool and approximately 40 cubic yards of contaminated soil were removed during the DSRP. The four TerraProbe boring locations were located in the area of the removal action. Confirmation soil samples collected during the removal action from the sidewalls and bottom of the excavation were submitted for off-site laboratory analysis. The results were all below the cleanup levels established for the removal action, as well as the USEPA Region IX residential soil PRGs.

The groundwater data set (see Table 5-3) consists of 8 groundwater samples collected from four monitoring wells. There are three rounds of data from CS-8/MW-2, one round of data from CS-8/MW-4, two rounds of data from FS-21/MW-1, one round of data from FS-21/MW-2, and one round of data from 04BH001 (a borehole advanced at the former location of FS-21/MW-2, which had been destroyed). Groundwater data from monitoring wells sampled during and after 1995 were obtained using low-flow sampling techniques. Therefore, only inorganics data from the groundwater sampling rounds at CS-8/MW-2 (March 1999), FS-21/MW-1 (March 1999), and 04BH001 (December 1999) were used in the PRE; inorganics data for non-low-flow sampling performed prior to 1995 were not evaluated in the PRE. Groundwater data for semivolatile organics obtained from the January 1988 sampling round at FS-21/MW-2 were not used in the PRE because the presence of these analytes was determined to be an artifact of suspended solids in the sample. The December 1999 sample round at this well location was analyzed for SVOCs, and none were detected. Therefore, the data for the December 1999 sampling round at 04BH001 were used in the PRE.

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| TABLE 5-1 |
|---|
| CONTAMINANTS OF POTENTIAL CONCERN IN SURFACE SOIL |

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| | | | RANGE OF | MAXIMUM | REGION IX | | |
|--------------------------------|-----------|---------|----------------|------------|-------------|----------|-------|
| | FREQUENCY | CRQL/ | DETECTED | BACKGROUND | RESIDENTIAL | SELECTED | 1 |
| | OF | CRDL | CONCENTRATIONS | CONC. | PRG | AS A | ł |
| CHEMICAL | DETECTION | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | CPC? | NOTES |
| SEMIVOLATILE ORGANIC COMPOUNDS | | | | | | | |
| Bis(2-ethylhexyl)phthalate | 1/2 | 0.33 | 0.064 - 0.064 | NA | 35 | NO | 2 |
| Butylbenzylphthalate | 1/2 | 0.33 | 0.14 - 0.14 | NA | 1,200 | NO | 2 |
| PCBs/PESTICIDES | | | | | | | |
| Aroclor-1260 | 1/2 | 0.16 | 0.65 - 0.66 | NA | 0.22 | YES | 1 |
| INORGANICS | | | | | | | |
| Aluminum | 2/2 | 40 | 1,300 - 3,220 | 8,930 | 7,600 | NO | 2 |
| Arsenic | 2/2 | 2 | 0.69 - 2.2 | 3.6 | 0.39 | YES | 1 |
| Barium | 2/2 | 40 | 9.6 - 11 | 10.4 | 540 | NO | 2 |
| Beryllium | 1/2 | 1 | 0.37 - 0.37 | 0.65 | 15 | NO | 2 |
| Calcium | 2/2 | 1,000 | 218 - 331 | 969 | NA | NO | 3 |
| Chromium | 2/2 | 2 | 2.1 - 5.3 | 6.8 | 30 | NO | 2 |
| Cobalt | 1/2 | 10 | 1.5 - 1.5 | 4.1 | 470 | NO | 2 |
| Copper | 1/2 | 5 | 9.2 - 9.2 | 5.2 | 290 | NO | 2 |
| Iron | 2/2 | 20 | 2,490 - 4,010 | 12,400 | 2,300 | NO | 3 |
| Lead | 2/2 | 3 | 5.6 - 6.9 | 12.05 | 400 | NO | 2 |
| Magnesium | 2/2 | 1,000 | 178 - 536 | 794.5 | NA | NO | 3 |
| Manganese | 2/2 | 3 | 31 - 59 | 108 | 180 | NO | 2 |
| Thallium | 1/2 | 2 | 0.4 - 0.4 | 0.25 | 0.63 | NO | 2 |
| Vapadium | 2/2 | 10 | 3.7 - 6.4 | 15.2 | 55 | NO | 2 |
| Zinc | 2/2 | 4 | 17 - 30 | 16 | 2,300 | NO | 2 |

NOTES:

NA = not applicable

CPC = contaminant of potential concern

CRQL = Contract Required Quantitation Limit

CRDL = Contract Required Detection Limit

mg/kg ≃ milligrams per kilogram

PRG - Preliminary Remediation Goal. Values are obtained from the Region IX PRG Table (USEPA, 1999). PRGs presented in this table are based on the lesser of a cancer risk of 1E-06 or a non-cancer bazard index of 0.1.

1 Chemical selected as a CPC because the maximum detected concentration exceeds the Region IX PRG.

2 Chemical not retained as a CPC because the maximum detected concentration is less than the Region IX PRG.

3 Chemical not retained as a CPC because it is an essential nutrient.

TABLE 5-2

CONTAMINANTS OF POTENTIAL CONCERN IN SUBSURFACE SOIL (2-10' bgs)

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| CHEMICAL | FREQUENCY OF DETECTION | CRQL/ CRDL (mg/kg) | RANGE OF DETECTED CONCENTRATIONS (wg/kg) | MAX. BKGRD CONC. (mg/kg) | REGION IX RESIDENTIAL PRG (mg/kg) | SELECTED AS A CPC? | NOTES |
|----------------------|------------------------------|--------------------------|---|-----------------------------------|--|--------------------------|-------|
| EMIVOLATILE | | | | | | | |
| RGANIC COMPOUNDS | | | | J J | | | |
| Butylbenzylphthalate | 1/2 | 0.34 | 0.34 | NA | 1200 | NO | 2 |
| NORGANICS | | | | | | | |
| Atuminum | 3/3 | - | 551 - 914 | 1,980 | 7,600 | NO | 2 |
| Arsenic | 1/3 | 2 | 0.55 J | 2.3 | 0.39 | YES | 1 |
| Barium | 3/3 | - | 1.7J - 4.0J | 14.7 | 540 | NO | 2 |
| Calcium | 3/3 | - | 27J - 81J | 933 | NA | NO | 3 |
| Chromium | 1/3 | 2 | 4.3 | 3.9 | 30 | NO | 2 |
| Copper | 2/3 | 5 | 0.94J - 4.1J | 4.3 | 290 | NO | 2 |
| Iron | 3/3 | - | 893J - 2020J | 2,600 | 2,300 | NO | 3 |
| Lead | 3/3 | - | 0.98J - 1.8 | 3.7 | 400 | NO | 2 |
| Magnesium | 3/3 | • | 129J - 354J | 742 | NA | NO | 3 |
| Manganese | 3/3 | - | 19J - 53J | 587 | 180 | NO | 2 |
| Thallium | 1/2 | 2 | 0.41J | { 1 | 0.63 | NO | 2 |
| Vanadium | 2/3 | 10 | 1.6J - 2.0J | 5.7 | 55 | NO | 2 |
| Zinc | 3/3 | - | 5.0 - 8.7J | 16 | 2,300 | NO | 2 |

NOTES:

NOTES: NA = not applicable CPC = contaminant of potential concern CRQL = Contract Required Quantitation Limit

CRDL = Contract Required Detection Limit

PRG - Prolingrams per kilogram PRG - Preliminary Remediation Goal. Values are obtained from the Region IX PRG Table (USEPA, 1999). PRGs presented in this table are based on the lesser of a cancer risk of 1E-06 or a non-cancer hazard index of 0.1.

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1 Chemical selected as a CPC because the maximum detected concentration exceeds the Region IX PRG. 2 Chemical not retained as a CPC because the maximum detected concentration is less than the Region IX PRG.

Chemical not retained as a CPC because it is an essential nutrient.
 Analyte was detected in all samples analyzed.

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TABLE 5-3 CONTAMINANTS OF POTENTIAL CONCERN IN GROUNDWATER

STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| CHEMICAL | FREQUENCY OF DETECTION | CRQL/CRDL (ug/L) | RANGE OF DETECTED CONCENTRATIONS (ug/L) | REGION IX TAPWATER PRG (ug/L) | USEPA MCL (ug/L) | MADEP MCL (ug/L) | SELECTED AS A CPC? | NOTES |
|---------------------------------|---------------------------|---------------------|---|-------------------------------------|------------------------|------------------------|-----------------------|-------|
| | | | | | | | | |
| Volatile Organic Compounds | | | | | | | | l |
| Ethylbenzene | 1/7 | 1 | 1.14 - 1.14 | 130 | 700 | 700 | NO | 2 |
| Xylene | 1/7 | 1 | 3.04 - 3.04 | 140 | 10000 | 10000 | NO | 2 |
| Semi-Volatile Organic Compounds | | | | | | | | |
| Bis(2-ethylhexyl)phthalate | 1/5 | 10 | 2 - 4 * | 4.8 | 6 | 6 | NO | 2 |
| Naphthalene | 1/5 | 10 | 2 - 2 | 0.62 | NA | NA | YES | 1 |
| Inorganic Compounds | | | | | | | | |
| Arsenic | 1/3 | 10 | 7.4 - 7.4 | 0.045 | 50 | 50 | YES | 1 1 |
| Aluminum | 1/2 | 80 | 1180 - 1180 | 3600 | NA | NA | NO | 2 |
| Barium | 2/2 | | 17.2 - 70.4 | 260 | 2000 | 2000 | NO | 2 |
| Beryllium | 1/3 | 5 | 0.94 - 0.94 | 7.3 | 4 | 4 | NO | 2 |
| Cadmium | 2/3 | 5 | 0.4 - 0.48 | 1.8 | 5 | 5 | NO | |
| Calcium | 2/2 | - | 2325 - 6030 | NA | NA | NA | NO | 3 |
| Magnesium | 2/2 | _ | 283 - 1200 | NA | NA | NA | NO | 3 |
| Manganese | 2/2 | - | 34.7 - 708 | 88 | NA | NA | YES | |
| Potassium | 2/2 | - | 1150 - 1300 | NA | NA | NA | NO | |
| Sodium | 2/2 | - | 18,900 - 21,700 | NA | NA | NA | NO | 3 |
| Zinc | 3/3 | 20 | 8.4 - 29.3 | 1100 | NA | NA | NO | 2 |

NOTES:

NA = Not available

CPC = contaminant of potential concern

CRQL = contract required quantitation limit

CRDL = contract required detection limit

ug/L = mircrograms per liter

* = range for original and duplicate results

PRG - Preliminary Remediation Goal. Values are obtained from the Region IX PRG Table (USEPA, 1999). PRGs presented in this table are based on the lesser of a cancer risk of 1E-06

or a non-cancer hazard index of 0.1.

MCL - Maximum Contaminant Level. USEPA primary drinking water standard (USEPA, 2000); Massachusetts primary drinking water standard (MADEP, 2000).

1 Chemical selected as a CPC because the maximum detected concentration exceeds the Region IX PRG or MCL.

2 Chemical not retained as a CPC because the maximum detected concentration is less than the Region IX PRG and MCL.

3 Chemical not retained as a CPC because it is an essential nutrient.

'-' = Analyte was detected in all samples analyzed.

5.2.2 Contaminants of Potential Concern

CPCs are analytes that are potentially related to contamination sources at the site and are present at concentrations that may pose a health risk of concern. CPCs were identified by comparing maximum detected site concentrations to health-based screening values. This methodology results in a conservative evaluation that does not overlook or dismiss analytes that could pose potentially substantial risks.

The procedure used to select CPCs is summarized as follows, and is consistent with USEPA Region I (1999) methodology:

- <u>Selected</u> as a CPC in **surface soils** if the maximum detected concentration exceeds the USEPA Region IX PRG for residential soils (USEPA, 1999b).
- <u>Selected</u> as a CPC in **subsurface soils** if the maximum detected concentration exceeds the USEPA Region IX PRG for residential soils (USEPA, 1999b).
- <u>Selected</u> as a CPC in **groundwater** if the maximum detected concentration exceeds the lesser of the USEPA Region IX PRG for tapwater (USEPA, 1999b), the USEPA MCLs (USEPA, 2000a) and Massachusetts MCLs (MADEP, 2000) for drinking water (the potentially applicable ARAR).
- <u>Eliminated</u> as a CPC if an essential nutrient. The following inorganic analytes are considered essential human nutrients: calcium, magnesium, iron, potassium, and sodium.

The PRGs are protective for direct contact (ingestion and dermal contact) exposures, as well as for inhalation of dusts and vapors. The PRGs are derived for a 1×10^{-6} cancer risk level or a non-cancer hazard quotient (HQ) of 1. Per USEPA Region I guidance (USEPA, 1995), the PRGs based on noncarcinogenic effects have been adjusted for a HQ of 0.1 for the purposes of CPC selection.

As shown in Table 5-1, aroclor-1260 and arsenic were selected as CPCs in surface soil. Arsenic was selected as a CPC in subsurface soil (see Table 5-2). In groundwater, naphthalene, arsenic, and manganese were selected as CPCs (see Table 5-3). The COPCs were further evaluated to assess whether they posed cumulative risks for unrestricted land use in excess of the USEPA and MADEP cumulative risk limits.

5.2.3 Cumulative Receptor Risks

Study Area CS-8/FS-21 is located outside the MMR Security Zone/Flightline Area. Therefore, it is assumed that future land use could include residential development. To provide a conservative assessment of potential future land uses and health risks associated with possible exposures to the CPCs in surface soil, subsurface soil, and groundwater, cumulative receptor risks are estimated for residential land use.

Cumulative receptor risks were calculated to assess whether the CPCs are present at concentrations that could pose a significant health risk for the future residential use of the site. This determination was made by comparing the calculated risks for a residential exposure scenario with the USEPA cumulative receptor risk limits of an excess lifetime cancer risk of 1×10^{-6} to 1×10^{-4} and a hazard index (HI) of 1 for noncancer risks, and the MADEP cumulative receptor risk limit of an excess lifetime cancer risk of 1×10^{-5} and a HI of 1 for noncancer risks.

Calculation of cumulative receptor risks involved four components:

- 1. Identification of CPC exposure point concentrations
- 2. Calculation of receptor CPC intakes
- 3. Quantification of CPC toxicity
- 4. Calculation of cancer and noncancer risks

5.2.3.1 Calculation of CPC Exposure Point Concentrations

In accordance with USEPA guidance, the EPCs for soil were based on the lesser of the 95% upper confidence limit (UCL) on the arithmetic mean concentration or the maximum detected concentration. Because the surface soil and subsurface soil data sets each have less than 10 samples, the 95% UCL statistic was not used (USEPA, 1992). Therefore, the maximum detected concentrations were used as the EPCs for soil. For groundwater, USEPA Region I requires that each monitoring well be treated as a separate exposure point. Therefore, the CPC concentrations reported in each monitoring well represent the EPCs. To streamline the risk calculation, the maximum detected concentration of each CPC from among all monitoring wells was used as the EPC. The health risks calculated using this approach are equal to or higher than the risks that would be associated with any single well.

5.2.3.2 Calculation of Receptor CPC Intakes

Intakes and risks for residential land use were quantified for child and adult resident scenarios in accordance with USEPA Region I risk characterization guidance (USEPA, 1994). It is assumed that adults and children who may live at the site in the future could be exposed to CPCs in surface soil through incidental ingestion, dermal contact, and inhalation of soil particulates with sorbed SVOCs or inorganic compounds. It is further assumed that subsurface soils may be relocated to the surface, thereby becoming accessible via these same exposure pathways. Groundwater is considered a potable water source. Therefore, exposures to groundwater via ingestion of the water as residential tapwater and dermal contact during showering are evaluated. Since VOCs were not selected as CPCs in groundwater, possible volatile inhalation exposures to groundwater are not evaluated (i.e., since VOCs were not selected as CPCs, they would not pose a risk of concern via this exposure pathway).

The exposure parameters used to quantify CPC intakes are provided in Table 5-4, and are based on USEPA Region I default exposure variables (USEPA, 1994). In summary, child and adult residents are assumed to contact soils 150 days per year, over a 30-year period.

To evaluate non-cancer effects to sensitive individuals, the 30-year exposure is evaluated separately for young children (exposures to children ages 1 through 6 over a six-year duration) and adults (exposures over a 24-year period). Use of groundwater as a potable source is evaluated for an adult resident receptor that is assumed to ingest 2 liters of water per day, 350 days per year, over a 30-year period.

Dermal exposures to soil were calculated using the recent USEPA RAGS Part E Supplemental Guidance for Dermal Risk Assessment (Interim) (USEPA, 2000c). As recommended in this guidance, the skin surface areas and soil adherence factors are 2,800 cm^2 and 0.2 mg/cm², respectively for children, and 5,700 cm² and 0.07 mg/cm², respectively, for adults. Dermal absorption factors for the CPCs in surface and subsurface soil (arsenic and arcolor-1260) are also published in this guidance (USEPA, 2000c). The dermal exposure assessment procedure described in the RAGS Part E interim guidance for the water pathway was also applied to the residential water use scenario. Based on these procedures, arsenic and manganese dermal exposures and risks are considered relatively insignificant (i.e., dermal dose < 10% ingestion dose) and are not quantified. Naphthalene dermal exposures and risks are quantified using the procedure described in the recent draft guidance.

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TABLE 5-4 **EXPOSURE PARAMETERS - RESIDENT**

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| PARAMETER | ADULT RESIDENT | CHILD RESIDENT | UNITS | SOURCE 1 |
|---|-------------------|-------------------|----------------------|-------------|
| Soil Ingestion Rate | 100 | 200 | mg/day | USEPA, 1994 |
| Fraction Ingested From Site | 100% | 100% | | Assumption |
| Drinking Water Ingestion Rate | 2 | NA | L/day | USEPA, 1994 |
| Exposure Time - outdoor ² | 8 | 8 | hours/day | Assumption |
| Exposure Time - bathing 7 | 0.58 | NA | hours/day | USEPA, 2000 |
| Exposure Frequency - outdoor ² | 150 | 150 | days/year | USEPA, 1994 |
| Exposure Frequency - indoor ³ | 350 | 350 | days/year | USEPA, 1994 |
| Exposure Duration | 24 | 6 | years | USEPA, 1994 |
| Body Weight | 70 | 15 | kg | USEPA, 1994 |
| Averaging Time | | | | |
| Cancer | 70 | 70 | years | USEPA, 1989 |
| Noncancer ⁵ | 24 | 6 | years | USEPA, 1989 |
| Surface Area Exposed - Soil ⁴ | 5700 | 2800 | cm ² /day | USEPA, 2000 |
| Surface Area Exposed - Groundwater ⁴ | 18000 | NA | cm²/day | USEPA, 2000 |
| Inhalation Rate ⁶ | 0.63 | 0.35 | m³/hour | USEPA, 1997 |
| Permeability Coefficient | chemical-specific | NA | cm/hour | USEPA, 2000 |
| Dermal Absorption Factor | chemical-specific | chemical-specific | ABSd | USEPA, 2000 |
| Soil Adherence Factor ⁴ | 0.07 | 0.2 | mg/cm ² | USEPA, 2000 |
| Particulate Emission Factor | 1.32E+09 | 1.32E+09 | m³/kg | USEPA, 1996 |

Notes:

1 - Exposure variables with source listed as "assumption" are site specific; the remainder are default values.

2 - Values used to calculate exposure to soil and soil-derived dust

- 3 Values used to calculate exposure to indoor air (vapors)
- 4 Values recommended for residential exposures
- 5 The AT for noncarcinogenic effects is equal to the exposure duration

6 - Adult value is recommended chronic value for adults (15.2 m³/24-hour period);

child value is recommended value for 3-5 year old children (8.3 m³/24-hour period). Values converted to rate/hour by dividing by 24-hours. 7 - RME value for time spent in a shower.

mg - milligrams

cm² - square centimeters

m³ - cubic meters

kg - kilograms

NA - Not Applicable; exposure not evaluated.

RME - Reasonable maximum exposure

USEPA, 1989. "Risk Assessment Guidance for Superfund, Volume 1, Human Health Evaluation Manual (Part A)"; Office of Emergency and Remedial Response; EPA-540/1-89/002 (interim final); Washington, D.C., December.

USEPA, 1991. "Risk Assessment Guidance for Superfund, Volume I: Human Health Evaluation Manual, Supplemental Guidance

Standard Default Exposure Factors"; Office of Emergency and Remedial Response, Toxics Integration Branch; OSWER Directive 9285.6-03

USEPA, 1994. "Risk Updates No. 2"; USEPA Region I, Waste Management Division; August. Values from "Attachment 2" to Risk Updates No. 2. USEPA, 1996. "Soil Screening Guidance: Technical Background Document' EPA/540/R-95/128. May.

USEPA, 1997. "Exposure Factors Handbook, Volume 1"; Office of Research and Development; EPA-600/P-95/002Fa; Washington, D.C.; August. USEPA, 2000. Risk Assessment Guidance for Superfund. Volume I: Human Health Evaluation Manual Supplemental Guidance (Part E, Dermal Risk Assessment) Interim Guidance

Tables 5-5 through 5-9 document the intake calculations and risk calculations.

5.2.3.3 Quantification of CPC Toxicity

Risks were calculated by combining CPC intakes with dose-response data that quantify the toxicity associated with each CPC. Cancer slope factors developed by USEPA (USEPA, 1997; 2000b) were used for evaluating carcinogenic effects. Chronic reference doses developed by USEPA (USEPA, 1997 and 2000b) were used for evaluating noncancer effects. For evaluating risks associated with dermal exposures to soil, dermal cancer slope factors and reference doses were calculated using the oral absorption factors and technical approach recommended in USEPA guidance (USEPA, 1999a and 2000c). In accordance with this guidance, manganese is the only CPC for which dermal doseresponse values are calculated. The dose-response values are presented in Tables 5-5 through 5-9.

5.2.3.4 Calculation of Cancer and Noncancer Risks

Table 5-10 provides a summary of the risk estimates for the resident receptor scenarios. The cumulative cancer risks for soil (i.e., sum of risks for ingestion, dermal, and inhalation exposures) are 2.6×10^{-6} for a child resident and 1.2×10^{-6} for an adult resident. The aggregate resident receptor cancer risk for soil is 3.8×10^{-6} . Non-cancer hazard index values are below 1 for soil (0.03 for the adult and 0.3 for the child). The cancer risk values for potential exposures to soil are within the USEPA cancer risk range of 1×10^{-6} to 1×10^{-4} and are below the MADEP cumulative cancer risk limit of 1×10^{-5} . The HI values for potential exposures to soil do not exceed the USEPA and MADEP threshold non-cancer risk limits of an HI of 1.

The non-cancer screening HI for groundwater is 1.5, primarily due to manganese and arsenic. The HQ for each of those analytes is below one. The HQ for naphthalene is very low and does not have a significant impact on the hazard index values. If analytes have different mechanisms of action, it is appropriate to segregate the HQs for the analytes because their effects would not be considered additive (USEPA, 1989).

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TABLE 5-5 INCIDENTAL INGESTION OF AND DERMAL CONTACT WITH SURFACE SOIL - RME UNRESTRICTED LAND USE - ADULT RESIDENT STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

EXPOSURE PARAMETERS

EQUATIONS

| | PARAMETER | | | |
|--|--------------------------|-------------------|-----------|--|
| CONCENTRATION SOIL | CS | See Below* | mg/kg | CANCER RISK - INTAKE (mg/kg-day) x CANCER SLOPE FACTOR (mg/kg-day)-1 |
| INGESTION RATE | IR | 100 | mg/day | |
| FRACTION INGESTED | FI | 100% | | HAZARD QUOTTENT = INTAKE (mg/kg-day) / REFERENCE DOSE (mg/kg-day) |
| SOIL ADHERENCE FACTOR | SAF | 0.07 | mg/cm² | |
| SURFACE AREA EXPOSED | SA | 5,700 | cm² | INTAKE = (INTAKE-INGESTION) + (INTAKE-DERMAL) |
| CONVERSION FACTOR | CF | 0.000001 | kg/mg | |
| BODY WEIGHT | BW | 70 | kg | $INTAKE-INGESTION = \frac{CS \times IR \times FI \times CF \times EF \times ED}{CS \times IR \times FI \times CF \times EF \times ED}$ |
| EXPOSURE FREQUENCY | EF | 150 | days/year | BW x AT x 365 days/yr |
| EXPOSURE DURATION | ED | 24 | years | |
| AVERAGING TIME | | | | INTAKE-DERMAL = $CS \times SA \times SAF \times AE \times CF \times EF \times ED$ |
| CANCER | AT | 70 | years | BW x AT x 365 days/yr |
| NONCANCER | AT | 24 | years | |
| DERMAL ABSORPTION EFFICIENCY | AE | Chemical-specific | unitless | |
| Notes: | | | | |
| For noncarcinogenic effects: AT = ED | | | | |
| All exposure variables are documented in Tal | ole 5-4. | | | |
| *The lesser of the 95 % upper confidence lim | it (UCL) & maximum con | centration. | | |
| ND = Value not determined | NE = Route not evaluated | | | |

CARCINOGENIC EFFECTS

| COMPOUND | SOIL CONCENTRATION (mg/kg) | INTAKE INGESTION (mg/kg-day) | DERMAL ABSORPTION EFFICIENCY | INTAKE DERMAL (mg/kg-day) | CANCER SLO ORAL (mg/kg-day)-1 | PE FACTOR DERMAL (mg/kg-day)-1 | CANCER RISK INGESTION | CANCER RISK DERMAL | TOTAL CANCER RISK | PERCENT TOTAL RISK |
|--------------|----------------------------------|------------------------------------|------------------------------------|---------------------------------|-------------------------------------|--------------------------------------|--------------------------|-----------------------|-------------------------|--------------------------|
| Arsenic | 2.2 | 4.4E-07 | 0.03 | 5.3E-08 | 1.5E+00 | 1.5E+00 | 6.6E-07 | 8.0E-08 | 7.4E-07 | 64.23% |
| Aroclor-1260 | 0.66 | 1.3 E-0 7 | 0.14 | 7.4E-08 | 2.0E+00 | 2.0E+00 | 2.7E-07 | 1.5E-07 | 4.1E-07 | 35.77% |
| | <u></u> | | | | SUMMARY CAN | NCER RISK | 9.3E-07 | 2.3E-07 | 1.2E-06 | |

NONCARCINOGENIC EFFECTS

| COMPOUND | SOIL CONCENTRATION (mg/kg) | INTAKE INGESTION (mg/kg-day) | DERMAL ABSORPTION EFFICIENCY | INTAKE DERMAL (mg/kg-day) | REFEREN(ORAL (mg/kg-day) | CE DOSE DERMAL (mg/kg-day) | HAZARD QUOTIENT INGESTION | HAZARD QUOTIENT DERMAL | TOTAL HAZARD QUOTIENT | PERCENT TOTAL RISK |
|--------------|---------------------------------------|------------------------------------|------------------------------------|---------------------------------|---------------------------------|----------------------------------|---------------------------------|------------------------------|-----------------------------|--------------------------|
| Arsenic | 2.2 | 1.3E-06 | 0.03 | 1.5E-07 | 3.0E-04 | 3.0E-04 | 4.3E-03 | 5.2E-04 | 4.8E-03 | 13.77% |
| Aroclor-1260 | 0.66 | 3.9E-07 | 0.14 | 2.2E-07 | 2.0E-05 | 2.0E-05 | 1.9E-02 | 1.1E-02 | 3.0E-02 | 86.23% |
| | · · · · · · · · · · · · · · · · · · · | | | SUMMARY | HAZARD INDE | x | 2.4E-02 | 1.1E-02 | 3.5E-02 | |

TABLE 5-6 INHALATION EXPOSURE TO PARTICULATES IN SURFACE SOIL - RME UNRESTRICTED LAND USE - ADULT RESIDENT STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

EXPOSURE PARAMETERS

EQUATIONS

| PARAMETER | SYMBOL | VALUE | UNITS | |
|--|-------------------------------|------------------------|-----------|--|
| CONCENTRATION SOIL* | CS | See below | mg/kg | CANCER RISK - INTAKE (mg/kg-day) x CANCER SLOPE FACTOR (mg/kg-day)-1 |
| CONCENTRATION AIR PARTICULATES | САр | Calculated | mg/m³ | |
| CONCENTRATION AIR VOLATILES | CAv | Calculated | mg/m³ | HAZARD QUOTIENT = INTAKE (mg/kg-day) / REFERENCE CONCENTRATION (mg/kg-day) |
| VOLATILIZATION FACTOR** | VF | Calculated | m²/kg | |
| PARTICULATE EMISSIONS FACTOR | PEF | 1.32E+09 | ug/m² | |
| INHALATION RATE | IhR | 0.63 | m²/hour | INTAKE - INHALATION = <u>(CAp + Cay) x RAF x INR x EF x EF x ED</u> |
| BODY WEIGHT | BW | 70 | kg | BW x AT x 365 days/yr |
| EXPOSURE TIME | ET | 8 | hours/day | |
| EXPOSURE FREQUENCY | EF | 150 | days/year | |
| EXPOSURE DURATION | ED | 24 | years | AIR CONCENTRATION PARTICULATES ~ CS x 1/PEF |
| RELATIVE ABSORPTION FACTOR | RAF | 100% | | |
| AVERAGING TIME | | | | AIR CONCENTRATION VOLATILES – CS x 1/VF |
| CANCER | AT | 70 | years | (VF not calculated because there are no VOCs selected as CPCs). |
| NONCANCER | AT | 24 | years | |
| Notes: * Soil concentration used is the lesser of the 95 % upp | er confidence limit (UCL) & | maximum concentratio | n | 7 |
| **Volatilization factor used only for volatile chemicals of pote | ntial concern. | | | |
| For noncarcinogenic effects: AT = ED | All exposure variables are do | ocumented in Table 5-4 | | |
| ND = Value not determined | | | | |

CARCINOGENIC EFFECTS

| | SOIL. | | AIR CONCI | ENTRATION | | CANCER SLOPE | | PERCENT |
|--------------|---------------|---------|-----------|--------------|-------------|---------------|---------|---------|
| COMPOUND | CONCENTRATION | VF | VOLATILES | PARTICULATES | INTAKE | FACTOR | CANCER | TOTAL |
| | (mg/kg) | (m³/kg) | (mg/m³) | (mg/m³) | (mg/kg-day) | (mg/kg-day)-1 | RISK | RISK |
| Arsenic | 2.2 | NA | | 1.7E-09 | 1.7E-11 | 1.5E+01 | 2.5E-10 | 96.15% |
| Aroclor-1260 | 0.66 | NA | | 5.0E-10 | 5,1E-12 | 2.0E+00 | 1.0E-11 | 3.85% |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | SUMMARY | CANCER RISK | 2.6E-10 | |

NONCARCINOGENIC EFFECTS

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| | SOIL | | AIR CONCE | NTRATION | | REFERENCE | | PERCENT |
|--------------|---------------|---------|-----------|--------------|-------------|--------------|----------|---------|
| COMPOUND | CONCENTRATION | VF | VOLATILES | PARTICULATES | INTAKE | DOSE | HAZARD | TOTAL |
| | (mg/kg) | (m³/kg) | (mg/m³) | (mg/m²) | (mg/kg-day) | (mg/kg-day) | QUOTIENT | RISK |
| Arsenic | 2.2 | NA | | 1.7E-09 | 4.9E-11 | ND | | |
| Aroclor-1260 | 0.66 | NA | | 5.0E-10 | 1.5E-11 | 2.0E-05 | 7.4E-07 | 100.00% |
| | | | | | { } | | | |
| | | | | | | | | |
| | | | | | SUMMARY | HAZARD INDEX | 7.4E-07 | |

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TABLE 5-7

INCIDENTAL INGESTION OF AND DERMAL CONTACT WITH SURFACE SOIL - RME UNRESTRICTED LAND USE - CHILD RESIDENT (1 TO 6 YEARS) STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

EXPOSURE PARAMETERS

EQUATIONS

| PARAMETER | SYMBOL | VALUE | UNITS | |
|--|--------------------------|-------------------|-----------------|---|
| CONCENTRATION SOIL | CS | See Below* | mg/kg | CANCER RISK = INTAKE (mg/kg-day) x CANCER SLOPE FACTOR (mg/kg-day)-1 |
| INGESTION RATE | IR | 200 | mg/day | |
| FRACTION INGESTED | FI | 100% | | HAZARD QUOTTENT = INTAKE (mg/kg-day) / REFERENCE DOSE (mg/kg-day) |
| SOIL ADHERENCE FACTOR | SAF | 0.2 | mg/cm² | |
| SURFACE AREA EXPOSED | SA | 2,800 | Cm ² | INTAKE = (INTAKE-INGESTION) + (INTAKE-DERMAL) |
| CONVERSION FACTOR | CF | 0.000001 | kg/mg | |
| BODY WEIGHT | BW | 15 | kg | INTAKE-INGESTION = $CS \times IR \times FI \times CF \times EF \times ED$ |
| EXPOSURE FREQUENCY | EF | 150 | days/year | BW x AT x 365 days/yr |
| EXPOSURE DURATION | ED | 6 | years | |
| AVERAGING TIME | | | | INTAKE-DERMAL – $CS \pm SA \pm SAF \pm AE \pm CF \pm EF \pm ED$ |
| CANCER | AT | 70 | years | BW x AT x 365 days∕yr |
| NONCANCER | AT | 6 | years | |
| DERMAL ABSORPTION | AE | Chemical-specific | unitless | |
| EFFICIENCY | | | | |
| Notes: | | | | |
| For noncarcinogenic effects: AT = ED | | | | |
| All exposure variables are documented in Tab | ble 5-4. | | | |
| | | | | |
| The lesser of the 95 % upper confidence lim | it (UCL) & maximum cond | entration. | | |
| ND = Value not determined | NE = Route not evaluated | | | |

CARCINOGENIC EFFECTS

| | SOIL | INTAKE | DERMAL | INTAKE | CANCER SLO | PE FACTOR | CANCER RISK | CANCER RISK | TOTAL | PERCENT |
|--------------|---------------|-------------|------------|-------------|---------------|---------------|-------------|-------------|---------|---------|
| COMPOUND | CONCENTRATION | INGESTION | ABSORPTION | DERMAL | ORAL | DERMAL | INGESTION | DERMAL | CANCER | TOTAL |
| | (mg/kg) | (mg/kg-day) | EFFICIENCY | (mg/kg-day) | (mg/kg-day)-1 | (mg/kg-day)-1 | | | RISK | RISK |
| Arsenic | 2.2 | 1.0E-06 | 0.03 | 8.7E-08 | 1.5E+00 | 1.5E+00 | 1.5E-06 | 1.3E-07 | 1.7E-06 | 66.07% |
| Aroclor-1260 | 0.66 | 3.1E-07 | 0.14 | 1.2E-07 | 2.0E+00 | 2.0E+00 | 6.2E-07 | 2.4E-07 | 8.6E-07 | 33.93% |
| | | | | | | | | | | |
| | | | | | SUMMARY CAN | NCER RISK | 2.2E-06 | 3.7E-07 | 2.5E-06 | |

NONCARCINOGENIC EFFECTS

| COMPOUND | SOIL CONCENTRATION (mg/kg) | INTAKE INGESTION (mg/kg-day) | DERMAL ABSORPTION EFFICIENCY | INTAKE DERMAL (mg/kg-day) | REFERENC ORAL (mg/kg-day) | CE DOSE DERMAL (mg/kg-day) | HAZARD QUOTIENT INGESTION | HAZARD QUOTIENT DERMAL | TOTAL HAZARD QUOTIENT | PERCENT TOTAL RISK |
|--------------|----------------------------------|------------------------------------|------------------------------------|---------------------------------|---------------------------------|----------------------------------|---------------------------------|------------------------------|-----------------------------|--------------------------|
| Arsenic | 2.2 | 1.2E-05 | 0.03 | 1.0E-06 | 3.0E-04 | 3.0E-04 | 4.0E-02 | 3.4E-03 | 4.4E-02 | 14.75% |
| Aroclor-1260 | 0.66 | 3.6E-06 | 0.14 | 1.4E-06 | 2.0E-05 | 2.0E-05 | 1.8E-01 | 7.1E-02 | 2.5E-01 | 85.25% |
| | | | | SUMMARY | HAZARD INDE | x | 2.2E-01 | 7.4E-02 | 3.0E-01 | |

TABLE 5-8 INHALATION EXPOSURE TO PARTICULATES IN SURFACE SOIL - RME UNRESTRICTED LAND USE - CHILD RESIDENT (1 TO 6 YEARS) STUDY AREA CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

EXPOSURE PARAMETERS

EQUATIONS

| PARAMETER | SYMBOL | VALUE | UNITS | |
|--|-------------------------------|------------------------|----------------------|---|
| CONCENTRATION SOIL* | CS | See below | mg/kg | CANCER RISK = INTAKE (mg/kg-day) x CANCER SLOPE FACTOR (mg/kg-day)-1 |
| CONCENTRATION AIR PARTICULATES | САр | Calculated | mg/m² | |
| CONCENTRATION AIR VOLATILES | CAv | Calculated | mg/m² | HAZARD QUOTIENT - INTAKE (mg/kg-day) / REFERENCE CONCENTRATION (mg/kg-day) |
| VOLATILIZATION FACTOR** | VF | Calculated | m³/kg | |
| PARTICULATE EMISSIONS FACTOR | PEF | 1.32E+09 | ug/m³ | |
| INHALATION RATE | IhR | 0.35 | m [*] /hour | INTAKE - INHALATION = $(CAp + Cav) \times RAF \times IhR \times ET \times EF \times ED$ |
| BODY WEIGHT | BW | 15 | kg | BW x AT x 365 days/yr |
| EXPOSURE TIME | ET | 8 | hours/day | 1 |
| EXPOSURE FREQUENCY | EF | 150 | days/year | |
| EXPOSURE DURATION | ED | 6 | years | AIR CONCENTRATION PARTICULATES = CS x 1/PEF |
| RELATIVE ABSORPTION FACTOR | RAF | 100% | | 1 |
| AVERAGING TIME | | | | AIR CONCENTRATION VOLATILES - CS x 1/VF |
| CANCER | AT | 70 | years | (VF not calculated because there are no VOCs selected as CPCs). |
| NONCANCER | AT | 6 | years | |
| Notes: * Soil concentration used is the lesser of the 95 % upp | er confidence limit (UCL) & | maximum concentration | n | |
| **Volatilization factor used only for volatile chemicals of pote | ntial concern. | | | |
| For noncarcinogenic effects: AT = ED | All exposure variables are de | cumented in Table 5-4. | | |
| ND = Value not determined | | | | |

CARCINOGENIC EFFECTS

| | SOIL | | AIR CONCE | NTRATION | _ | CANCER SLOPE | | PERCENT |
|--------------|---------------|---------|-----------|--------------|-------------|---------------|---------|---------|
| COMPOUND | CONCENTRATION | VF | VOLATILES | PARTICULATES | INTAKE | FACTOR | CANCER | TOTAL |
| | (mg/kg) | (m³/kg) | (mg/m³) | (mg/m³) | (mg/kg-day) | (mg/kg-day)-1 | RISK | RISK |
| Arsenic | 2.2 | NA | | 1.7E-09 | 1.1E-11 | 1.5E+01 | 1.6E-10 | 96.15% |
| Aroclor-1260 | 0.66 | NA | | 5.0E-10 | 3.3E-12 | 2.0E+00 | 6.6E-12 | 3.85% |
| | 1 1 | | | | 1 1 | | | |
| | | | | <u> </u> | | | | |
| | | | | | SUMMARY | CANCER RISK | 1.7E-10 | |

NONCARCINOGENIC EFFECTS

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| | SOIL | | AIR CONCE | NTRATION | | REFERENCE | | PERCENT |
|--------------|---------------|----------------------|-----------|--------------|-------------|--------------|----------|---------|
| COMPOUND | CONCENTRATION | VF | VOLATILES | PARTICULATES | INTAKE | DOSE | HAZARD | TOTAL |
| | (mg/kg) | (m ¹ /kg) | (mg/m³) | (mg/m²) | (mg/kg-day) | (mg/kg-day) | QUOTIENT | RISK |
| Arsenic | 2.2 | ŇA | | 1.7E-09 | 1.3E-10 | ND | | |
| Aroclor-1260 | 0.66 | NA | | 5.0E-10 | 3.8E-11 | 2.0E-05 | 1.9E-06 | 100,00% |
| | | | | | | | | |
| | | | | | SUMMARY | HAZARD INDEX | 1.9E-06 | |

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TABLE 5-9 INGESTION OF AND DERMAL CONTACT WITH GROUND WATER - RME UNRESTRICTED LAND USE - ADULT RESIDENT STUDY AREA CS-8/FS-21DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION EQUATIONS

EXPOSURE PARAMETERS

| PARAMETER | SYMBOL | VALUE | UNITS | |
|--------------------------------------|----------------------------|---------------------|------------------------|---|
| CONCENTRATION WATER | CW | maximum | mg/liter | CANCER RISK = INTAKE (mg/kg-day) x CANCER SLOPE FACTOR (mg/kg-day)" |
| ABSORBED DOSE PER EVENT | DAcvent | calculated | mg/cm² per event | |
| INGESTION RATE | IR | 2 | liters/day | |
| FRACTION ABSORBED | FA | Chemical-specific | dimensionless [a] | |
| SURFACE AREA EXPOSED | SA | 18,000 | cm ² | HAZARD QUOTIENT = INTAKE (mg/kg-day) / REFERENCE DOSE (mg/kg-day) |
| CONVERSION FACTOR | CF | 0.001 | liters/cm ¹ | |
| BODY WEIGHT | BW | 70 | kg | 1 |
| EVENT TIME | ET | 0.58 | hours/event | |
| LAG TIME PER EVENT | LT | Chemical-specific | hours/event [a] | |
| EVENT FREQUENCY | EV | t t | event/day | |
| EXPOSURE FREQUENCY | EF | 350 | days/year | INTAKE = (INTAKE-INGESTION) + (INTAKE-DERMAL) |
| EXPOSURE DURATION | ED | 30 | years | |
| AVERAGING TIME | | | | INTAKE-INGESTION - <u>CWxIRxEFxED</u> |
| CANCER | AT | 70 | years | BW x AT x 365 days/yr |
| NONCANCER | AT | 30 | years | |
| PERMEABILITY COEFFICIENT | Kpevent | Chemical-specific | cm/hour [b] | INTAKE-DERMAL - DAEvent ISA I EV I EF I ED |
| | | | L <u> </u> | BW x AT x 365 days/yr |
| Notes: | | | | 1 |
| For noncarcinogenic effects: AT = ED | All exposure variables are | documented in Table | · 5-4. | DAevent (organics) = 2 * FA * CW * Kpevent * SQRT(6 * LT x ET/Pi) * CF; Pi ~ 3.1416 |
| | | | | applicable where ET <= time to reach steady state (hr) = 2.4 * LT |
| 1 | | | | DAevent (organics) = FA * CW * Kpevent * [ET/(1+B) + (2 * LT) * ((1 + 3 * B + 3 * B ²)/(1 + B) ³)] * CF |
| | | | | applicable where ET > time to reach steady state (hr) = 2.4 * LT |
| | | | | |
| | | | | DAevent (inorganics) = CW * Kpevent * ET * CF |
| | | | | Source: RAGS Part E Supplemental Guidance for Dermal Risk Assessment; Interim Guidance (EPA, 2000). |
| | | - | | 1 |

CARCINOGENIC EFFECTS

| COMPOUND | WATER CONCENTRATION (mg/L) | INTAKE INGESTION (mg/kg-day) | PERMEABILITY COEFFICIENT (cm/kour) | DAgyport mg/cm ^{2-mga} | INTAKE DERMAL (mg/kg-day) | CANCER SLO ORAL (mg/kg-day) ⁻¹ | DPE FACTOR DERMAL (mg/kg-day) ^{**} | CANCER RISK INGESTION | CANCER RISK DERMAL | TOTAL CANCER RISK | PERCENT TOTAL RISK |
|----------|----------------------------------|------------------------------------|--|------------------------------------|---------------------------------|---|---|-----------------------------|--------------------------|-------------------------|--------------------------|
| Arsenic | 0.0074 | 8.7E-05 | 1.0E-03 | NE | NE | 1.5E+00 | 1.5E+00 | 1.3E-04 | NE | 1.3E-04 | 100.00% |
| | | | ····· | | | SUMMARY CAL | NCER RISK | 1.3E-04 | 0.0E+00 | 1.3E-04 | |

NONCARCINOGENIC EFFECTS

| COMPOUND | WATER CONCENTRATION (mg/L) | INTAKE INGESTION (mg/kg-day) | PERMEABILITY COEFFICIENT (cm/hour) | DA _{EVENT} mg/cm ⁴ | INTAKE DERMAL (mg/kg-day) | REFEREN ORAL (mg/kg-day) | CE DOSE DERMAL (mg/kg-day) | HAZARD QUOTIENT INGESTION | HAZARD QUOTIENT DERMAL | TOTAL Hazard Quotient | PERCENT TOTAL RISK |
|-------------------------------------|----------------------------------|------------------------------------|--|---|---------------------------------|--------------------------------|----------------------------------|---------------------------------|------------------------------|-----------------------------|---------------------------|
| Arsenic Naphthalene Manganese | 0.0074 0.002 0.708 | | 1.0E-03 4.9E-02 1.0E-03 | NE 1.5E-07 NE | NE 3.8E-05 NE | 3.0E-04 2.0E-02 2.4E-02 | 3.0E-04 2.0E-02 9.6E-04 | 6.8E-01 2.7E-03 8.1E-01 | NI 1.9E-03 NI | 4.6E-03 | 45.40% 0.31% 54.29% |
| | <u></u> | | | | ł | SUMMARY HAZ | ARD INDEX | 1.5E+00 | 1.9E-0 | 1.5E+00 | |

NE = not evaluated for dermal exposure. Per Appendix B, Exhibit B.4 of RAGS Part E Supplemental Guidance for Dermal Risk Assessment (Interim Guidance), arsenic and manganese are not included as chemicals to be assessed for dermal exposure. EPA considers dermal exposure for these compounds to be relatively insignificant compared to ingestion exposures (i.e., dermal dose < 10% ingestion dose).

[a] For naphthalene, FA = 1 and LT = 0.55 hours/event (EPA, 2000 [Appendix B])

[b] Kp values obtained from EPA, 2000 (Exhibit 3.1 for inorganics and Appendix B for organics)

TABLE 5-10 RISK SUMMARY TABLE - POTENTIAL FUTURE LAND USE

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

| Exposure Scenario | Population | Medium | Exposure Route | HI | ELCR | - |
|------------------------|---------------------------|--------------------|--|--|---|---|
| Resident - 0-10 ft bgs | Child Resident | Soil 0-10 feet bgs | Ingestion Dermal Dust Inhalation | 0.2 0.1 <u>0.000002</u> | 2.2E-06 3.7E-07 <u>1.7E-10</u> | - |
| | | | | 0.3 | 2.6E-06 | |
| | Adult Resident | Soil 0-10 feet bgs | Ingestion Dermal Dust Inhalation Soil 0-10 ft bgs | 0.02 0.01 <u>0.0000007</u> 0.03 | 9.3E-07 2.3E-07 <u>2.6E-10</u> 1.2E-06 | |
| | | Groundwater* | Ingestion Dermal contact Groundwater | 1.5 <u>0.002</u> 1.5 | 1.3E-04 <u>NE</u> 1.3E-04 | - |
| | Child & Adult Resident | Soil 0-10 feet bgs | Total Exposure | 0.33 | 3.8E-06 | |

Notes:

HI - Hazard Index

ELCR - Excess Lifetime Cancer Risk

NE - not evaluated

* Groundwater hazard index is a screening hazard index. Segregated hazard quotients for specific endpoints are all below one.

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In this case, the HQs should be segregated. For arsenic, the critical effect is hyperpigmentation, keratosis, and possible vascular complications (IRIS, 2000). For manganese, the critical effect is central nervous system (neurological effects) as reported in IRIS. Because the HQs for arsenic and manganese are each below 1, the risk assessment results indicate that neither the skin-related effects associated with arsenic or the neurological effects associated with manganese are likely. Therefore, the screening HI overestimates the non-cancer risks. The segregated HQs for CPCs in groundwater are appropriate to consider here and they are below 1.

The cumulative cancer risk for groundwater used as potable water is 1.3×10^{-4} for the adult resident, which is just above the USEPA cancer risk range of 1×10^{-6} to 1×10^{-4} , and above the MADEP cumulative cancer risk limit of 1×10^{-5} . However, the risk associated with groundwater is attributable to arsenic. Arsenic was detected in one groundwater well at a concentration (7.4 µg/L) below the Federal and Massachusetts MCL of 50 µg/L, and just above the proposed USEPA MCL of 5 µg/L.

Based on this evaluation, the risks associated with residential use of the surface soil, subsurface soil, and groundwater at Study Area CS-8/FS-21 meet risk management criteria established by USEPA. Cancer risks associated with groundwater exceed the USEPA and MADEP cumulative cancer risk limit due to arsenic; however, the detection of arsenic in groundwater is below the current Federal and Massachusetts drinking water standards.

5.2.4 Uncertainty Assessment

Study-area-specific uncertainties associated with the PRE were identified as follows:

- 1. Arsenic was not detected at concentrations in surface soil greater than maximum background levels. It was retained as a CPC only because Study Area CS-8/FS-21 maximum surface-soil concentrations exceeded the Region IX residential soil PRG. The residential cancer risk associated with exposure to the maximum background concentration of arsenic $(3x10^{-6})$ is greater than the risk associated with exposure to the maximum concentration detected at the site $(2x10^{-6})$.
- 2. The conservative nature of the PRE methodology may overestimate potential health risks. For example, arsenic in groundwater was detected

above laboratory detection in just one groundwater sample (04BH001). Groundwater samples collected upgradient (CS-8/MW-2) and downgradient (FS-21/MW-1) of this sample were both below detection limits, yet the arsenic detection in groundwater from borehole 04BH001 (7.4 μ g/L) was used to represent the arsenic concentration in groundwater throughout the site in terms of risk.

3. The PRE was conducted using an SI data set that is less comprehensive than an RI data set. Therefore, the human-health PRE may potentially underestimate risk. An SI data set, however, is designed to evaluate the presence of contaminants (worst case scenario). Therefore, it is unlikely that contaminants are present in concentrations that would pose an unacceptable risk.

5.3 ECOLOGICAL PRELIMINARY RISK EVALUATION

The ecological PRE was completed to provide a screening-level evaluation of potential risks that CPCs in surface soil may pose to ecological receptors at Study Areas CS-8/FS-21. PRE methodology is discussed in detail in the RAH (ASG, 1994). The following subsections discuss the PRE and its results.

5.3.1 Data Evaluation

Table 5-11 summarizes the frequencies of detection, the ranges of CRQLs and CRDLs, and the range of detected concentrations of CPCs in surface soil. Ecological CPCs were selected in accordance with steps described in the RAH. Inorganic analytes detected in surface soil at concentrations below MMR background and also below the lowest Tier I and Tier II HECs presented in the RAH were not considered CPCs. In addition, the essential nutrients calcium, iron, magnesium, potassium, and sodium were not retained as CPCs because they have low toxicity and are present at concentrations below MMR background concentrations. As shown in Table 5-11, only zinc was selected as a CPC. The maximum detected concentration of zinc was used in the ecological PRE.

5.3.2 Ecological Exposure Assessment

Study Areas CS-8/FS-21 cover approximately 3 acres and include three buildings for vehicle and equipment maintenance, two concrete wash pads, and two fuel pump islands.

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Most of the northern section of the study areas is paved. The unpaved section consists of mowed grass toward the southern end and loose gravel to the west. A chain-link fence borders the site to the north and south. The grassy area contains species typical of a ruderal habitat, such as grasses (Graminae) and herbaceous perennials (e.g., goldenrod [Solidago sp.] and bluestem grass [Andropogon scoparius]). The majority of the site is paved or covered with crushed rock, and there is a lack of suitable vegetative cover or forage to support most species. Moreover, the frequent human activity at and in the vicinity of the site discourages the presence of wildlife, and the chain-link fence deters larger ecological receptors from entering the site. A pitch-pine/oak woodland, containing characteristic species described in the RAH, is located southwest of the site, beyond the perimeter fence. This wooded area may provide food and cover for small mammals and birds.

The following receptor species, representing a range of taxonomic groups and trophic levels, were used to evaluate risk from surface-soil contamination in the PRE: white-footed mouse (omnivorous small mammal), cardinal (insectivorous small bird), and red fox (omnivorous predatory mammal). These three species were chosen to represent other species within the same trophic level at Study Areas CS-8/FS-21 (ASG, 1994). These three receptors could potentially occur in the nearby woodland and could occasionally use the limited habitat at the study areas. Species-specific ecological exposure parameters used to estimate potential dietary exposure to surface-soil contaminants are in Appendix P of the RAH.

5.3.3 Ecological Effects Assessments

Development of HECs for animal and plant receptors is described in the RAH. HECs for receptor species and Soil Critical Concentration values for phytotoxicity are in Appendices F and O of the RAH, respectively.

5.3.4 Risk Screening

PREs characterize potential ecological risks by comparing maximum site concentrations to pre-calculated, medium-specific HECs or phytotoxicity data. For the ecological PRE, Tier I HECs (based on current exposures) and Tier II HECs (based on future exposures) are the same because exposures of ecological receptors are likely to be similar under current and future land-use conditions (ASG, 1994). Risks to ecological receptors were evaluated by calculating HQs and HIs described in the RAH. The ecological HQ is a receptor- and chemical-specific value equivalent to the ratio of the maximum detected site concentration

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TABLE 5-11 ECOLOGICAL CONTAMINANTS OF POTENTIAL CONCERN IN SURFACE SOIL

| ····· | T | г | MAXIMUM | HAZARD FOULV | ALENT CONCENT | RATIONS (mg/kg) | SOIL | RANGEOF | | |
|------------------------------|-----------|---------|------------|--------------|---------------|-----------------|---------------|----------------|----------|-------|
| | FREQUENCY | CROL | BACKGROUND | | | | CRITICAL | DETECTED | SELECTED | |
| | | | | WHITE- | | | CONCENTRATION | | | |
| | OF | CRDL | CONC. | FOOTED | RED FOX | CARDINAL | FOR PLANTS | CONCENTRATIONS | AS A | |
| CHEMICAL | DETECTION | (mg/kg) | (mg/kg) | | | · | (mg/kg) | (mg/kg) | CPC? | NOTES |
| SEMIVOLATILE ORGANIC COMPOUN | //// | ! | | | | | | | | |
| Bis(2-ethylhexyl)phthalate | 1/2 | 0.33 | NA | 94.7 | 3,220 | 11.2 | 100 | 0.064 - 0.064 | NO | 2 |
| Butylbenzylphthalate | 1/2 | 0.33 | NA | 11,100 | 223,000 | 1,340 | 11.2 | 0.14 - 0.14 | NO | 2 |
| PCBs/PESTICIDES | | 1 | 1 | | | | | | | |
| Aroclor-1260 | 1/2 | 0.16 | NA | 17.8 | 29.9 | 17.6 | 10 | 0.66 - 0.66 | NO | 2 |
| INORGANICS | | | | | | | | | | |
| Aluminum | 2/2 | 40 | 8,930 | 24,200 | 317,000 | 54,500 | NA | 1,300 - 3,220 | NO | 2,4 |
| Arsenic | 2/2 | 2 | 3.6 | 25.3 | 762 | 33.5 | 5 | 0.69 - 2.2 | NO | 4 |
| Barium | 2/2 | 40 | 10.4 | 3,900 | 73,000 | 1,050 | NA | 9.6 - 11 | NO | 2 |
| Beryllium | 1/2 | 1 | 0.65 | 34,100 | 152,000 | 35,500 | NA | 0.37 - 0.37 | NO | 2, 4 |
| Calcium | 2/2 | 1,000 | 969 | NA | NA | NA | NA | 218 - 331 | NO | 3, 4 |
| Chromium | 2/2 | 2 | 6.8 | 7.2 | 185 | 707 | 5 | 2.1 - 5.3 | NO | 4 |
| Cobalt | 1/2 | 10 | 4.1 | 461 | 2,640 | 287 | 15 | 1.5 - 1.5 | NO | 2, 4 |
| Copper | 1/2 | 5 | 5.2 | 50.8 | 4,610 | 12 | 20 | 9.2 - 9.2 | NO | 2 |
| Iron | 2/2 | 20 | 12,400 | NA | NA | NA | NA | 2,490 - 4,010 | NO | 3,4 |
| Lead | 2/2 | 3 | 12.05 | 169 | 66,100 | 222 | 30 | 5.6 - 6.9 | NO | 2,4 |
| Magnesium | 2/2 | 1,000 | 794.5 | NA | NA | 222 | NA | 178 - 536 | NO | 3,4 |
| Manganese | 2/2 | 3 | 108 | 239 | 13,200 | 59.6 | 300 | 31 - 59 | NO | 2, 4 |
| Thallium | 1/2 | 2 | 0.25 | 1.28 | 5.29 | 2.87 | 5 | 0.4 - 0.4 | NO | 2 |
| Vanadium | 2/2 | 10 | 15.2 | 25.9 | 111 | 43.8 | NA | 3.7 - 6.4 | NO | 2,4 |
| Zinc | 2/2 | 4 | 16 | 32.7 | 7,720 | 39.5 | 100 | 17 - 30 | NO | 2 |

STUDY AREAS CS-8/FS-21 DECISION DOCUMENT MASSACHUSETTS MILITARY RESERVATION

NOTES:

NA = not applicable

CPC = contaminant of potential concern

CRQL = Contract Required Quantitation Limit

CRDL = Contract Required Detection Limit

mg/kg = milligrams per kilogram

Hazard Equivalent Concentration from Appendix F-2 and Soil Phytotoxicity Value from Appendices O-3 and O-4 of the MMR Risk Assessment Handbook (ASG, 1994).

1 Chemical selected as a CPC because maximum detected concentration exceeds HEC and background.

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2 Chemical not retained as a CPC because the maximum detected concentration was less than the HEC.

3 Chemical not retained as a CPC because it is an essential nutrient.

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4 Chemical not retained as a CPC because the maximum detected concentration was less than the maximum background concentration.

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of a particular CPC to the receptor-specific HEC for that particular CPC. The ecological HI is a receptor-specific value equivalent to the sum of all chemical-specific HQs for a given receptor. Separate HIs are calculated for organic and inorganic CPCs. If an HI for a receptor for organics exceeds 1, or for inorganics exceeds 10, the RAH recommends an RI (ASG, 1994). Otherwise, a no-further-action decision can be made.

Surface Soil. A screening-level evaluation of potential risk from surface soil was conducted by using the maximum concentration of the surface-soil analytes detected with species-specific HECs. No analyte concentrations exceeded the HECs.

5.3.5 Ecological Uncertainty Assessment

Several uncertainties are associated with the ecological PRE at Study Areas CS-8/FS-21. Additional general uncertainties inherent in the ecological PRE process are outlined in the RAH. Study-area-specific uncertainties were identified as follows:

- 1. There is uncertainty associated with the HECs. Several HECs are lower than MMR maximum background concentrations; therefore, some of the risks at the study areas may be attributed to background levels (i.e., natural content of inorganic analytes in soil).
- 2. Concentrations of inorganics in surface soil were compared to Soil Critical Concentration values for plants. Soil Critical Concentrations were not available for barium. Insufficient data for this analyte adds an uncertainty to the phytotoxicity evaluation for Study Areas CS-8/FS-21.
- 3. As described in Subsection 5.3.2, Study Areas CS-8/FS-21 contain limited ecological habitat in a semi-urbanized environment. It is unlikely that significant use of these study areas by all ecological receptors evaluated in this PRE will occur.

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5.4 SUMMARY OF PRELIMINARY RISK EVALUATION RESULTS

Results of the human-health PRE indicated that arsenic and aroclor-1260 were selected as CPCs in surface soil, arsenic was selected as a CPC in subsurface soil, and naphthalene, arsenic, and manganese were selected as CPCs in groundwater. Health risk calculations for potential resident receptor exposures to the CPCs indicate that risks associated with residential use of the surface soil, subsurface soil, and groundwater at Study Area CS-8/FS-21 meet risk management criteria established by USEPA. Cancer risks associated with groundwater exceed the MADEP cumulative cancer risk limit due to arsenic; however, the detection of arsenic in groundwater is below the current Federal and Massachusetts drinking water standards.

Tiers I and II of the ecological PRE showed that no maximum surface-soil concentrations exceeded the lowest HECs. HIs for all three evaluated receptors did not exceed 1 for organics or 10 for inorganics. Both organic and inorganic HIs for plants were less than 1.

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6.0 DRAINAGE STRUCTURE REMOVAL ACTIVITIES

One drainage structure, a cesspool (i.e., 04CDXX1), was removed in April 1996 at Study Areas CS-8/FS-21 as part of the MMR Drainage Structure Removal Program (DSRP) (Jacobs, 1996). The cesspool was located outside the flightline to the west of Building OMS-22, which is located in the Army National Guard Motor Pool area off of South Truck Road. The cesspool walls were constructed of concrete barrel block, with an open dirt bottom and a 4-inch-diameter ceramic inlet pipe. No outlet pipe was present. At the time of removal, the structure had no cover and the walls had collapsed.

No sludge or liquid was removed from the cesspool. The structure and approximately 40 cubic yards of soil and debris were excavated and removed. PI detector screening, sampling and analysis of soils from the open excavation, and backfilling of the excavation were completed following DSRP protocols (Jacobs, 1996). None of the soil analytical results exceeded soil target clean-up levels established for the DSRP; therefore, the cesspool was removed with a clean closure. The Closure Report for the cesspool is included as Appendix D.

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7.0 SUMMARY AND CONCLUSIONS

SI efforts at Study Area CS-8/FS-21, which included a soil-gas survey, 10 test pits, six soil borings, four monitoring wells, field-screening of soil samples, and laboratory analysis of soil and groundwater samples, showed no significant contamination of soil or groundwater. Chemical source areas were not found through the soil-gas, test-pit, or soil-boring investigations, and Cesspool 04CDXX1 was removed with a clean closure. Additionally, the former 12,500-gallon diesel fuel UST and the former 5,000-gallon MOGAS UST (i.e., Current Product Tank No. 90) were removed with no evidence of soil contamination during the excavations.

Results of the human-health PRE indicated that arsenic and aroclor-1260 were selected as CPCs in surface soil, arsenic was selected as a CPC in subsurface soil, and naphthalene, arsenic, and manganese were selected as CPCs in groundwater. Health risk calculations for potential resident receptor exposures to the CPCs indicate that risks associated with residential use of the surface soil, subsurface soil, and groundwater at Study Area CS-8/FS-21 meet risk management criteria established by USEPA. Cancer risks associated with groundwater exceed the USEPA and MADEP cumulative cancer risk limit due to arsenic; however, the detection of arsenic in groundwater is below the current Federal and Massachusetts drinking water standards.

Tier I and Tier II of the ecological PRE showed that no maximum surface-soil concentrations exceeded the lowest HECs. HIs for all three evaluated receptors did not exceed 1 for organics or 10 for inorganics. Both organic and inorganic HIs for plants were less than 1.

Therefore, since human-health and ecological risk assessments suggest that unacceptable levels of risk are not anticipated and the potential contaminant source areas have been removed, AFCEE concludes that Study Areas CS-8/FS-21 do not require additional characterization or remediation efforts, and that no further action within the IRP is appropriate.

8.0 DECISION

On the basis of these findings, there is no evidence of significant environmental contamination or of human health or ecological risks at Study Areas CS-8/FS-21. The decision has been made to remove these study areas from further consideration in the IRP process.

GARY M. ERI Director Air Force Center for Environmental Excellence

19 Oct 00

Date

CS-8/FS 21 DECISION DOCUMENT

U.S. Environmental Protection Agency New England Division

M Concur

RICHARD CAVAGNER® Deputy Director Office of Site Remediation and Restoration

W 1,2000 Date

Date

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[] Non-concur (please provide reasons

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CS-8/FS-21 DECISION DOCUMENT

Commonwealth of Massachusetts Department of Environmental Protection

FOR

Concur

[1]

PAUL A. TAURASI, P.E. Regional Director Southeast Region

11-22-00

Date

[] Non-concur (please provide reasons): _____

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ARGEO PAUL CELLUCCI Governor

JANE SWIFT Lieutenant Governor COMMONWEALTH OF MASSACHUSETTS EXECUTIVE OFFICE OF ENVIRONMENTAL AFFAIRS **DEPARTMENT OF ENVIRONMENTAL PROTECTION** 20 RIVERSIDE DRIVE, LAKEVILLE, MA 02347 508-946-2700

> BOB DURAND Secretary

LAUREN A. LISS Commissioner

November 22, 2000

Mr. Robert M. Gill Remediation Program Manager HQ AFCEE/MMR 322 East Inner Road Otis ANG Base, Massachusetts 02542 RE: BOURNE--BWSC-4-0037 Massachusetts Military Reservation (MMR), Study Area CS-8/FS-21, Final Decision Document, Concurrence

Dear Mr. Gill:

The Department of Environmental Protection (the "Department") has reviewed a document titled "DECISION DOCUMENT STUDY AREA CS-8/FS-21, FINAL" dated October 2000 and prepared by Harding Lawson Associates of Portland, Maine for the Air Force Center for Environmental Excellence (AFCEE).

Study Area CS-8/FS-21 is located on the south side of the intersection of South Outer Road and East Truck Road at the MMR. The Study Area has been operating as the Army National Guard Vehicle Repair Shop since 1950. All underground fuel storage tanks and pump islands have been removed and no contaminants were detected during the removals. A cesspool, the only underground drainage structure, was removed in 1996. Soil and groundwater were tested for all contaminants of concern. Based on current site conditions, the site poses no significant risk of harm to human health or the environment. The DD proposes no further action at Study Area CS-8/FS-21.

The Department concurs with the DD. The Department's concurrence with this DD is based upon representations made to the Department by the AFCEE and assumes that all information provided is substantially complete and accurate. Without limitation, if the Department determines that any material omissions or misstatements exist, if new information becomes available, or if conditions at the Study Area change, resulting in potential or actual human exposure or threats to the environment, the Department reserves its authority under M.G.L. c. 21E, and the MCP, 310 CMR 40.0000 <u>et seq</u>., and any other applicable law or regulation to require further response actions.

> This information is available in alternate format by calling our ADA Coordinator at (617) 574-6872. DEP on the World Wide Web: http://www.magnet_state.ma.us/dep Printed on Recycled Paper

Please incorporate this letter into the Administrative Record for the Study Area CS-8/FS-21. If you have any questions regarding this letter, please contact Leonard J. Pinaud at (508) 946-2871.

Sincerely, FCR A. Taurasi, P.E.,

Regional Director

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DEP-SERO Cc: ATTN: Mildred Garcia-Surette, Deputy Regional Director Leonard J. Pinaud, Chief, Federal Facilities Remediation Section

Distributions: SERO SMB Plume Containment Team Public Information Team Long Range Water Supply PAT Boards of Selectmen Boards of Health

GLOSSARY OF ACRONYMS AND ABBREVIATIONS

| ABB-ES | ABB Environmental Services, Inc. |
|--------|--------------------------------------|
| ASG | Automated Sciences Group |
| BAP | benzo(a)pyrene |
| BEHP | bis(2-ethylhexyl)phthalate |
| bgs | below ground surface |
| CS | Chemical Spill |
| CLP | Contract Laboratory Program |
| CPC | contaminant of potential concern |
| CRDL | Contract Required Detection Limit |
| CRQL | Contract Required Quantitation Limit |
| D-BHC | delta-benzene hexachloride |
| DCE | dichloroethene |
| DDT | dichlorodiphenyltrichloroethane |
| DSRP | Drainage Structure Removal Program |
| EDB | ethylene dibromide |
| EPC | Exposure Point Concentration |
| FS | fuel spill |
| gal/yr | gallons per year |
| GC | gas chromatograph |
| HARM | Hazard Assessment Rating Methodology |
| HEC | hazard equivalent concentration |
| HI | hazard index |
| HQ | hazard quotient |
| IDL | instrument detection limit |
| IRP | Installation Restoration Program |
| MCL | Maximum Contaminant Level |
| mg/kg | milligrams per kilogram |
| MMR | Massachusetts Military Reservation |
| MOGAS | motor vehicle gasoline |
| MSL | mean sea level |

GLOSSARY OF ACRONYMS AND ABBREVIATIONS

| MTBE | methyl-tert-butyl ether |
|-------|---|
| PA | preliminary assessment |
| PAH | polycyclic aromatic hydrocarbon |
| PCB | polychlorinated biphenyl |
| PCE | tetrachloroethene |
| PI | photoionization |
| ppm | parts per million |
| PRG | preliminary remediation goal |
| PRE | preliminary risk evaluation |
| QC | quality control |
| RAH | Risk Assessment Handbook |
| RI | remedial investigation |
| SI | site investigation |
| SMCL | Secondary Maximum Contaminant Level |
| SQL | sample quantitation limit |
| SVOC | semivolatile organic compound |
| TAL | Target Analyte List |
| TCE | trichloroethene |
| TCL | Target Compound List |
| TPH | total petroleum hydrocarbons |
| TSDF | Treatment, Storage, and Disposal Facility |
| UCL | upper confidence limit |
| USEPA | U.S. Environmental Protection Agency |
| UST | underground storage tank |
| UTES | Unit Training Equipment Storage |
| VOC | volatile organic compound |
| µg/kg | micrograms per kilogram |
| μg/L | micrograms per liter |

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APPENDIX A

APPENDIX A SUMMARY OF CLP ANALYTICAL DATA

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DATA QUALITY ASSESSMENT

The CLP analytical data from Study Areas CS-8/FS-21 were reviewed and evaluated for suitability of use in a risk assessment or as support for engineering studies. Twenty soil samples, all collected in September 1987 as part of the Task 2-3A SI, were analyzed for TCL VOCs, TCL SVOCs, TCL pesticides and PCBs, TAL inorganic analytes, and cyanide. Five of the 20 samples were taken from depths of 10 feet bgs or less. In January 1988, as part of the Task 2-3B SI, samples were collected from four monitoring wells (i.e., CS-8/MW-2, CS-8/MW-4, FS-21/MW-1, and FS-21/MW-2) and analyzed for TCL VOCs and SVOCs, TAL inorganics, and nitrates/nitrites. Duplicate samples were collected from CS-8/MW-2 and CS-8/MW-4. In June 1989, as part of the Task 2-3C SI, one groundwater sample and a duplicate were collected from CS-8/MW-2 and analyzed for TCL VOCs and SVOCs, TAL inorganic analytes, and cyanide.

All Tasks 2-3A, 2-3B, and 2-3C analytical data were CLP data. In preparation for this report, the data were reviewed for holding times, correct methodology, and presence of chemicals in laboratory and field blanks. Corrections for concentrations of chemicals found in any of the associated blanks were applied following USEPA validation guidelines and other guidance (USEPA, 1988, 1989a, and 1989b). For those samples for which associated method blanks or field blanks were not clearly labeled, corrections were applied as determined by the time frame of the sampling and analysis (USEPA, 1989b). The data were not fully validated following USEPA guidance.

The evaluated data were reported with one of four qualifications that affect the suitability for use in a risk assessment:

| an unqualified value | use in risk assessment |
|--|--------------------------------------|
| • a value with a U; not detect | ed at the SQL use in risk assessment |
| • a value with a J; estimated | value use in risk assessment |
| • R; rejected data | do not use in risk assessment |
| • a value with a J; estimated | value use in risk assessmen |

Based on the evaluation, the data are considered adequate for use in a risk assessment. None of the data were rejected. Much of the data, especially the inorganic analyses, was outside the laboratory method criteria and data validation guidelines, but not so far as to cause the data to be rejected. The identities of these chemicals are not considered uncertain, but the concentrations are considered estimated values and are therefore qualified with a J. Some of the concentrations of the chemicals are estimated (J) because they are below the sample quantitation limit (SQL); that is, the sample-specific instrument detection limit (IDL) for inorganic compounds or the CRQL for organic compounds. As before, the identities of

APPENDIX A

these chemicals are not considered uncertain but the reported concentrations are. Estimated data (J) are suitable for use in a risk assessment (USEPA, 1989b).

The inorganic data tables available for review from 1987 and 1988 did not report values for detection limits; dashes were used instead. In reviewing the data, the undetected analytes were reported using the SQL based on the CRDL rather than the IDL, qualified with a U. The current validation procedure for reporting undetected inorganics uses the IDL for a value, qualified with a U. Because the CRDL is always higher than the IDL, non-detect values in the older data sets are higher than would have been reported if the IDLs had been known.

As part of the data evaluation, the identities and concentrations of chemicals reported in laboratory method blanks and field blanks (i.e., trip blanks, sampler blanks, equipment rinseates, and source waters) and the effect on the sample results were evaluated. Following USEPA guidance (CLP functional guidelines and USEPA, 1989b), action levels were set for those chemicals detected in the blanks. Only those chemicals whose concentrations exceeded the action levels were considered site-related and not a laboratory or sampling artifact. If the chemical was a common laboratory contaminant as defined by USEPA (i.e., methylene chloride, acetone, 2-butanone, toluene, and phthalates), 10 times the concentration was used for the action level. If the chemical was not a common laboratory contaminant, five times the blank concentration was used for the action level. The individual samples and the laboratory method blanks associated with each are reviewed during evaluation. As part of the data evaluation, blank concentrations over similar time periods were grouped together and concentrations for any samples within that same time period were adjusted following USEPA guidelines (USEPA, 1989b).

VOC ANALYSES

Methylene chloride, acetone, and 2-butanone were the only TCL VOCs detected in either the soil or water samples at Study Areas CS-8/FS-21. All methylene chloride and acetone in the soil samples collected in January 1988 and September/October 1987 are below the action levels of 130 and 500 μ g/kg, respectively, based on the maximum concentrations observed in the method blanks (13 μ g/kg and 50 μ g/L, respectively). No 2-butanone was detected in the soil samples. A single 2-butanone detection at 19 μ g/L in a groundwater sample was below the action level of 85 μ g/L (based on a blank detection of 8.5 μ g/L); it therefore was considered a sampling or analytical artifact and was not reported in the data tables.

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Similarly, methylene chloride and acetone were detected but at concentrations below the action levels of 30 and 110 μ g/L, respectively, in groundwater samples collected during the June 1989 sampling event. As a result of this evaluation, neither methylene chloride nor acetone is considered a site-related contaminant at Study Areas CS-8/FS-21. They are considered laboratory and sampling artifacts and therefore are not included in the tables.

Detections of 2-butanone during the Task 2-3C SI were above the action levels and do appear in the data tables. Although reported in the data tables, the presence of 2-butanone is considered a sampling artifact resulting from incomplete rinsing of the interior of sampling equipment with distilled water after decontamination with methyl hydrate, which contains approximately 3 percent 2-butanone. The 2-butanone concentrations reported in groundwater samples are not considered representative of those in the aquifer, as was demonstrated previously (E. C. Jordan Co., 1990b and 1992). The 2-butanone is not considered a site-related contaminant.

SVOC ANALYSES

SVOCs detected in samples from Study Areas CS-8/FS-21 are PAHs, bis(2-ethylhexyl)phthalate (BEHP), and butylbenzylphthalate. BEHP and butylbenzylphthalate were detected in one soil sample below the SQL in the Fall 1987 sampling event. Phthalates were not detected in either laboratory or field quality control (QC) samples, according to the Task 2-3A SI report. Consequently, there is no evidence that the phthalates may be considered laboratory or sampling artifacts for this sample.

PAHs were detected in two groundwater samples. Values for the PAH concentrations were estimated and qualified with a J because they were below the SQLs. Naphthalene was detected at 2J μ g/L in duplicate samples collected from CS-8/MW-2 in June 1989. However, naphthalene was not detected in duplicate samples collected from the same well in October 1987.

Chrysene, fluoranthene, pyrene, benzo(b)fluoranthene, and benzo(k)fluoranthene were detected at concentrations below the SQL in one groundwater sample, JFS21MW202, collected in January 1988. The laboratory was unable to distinguish between benzo(b)fluoranthene and benzo(k)fluoranthene; the value reported represents the sum of the two chemicals. This sample was noted in the sampling log and reported in the SI report to be light brown and contain suspended solids. Monitoring well FS-21/MW-2 is located directly below the MOGAS tank. SVOCs were not detected in the corresponding soil samples from FS-21/TB-2 at 59 to 61 feet. Field-screening for VOCs indicated the possible

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presence of benzene, TCE, and others; however, no TCL VOCs, other than known laboratory artifacts, were detected. It is possible that the PAHs detected are laboratory or sampling artifacts and not representative of the groundwater; however, they have been retained for consideration as CPCs.

PESTICIDES AND PCBS

4,4'-DDT and Aroclor-1260 were the only TCL pesticide or PCB detected in any of the soil samples. The pesticide 4,4'-DDT was detected in three soil samples, at 45 μ g/kg in FS21TB0259 (at 59 to 61 feet bgs), at 18 μ g/kg in CS8XTB0219 (at 19 to 21 feet bgs), and at 24 μ g/kg in CS8XTB0254 (at 54 to 56 feet bgs). Aroclor-1260 was detected in one sample at 660 μ g/kg in CS8TP0101 (at 1 to 2 feet). Neither 4, 4'-DDT nor Aroclor-1260 was detected in any associated laboratory or field blank; therefore, the 4,4'-DDT and PCB are considered CPCs. None of the groundwater samples were analyzed for TCL pesticides and PCBs.

INORGANIC ANALYSES

Much of the inorganic soil data is qualified as estimated (J) as a result of analyses that are out of criteria for the methods or for results that are below the SQL. The reasons for the estimated qualifier vary and were summarized in the SI and RI reports. None of the inorganic soil data were rejected. Laboratory reports of analyses in the SI and RI reports indicate the type of QC issues. The most common QC issues in the inorganic analyses for Study Areas CS-8/FS-21 data are as follows:

- spiked sample recovery out of control limits (laboratory qualifier N)
- value between the SQL and the IDL (laboratory qualifier [])
- matrix interferences, especially for the zinc data (laboratory qualifier E)
- duplicate analyses do not agree within control limits (laboratory qualifier *)

These laboratory qualifiers were changed to J's during the data evaluation. These qualified data are adequate for use in risk assessment.

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Zinc was detected in the method blanks associated with groundwater collected in June 1989; none of the zinc in the groundwater samples was above the action limit (i.e., five times the blank concentration). There were no other QC issues reported in the SI reports for the soil or water samples collected in the fall of 1987 and January 1988 (E.C. Jordan Co., 1989a and 1990a).

| FS21 | TBR-/ | APPE | NDIX |
|------|-------|------|------|
|------|-------|------|------|

| | R | evi | sed | 5/ | 6/ | 93 |
|--|---|-----|-----|----|----|----|
|--|---|-----|-----|----|----|----|

| | SITE DEPTH ID NUMBER DATE MATRIX | FS21-TB1 13-15 FS21TB0113 09/03/87 S01L | FS21-TB1 18-20 FS21TB0118 09/03/87 S01L | FS21-TB1 28-30 FS21TB0128 09/03/87 S0IL | FS21-TB1 33-35 FS21TB0133 09/03/87 SOIL | FS21-TB1 63-65 FS21TB0163 09/03/87 S01L | FS21-TB1 63-65 FS21TB1D63 09/03/87 S01L | FS21-TB2 0-2 FS21TB02X0 09/10/87 S0IL |
|---|--|--|---|---|---|---|---|--|
| ANALYTE | | | | | | | | |
| INORGANIC ANALYTES (mg/kg) | - CRDL | | | | | | | |
| ARSENIC BERYLLIUM CHROMIUM COPPER LEAD SELENIUM ZINC BARIUM IRON MANGANESE VANADIUM ALUMINUM MAGNESIUM CALCIUM | 2 1 2 5 3 1 4 40 20 3 0 100 1000 1000 | NR NR NR NR NR NR NR NR NR NR NR NR NR | NR NR NR NR NR NR NR NR NR NR NR NR NR N | 0.66 J 1 U 2 U 5 U 1.1 0.35 J 6.0 3.2 J 1300 41 J 1.9 J 677 251 J 85 J | NR NR NR NR NR NR NR NR NR NR NR NR NR N | 2 U 1 U 2 U 5 U 0.72 J 6.6 1.2 J 825 8.3 J 12 U 352 1200 U 68 J | 0.65 J 1.3 2 U 6 U 0.79 J 1 U 4.1 J 50 U 573 5.20 J 12 U 316 J 1200 U 75 J | 2.2 1 U 5.3 5 U 5.6 1 U 30 J 9.6 J 4010 J 31 J 6.4 J 3220 178 J 218 J |
| COBALT THALLIUM | 10 2 | | NR NR | : | NR NR | - | - | 11 U 2 U |
| VOLATILE ORGANIC (ug/kg) | - | - | - | - | | - | – | - |
| SEMI-VOLATILE ORGANIC (Ug/kg) BIS(2-ETHYLHEXYL)PHTHALATE BUTYLBENZYLPHTHALATE | - CRQL - 330 330 | NR | NR NR | 340 U 340 U | NR NR | 400 U 400 U | 400 U 400 U | 64 J 140 J |
| PEST/PCBs (ug/kg) | CRQL | | | | | | | |
| 4,4'-DDT Aroclor 1260 | - 16 160 | | NR NR | 16 ປ - | NR NR | 20 U | 20 U | 18 U 180 U |
| PH | | NR | NR | 6.37 | NR | 6.22 | 6.09 | 6.3 |
| PERCENT SOLIDS | | 98 | 98 | 97 | 97 | 81 | 81 | 91 |

Notes:

For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

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| ANALYTE | SITE DEPTH ID NUMBER DATE MATRIX | FS21-TB2 10-12 FS21TB0210 09/10/87 S01L | FS21-TB2 39-41 FS21B0239 09/10/87 S01L | FS21-TB2 59-61 FS21TB0259 09/10/87 SOIL |
|---|---|---|---|--|
| INORGANIC ANALYTES (mg/kg) | CRDI. | | | |
| ARSENIC BERYLLIUM CHROMIUM COPPER LEAD SELENIUM ZINC BARIUM IRON MANGANESE VANADIUM ALUMINUM MAGNESIUM CALCIUM COBALT THALLIUM | 2 1 2 5 3 1 4 40 20 3 10 1000 1000 1000 1000 1000 2 | 1 U 4.3 0.94 J 1.8 1 U 8.7 J 4.0 J 2020 J 53 J 2.0 J 914 354 J 76 J | NR NR NR NR NR NR NR NR NR NR NR NR NR N | 2 U 1 U 2 U 6 U 1.1 J 0.39 J 9.9 J 3.4 J 865 J 56 J 12 U 432 95 J 136 J |
| VOLATI'S ORGANIC (ug/kg) | | | | |
| SEMI-VOLATILE ORGANIC (ug/kg) | CRQL | - | - | - |
| BIS(2-ETHYLHEXYL)PHTHALATE BUTYLBENZYL#WTHALATE | 330 330 | | NR NR | 380 U 380 U |
| P2S7/PCBs (ug/kg) | CRQL | | | |
| 4,4'-DDT Aroclor 1260 | 16 160 | | NR NR | 45 190 U |
| РН | | 8.1 | NR | 6.99 |
| PERCENT SOL IUS | | 98 | | 86 |

Notes:

- For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), indicates that compound was undetected.

CS8XTBR-APPENDIX

| ANALYTE | SITE DEPTH ID NUMBER DATE MATRIX CRDL | CS8-TB2 54-56 CS8XTB0254 09/01/87 S0IL | CS8-TB2 54-56 CS8XTBD254 09/01/87 SOIL | CS8-TB3 8-10 CS8XTB0308 09/01/87 SOIL | CS8-TB3 18-20 CS8XTB0318 09/01/87 S0IL | CS8-TB3 58-60 CS8XTB0358 09/01/87 S01L | CS8-TB4 0-2 CS8XTB04X0 09/04/87 S01L | CS8-TB4 5-7 CS8XTB0405 09/04/87 SOIL |
|---|--|--|--|---|---|---|---|---|
| INORGANIC ANALYTES (mg/kg) | CRUE | | | | | | | |
| ARSENIC BERYLLIUM CHROMIUM COPPER LEAD THALLIUM ZINC BARIUM IRON MANGANESE VANADIUM ALUMINUM COBALT MAGNESIUM CALCIUM | 2 1 5 5 2 4 40 20 3 10 100 1000 1000 | 1 U 2 U 4.3 J 0.9 J 0.47 J 7.1 J 1.4 J 822 9.5 J 11 U 409 10 U 103 J | 2 U 1 U 2 U 4.3 J 0.81 J 0.52 J 8.8 J 1.3 J 1060 24 J 12 U 469 12 U 147 J 29 J | NR NR NR NR NR NR NR NR NR NR NR NR NR N | 2 U 1 U 2 U 3.5 J 2.1 J 0.52 J 7.1 J 1.7 J 821 J 27 J 10 U 429 10 U 83 J 23 J | 2 U 1 U 2 U 3.6 J 0.92 J 0.54 J 5.3 1.4 J 416 5.4 J 12 U 225 12 U 1160 U 23 J | NR NR NR NR NR NR NR NR NR NR NR NR NR N | NR NR NR NR NR NR NR NR NR NR NR NR NR N |
| SELENIUM | 1000 | 1 U | 1 U | NR | 25 J 1 U | 1 U | NR | · NR |
| VOLATILE ORGANIC (ug/kg) | CRQL | - | - | - | • | - | - | - |
| SEMI-VOLATILE ORGANIC (ug/kg) | CRQL | | | | | | | |
| BIS(2-ETHYLHEXYL)PHTHALATE BUTYLBENZYLPHTHALATE | 330 330 | | : | NR NR | - | - | NR NR | NR NR |
| PEST/PCBs (ug/kg) | CRQL | | | | | | | |
| 4,4'-DDT PCB-1260 | 16 160 | | 24 190 U | NR NR | 16 U 160 U | 19 U 190 U | NR NR | NR NR |
| PH | | 6.66 | 6.68 | NR | 6.7 | 6.84 | NR | NR |
| PERCENT SOLIDS | | 82 | 86 | 98 | 95 | 86 | 91 | 97 |

Notes:

1. For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.

Was analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

CS8XTBR-APPENDIX

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| ANALYTE INORGANIC ANALYTES (mg/kg) | SITE DEPTH ID NUMBER DATE MATRIX CRDL | CS8-TB4 8-10 CS8XTB0408 09/04/87 SOIL | CS8-TB4 10-12 CS8XTB0410 09/04/87 SOIL | CS8-TB4 34-36 CS8XTB0434 09/04/87 S0IL | CS8-TB4 54-56 CS8XTB0454 09/04/87 SOIL |
|---|--|--|---|--|--|
| ARSENIC BERYLLIUM CHROMIUM COPPER LEAD THALLIUM ZINC BARIUM IRON MANGANESE VANADIUM ALUMINUM COBALT MAGNESIUM CALCIUM SELENIUM | 2 1 2 5 3 4 40 20 3 10 100 1000 1000 | NR NR NR NR NR NR NR NR NR NR NR NR | 0.55 J 1 U 2 U 5 U 0.98 J 2 U 5.0 2.1 J 893 J 46 J 1.6 J 551 10 U 171 J 81 J 1 U | 0.55 J 1 U 2 U 5 U 0.52 J 2.1 J 5.0 2.1 J 583 16 J 10 U 373 10 U 82 J 63 J 0.33 J | 0.63 J 1 U 2 U 5 U 0.67 J 2 U 6.3 3.6 J 22 U 11 U 612 11 U 148 J 93 J 11 U |
| VOLATILE ORGANIC (ug/kg) | CRQL | - | - | - | - |
| SEMI-VOLATILE ORGANIC (ug/kg) BIS(2-ETHYLHEXYL)PHTHALATE BUTYLBENZYLPHTHALATE | CRQL 330 330 | | : | - | - |
| PEST/PCBs (ug/kg) | CRQL | | | | |
| 4,4'-DDT PCB-1260 | 16 160 | | 16 U 160 U | 16 U 160 U | 16 U 160 U |
| PH | | NR | 6.13 | 6.37 | 5.87 |
| PERCENT SOLIDS | | 97 | 97 | 97 | 90 |

Notes:

- For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), indicates that compound
- was undetected.

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CS8XTBR-APPENDIX

| ANALYTE | SITE DEPTH ID NUMBER DATE MATRIX | CS8-TP1 1-2 CS8TP0101 09/02/87 S0IL | CS8-TP6 5-6 CS8TP0605 09/02/87 S0IL | CS8-TB1 23-25 CS8XTB0123 09/01/87 S0IL | CS8-TB1 43-45 CS8XTB0143 09/01/87 S0IL | CS8-TB1 43-45 CS8XTBD143 09/01/87 S0IL | CS8-TB1 58-60 CS8XTB0158 09/01/87 S01L | CS8-TB2 19-21 CS8XTB0219 09/01/87 S01L |
|--|--|---|--|--|--|---|---|--|
| INORGANIC ANALYTES (mg/kg) | CRDL | | | | | | | |
| ARSENIC BERYLLIUM CHROMIUM COPPER LEAD THALLIUM ZINC BARIUM IRON MANGANESE VANADIUM ALUMINUM COBALT MAGNESIUM | 2 1 2 5 3 2 4 4 40 20 3 10 100 1000 1000 | 0.69 J 0.37 J 2.1 J 9.2 J 6.9 0.40 J 17 J 11 J 2490 59 J 3.7 1300 1.5 J 536 J 331 J | 2 U 1 U 2 U 4.1 J 1.0 J 0.41 J 6.6 J 1.7 J 961 19 J 647 10 U 129 J | 2 U 1 U 2 U 3.1 J 0.84 J 6.7 J 1.5 J 418 10 J 10 U 259 10 U 1000 U | 2 U 1 U 2 J 3.2 J 0.82 J 0.45 J 8.1 J 2.6 J 718 15 J 10 U 401 10 U 95 J 36 J | 2 U 0.40 J 3.1 1.2 0.44 J 9.2 J 4.1 J 2170 23 J 2.9 J 925 1.30 J 56 J | 2 U 1 U 2 U 4.7 J 0.72 J 0.75 J 12 J 2.8 J 979 53 J 11 U 483 1.5 J 196 J | 2 U 1 U 2 U 3.6 J 0.78 J 0.41 J 7.6 J 1.5 J 690 22 J 10 U 420 10 U 1000 U |
| CALCIUM SELENIUM | 1 | 331 J 1 U | 27 J 0.3 U | 27 J 1 U | 30 J 1 U | 50 J 1 U | 42 J 1 U | 27 J 1 U |
| VOLATILE ORGANIC (ug/kg) | CRQL | - | - | • | - | - | - | - |
| SEMI-VOLATILE ORGANIC (ug/kg) | CRQL | | | | | | | |
| BIS(2-ETHYLHEXYL)PHTHALATE BUTYLBENZYLPHTHALATE | 330 330 | 350 U 350 U | 340 U 340 | 330 U 330 U | 340 U 340 U | 340 U 340 U | 390 U 390 U | 92 U 340 U |
| PEST/PCBs (ug/kg) | CRQL | | | | | | | |
| 4,4'-DDT PCB-1260 | 16 160 | 16 U 660 | 16 U 160 U | 16 U 160 U | 16 U 160 U | 16 U 160 U | 19 U 190 U | 18 160 U |
| PH | | 6.52 | 6.94 | 6.53 | 6.56 | 6.63 | 6_61 | 6.73 |
| PERCENT SOLIDS | | 94 | 98 | 99 | 97 | 98 | 84 | 97 |

Notes:

For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

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| | SIT | Έ; | CS8-MW2 | | CS8-MW2 DUP | | CS8-MW4 | CS8-MW4 DUP | | FS21-MW1 | F | S21-MW2 |
|---|---|--|--|-----------------------|---|----------------------------|--|---------------------------------|----------------------------|---|---|--|
| | ID NUMBE DAT MATRI | Έ: | JCS8MW2002 01/07/88 WATER | | JDUP002X02 01/07/88 WATER | | JCS8MW4001 10/28/87 WATER | JDUP003X01 10/28/87 WATER | | JFS21MW001 10/28/87 WATER | | 21MW202 01/07/88 WATER |
| INORGANIC COMPOUNDS | ANALYTICAL METHOD | CRDL ug/L | | | | | | | | | | |
| Alumninum Arsenic Barium Beryllium Calcium Iron Lead Magnesium Manganese Potassium Sodium Cobalt | P F P P P P P P P P P P P | 200 10 200 5 5000 100 5 5000 5000 5000 5 | 44 2.5 37 3.6 5220 89 5.2 1200 69 2690 12100 50 | 1 1 1 1 1 | 200 2.7 35 5020 78 2.7 915 66 5000 10900 50 | 0 1 0 1 1 0 | 1980 3.5 J 25 J 3960 J 654 5 U 1250 J 181 1890 J 25000 J 7 U | 5000 607 | 1 1 1 1 1 1 | 1670 10 U 67 J 5 U 4600 J 21 U 9.7 J 1540 J 281 2110 J 12000 J 6.4 U | · | 200 U 2.9 J 44 J 2 J 10200 3.6 J 1080 J 2700 3250 J 10800 50 U |
| VOLATILE ORGANIC CON | | 50 | 50 | U | 50 | U | 70 | 0.4 | U | 0.4 U | | 50 0 |
| 2-Butanone | | 10 | 10 | U | 10 | U | 10 (| J 10 | U | 19 U | | NR |
| SEMI-VOLATILE ORGANIC | COMPOUNDS | CRQL ug/L | | | | | | | | | | |
| Benzo(b)Fluoranthene/Benzo(k) Chrysene Fluoranthene Pyrene Naphthalene | Fluoranthene | e 10 10 10 10 10 | 20 20 20 20 20 20 | บ บ บ บ บ | 20 20 20 20 20 | บ บ บ บ | 10 t 10 t 10 t |) 10 J 10 J 10 | U U | 10 10 10 | | 3 J 2 J 3 J 3 J 20 U |
| DILUTION FACTOR: OTHER | | RL mg/L | 2 | | 2 | | | | | | | 2 |
| Nitrate-N Nitrite-N | | 0.040 | _ _ | | - | | - | • - | | - | | - |

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- For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), indicates that compound was undetected.

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| PROJECT: MMR 2- | | | | Volatile Aqueous A | nalysis (ug/L) |
|---------------------------------|--|---|---|---|----------------|
| Tabl Validation / | e 2 Summary Table | | | | |
| | SAMPLE LOCATION: LAB NUMBER: DATE SAMPLED: DATE ANALYZED: | CS8MW2003 270836 06/28/89 07/02/89 | DUP001X03 270837 06/28/89 07/02/89 | DUP002X03 270841 06/28/89 07/02/89 | |
| ANALYTE | CRQL | | | | |
| Volatile Organi | c Compounds | - | | | |
| 2-Butanone | 10 | 6400 | 3800 | 1200 | |
| =============================== | Dilution Factor: | 42 | 29 | | |
| Associat | iated Method Blank: ed Equipment Blank: ociated Trip Blank: | CB890702B23 SB002X003 TB002X003 | CB890702B23 SB002X003 TB002X003 | | |
| | | Notes: | | | |

For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

07-Sep-93

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PROJECT: MMR 2-3C

Semivolatile Aqueous Analysis (ug/L)

14-Jun-93

MR215S2R

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Table 2 Validation / Summary Table

| L/ DATE DATE I | LOCATION: AB NUMBER: SAMPLED: EXTRACTED: ANALYZED: | CS8MW2003 270836 06/28/89 06/30/89 07/02/89 | DUP001X03 270837 06/28/89 06/30/89 07/02/89 | |
|--|--|---|---|------|
| ANALYTE | CRQL | | | |
| Naphthalene 2-Methylnaphthalene Acenaphthylene Acenaphthene Fluorene Phenanthrene Anthracene Fluoranthene Pyrene Benzo(a)Anthracene Chrysene Benzo(b)Fluoranthene Benzo(b)Fluoranthene Benzo(a)Pyrene Indeno(1,2,3-cd)pyrene Dibenz(a,h)anthracene Benzo(g,h,i,)perylene | 10 10 10 10 10 10 10 10 10 10 10 10 | 2 10 10 10 10 10 10 10 10 10 10 10 10 | U 10 U 10 | |
| PRESERVICE Diluti | | *====================================== | | 2222 |
| Associated Met Associated Equipm | hod Blank: | | G3J71066A08 SB002X003 | |

Notes:

For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

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PROJECT: MMR 2-3C

Inorganic Aqueous Analysis (ug/L)

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14-Jun-93

Table 2 Validation / Summary Table

| | SAMPLE LOCATION: | CS8MW2003 | DUP001x03 |
|---|---|--|---|
| | LAB NUMBER: | 270863 | 270871 |
| | DATE SAMPLED: | 06/28/89 | 06/28/89 |
| ANALYTE | CRDL | | |
| Aluminum Arsenic Barium Calcium Cobalt Iron Lead Magnesium Manganese Potassium Sodium | 200 10 200 5 5000 50 100 5 5000 15 5000 5000 | 180 J 1.2 U 53.4 J 0.5 UJ 4800 J 12.6 J 2070 1.1 J 994 U 198 1740 U 10200 | 146 J 1.2 U 39.3 J 0.5 UJ 3670 J 13.8 J 2080 2.2 UJ 996 U 198 1740 U 10800 |
| | ated Method Blank: | 17094C | 17094C |
| | d Equipment Blank: | SB002X003 | SB002X003 |

Notes:

For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

PROJECT: MMR 2-3C

TPH Aqueous Analysis (mg/L)

08-Sep-93

MR215M2R

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Table 2 Validation / Summary Table

| | SAMPLE LOCATION: LAB NUMBER: DATE SAMPLED: DATE EXTRACTED: DATE ANALYZED: | CS8MW2003 270880 06/28/89 07/06/89 07/08/89 | DUP001x03 270881 06/28/89 07/06/89 07/08/89 | |
|----------------------------|---|---|---|---|
| ANALYTE | RL | | | |
| Petroleum Hydrocar | bons, Total 1 | 1 | บ 1 บ | |
| 2222#222# 2 2#22#22 | ************* | | | = |
| | ed Method Blank: Equipment Blank: | 271849 SB002X003 | 271849 SB002X003 | |

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Notes:

For data in 1988 and earlier, undetected values for inorganics are indicated by CRDL, qualified with a U.
 NR = Analysis not requested.
 In samples not used in the preliminary risk evaluation (PRE), - indicates that compound was undetected.

Page 1

| Location | Sample ID | Date | Depth | Type | Matrix | Test | Prep | Analyte | Result | DL | RL | Units | Qual |
|------------|-------------|---------|-------|------|--------|--------|--------|---------------------------------------|--------|-------|------|-------|----------|
| | | - | | | | | | · · · · · · · · · · · · · · · · · · · | 1 | | | | |
| FS-21 MW-1 | | | | | | | | | | | | | <u> </u> |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | ALUMINUM (TOTAL) | 1180 | 22 | 100 | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | ANTIMONY (TOTAL) | ND | 1.9 | 5 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | ARSENIC (TOTAL) | ND | 1.9 | 5 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | BARIUM (TOTAL) | 70.4 | 0.3 | 20 | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | BERYLLIUM (TOTAL) | 0.94 | 0.4 | 1 | UG/L | J |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | CADMIUM (TOTAL) | 0.4 | 0.4 | 1 | UG/L | J |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | CALCIUM (TOTAL) | 6030 | 28.1 | 500 | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | CHROMIUM (TOTAL) | ND | 1.2 | 5 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | COBALT (TOTAL) | ND | 2.6 | 5 | UG/L | υ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | COPPER (TOTAL) | ND | 0.8 | 5 | UG/L | UJ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | IRON (TOTAL) | ND | 15.6 | 100 | UG/L | UJ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | LEAD (TOTAL) | ND | 1.1 | 2 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | MAGNESIUM (TOTAL) | 1200 | 21.8 | 500 | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | MANGANESE (TOTAL) | 708 | 0.3 | 10 | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | NICKEL (TOTAL) | ND | 3.8 | 20 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | POTASSIUM (TOTAL) | 1300 | 46.3 | 1500 | UG/L | J |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | SELENIUM (TOTAL) | ND | 2.7 | 5 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | SILVER (TOTAL) | ND | 0.95 | 10 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | SODIUM (TOTAL) | 21700 | 98.8 | 500 | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | THALLIUM (TOTAL) | ND | 2.5 | 10 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | VANADIUM (TOTAL) | ND | 1 | 10 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C200.7 | TOTAL | ZINC (TOTAL) | 19.3 | 1.9 | 20 | UG/L | J |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | C245.2 | TOTAL | MERCURY (TOTAL) | ND | 0.012 | 0.2 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,1,1-TRICHLOROETHANE | ND | 0.21 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | Ñ1 | WG | CVOL | METHOD | 1,1,2,2-TETRACHLOROETHANE | ND | 0.18 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,1,2-TRICHLOROETHANE | ND | 0.23 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,1-DICHLOROETHANE | ND | 0.19 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,1-DICHLOROETHENE | ND | 0.21 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | METHOD | 1,2,4-TRICHLOROBENZENE | ND | 0.31 | 1 | UG/L | υ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.37 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,2-DIBROMOETHANE (EDB) | ND | 0.22 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | 1,2-DICHLOROBENZENE | ND | 0.26 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,2-DICHLOROETHANE | ND | 0.18 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | | N1 | WG | CVOL | METHOD | 1,2-DICHLOROPROPANE | ND | 0.15 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 1,3-DICHLOROBENZENE | ND | 0.24 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | | | WG | CVOL | METHOD | 1,4-DICHLOROBENZENE | ND | 0.2 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | 2-HEXANONE | ND | 0.87 | 5 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | | N1 | WG | CVOL | METHOD | ACETONE | ND | 0.82 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | BENZENE | ND | 0.19 | | UG/L | |

Appendix A - Additional Groundwater Sampling CS-8/FS-21 Decision Document

Appendix A - Additional Groundwater Sampling

CS-8/FS-21 Decision Document

| Location | Sample ID | Date | Depth | Туре | Matrix | Test | Prep | Analyte | Result | DL | RL | Units | Qua |
|------------|--|---------|-------|------|--------|--------|--|-------------------------------------|--------|--------|------|-------|-----------|
| 04MW0001 | 04MW0001-02 | 3/10/99 | | | WG | CVOL | METHOD | BROMOCHLOROMETHANE | ND | 0.23 | | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | | | WG | CVOL | METHOD | BROMODICHLOROMETHANE | ND | 0.19 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | BROMOFORM | ND | 0.27 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | BROMOMETHANE | ND | 0.16 | 1 | UG/L | <u>lu</u> |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | METHOD | CARBON DISULFIDE | ND | 0.21 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | CARBON TETRACHLORIDE | ND | 0.16 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | N1 | WG | CVOL | METHOD | CHLOROBENZENE | ND | 0.19 | 1 | UG/L | U |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | METHOD | CHLOROETHANE | ND | 0.19 | 1 | UG/L | tu - |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | METHOD | CHLOROFORM | ND | 0.16 | 1 | UG/L | lu – |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | METHOD | CHLOROMETHANE | ND | 0.18 | | UG/L | _ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | METHOD | CIS-1,2-DICHLOROETHYLENE | ND | 0.2 | | UG/L | _ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | CIS-1,3-DICHLOROPROPENE | ND | 0.14 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | DIBROMOCHLOROMETHANE | ND | 0.24 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | ETHYLBENZENE | ND | 0.18 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | wğ | CVOL | | METHYL ETHYL KETONE (2-BUTANONE) | ND | 0.97 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | METHYL ISOBUTYL KETONE (4-METHYL-2- | ND | 0.81 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | METHYLENE CHLORIDE | ND | 0.19 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | STYRENE | ND | 0.17 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | TERT-BUTYL METHYL ETHER | ND | 0.17 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | 4 | TETRACHLOROETHYLENE(PCE) | ND | 0.35 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | TOLUENE | ND | 0.00 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | TRANS-1,2-DICHLOROETHENE | ND | 0.18 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | TRANS-1,3-DICHLOROPROPENE | ND | 0.10 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | | | WG | CVOL | | TRICHLOROETHYLENE (TCE) | ND | 0.14 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | VINYL CHLORIDE | ND | 0.10 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | CVOL | | XYLENES, TOTAL | ND | 0.14 | | UG/L | |
| 04MW0001 | 04MW0001-02 | 3/10/99 | 60.1 | | WG | E504 | design of the second se | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.0045 | | UG/L | _ |
| 04MW0001 | 04MW0001-02 | 3/10/99 | | | WG | E504 | | 1,2-DIBROMOETHANE (EDB) | ND | 0.0043 | | UG/L | |
| | | | | 141 | 100 | L-004 | | | | 0.0047 | 0.01 | 100/2 | <u>++</u> |
| CS-8 MW-2 | —————————————————————————————————————— | | | | ┣─── | | | | | | | | + |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | C200.7 | TOTAL | ALUMINUM (TOTAL) | ND | 80.1 | 100 | UG/L | ┼╓╌╴ |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | ANTIMONY (TOTAL) | ND | 1.9 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | ARSENIC (TOTAL) | ND | 1.9 | | | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | BARIUM (TOTAL) | 17.2 | | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | BERYLLIUM (TOTAL) | ND | 0.3 | | UG/L | |
| | | | | | | | | | | | | | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | | 0.35 | 0.4 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | | 2590 | 28.1 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | | WG | C200.7 | TOTAL | | ND | 1.2 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | | ND | 0.6 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | N1 | WG | C200.7 | TOTAL | COPPER (TOTAL) | ND | 0.8 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | C200.7 | TOTAL | IRON (TOTAL) | ND | 83.5 | 100 | UG/L | . [U |
| CS8 mod1da | ata.xls | | | | | | Page 2 of | 5 | | | | 8/5/9 | 19 |
| - | _ <i>_</i> _ | - | - | | - | - | - | | - | 3 | | ÷ | |
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Appendix A - Additional Groundwater Sampling CS-8/FS-21 Decision Document

| Location | Sample ID | Date | Depth | Туре | Matrix | Test | Prep | Analyte | Result | DL | RL. | Units | Qual |
|----------|-------------|---------|-------|------|--------|--------|--------|-----------------------------|--------|-------|------|-------|-----------|
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | C200.7 | TOTAL | LEAD (TOTAL) | ND | 1.1 | 2 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | WG | C200.7 | TOTAL | MAGNESIUM (TOTAL) | 287 | 21.8 | 500 | UG/L | J |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | | C200.7 | TOTAL | MANGANESE (TOTAL) | 34.1 | 0.3 | 10 | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | | C200.7 | TOTAL | NICKEL (TOTAL) | ND | 1.9 | 20 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | | C200.7 | TOTAL | POTASSIUM (TOTAL) | 1120 | 46.3 | 1500 | UG/L | J |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | WG | C200.7 | TOTAL | SELENIUM (TOTAL) | ND | 2.7 | 5 | UG/L | υ |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | WG | C200.7 | TOTAL | SILVER (TOTAL) | ND | 0.59 | 10 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | C200.7 | TOTAL | SODIUM (TOTAL) | 18800 | 98.8 | 500 | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | | C200.7 | TOTAL | THALLIUM (TOTAL) | ND | 2.5 | 10 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | WG | C200.7 | TOTAL | VANADIUM (TOTAL) | ND | 1 | 10 | UG/L | υ |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | C200.7 | TOTAL | ZINC (TOTAL) | ND | 4.2 | 20 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | | C245.2 | TOTAL | MERCURY (TOTAL) | ND | 0.025 | 0.2 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | | | CVOL | METHOD | 1,1,1-TRICHLOROETHANE | ND | 0.21 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,1,2,2-TETRACHLOROETHANE | ND | 0.18 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,1,2-TRICHLOROETHANE | ND | 0.23 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | | CVOL | METHOD | 1,1-DICHLOROETHANE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,1-DICHLOROETHENE | ND | 0.21 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,2,4-TRICHLOROBENZENE | ND | 0.31 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | | WG | CVOL | METHOD | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.37 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,2-DIBROMOETHANE (EDB) | ND | 0.22 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,2-DICHLOROBENZENE | ND | 0.26 | 1 | UG/L | <u>lu</u> |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WĠ | CVOL | METHOD | 1,2-DICHLOROETHANE | ND | 0.18 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,2-DICHLOROPROPANE | ND | 0.15 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,3-DICHLOROBENZENE | ND | 0.24 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 1,4-DICHLOROBENZENE | ND | 0.2 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | 2-HEXANONE | ND | 0.87 | 5 | UG/L | . U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | ACETONE | ND | 0.82 | 5 | UG/L | . U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | BENZENE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | BROMOCHLOROMETHANE | ND | 0.23 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | BROMODICHLOROMETHANE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | | CVOL | METHOD | BROMOFORM | ND | 0.27 | 1 | UG/L | . U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | BROMOMETHANE | ND | 0.16 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | CARBON DISULFIDE | ND | 0.21 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | CARBON TETRACHLORIDE | ND | 0.16 | 1 | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | CHLOROBENZENE | ND | 0.19 | 1 | ŪG/L | U I |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WĞ | CVOL | METHOD | CHLOROETHANE | ND | 0.19 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | | CVOL | METHOD | CHLOROFORM | ND | 0.16 | 1 | UG/L | TU - |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | | CHLOROMETHANE | ND | 0.18 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | CIS-1,2-DICHLOROETHYLENE | ND | 0.2 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | | CIS-1,3-DICHLOROPROPENE | ND | 0.14 | | UG/L | |
| 04MW0002 | 04MW0002-03 | 3/10/99 | | | | CVOL | | DIBROMOCHLOROMETHANE | ND | 0.24 | | UG/L | |

Appendix A - Additional Groundwater Sampling

CS-8/FS-21 Decision Document

| Location | Sample ID | Date | Depth | Type | Matrix | Test | Prep | Analyte | Result | DL | RL | Units | Qual |
|------------------|---------------|---------|-------|------|--------|--------|-------------|-------------------------------------|--------|--------|------|--------|------|
| | 04MW0002-03 | 3/10/99 | 60 | | WG | | | | ND | 0.18 | | UG/L | |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | METHYL ETHYL KETONE (2-BUTANONE) | ND | 0.97 | 5 | UG/L | U I |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | METHYL ISOBUTYL KETONE (4-METHYL-2- | ND | 0.81 | | UG/L | |
| | 04MW0002-03 | 3/10/99 | 60 | | WG | CVOL | | | ND | 0.19 | 2 | UG/L | U |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | STYRENE | ND | 0.17 | | UG/L | |
| | 04MW0002-03 | 3/10/99 | 60 | | WG | CVOL | | | ND | 0.17 | | UG/L | |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | TETRACHLOROETHYLENE(PCE) | ND | 0.18 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | TOLUENE | ND | 0.19 | 1 | ŨG/L | U |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | TRANS-1,2-DICHLOROETHENE | ND | 0.18 | 1 | UG/L | U |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | TRANS-1,3-DICHLOROPROPENE | ND | 0.14 | 1 | UG/L | υ |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | TRICHLOROETHYLENE (TCE) | ND | 0.16 | 1 | ŪG/L | U |
| | 04MW0002-03 | 3/10/99 | 60 | | WG | CVOL | | | ND | 0.14 | | | U |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | CVOL | METHOD | XYLENES, TOTAL | ND | 0.2 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | E504 | METHOD | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.0045 | 0.01 | UG/L | U |
| | 04MW0002-03 | 3/10/99 | 60 | N1 | WG | E504 | | 1,2-DIBROMOETHANE (EDB) | ND | 0.0047 | 0.01 | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | ALUMINUM (TOTAL) | ND | 65.1 | 100 | UG/L | ΰ |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | | ANTIMONY (TOTAL) | ND | 1.9 | 5 | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | | ARSENIC (TOTAL) | ND | 1.9 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | BARIUM (TOTAL) | 17.1 | 0.3 | | UG/L | J |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | BERYLLIUM (TOTAL) | ND | 0.4 | | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | | CADMIUM (TOTAL) | 0.62 | 0.4 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | CALCIUM (TOTAL) | 2620 | 28.1 | 500 | UG/L | |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | CHROMIUM (TOTAL) | ND | 2.01 | | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | COBALT (TOTAL) | ND | 0.6 | 5 | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | COPPER (TOTAL) | ND | 0.8 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | IRON (TOTAL) | ND | 67.3 | 100 | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | TOTAL | LEAD (TOTAL) | ND | 1.1 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | MAGNESIUM (TOTAL) | 280 | 21.8 | 500 | UG/L | J |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | MANGANESE (TOTAL) | 35.2 | 0.3 | 10 | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | TOTAL | NICKEL (TOTAL) | ND | 1.9 | | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | POTASSIUM (TOTAL) | 1180 | 46.3 | 1500 | UG/L | J |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | SELENIUM (TOTAL) | ND | 2.7 | 5 | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | TOTAL | SILVER (TOTAL) | ND | 1.9 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | C200.7 | TOTAL | SODIUM (TOTAL) | 19000 | 98.8 | 500 | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | TOTAL | THALLIUM (TOTAL) | ND | 2.5 | | UG/L | U |
| | 04MW0002-03FD | 3/10/99 | | | WĠ | C200.7 | TOTAL | VANADIUM (TOTAL) | ND | 1 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | C200.7 | TOTAL | ZINC (TOTAL) | 8.4 | 1.9 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | C245.2 | TOTAL | MERCURY (TOTAL) | ND | 0.026 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | CVOL | | 1,1,1-TRICHLOROETHANE | ND | 0.21 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | CVOL | METHOD | | ND | 0.18 | | UG/L | |
| | 04MW0002-03FD | 3/10/99 | | | WG | CVOL | METHOD | | ND | 0.23 | | UG/L | |
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Appendix A - Additional Groundwater Sampling

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CS-8/FS-21 Decision Document

| Location | Sample ID | Date | Depth | Type | Matrix | Test | Prep | Analyte | Result | DL | RL | Units | Qual |
|----------|---------------|---------|-------|------|--------|------|--|-------------------------------------|--------|--------|----|-------|------|
| 04MW0002 | 04MW0002-03FD | 3/10/99 | | | | CVOL | METHOD | 1,1-DICHLOROETHANE | ND | 0.19 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,1-DICHLOROETHENE | ND | 0.21 | 1 | UG/L | υ |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,2,4-TRICHLOROBENZENE | ND | 0.31 | 1 | UG/L | υ |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.37 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/*0/99 | 60 | FD1 | WG | CVOL | METHOD | 1,2-DIBROMOETHANE (EDB) | ND | 0.22 | 1 | UG/L | TU T |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,2-DICHLOROBENZENE | ND | 0.26 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,2-DICHLOROETHANE | ND | 0.18 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,2-DICHLOROPROPANE | ND | 0.15 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,3-DICHLOROBENZENE | ND | 0.24 | 1 | UG/L | TU 1 |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 1,4-DICHLOROBENZENE | ND | 0.2 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | 2-HEXANONE | ND | 0.87 | 5 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | ACETONE | ND | 0.82 | 5 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | BENZENE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | BROMOCHLOROMETHANE | ND | 0.23 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | BROMODICHLOROMETHANE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | BROMOFORM | ND | 0.27 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | BROMOMETHANE | ND | 0.16 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CARBON DISULFIDE | ND | 0.21 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CARBON TETRACHLORIDE | ND | 0.16 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | | CHLOROBENZENE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CHLOROETHANE | ND | 0.19 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CHLOROFORM | ND | 0.16 | 1 | UG/L | tu 1 |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CHLOROMETHANE | ND | 0.18 | 1 | UG/L | tu l |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CIS-1,2-DICHLOROETHYLENE | ND | 0.2 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | CIS-1,3-DICHLOROPROPENE | ND | 0.14 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | DIBROMOCHLOROMETHANE | ND | 0.24 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | ETHYLBENZENE | ND | 0.18 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | METHYL ETHYL KETONE (2-BUTANONE) | ND | 0.97 | 5 | UG/L | JU I |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | METHYL ISOBUTYL KETONE (4-METHYL-2- | ND | 0.81 | 5 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | METHYLENE CHLORIDE | ND | 0.19 | 2 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | STYRENE | ND | 0.17 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | TERT-BUTYL METHYL ETHER | ND | 0.17 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | TETRACHLOROETHYLENE(PCE) | ND | 0.18 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | TOLUENE | ND | 0.19 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | TRANS-1,2-DICHLOROETHENE | ND | 0.18 | 1 | UG/L | U |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | | TRANS-1,3-DICHLOROPROPENE | ND | 0.14 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | | CVOL | METHOD | TRICHLOROETHYLENE (TCE) | ND | 0.16 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | CVOL | METHOD | VINYL CHLORIDE | ND | 0.14 | | UG/L | - |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | | FD1 | | CVOL | | XYLENES, TOTAL | ND | 0.2 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | E504 | | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.0045 | | UG/L | |
| 04MW0002 | 04MW0002-03FD | 3/10/99 | 60 | FD1 | WG | E504 | and the second | | ND | 0.0047 | | UG/L | |

Allachment A CS-8/FS-21 Groundwater Monitoring Results

| Location | Date | Analyle | Result | ĎL. | RL | Units | Qualilier |
|------------|----------|------------------------------|-----------|--------|------|---------|-----------|
| 04B/10001 | 12/14/99 | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.0017 | 0.01 | jug/L | U |
| 04BH0001 | 12/14/99 | 1,2-DIBROMOETHANE (EDB) | ND | 0.0083 | 0.01 | 119/1 | U |
| 04BH0001 | 12/14/99 | ANTIMONY (TOTAL) | ND | 4.5 | 60 | μg/L | U |
| 04BH0001 | 12/14/99 | ARSENIC (TOTAL) | 7.4 | 3.9 | 10 | ug/L | J |
| 04BH0001 | 12/14/99 | BERYLLIUM (TOTAL) | ND | 0.7 | 5 | µg/L | ŪJ |
| 04BH0001 | 12/14/99 | CADMIUM (TOTAL) | ND | 0.4 | 5 | IIg/L | ŪJ |
| 048H0001 | 12/14/99 | CHROMIUM (TOTAL) | ND | 0.7 | 10 | ug/L | U |
| 04BH0001 | 12/14/99 | COPPER (TOTAL) | ND | 2 | 25 | µg/L | U |
| 048110001 | 12/14/99 | NICKEL (TOTAL) | ND. | 1.4 | 40 | µg/L. | UJ. |
| 048H0001 | 12/14/99 | SELENIUM (TOTAL) | ND | 3.6 | 5 | µg/L | U U |
| 048H0001 | 12/14/99 | SILVER (TOTAL) | ND | 0.8 | 10 | 199/1 | U |
| 04BH0001 | 12/14/99 | ZINC (TOTAL) | 29.6 | 0.4 | 20 | µg/L | |
| 04BH0001 | 12/14/99 | LEAD (TOTAL) | ND | 6 | 15 | µg/L | U |
| 048H0001 | 12/14/99 | LEAD (TOTAL) | ND | 6 | 15 | rig/L | U |
| 04BH0001 | 12/14/99 | MERCURY (TOTAL) | ND | 0.1 | 0.2 | µg/L | Ū |
| 04BH0001 | 12/14/99 | THALLIUM (TOTAL) | ND | 1.7 | 2 | 119/1 | Ū |
| 04BH0001 | 12/14/99 | THALLIUM (TOTAL) | ND | . 1.7 | 2 | Lig/L | U |
| D4BH0001 | 12/14/99 | BENZO(B)FLUORANTHENE | ND | 0.023 | 0.1 | -11g/L | U |
| 04BH0001 | 12/14/99 | BENZO(K)FLUORANTHENE | ND | 0.023 | 0.1 | hâyr | U |
| 04BH0001 | 12/14/99 | 2,2'-OXYBIS(1-CHLORO)PROPANE | ND | 3.27 | 4.81 | 119/1. | ŧJ |
| 04BH0001 | 12/14/99 | 2,4,5-TRICHLOROPHENOL | ND | 2,02 | 19.2 | 119/1. | U |
| 048110001 | 12/14/99 | 2,4,6-TRICHLOROPHENOL | ND | 2.21 | 4.81 | 119/1. | U |
| 04BH0001 | 12/14/99 | 2,4-DICHLOROPHENOL | ND | 2.6 | 4.81 | µg/L | U |
| 04BI-10001 | 12/14/99 | 2,4-DIMETHYLPHENOL | ND | 2.4 | 4.81 | µg/L | <u> </u> |
| 04BH0001 | 12/14/99 | 2,4-DINITROPHENOL | ND | 2.5 | 19.2 | 11g/L | U |
| 04BH0001 | 12/14/99 | 2,4-DINITROTOLUENE | ND | 1.92 | 4.81 | | <u> </u> |
| 04BH0001 | 12/14/99 | 2,G-DINITROTOLUENE | ND | 1.44 | 1.81 | 119/1_ | <u> </u> |
| 04BH0001 | 12/14/99 | 2-CHLORONAPHTHALENE | ND | 2.79 | 4.81 | 119/1. | <u> </u> |
| 04BH0001 | 12/14/99 | 2-CHLOROPHENOL | ND | 3.08 | 4.81 | µg/1 | <u> </u> |
| 04BH0001 | 12/14/09 | 2-METHYLNAPHTHALENE | ND | 2.98 | 4.81 | <u></u> | <u> </u> |
| 04BH0001 | 12/14/99 | 2-METHYLPHENOL (O-CRESOL) | ND | 2.88 | 4.81 | <u></u> | U |
| 048110001 | 12/14/99 | 2-NITROANILINE | ND | 1.64 | 19.2 | H11/L | U |
| 04BH0001 | 12/14/99 | 2-NITROPHENOL | ND | 2.98 | 4.81 | pg/l. | U |
| 048H0001 | 12/14/99 | 3,3'-DICHLOROBENZIDINE | ND | 2.88 | 4.81 | µg/L | U |
| 04BH00D1 | 12/14/99 | 3-NITROANILINE | <u>ND</u> | 2.02 | 19.2 | Lig/L | U |
| 048H0001 | 12/14/99 | 4,6-DINITRO-2-METHYLPHENOL | ND | 2.6 | 19,2 | 111/1. | U |

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|------------|-----------|-----------------------------|-----------|--------|-------|---------------|--------------|-------------|
| Location | Date | Analyle | L. Result | DL. | RL | Units | Qualifier | |
| 04B10001 | 12/14/99. | 4-JROMOPHENYL PHENYL ETHER | ND | 2.21 | 4.81 | pg/L | U. | |
| -048H0001 | 12/14/99 | 4-CHLORO-3-METHYLPHENOL | ND | 2.12 | 4.81 | -ug/L | <u> </u> | |
| 04BH0001 | 12/14/99 | 4-CHLOROANILINE | ND | 2.5 | 4.81 | Jug/L | UJ | ł. |
| 04BH0001 | 12/14/99 | 4-CHLOROPHENYL PHENYL ETHER | ND · | 1.64 | 4.81 | µg/L | U | Í |
| 048H0001 | 12/14/99 | 4-METHYLPHENOL (P-CRESOL) | ND | 2.69 | 4.81 | µg/L | U | |
| 04BH0001 | 12/14/99 | 4-NITROANILINE | ND | 4.14 | 19.2 | 119/L | Ū | 4 |
| 04BH0001 | 12/14/99 | 4-NITROPHENOL | ND | 4.04 | 19.2 | 119/L | U | 1 |
| 04BH0001 | 12/14/99 | ACENAPHTHENE | ND | 2.12 | 4.81 | µg/L | U . | 1 |
| 04BH0001 | 12/14/99 | ACENAPHTHYLENE | ND | 2.21 | 4.81 | µg/L | U | |
| 04BH0001 | 12/14/99 | ANTHRAGENE | ND | 1.64 | | -ug/L | . <u>U</u> | _ |
| 048110001 | 12/14/99 | BENZO(A)ANTHRACENE | ND | 0.865 | 4.81 | µg/L | <u>U</u> · : | } . |
| 04BH0001 | 12/14/99 | BENZO(A)PYRENE | - | - | - | µg/L | R | ł |
| 04BH0001 | 12/14/99 | BENZO(B)FLUORANTHENE | ND | 1.35 | 4.81 | Ug/L | U | |
| 048110001 | 12/14/99 | BENZO(G,H,I)PERYLENE | ND | 3.08 | 4.81 | µg/L | ป | |
| 048110001 | 12/14/99 | BENZO(K)FLUORANTHENE | ND | 1.25 | 4.81 | jug/L | u | ł |
| 048110001 | 12/14/99 | BENZYL BUTYL PHTHALATE | ND | 1.25 | 4.81 | 119/L | U | 8 |
| 048/10001 | 12/14/99 | BIS(2-CHLOROETHOXY) METHANE | ND | · 3.17 | 4.81 | ug/L | U | { |
| 04BH0001 | 12/14/99 | BIS(2-CHLOROETHYL) ETHER | ND | 3.36 | 4.81 | µg/1_ | <u> </u> | - |
| 048H0001 | 12/14/99 | BIS(2-ETHYLHEXYL) PHTHALATE | ND | 3,75 | 4.81 | µg/L | U | [[|
| 04BH0001 | 12/14/99 | CHRYSENE | ND | 1.06 | 4.81 | 11g/L | U | 1 |
| 04BH0001 | 12/14/09 | DI-N-BUTYL PHITHALATE | ND | 1.92 | 4.81 | Light | Ű | 1 |
| 04BH0001 | 12/14/99 | DI-N-OCTYLPHTHALATE | ND | 3.17 | 4.81 | Ig/l. | U | fi - |
| 04BH0001 | 12/14/99 | DIBENZ(A,H)ANTHRACENE | ND | 2.6 | 4.81 | JIG/L | U | J. |
| 04BH0001 | 12/14/99 | DIBENZOFURAN | ND | 1.92 | 4.81 | µg/L | U | X |
| 048H0001 | 12/14/99 | DIETHYL PHTHALATE | ND . | 1.25 | 4.81 | /L | | |
| 04BH0001 | 12/14/99 | DIMETHYL PHITHALATE | ND | 1,83 | 1.81 | L | IJ | |
| 04BH0001 | 12/14/99 | FLUORANTIENE | ND | 1.44 | 4.81 | ! <u>19/1</u> | U | 1 |
| 04BH0001 | 12/14/99 | FLUORENE | ND | 1.64 | 4.81 | <u>_ug/L</u> | 0 | 8 |
| 04BH0001 | 1/2/14/99 | HEXACHLOROBENZENE | ND | 2.02 | 4.81 | <u> </u> | <u> </u> | |
| 0481-10001 | 12/14/99 | HEXACHLOROBUTADIENE | ND | 3.27 | 4.81 | ING/L | 11 | |
| 04BH0001 | 12/14/99 | HEXACHLOROCYCLOPENTADIENE | ND | 1.06 | 4.81 | Ing/L | <u> </u> | |
| 04BH0001 | 12/14/99 | HEXACHLOROETHANE | NĎ | 3.36 | 4.81 | 1141. | (j | .1 |
| 048110001 | 12/14/99 | INDENO(1,2,3-C,D)PYRENE | ND | 2.6 | 4.81 | p./L | U | J - |
| 04BH0001 | 12/14/99 | ISOPHORONE | ND | 2.80 | 4.81 | 119/1- | U | |
| 04BH0001 | 12/14/99 | N-NITROSODI-N-PROPYLAMINE | ND | 3.17 | 4.81 | LIGA | U.I | .# |
| 04BH0001 | 12/14/99 | N-NITROSODIPHENYLAMINE | ND | 1.54 | 4.81 | uga. | UJ . | |

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Atlachment A CS-8/FS-21 Groundwaler Monitoring Results

| Location | Dale 1 | Analyte | Result | DL I | RL | Units | Qualifier |
|------------|----------|-----------------------------|--------|------|-------|--------------|-----------|
| 04BH0001 | 12/14/99 | NAPHTHALENE | ND ND | 2.98 | 4.81 | µg/L | U |
| 048H0001 | 12/14/99 | NITROBENZENE | ND | 3.17 | 4.81 | 119/4. | Ū |
| 048140001 | 12/14/99 | PENTACHLOROPHENOL | ND | 2.31 | 19.2 | μg/L | ŭ |
| 048110001 | 12/14/99 | PHENANTI IRENE | ND | 1.64 | 4.81 | µg/L | U |
| 0481-10001 | 12/14/99 | PHENOL | ND | 3.08 | 4.81 | . <u>[</u> | <u> </u> |
| 04BH0001 | 12/14/99 | PYRENE | ND | 1.83 | 4.81 | ug/L ug/L | <u>u</u> |
| 04BH0001 | 12/14/99 | 1,1,1-TRICHLOROETHANE | ND | 0.23 | 1 | µg/L | <u>U</u> |
| -04BH0001 | 12/14/99 | 1,1,2,2-TETRACHLOROETHANE | ND | 0.32 | | µg/L | <u> </u> |
| 048110001 | 12/14/99 | 1, 1, 2-TRICHLOROETHANE | ND | 0.32 | 1 | ug/l. | U U |
| 04BH0001 | 12/14/99 | 1,1-DICHLOROETHANE | NÐ | 0.29 | | l lig/L | <u> </u> |
| 04BH0001 | 12/14/99 | 1,1-DICHLOROETHENE | ND ND | 0.25 | l | ug/l. | <u>U</u> |
| 0481-10001 | 12/14/99 | 1,2,4-TRICHLOROBENZENE | ND | 0.31 | | 119/L | <u>U</u> |
| 048H0001 | 12/14/99 | 1,2-DIBROMO-3-CHLOROPROPANE | ND | 0.43 | 1 | Ing/L | <u> </u> |
| 04BH0001 | 12/14/99 | 1.2-DIBROMOETHANE (EDB) | ND | 0.28 | 1 | L have | |
| 04BH0001 | 12/14/99 | 1.2-DICHLOROBENZENE | ND | 0.24 | 1 | lig/L | <u> </u> |
| 04BH0001 | 12/14/99 | 1.2-DICHLOROETHANE | ND | 0.3 | 1. | µg/L | Ű |
| 04BH0001 | 12/14/99 | 1.2-DICHLOROPROPANE | ND | 0,31 | 1 | µg/L | U U |
| 04BH0001 | 12/14/99 | 1,3-DICHLOROBENZENE | ND | 0.25 | 1 | 11g/L | <u> </u> |
| 04BI-10001 | 12/14/99 | · 1.4-DICHLOROBENZENE | ND | 0.26 | 1 | 11g/L | <u> </u> |
| 0481-0001 | 12/14/99 | 2-HEXANONE | | - | | µg/L | <u> </u> |
| 048H0001 | 12/14/99 | ACETONE | | - | | ug/L | R |
| 04BH0001 | 12/14/99 | BENZENE | ND | 0,28 | 1 | 110/L | |
| 04BH0001 | 12/14/99 | BROMOCHLOROMETHANE | ND | 0.3 | 1 | Jig/L | Ū |
| 04BH0001 | 12/14/99 | BROMODICHLOROMETHANE | ND | 0,25 | 1 | µg/L | U |
| 04BH0001 | 12/14/99 | BROMOFORM | ND | 0,26 | 1 | JIGA_ | Ų |
| 04BI-10001 | 12/14/99 | BROMOMETHANE | ND | 0.28 | 1 | 11g/1_ | U |
| 04BH0001 | 12/14/99 | CARBON DISULFIDE | ND | 0.29 | 1 | µg/L | U |
| 04BH0001 | 12/14/99 | CARBON TETRACHLORIDE | ND | 0.27 | 1 | 11g/L_ | U |
| 04BH0001 | 12/14/99 | CHLOROBENZENE | ND | 0.25 | 1 | ug/l_ | U |
| 04BH0001 | 12/14/99 | CHLOROETHANE | ND | 0.27 | 1 | lig/L | U |
| 04BH0001 | 12/14/99 | CHLOROFORM | ND | 0,29 | 1 | LIG/L | U |
| 04BI-10001 | 12/14/99 | CHLOROMETHANE | ND | 0.28 | 1 | 119/1. | U |
| 04BH0001 | 12/14/99 | CIS-1,2-DICHLOROETHENE | ND | 0.24 | 1 | Hg/L_ | U |
| 04BH0001 | 12/14/99 | CIS-1,3-DICHLOROPROPENE | ND | 0.32 | 1 | fajlt. | 1 |
| 04BH0001 | 12/14/99 | DIBROMOCHLOROMETHANE | ND | 0.28 | 1 | HIJA | 1 |
| 04BH0001 | 12/14/99 | ETHYLBENZENE | 1.14 | 0.21 | 1 | 1111/L. | |

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CDD CS-8/FS-21 Groundwater Monitoring Results

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| Location | Date | Analyle | Result | DL | RL | Units | Qualifier- |
|------------------------|----------|----------------------------------|--------|------|-----|--------|---|
| 04BI10001 | 12/14/99 | METHYL ETHYL KETONE (2-BUTANONE) | - | - | - | µg/L | R |
| 04BH0001 | 12/14/99 | METHYL ISOBUTYL KETONE | ND | 1.42 | | µg/L | U |
| - 04 0H0901 | | METHYL-TERT-BUTYL-ETHER (MTBE) | ND | 0.45 | 1 | /L | . U . |
| -04BH0001 | 12/14/99 | METHYLENE CHLORIDE | ND | 0.28 | . 2 | · µg/L | ······································· |
| 04BH0001 | 12/14/99 | STYRENE | ND | 0.26 | 1 | 11g/L | U |
| 048H0001 | 12/14/99 | TETRACHLOROETHENE(PCE) | ND | 0.22 | 1 | µg/L | U |
| 04BH0001 | 12/14/99 | TOLUENE | ND | 0,29 | 1 | µg/L | U |
| 04BH0001 | 12/14/99 | TRANS-1,2-DICHLOROETHENE | ND | 0.24 | 1 | 119/1_ | U |
| 04BI 10001 | 12/14/99 | TRANS-1,3-DICHLOROPROPENE | ND | 0.44 | 1 | pg/L | U |
| 048H0001 | 12/14/99 | TRICHLOROETHENE(TCE) | NO | | | | U- |
| 04BH0001 | 12/14/99 | VINYL CHLORIDE | ND | 0.27 | 1 | 119/1 | U |
| 04BH0001 | 12/14/99 | XYLENES, TOTAL | 3.04 | 0.79 | 1 | µg/L_ | |

DL = detection limit

RL = reporting limit

ND = nondelect

 $\mu g/L \approx micrograms per liter$ U = concentration of analyte is below detection fimit

J = estimated concentration

Us = concentration of analyte is estimated to be below detection limit

R = rejected

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Attachment A CS-8/FS-21 Groundwater Monitoring Results

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| Location | Sample Date | Témpérature (⁰ C) | Dissolved Oxygen (ng/L) | Turbidily (NTU) | Oxidallon/Reduction Potentiat (mV) | Specific Conductance (IIS/cm) | pH (SV) | |
|---|-----------------------------------|----------------------------------|-------------------------------|--|--|---------------------------------------|---------------------------------------|------------|
| 048110001 | 12/14/99 | 11.48 | 0.41 | 7.90 | 155.70 | 124.00 | 6,18 | • |
| C = degrees centigr ng/L = milligrams pe VTU = nephelometri nV = millivolis IS/cm = microsieme SU = standerd units | ade ar liter c lurbidity ur | bils | <u> </u> | | | · · · · · · · · · · · · · · · · · · · | · · · · · · · · · · · · · · · · · · · | |
| | | | · · · · · · | · · · · | | | | |
| •••• | | | | | | | • | |
| • • • • • • • • • • • | | | · · · · · · · · · · · · · · · | •••••••••••••••••••••••••••••••••••••• | ···· ·· ·· ·· ·· ·· ·· ·· ·· ·· ·· ·· · | · · ~ · · | | . . |
| • | | | | · · · | | | - | |
| · · · · · · · · · · · · · · · · · · · | | · · | · . | • | | ••• | | • |
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CS-8/FS-21 Sample Information

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| - | | | Ground | | | Northing | Easling |
|---|--|--------------|-----------|---------------|-------------|------------|------------|
| · | ······································ | | Surface | Sample | Groundwater | Coordinate | Coordinate |
| | Well | | Elevation | Elevation | Elevation | | |
| | Identification | Date Sampled | (ft msl) | (ft msl) | (ft msl) | (feet) | (feet) |
| | 04BH0001 | 14-Dec-99 | 107.21 | 35.01 - 40.01 | 50.21 | 237008.12 | 862883.32 |

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Elevation is measured in feet mean sea level (ft msl).

Elevatori is medsured is loet mean sed level (it mar).

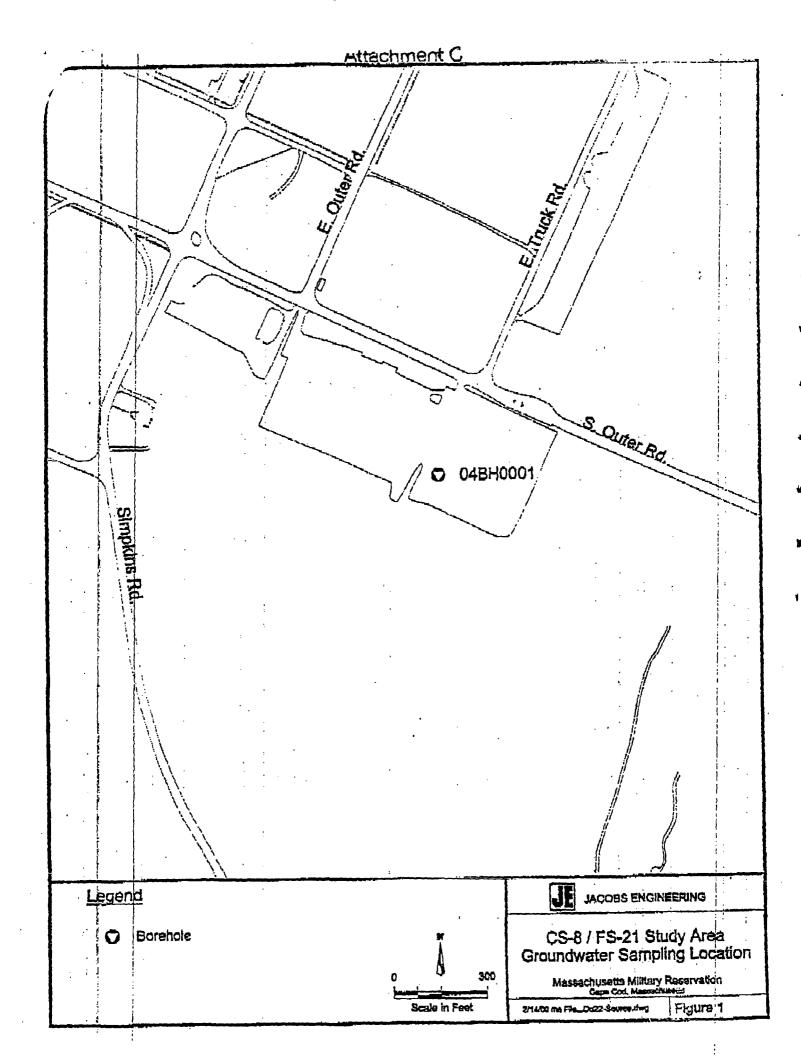
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APPENDIX B

APPENDIX B SUMMARY OF FIELD LABORATORY ANALYTICAL DATA FOR TERRAPROBE BORINGS

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| Parameter | PHASE I Site: CS-8/CD-1 Matr: SOIL Dpth: 4 ISIS: 04TRX11XX4X1XF Date: 28-FEB-92 | SUMP REMOVAL AC CS-8/CD-1 SOIL 8 04TRX11XX8X1XF 28-FEB-92 | TION PROGRAM - M CS-8/CD-1 SOIL 12 04TRX11X12X1XF 28-FEB-92 | MASSACHUSETTS MIL CS-8/CD-1 SOIL 18 04TRX11X18X1XF 28-FEB-92 | ITARY RESERVATIO CS-8/CD-1 SOIL 24 04TRX11X24X1XF 28-FEB-92 | N CS-8/CD-1 SOIL 4 04TRX12XX4X1XF 28-FEB-92 | CS-8/CD-1 SOIL 8 04TRX12XX8X1DF 28-FEB-92 | CS-8/CD-1 SOIL 8 04TRX12XX8X1XF 28-FEB-92 |
|--------------------------|--|--|--|---|--|--|---|---|
| PI meter reading (ppm) | 1.3 | 0.0 | 0.0 | 0.6 | 0.0 | 0.0 | 1.2 | 1.2 |
| VOCs (ug/kg) Toluene | ND | ND | ND | ND | ND | ND | 130 | ND |
| Ethylbenzene | ND | ND | ND | ND | ND | ND | 190 | ND |
| o-Xylene | ND | ND | ND | ND | ND | ND | 67 | ND |
| m/p-Xylene | ND | ND | ND | ND | ND | ND | 250 | ND |
| SVOCs (ug/kg) | ND | ND | . ND | ND | ND | ND | ND | ND |
| Pest/PCBs (ug/kg) | | | | | | | | |
| Delta-BHC | ND | ND | ND | ND | ND | ND | 330 | 530 |
| Heptachlor epoxide | ND | ND | ND | ND | ND | 140 | 1700 | 1800 |
| Inorganics (mg/kg) | | | | · | | | | |
| Lead | ND | ND | ND | ND | ND | ND | ND | ND |
| Copper | ND | ND | ND | ND | ND | ND | 270 | ND |
| Zinc | ND | ND | ND | ND | ND | ND | 380 | ND |
| Arsenic | ND "ND | ND | ND | ND | ND ND | NÐ | ND | ND · |
| Chromium TRK (market) | 130 | ND 83 | ND 84 | ND 120 | 120 | ND 130 | ND 980 | 820 |
| TPH (mg/kg) | 100 | 60 | 04 | 120 | 120 | 120 | 900 | 820 |

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|------------------------------------|---|-----------------------------|----------------------|-----------------------------|---------------------|-----------------------------|-----------------------------|-----------------------------|
| | PHASE I | SUMP REMOVAL AC | | | LITARY RESERVATIO | | | |
| | Site: CS-8/CD-1 | CS-8/CD-1 | CS-8/CD-1 | CS-8/CD-1 | CS-8/CD-1 | CS-8/CD-1 | CS-8/CD-1 | CS-8/CD-1 |
| | Matr: SOIL | SOIL | SOIL | SOIL | SOIL | SOIL | SOIL | SOIL |
| | Dpth: 12 | 18 0/ TOV12V18V1V5 | 24 04TRX12X24X1XF | 4 | 8 04TRX13XX8X1DF | 8 | 12 | 18 |
| Parameter | ISIS: 04TRX12X12X1XF Date: 28-FEB-92 | 04TRX12X18X1XF 28-FEB-92 | 28-FEB-92 | 04TRX13XX4X1XF 04-MAR-92 | 04-MAR-92 | 04TRX13XX8X1XF 04-MAR-92 | 04TRX13X12X1XF 04-MAR-92 | 04TRX13X18X1XF 04-MAR-92 |
| *arameter | Date: 20-110 72 | 20-100-15 | 20-FED-72 | 04-MAK-76 | U4-19AK-76 | U4 * MAK * 76 | U4*MAX*76 | 04-mak-72 |
| ol meter reading (ppm) | 5.2 | 1.9 | 1.9 | 0.0 | 0.0 | 0.0 | 0.0 | 0.0 |
| /OCs (ug/kg) | | | | | | _ | | |
| Toluene | ND | ND | ND | | | ND | ND | ND |
| Ethylbenzene | ND | ND [·] | ND | | | ND | ND | ND |
| o-Xylene | ND ND | ND ND | ND ND | | ND | ND | ND | ND |
| n/p-Xylene SVOCs (ug/kg) | ND ND | ND ND | NU ND | UM ND | ND ND | ND ND | ND ND | ND ND |
| Pest/PCBs (ug/kg) | ND ND | 10 | 20 | UI UI | 10 | NU | hν | ND |
| Delta-BHC | 150 | ND | ND | ND | ND | ND | ND | ND |
| leptachlor epoxide | 1300 | ND | ND | | | ND | ND | ND |
| Inorganics (mg/kg) | | | | | | | | |
| ead | ND | ND | ND | ND | ND | ND | ND | ND |
| opper | ND | ND | ND | | ND | ND | ND | ND |
| linc | ND | ND | ND | ND | ND | ND | ND | ND |
| Arsenic | ND | ND | ND | ND | ND | ND | ND | ND |
| Chromium | ND 83 | ND 120 | ND 40 | ND 170 | ND 330 | ND 320 | ND 290 | ND |
| TPH (mg/kg) | CO | 120 | 40 | 170 | 066 | 320 | 290 | 410 |

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|--|--|-----------------------------|-----------------------------|-----------------------------|-----------------------------|-----------------------------|--|--|
| | PHASE I SUMP REMOVAL ACTION PROGRAM - MASSACHUSETTS MILITARY RESERVATION | | | | | | | |
| | Site: CS-8/CD-1 Matr: SOIL Dpth: 24 | CS-8/CD-1 SOIL | CS-8/CD-1 SOIL 8 | CS-8/CD-1 SOIL 12 | CS-8/CD-1 SOIL 18 | CS-8/CD-1 SOIL 24 | | |
| Parameter | ISIS: D4TRX13X24X1XF Date: 04-MAR-92 | 04TRX14XX4X1XF 04-mar-92 | 04TRX14XX8X1XF 04-MAR-92 | 04TRX14X12X1XF 04-MAR-92 | 04TRX14X18X1XF 04-MAR-92 | 04TRX14X24X1XF 04-MAR-92 | | |
| PI meter reading (ppm) VOCs (ug/kg) | 0.0 | 0.0 | 0.0 | 2.9 | 0.0 | 0.0 | | |
| Toluene | ND | ND | ND | ND | ND | ND | | |
| Ethylbenzene | ND | ND | ND | ND | ND | ND | | |
| -Xylene | ND | ND | ND | ND | ND | ND | | |
| p-Xylene | ND | 6.1 | ND | ND | ND | ND | | |
| OCs (ug/kg) st/PCBs (ug/kg) | ND | ND | ND | ND | ND | ND | | |
| elta-BHC | NÐ | ND | ND | ND | ND | ND | | |
| eptachlor epoxide norganics (mg/kg) | ND | ND | ND | ND | ND | ND | | |
| ead | ND | ND | ND | ND | ND | ND | | |
| opper | ND | ND | ND | ND | ND | ND | | |
| inc | ND | ND | ND | ND | ND | ND | | |
| rsenic | ND | ND | ND | ND | ND | ND | | |
| hromium | ND | ND | ND | ND | ND | ND | | |
| TPH (mg/kg) | 410 | . 180 | 160 | 120 | 210 | 410 | | |

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APPENDIX C

APPENDIX C UNDERGROUND STORAGE TANK REMOVAL DET AILS

mmr/docs/cs8dd/text/cs8-fs21 Final DD2. doc

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March 5, 1996 Reference #9509874

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Colonel Louis Volpe State Quartermaster The Commonwealth of Massachusetts Office of the State Quartermaster Camp Curtis Guild, Reading, Massachusetts 01867-1999

Re: UST Removal Report 5000 Gallon Gasoline Tank Office and Maintenance #22, Building #9002 Outer Road Camp Edwards, Massachusetts 02542

Dear Colonel Volpe:

On February 13, 1996, Simmons Environmental Services, Inc. (SIMMONS) observed the removal of a 6000 gallon gasoline underground storage tank (UST) at the area designated as Office and Maintenance #22, Building #9002, Outer Road in Camp Edwards, Massachusetts. (See Site Plan).

The UST removal, cleaning and disposal was conducted by personnel from ZECCO, INC. of Northboro, Massachusetts. Philip Motta, Assistant Chief Fire Inspector for Otis Air Force Base was onsite for inspection of the UST and its removal. Stephan H. Landry, Hydrogeologist of SIMMONS, was onsite to observe the subsurface conditions, field screen the UST excavation soils for headspace volatiles, and collect UST excavation floor soil samples for laboratory analyses. Mr. John Callan was onsite as the site Project Manager.

The UST was situated underneath a 10" thick concrete pad and connected to an adjacent fuel pump. The UST area was approximately 50' east of the OMS #22 building. There was no surficial staining observed on the concrete pavement or around the UST fill pipe. According to the ZECCO vacuum truck operator, prior to excavation, approximately 1600 gallons of gasoline product was removed from the UST. The top of the UST was approximately 2.5' below grade. The UST was measured to be 8'1" in diameter and 13'8" in length. Upon removal, the UST was noted to be in excellent condition with no evidence of holes, pitting or staining. According to Mr. Callan, the UST was approximately 6-7 years old.

> 375 Elm Street Salisbury, MA 01952 Telephone 508-463-6669 Fax 508-463-6679



There was no overt evidence of soil staining or odors from the UST excavation soils. As specified by Mr. Callan, representative soils from the excavation north end and south end bottoms and east sidewall were collected and field screened in accordance with the Massachusetts Department of Environmental Protection (MDEP) jar headspace methodology using a precalibrated HNu Photoionization Detector (PID) model 101 (10.2 eV). The results of the headspace field screening is shown in the table below expressed in parts per million (ppm). Each of the soil samples showed less than 10 ppm headspace, the samples were then transported under a Chain of Custody to a state certified laboratory for analysis of volatile organic compounds (VOCs) by EPA Method 8240 with Ethylene Dibromide, Semi-volatile organic compounds (TPHs) by EPA Method 8100 modified and total lead. The results of these analyses are also shown in the table below.

| LOCATION | PID. (ppm) | TPHs (mg/Kg) | VOCs (ug/Kg) | -SVOCs- (ug/Kg) | TotaFlead (mg/Kg) |
|---------------------|---------------|-----------------|-----------------|--------------------|----------------------|
| North End Bottom | 0 | ND | ND | ND | 5.2 |
| South End Bottom | 2.2 | ND | ND | ND | 3.9 |
| East Wall | 0 | ND | ND | ND | <2.6 |

ND = No Detectable concentrations above laboratory method's detection limit

There were no exceedances of the MDEP's Reportable Concentrations as listed in 310 CMR 40.0000, the Massachusetts Contingency Plan (MCP), for the compounds analyzed. Therefore, no further actions are warranted. The analytical data and site diagram of the work to date are enclosed. If there are any questions, please call either of the undersigned at 508-463-6669.

Very truly yours,

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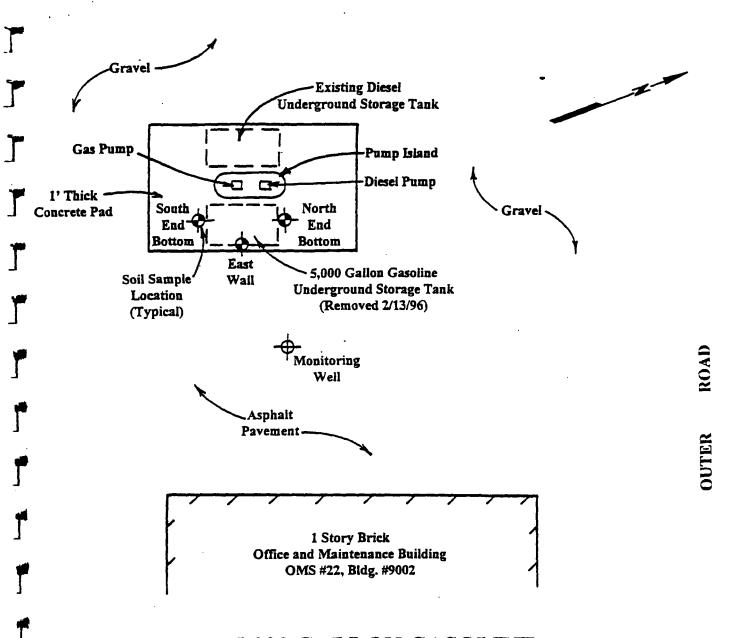
Rohyn Chausson

ジー Stephan H. Landry Hydrogeologist

Id ASin

William A. Simmons, LSP Principal

WAS:rac



5,000 GALLON GASOLINE UNDERGROUND STORAGE TANK CLOSURE &

SOIL SAMPLING PLAN

FOR PROPERTY AT: OMS #22, BLDG. #9002 OUTER ROAD CAMP EDWARDS, MA

PREPARED BY: SIMMONS ENVIRONMENTAL SERVICES, INC. 375 ELM STREET SALISBURY, MA 01952 (508) 463-6669

| 0' | 10' | 20' | 40' |
|----|-----|-----|-----|
| | | | |

SCALE: 1"=20' PROJECT #: 9509874 DATE: FEBRUARY 13, 1996 DRAFTED: SHL/AK

AMRO Environmental Laboratories Corporation



111 Herrick Street, Merrimack, NH 03054 TEL: (603) 424-2022 · FAX: (603) 429-8496

February 23, 1996

Mr. William Simmons Simmons Environmental Services, Inc. 375 Elm St. Crossroads Plaza Salisbury, MA 01952

RE Your project: 9509874 Mass Guard

Dear William:

Enclosed please find the results for the above-referenced project, received on February 14, 1996. AMRO operates a Quality Control Program which meets or exceeds EPA and state requirements. A copy of the appropriate State Certificate is attached. No quality control deviations were noted during the analyses associated with this project. This project was assigned AMRO Project Number 12257. If you have any questions regarding this project in the future, please refer to this number.

Please be advised that any unused sample volume and sample extracts will be stored for a period of thirty (30) days from this report date. After this time, AMRO will properly dispose of the remaining sample. If you require further analysis, or need the samples held for a longer period, please contact us immediately.

This letter is an integral part of your data report.

Please do not hesitate to call if you have any questions.

Sincerely,

Maria N. Borduz, Ch.E. President

Encl.

| client: | | ironmental Laborato | Clie | nt Design | | | |
|--|---|---|--|---|---|----------------|----------------------|
| 375 Elm Crossroa Salisbur | St. Nds Plaza | tal Services, Inc. 01952 nmons | 95 | 09874 Mas | s Guard | | |
| Samples Qt | y/Type: : | 3/Solid | Date Sa Date Re | esignatic ampled: ec'vd: omplete: 14071 | on: 12257 02/13/ 02/14/ 02/22/ | 96 | |
| Sample Identity | AMRO Identity | Test Parameter | Results | Units | Date of Analysis | Run by | EPA Metho |
| SK UST, East Wall | 12257-01 | Total Solids Digestion Lead, Total | 96.5 | % mg∕Kg | 02/15/96 02/16/96 02/22/96 | RR TM TM | 2540 3050 6010 |
| SK UST, South End Bottom | 12257-02 | Total Solids Digestion Lead, Total | 97.1 3.9 | ۶, mg/Kg | 02/15/96 02/16/96 02/22/96 | RR TM TM | 2540 3050 6010 |
| SK UST, North End Bottom | 12257-03 | Total Solids Digestion Lead, Total | 96.8 <2.6 | ¥ mg∕Kg | 02/15/96 02/16/96 02/22/96 | RR TM TM | 2540 3050 6010 |
| All ana US St Ed The fol all section < = | EPA Methods andard Meth ition, 1992 lowing star ns: | ormed in accordance of Chemical Analy ods for the Examin and USEPA S dard abbreviations of followed by the | sis for W ation of W846 Manu and conv | Water an Mal, 3rd. Ventions | d Wastewat ed. | | |
| | | | Certified | i by: <u> </u> | haria N. Bor | . Bo duz, | ch.E. |

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LABORATORY REPORT

Petroleum Hydrocarbons by Gas Chromatography EPA Method 8100 (Modified)

Client: <u>Simmons Environmental Services</u>, Inc. Client I.D.: <u>9509874 Mass Guard</u>

<u>SK UST, East Wall</u> AMRO I.D.: <u>12257-01</u> Date sampled: <u>02/13/96</u> Date received: <u>02/14/96</u> Date prepared: <u>02/19/96</u> Date analyzed: <u>02/20/96</u> Sample Qty/Type: <u>1/Solid</u>

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-------------------------|--------------------|---------------------------|
| Gasoline | ND | 51 |
| Kerosene | ND | 51 |
| Mineral Spirits | ND | 51 |
| Fuel Oil #2/Diesel | ND | 51 |
| Fuel Oil #4 | ND | 51 |
| Fuel Oil #6 | ND | 100 |
| Motor Oil/Hydraulic Oil | ND | 51 |

Gasoline results are provided for screening purposes only. The recommended procedure for gasoline analysis is a modified EPA 8015 or 8240 (purge and trap). Solid Content = 96.5%. Results are in dry weight. Comments:

ND = Not Detected at or above the reporting limit.

Analyzed By: JK

Approved by <u>Jamen Mandone</u> for Nancy Stewart

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LABORATORY REPORT

Petroleum Hydrocarbons by Gas Chromatography EPA Method 8100 (Modified)

Client: <u>Simmons Environmental Services</u>, <u>Inc.</u> Client I.D.: <u>9509874 Mass Guard</u>

<u>SK UST, South End Bottom</u> AMRO I.D.: <u>12257-02</u> Date sampled: <u>02/13/96</u> Date received: <u>02/14/96</u> Date prepared: <u>02/19/96</u> Date analyzed: <u>02/20/96</u> Sample Qty/Type: <u>1/Solid</u>

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-------------------------|--------------------|---------------------------|
| Gasoline | ND | 50 |
| Kerosene | ND | 50 |
| Mineral Spirits | ND | 50 |
| Fuel Oil #2/Diesel | ND | 50 |
| Fuel Oil #4 | ND | 50 |
| Fuel Oil #6 | ND | 99 |
| Motor Oil/Hydraulic Oil | ND | 50 |

Gasoline results are provided for screening purposes only. The recommended procedure for gasoline analysis is a modified EPA 8015 or 8240 (purge and trap). Solid Content = <u>97.1%</u>. Results are in dry weight. Comments:

ND = Not Detected at or above the reporting limit.

Analyzed By: JK

Y Lauren Mardone for Nancy Stewart Approved by Jac

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LABORATORY REPORT

Petroleum Hydrocarbons by Gas Chromatography EPA Method 8100 (Modified)

Client: <u>Simmons Environmental Services</u>, Inc. Client I.D.: <u>9509874 Mass Guard</u> <u>SK UST</u>, North End Bottom AMRO I.D.: <u>12257-03</u>

Date sampled: 02/13/96 Date received: 02/14/96 Date prepared: 02/19/96 Date analyzed: 02/20/96 Sample Qty/Type: 1/Solid

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-------------------------|--------------------|---------------------------|
| Gasoline | ND | 50 |
| Kerosene | ND | 50 |
| Mineral Spirits | ND | 50 |
| Fuel Oil #2/Diesel | ND | 50 |
| Fuel Oil #4 | ND | 50 |
| Fuel Oil #6 | ND | 99 |
| Motor Oil/Hydraulic Oil | ND | 50 |

Gasoline results are provided for screening purposes only. The recommended procedure for gasoline analysis is a modified EPA 8015 or 8240 (purge and trap). Solid Content = <u>96.8%</u>. Results are in dry weight. Comments:

ND = Not Detected at or above the reporting limit.

Analyzed By: JK

Approved by runen Mardane for Nancy Stewart

LABORATORY REPORT EPA Method 8240B Volatile Organic Compounds

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| Client: <u>Simmons Environmental Services</u> , Inc. |
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| Client I.D.: <u>9509874 Mass Guard</u> |
| SK UST, East Wall |
| AMRO I.D.: 12257-01 |
| Date sampled: 02/13/96 Date received: 02/14/96 |
| Date prepared: 02/15/96 Date analyzed: 02/16/96 |
| Sample Qty/Type: 1/Solid |

| Test Parameter | Results (ug/kg) | Reporting Limit(ug/kg) |
|--------------------------------|--------------------|---------------------------|
| Chloromethane | ND | 67 |
| Bromomethane | ND | 67 |
| Vinyl Chloride | ND | 67 |
| Chloroethane | ND | 67 |
| Methylene Chloride | ND | 27 |
| Acetone | ND | 130 |
| Carbon Disulfide | ND | 27 |
| 1,1-Dichloroethene | ND | 27 |
| 1,1-Dichloroethane | ND | 27 |
| 1,2-Dichloroethene (trans) | ND | 27 |
| 1,2-Dichloroethene (cis) | ND | 27 |
| Chloroform | ND | 27 |
| 1,2-Dichloroethane | ND | 27 |
| 2-Butanone (MEK) | ND | 130 |
| 1,1,1-Trichloroethane | ND | 27 |
| Carbon Tetrachloride | ND | 27 |
| Vinyl Acetate | ND | 130 |
| Bromodichloromethane | ND | 27 |
| 1,2-Dichloropropane | ND | 27 |
| cis-1,3-Dichloropropene | ND | 27 |
| Trichloroethene | ND | 27 |
| Dibromochloromethane | ND | 27 |
| 1,1,2-Trichloroethane | ND | 27 |
| Benzene | ND | 27 |
| trans-1,3-Dichloropropene | ND | 27 |
| Bronoform | ND | 27 |
| 4-Methyl-2-Pentanone (MIBK) | ND | 130 |
| 2-Hexanone | ND | 130 |
| Tetrachloroethene | ND | 27 |
| 1,1,2,2-Tetrachloroethane | ND | 27 |
| Toluene | ND | 27 |
| Chlorobenzene | ND | 27 |
| Ethylbenzene | ND | 27 |
| Styrene | ND | 27 |
| Xylene (total) | ND | 27 |
| Methyl-tert-butyl ether (MTBE) | ND | 27 |
| 1,2-Dibromoethane (EDB) | ND | 27 |

Solid Content = 96.5%. Results are in dry weight. ND = Not Detected at or above the reporting limit

Analyzed By: <u>SK</u>

Approved by <u>Kamen Mandone</u> for Nancy Stewart

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LABORATORY REPORT

EPA Method 8240B Volatile Organic Compounds

| Client: <u>Simmons Environmental Services</u> , Inc. |
|--|
| Client I.D.: <u>9509874 Mass Guard</u> |
| SK UST, South End Bottom |
| AMRO I.D.: <u>12257-02</u> |
| Date sampled: <u>02/13/96</u> Date received: <u>02/14/96</u> |
| Date prepared: 02/15/96 Date analyzed: 02/16/96 |
| Sample Qty/Type: 1/Solid |

| Test Parameter | Results (ug/kg) | Reporting Limit(ug/kg) |
|--------------------------------|--------------------|---------------------------|
| Chloromethane | ND | 65 |
| Bromomethane | ND | 65 |
| Vinyl Chloride | ND | 65 |
| Chloroethane | ND | 65 |
| Methylene Chloride | ND | 26 |
| Acetone | ND | 130 |
| Carbon Disulfide | ND | 26 |
| 1,1-Dichloroethene | ND | 26 |
| 1,1-Dichloroethane | ND | 26 |
| 1,2-Dichloroethene (trans) | ND | 26 |
| 1,2-Dichloroethene (cis) | ND | 26 |
| Chloroform | ND | 26 |
| 1,2-Dichloroethane | ND | 26 |
| 2-Butanone (MEK) | ND | 130 |
| 1,1,1-Trichloroethane | ND | 26 |
| Carbon Tetrachloride | ND | 26 |
| Vinyl Acetate | ND | 130 |
| Bromodichloromethane | ND | 26 |
| 1,2-Dichloropropane | ND | 26 |
| cis-1,3-Dichloropropene | ND | 26 |
| Trichloroethene | ND | 26 |
| Dibromochloromethane | ND | 26 |
| 1,1,2-Trichloroethane | ND | 26 |
| Benzene | ND | 26 |
| trans-1,3-Dichloropropene | ND | 26 |
| Bromoform | ND | 26 |
| 4-Methyl-2-Pentanone (MIBK) | ND | 130 |
| 2-Hexanone | ND | 130 |
| Tetrachloroethene | ND | 26 |
| 1,1,2,2-Tetrachloroethane | ND | 26 |
| Toluene | ND | 26 |
| Chlorobenzene | ND | 26 |
| Ethylbenzene | ND | 26 |
| Styrene | ND | 26 |
| Xylene (total) | ND | 26 |
| Methyl-tert-butyl ether (MTBE) | ND | 26 |
| 1,2-Dibromoethane (EDB) | ND | 26 |

Solid Content = 97.1%. Results are in dry weight. ND = Not Detected at or above the reporting limit

Analyzed By: <u>SK</u>

Approved by <u>Lauren</u> Mardane for Nancy Stewart

| EPA Method | | |
|--|-----------------------|-----------------|
| Volatile Organic | Compounds | |
| Client: Simmons Environment | | nc. |
| Client I.D.: <u>9509874 Mass Gr</u> SK UST, North | | |
| AMRO I.D.: 12257-03 | | |
| Date sampled: 02/13/96 Date | | |
| Date prepared: <u>02/15/96</u> Date Sample Qty/Type: <u>1/Solid</u> | e analyzed: <u>02</u> | /16/96 |
| Test | Results | Reporting |
| Parameter | (ug/kg) | Limit(ug/k |
| Chloromethane | ND | 66 |
| Bromomethane | ND | 66 |
| Vinyl Chloride | ND | 66 |
| Chloroethane | ND | 66 |
| Methylene Chloride | ND | 26 |
| Acetone Carbon Disulfide | ND | 130 |
| | ND | 26 |
| 1,1-Dichloroethene | ND ND | 26 |
| 1,1-Dichloroethane 1,2-Dichloroethene (trans) | ND | 26 26 |
| 1,2-Dichloroethene (cialis) | ND | 26 |
| Chloroform | ND | 26 |
| 1,2-Dichloroethane | ND | 26 |
| 2-Butanone (MEK) | ND | 130 |
| 1,1,1-Trichloroethane | ND | 26 |
| Carbon Tetrachloride | ND | 26 |
| Vinyl Acetate | ND | 130 |
| Bromodichloromethane | ND | 26 |
| 1,2-Dichloropropane | ND | 26 |
| cis-1,3-Dichloropropene | ND | 26 |
| Trichloroethene | ND | 26 |
| Dibromochloromethane | ND | 26 |
| 1,1,2-Trichloroethane | ND | 26 |
| Benzene | ND | 26 |
| trans-1,3-Dichloropropene | ND | 26 |
| Bromoform | ND | 26 |
| 4-Methyl-2-Pentanone (MIBK) | ND | 130 |
| 2-Hexanone | ND | 130 |
| Tetrachloroethene | ND | 26 [°] |
| 1,1,2,2-Tetrachloroethane | ND | 26 |
| Toluene | ND | 26 |
| Chlorobenzene | ND | 26 |
| Ethylbenzene | ND | 26 |
| Styrene | ND | 26 |
| Xylene (total) | ND | 26 |
| Methyl-tert-butyl ether (MTBE) | ND | 26 |
| 1,2-Dibromoethane (EDB) | ND | 26 |

Analyzed By: <u>SK</u>

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Approved by <u>Lauren Mandone</u> for Nancy Stewart

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QC REPORT FOR AMRO PROJECT #12257

VOLATILE ORGANICS

| 1,1-Dichloroethene | 99\$ |
|--------------------|------|
| Trichloroethene | 978 |
| Benzene | 102% |
| Toluene | 99% |
| Chlorobenzene | 97% |

ANALYST: SK

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LABORATORY REPORT

EPA Method 8270 Semivolatile Organic Compounds Base/Neutral Extractables Page 1 of 2

| Client: Simmons Environmental Services, Inc. |
|---|
| Client I.D.: <u>9509874 Mass Guard</u> |
| SK UST, East Wall |
| AMRO I.D.: 12257-01 |
| Date sampled: 02/13/96 Date received: 02/14/96 |
| Date prepared: 02/15/96 Date analyzed: 02/15/96 |
| Sample Qty/Type: 1/Solid |

| 1,3-DichlorobenzeneND0.261,4-DichlorobenzeneND0.26Bis(2-Chloroethyl)EtherND0.26HexachloroethaneND0.261,2-DichlorobenzeneND0.26Bis(2-Chloroisopropyl)EtherND0.26Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.262,4-DinitrotolueneND0.26JitrotolueneND0.260 iethylphthalateND0.260 iethylphthalateND0.260 iethylphthalateND0.260 iethylphthalateND0.260 iethylphthalateND0.261 - NitrosodiphenylamineND0.261 - NitrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - No itrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - NitrosodiphenyletherND0.261 - No itrosodiphenylether< | ter | Results (mg/kg) | Reporting Limit(mg/kg) |
|--|---------------|--------------------|---------------------------|
| 1,4-DichlorobenzeneND0.26Bis(2-Chloroethyl)EtherND0.26HexachloroethaneND0.261,2-DichlorobenzeneND0.26Bis(2-Chloroisopropyl)EtherND0.26N-Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.26Dimethyl PhthalateND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26N-NitrosodiphenylamineND0.26 | chlorobenzene | ND | 0.26 |
| Bis(2-Chloroethyl)EtherND0.26HexachloroethaneND0.261,2-DichlorobenzeneND0.26Bis(2-Chloroisopropyl)EtherND0.26N-Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.260.260.260.261.2,2,4-DinitrotolueneND0.261.2,2,4-Trichlorophenyl-PhenyletherND0.261.2,2,4-Trichlorophenyl-PhenyletherND0.261.2,4-Trichlorophenyl-PhenyletherND0.262,6-DinitrotolueneND0.262,6-DinitrotolueneND0.261.4-Chlorophenyl-PhenyletherND0.261.5-DinitrotolueneND0.261.5-DinitrotolueneND0.261.5-DinitrotolueneND0.261.5-DinitrotolueneND0.261.5-DinitrotolueneND0.261.5-DinitrotolueneND0.261.5-DinitrotolueneND0.261.5-Dinitrotoluene <td></td> <td></td> <td></td> | | | |
| HexachloroethaneND0.261,2-DichlorobenzeneND0.26Bis(2-Chloroisopropyl)EtherND0.26N-Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| 1,2-DichlorobenzeneND0.26Bis(2-Chloroisopropyl)EtherND0.26N-Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.26Dimethyl PhthalateND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26DiethylphthalateND0.26DiethylphthalateND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| Bis(2-Chloroisopropyl)EtherND0.26N-Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.26Jimethyl PhthalateND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| N-Nitroso-Di-n-PropylamineND0.26NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26DiethylphthalateND0.26DiethylphthalateND0.26DiethylphthalateND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| NitrobenzeneND0.26HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| HexachlorobutadieneND0.261,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26ND0.260.260.16ND0.260.26ND0.260.26ND0.260.26ND0.260.26ND0.260.26ND0.26 | | | |
| 1,2,4-TrichlorobenzeneND0.26IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| IsophoroneND0.26Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| Bis(2-Chloroethoxy)MethaneND0.26HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26DiethylphthalateND0.260.26ND0.26 | | _ | |
| HexachlorocyclopentadieneND0.262-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | _ | |
| 2-ChloronaphthaleneND0.26Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| Dimethyl PhthalateND0.262,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| 2,6-DinitrotolueneND0.264-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| 4-Chlorophenyl-PhenyletherND0.262,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| 2,4-DinitrotolueneND0.26DiethylphthalateND0.26N-NitrosodiphenylamineND0.26 | | | |
| Diethylphthalate ND 0.26 N-Nitrosodiphenylamine ND 0.26 | | | |
| N-Nitrosodiphenylamine ND 0.26 | | ND | 0.26 |
| | | | 0.26 |
| | | ND | 0.26 |
| Hexachlorobenzene ND 0.26 | | ND | 0.26 |
| Di-n-Butylphthalate ND 0.26 | utylphthalate | ND | 0.26 |
| Butylbenzylphthalate ND 0.26 | | ND | 0.26 |
| Bis(2-Ethylhexyl)Phthalate ND 0.26 | | ND | |
| 3,3'-Dichlorobenzidine ND 0.26 | | ND | 0.26 |
| Di-n-Octyl Phthalate ND 0.26 | | | 0.26 |
| N-Nitrosodimethylamine ND 0.26 | | | |
| Pyrene ND 0.26 | • | | 0.26 |
| Benzo(a)Anthracene ND 0.26 | a)Anthracene | ND | 0.26 |
| Chrysene ND 0.26 | - | - | |
| Benzo(b)Fluoranthene ND 0.26 | | | |
| Benzo(k) Fluoranthene ND 0.26 | • | | |
| Benzo(a) Pyrene ND 0.26 | • | | |

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LABORATORY REPORT

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EPA Method 8270/Solid Sample Semivolatile Organic Compounds Base/Neutral Extractables Page 2 of 2

Client: <u>Simmons Environmental Services, Inc.</u> Client I.D.: <u>SK UST, East Wall</u> AMRO I.D.: <u>12257-01</u>

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-------------------------|--------------------|---------------------------|
| Indeno(1,2,3-c,d)Pyrene | ND | 0.26 |
| Dibenzo(a,h)Anthracene | ND | 0.26 |
| Benzo(g,h,i)Perylene | ND | 0.26 |
| Naphthalene | ND | 0.26 |
| Acenaphthylene | ND | 0.26 |
| Acenaphthene | ND | 0.26 |
| Fluorene | ND | 0.26 |
| Phenanthrene | ND | 0.26 |
| Anthracene | ND | 0.26 |
| Fluoranthene | ND | 0.26 |
| 2-Methylnaphthalene | ND | 0.26 |
| Dibenzofuran | ND | 0.64 |

Solid Content = 96.5%. Results are in dry weight. ND = Not Detected at or above the reporting limit.

Analyzed By: NM

Approved by Lamen Mardone for Nancy Stewart

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LABORATORY REPORT

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EPA Method 8270 Semivolatile Organic Compounds Base/Neutral Extractables Page 1 of 2

| Client: Simmons Environmental Services, Inc. |
|---|
| Client I.D.: 9509874 Mass Guard |
| SK UST, South End Bottom |
| AMRO I.D.: 12257-02 |
| Date sampled: 02/13/96 Date received: 02/14/96 |
| Date prepared: 02/15/96 Date analyzed: 02/15/96 |
| Sample Qty/Type: 1/Solid |

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-----------------------------|--------------------|---------------------------|
| 1,3-Dichlorobenzene | ND | 0.25 |
| 1,4-Dichlorobenzene | ND | 0.25 |
| Bis(2-Chloroethyl)Ether | ND | 0.25 |
| Hexachloroethane | ND | 0.25 |
| 1,2-Dichlorobenzene | ND | 0.25 |
| Bis(2-Chloroisopropyl)Ether | ND | 0.25 |
| N-Nitroso-Di-n-Propylamine | ND | 0.25 |
| Nitrobenzene | ND | 0.25 |
| Hexachlorobutadiene | ND | 0.25 |
| 1,2,4-Trichlorobenzene | ND | 0.25 |
| Isophorone | ND | 0.25 |
| Bis(2-Chloroethoxy)Methane | ND | 0.25 |
| Hexachlorocyclopentadiene | ND | 0.25 |
| 2-Chloronaphthalene | ND | 0.25 |
| Dimethyl Phthalate | ND | 0.25 |
| 2,6-Dinitrotoluene | ND | 0.25 |
| 4-Chlorophenyl-Phenylether | ND | 0.25 |
| 2,4-Dinitrotoluene | ND | 0.25 |
| Diethylphthalate | ND | 0.25 |
| N-Nitrosodiphenylamine | ND | 0.25 |
| 4-Bromophenylphenyl Ether | ND | 0.25 |
| Hexachlorobenzene | ND | 0.25 |
| Di-n-Butylphthalate | ND | 0.25 |
| Butylbenzylphthalate | ND | 0.25 |
| Bis(2-Ethylhexyl)Phthalate | ND | 0.25 |
| 3,3'-Dichlorobenzidine | ND | 0.25 |
| Di-n-Octyl Phthalate | , ND | 0.25 |
| N-Nitrosodimethylamine | ND | 0.25 |
| Pyrene | ND | 0.25 |
| Benzo(a)Anthracene | ND | 0.25 |
| Chrysene | ND | 0.25 |
| Benzo(b)Fluoranthene | ND | 0.25 |
| Benzo(k) Fluoranthene | ND | 0.25 |
| Benzo(a) Pyrene | ND | 0.25 |
| | | |

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LABORATORY REPORT

EPA Method 8270/Solid Sample Semivolatile Organic Compounds Base/Neutral Extractables Page 2 of 2

Client: <u>Simmons Environmental Services, Inc.</u> Client I.D.: <u>SK UST, South End Bottom</u> AMRO I.D.: <u>12257-02</u>

| Test ParameterResults (mg/kg)Reporting Limit(mg/kg)Indeno(1,2,3-c,d)PyreneND0.25Dibenzo(a,h)AnthraceneND0.25Benzo(g,h,i)PeryleneND0.25NaphthaleneND0.25AcenaphthyleneND0.25FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.25FluorantheneND0.25FluorantheneND0.25FluorantheneND0.25JobenzofuranND0.25OibenzofuranND0.25OibenzofuranND0.25OibenzofuranND0.25OibenzofuranND0.63 | | ال الحديث الذي يود عن الذي يود بعن الذي العرب الحديث الحديث الحديث الحديث الحديث الحديث الحديث العرب | | |
|--|--------|--|----|------|
| Dibenzo(a,h)AnthraceneND0.25Benzo(g,h,i)PeryleneND0.25NaphthaleneND0.25AcenaphthyleneND0.25AcenaphtheneND0.25FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.25FluorantheneND0.25ServiceND0.25AnthraceneND0.25ServiceND0.25ServiceND0.25ServiceND0.25ServiceND0.25ServiceND0.25ServiceND0.25ServiceND0.25ServiceND0.25 | | ter | | |
| Benzo (g,h,i) PeryleneND0.25NaphthaleneND0.25AcenaphthyleneND0.25AcenaphtheneND0.25FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.25StructureND0.25OutputND0.25AnthraceneND0.25StructureND0.25StructureND0.25StructureND0.25StructureND0.25StructureND0.25StructureND0.25 | Indeno | (1,2,3-c,d) Pyrene | ND | 0.25 |
| NaphthaleneND0.25AcenaphthyleneND0.25AcenaphtheneND0.25FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.252-MethylnaphthaleneND0.25 | Dibenz | o(a,h)Anthracene | ND | 0.25 |
| AcenaphthyleneND0.25AcenaphtheneND0.25FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.252-MethylnaphthaleneND0.25 | Benzo(| g,h,i)Perylene | ND | 0.25 |
| AcenaphtheneND0.25FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.252-MethylnaphthaleneND0.25 | | | ND | 0.25 |
| FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.252-MethylnaphthaleneND0.25 | Acenap | hthylene | ND | 0.25 |
| FluoreneND0.25PhenanthreneND0.25AnthraceneND0.25FluorantheneND0.252-MethylnaphthaleneND0.25 | Acenap | hthene | ND | 0.25 |
| AnthraceneND0.25FluorantheneND0.252-MethylnaphthaleneND0.25 | | | ND | 0.25 |
| FluorantheneND0.252-MethylnaphthaleneND0.25 | Phenan | threne | ND | 0.25 |
| 2-Methylnaphthalene ND 0.25 | Anthra | cene | ND | 0.25 |
| | Fluora | nthene | ND | 0.25 |
| Dibenzofuran ND 0.63 | 2-Meth | ylnaphthalene | ND | 0.25 |
| | | | ND | 0.63 |

Solid Content = 97.1%. Results are in dry weight. ND = Not Detected at or above the reporting limit.

Analyzed By: <u>NM</u>

y <u>Lauren Mardone</u> for Nancy Stewart Approved by

LABORATORY REPORT

EPA Method 8270 Semivolatile Organic Compounds Base/Neutral Extractables Page 1 of 2

Client: <u>Simmons Environmental Services</u>, Inc. Client I.D.: <u>9509874 Mass Guard</u>

SK UST, North End Bottom

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AMRO I.D.: <u>12257-03</u> Date sampled: <u>02/13/96</u> Date received: <u>02/14/96</u> Date prepared: <u>02/19/96</u> Date analyzed: <u>02/19/96</u> Sample Qty/Type: <u>1/Solid</u>

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-----------------------------|--------------------|---------------------------|
| 1,3-Dichlorobenzene | ND | 0.25 |
| 1,4-Dichlorobenzene | ND | 0.25 |
| Bis(2-Chloroethyl)Ether | ND | 0.25 |
| Hexachloroethane | ND | 0.25 |
| 1,2-Dichlorobenzene | ND | 0.25 |
| Bis(2-Chloroisopropyl)Ether | ND | 0.25 |
| N-Nitroso-Di-n-Propylamine | ND | 0.25 |
| Nitrobenzene | ND | 0.25 |
| Hexachlorobutadiene | ND | 0.25 |
| 1,2,4-Trichlorobenzene | ND | 0.25 |
| Isophorone | ND | 0.25 |
| Bis(2-Chloroethoxy)Methane | ND | 0.25 |
| Hexachlorocyclopentadiene | ND | 0.25 |
| 2-Chloronaphthalene | ND | 0.25 |
| Dimethyl Phthalate | ND | 0.25 |
| 2,6-Dinitrotoluene | ND | 0.25 |
| 4-Chlorophenyl-Phenylether | ND | 0.25 |
| 2,4-Dinitrotoluene | ND | 0.25 |
| Diethylphthalate | ND | 0.25 |
| N-Nitrosodiphenylamine | ND | 0.25 |
| 4-Bromophenylphenyl Ether | ND | 0.25 |
| Hexachlorobenzene | ND | 0.25 |
| Di-n-Butylphthalate | ND | 0.25 |
| Butylbenzylphthalate | ND | 0.25 |
| Bis(2-Ethylhexyl)Phthalate | ND | 0.25 |
| 3,3'-Dichlorobenzidine | ND | 0.25 |
| Di-n-Octyl Phthalate | ND | 0.25 |
| N-Nitrosodimethylamine | ND | 0.25 |
| Pyrene | ND | 0.25 |
| Benzo(a)Anthracene | ND | 0.25 |
| Chrysene | ND | 0.25 |
| Benzo(b)Fluoranthene | ND | 0.25 |
| Benzo(k) Fluoranthene | ND | 0.25 |
| Benzo(a)Pyrene | ND | 0.25 |
| | | |

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LABORATORY REPORT

EPA Method 8270/Solid Sample Semivolatile Organic Compounds Base/Neutral Extractables Page 2 of 2

Client: <u>Simmons Environmental Services</u>, Inc. Client I.D.: <u>SK UST</u>, North End Bottom AMRO I.D.: <u>12257-03</u>

| Test Parameter | Results (mg/kg) | Reporting Limit(mg/kg) |
|-------------------------|--------------------|---------------------------|
| Indeno(1,2,3-c,d)Pyrene | ND | 0.25 |
| Dibenzo(a,h)Anthracene | ND | 0.25 |
| Benzo(g,h,i)Perylene | ND | 0.25 |
| Naphthalene | ND | 0.25 |
| Acenaphthylene | ND | 0.25 |
| Acenaphthene | ND | 0.25 |
| Fluorene | ND | 0.25 |
| Phenanthrene | ND | 0.25 |
| Anthracene | ND | 0.25 |
| Fluoranthene | ND | 0.25 |
| 2-Methylnaphthalene | ND | 0.25 |
| Dibenzofuran | ND | 0.64 |

Solid Content = 96.8. Results are in dry weight. ND = Not Detected at or above the reporting limit.

Analyzed By: <u>NM</u>

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| I T | . I I | I AIP | 2.1 | 196 | | 112. | Kunt | ~ | Yes | No | | V/A | | | | • | | | | | | | |

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The Commonwealth of Massachusetts



Department of Environmental Protection Division of Environmental Analysis

Certifies

Laboratory ID #: M-NH012

Amro Environmental Lab Tl Herrick St. Merrimack, NH 03054

for the Chemical Analysis of Potable and Non-Potable Water

pursuant to 310 CMR 42.00

Laboratory Director: Nancy Stewart Expiration Date:

06/30/96

This certificate supercedes all previous Massachusetts certificates issued to this laboratory. The laboratory is regulated by and shall be responsible for being in compliance with Massachusetts regulations at 310 CMR 42.00.

This certificate is valid only when accompanied by the latest dated Certified Parameter List as issued by the Massachusetts D.E.P.

Certification is no guarantee of the validity of the data. This certification is subject to unannounced laboratory inspections.

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Director, Division of Environmental Analysis

01/01/95 Issued

Soil Analysis

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SQM #34-98 Eldg. 5218 & OMS #22 Camp Edwards Bourne, Massachusetts

Prepared for

Office of the State Quartermaster 50 Maple Street Milford, MA 0175⁻

By

TMC Services, Inc. P O Box 481 Bellingham, MA 02019

March 6, 1998



Post Office Box 451 Beilingham, MA 02019

TEL. (508) 960-3757 FAX (508) 960-4801

March 6, 1998

Mr. David Tessicini The Commonwealth of Massachusetts Military Division Office of the State Quartermaster 50 Mapie Street Milford, MA 01757

> RE: SQM# 45-98 Removal of Two (2) Underground Storage Tanks & Soil Analysis Bldg. 5218 & OMS# 22 Camp Edwards

Dear Mr Tessicini:

The following information and enclosed documents have been prepared to satisfy the Underground Storage Tank (UST) Closure Assessment Report requirement listed in the project specifications.

SCOPE OF WORK

This project entailed the removal of USTs from the following locations:

Building# 5218 Turpentine Road Camp Edwards Tank Size: 500 gallons Contents: 500 gallons No. 2 Fuel Oil

OMS 22 South Outer Road Camp Edwards Tank Size: 12,500 gallons Contents: 200 gallons Diesel Fuel



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NARRATIVE

Building= 5218, Turpentine Road

Tank removal was initiated on February 12, 1998. The tank was discovered approximately two feet below grade at the rear of the building on the south west corner. Tank contents were emptied into drums pending removal for disposal. Drv ice was placed in the emptied tank and allowed to sit. Under observation of the Fire Marshal from Otis Air National Guard, the tank was then extracted from the ground using a Backhoe Loader. Grab samples were then taken from all four walls of the tank grave and one from its base. A head space was performed on each of these sample points using a Thermal Environmental Photo Ionization Detector (PID) Model 580E [results are listed in Sampling Data Sheet]. Two additional samples were collected from each of these points. Fifteen grams of soil was collected for VPH in a 40mL VOC vial with 15 grams of methanol. The second sample was placed in a wide mouth glass jar with Teflon lined screw cap for analysis of EPH [see laboratory data]. Samples were collected using a stainless steel spoon from a point approximately 12" from the surface of the excavation. The sample spoon was cleaned between sample points.

OMS 22. South Outer Road

The tank was removed on February 12, 1998. The tank was located on the south side on the building. Two observation wells were located at the ends of the tank. Both wells were opened and examined. The wells did not contain ground water. The contents of the tank was emptied into drums. Dry ice was then added to the tank. The concrete apron above the tank was broken up and removed along with the fuel dispensing unit. The tank was then exposed using a track excavator. The emptied tank was then removed from the excavation under observation of the Fire Marshal from Otis Air National Guard. Grab samples were then taken from all four walls of the tank grave and one from its base. A head space was performed on each of these sample points using a Thermal Environmental Photo Ionization Detector (PID) Model 580E [results are listed in Sampling Data Sheet]. Two additional samples were collected from each of these points. Fifteen grams of soil was collected for VPH in a 40mL VOC vial with 15 grams of methanol. The second sample was placed in



a wide mouth glass jar with Terlon lined screw cap for analysis of EPH [see laboratory data]. Samples were collected using a stainless steel spoon from a point approximately 12" from the surface of the excavation. The sample spoon was cleaned between sample points.

The emptied and clean tanks were removed and disposed of in accordance with Massachusetts requirements at a State certified tank disposal yard [see tank disposal permit and cards]. The removed oil and diesel from the two tanks was transported to United Oil Recovery Inc. Meriden. Connecticut, on February 19, 1998 [see copies of Manifests]

SUMMARY AND CONCLUSIONS

Building# 5218, Turpentine Road

PID results from the base of the tank indicated residual impact of petroleum product. These results were not duplicated in the EPH/VPH analysis of the soils taken from the same point [Sample CE05]. There were no other readings of petroleum contaminants from this tank grave.

OMS 22, South Outer Road

There were no PID readings or indication of petroleum contamination during the removal of the tank at OMS \ddagger 22. Petroleum contamination was not detected from samples taken from the tank grave at this site.

Should you have any further questions, please contact me at 508-966-3737.

Sincere TMC Services, Inc.

Matthew Clark President

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SAMPLING DATA SHEET

Project Name: CAMP EDWARDS Site Name: OM S 22 Address: SOUTH OUTER ROAD NA BUSRNE

Date: 2/12/92

Weather Conditions:

Clasoy RAIN - 40'S HIGH WINO,

| Sample No | Grab or Composite | PID Results | Comments |
|-----------|-------------------|-------------|----------------------------|
| 06 | GRAG | 0.0 | 8' BELOW GRADE |
| 07 | SLAN | 0.0 | 8' BEWSGAMPE |
| 03 | GRAN | 0.0 | 9' BELL GRADE |
| 09 | Shan | 0.0 | 3' BELOW GRAME |
| 10 | 5.2000 | 0.0 | 13' BELON GRADE TONK. BASE |
| | | | |
| | | | |
| | | | |

13' BEWE GRADE SANNIT WATER WAS DETECTED AT ~ ¥

8 09 Ċ Ø 06 Ю 3 07 TANK GRAJE

Sketch stock pile area and sample points. Note any outstanding features or information.

Sampler's Certification: I hear by certify that the above information is true and accurate.

2-12-93 N. A.2, 2, 7, A al Dire Innai Company Date Name smpdoc/rac

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

PROJECT NARRATIVE

CLIENT: TMC Services CLIENT PROJECT ID: Camp Edwards ESS PROJECT ID: 980408

Sample Receipt

One liquid and ten solid samples were received on February 13, 1998 for the analyses specified on the enclosed Chain of Custody Record.

Analytical Summary

The project as described above has been analyzed in accordance with the ESS Quality Assurance Plan. This plan utilizes the following methodologies: US EPA SW-846, US EPA Methods for Chemical Analysis of Water and Wastes per 40 CFR Part 136, APHA Standard Methods for the Examination of Water and Wastewater, American Society for Testing and Materials (ASTM), and other recognized methodologies. These analyses with these noted observations are in conformance to the Quality Assurance Plan.

Volatile Organics Analysis

Surrogate recoveries were outside of the recommended ranges for samples 980408-02. -05. -07. -08 and -10 due to matrix interferences.

No other observations noted.

This signed Certificate of Analysis is our approved release of your analytical results. Beginning with this Project Narrative, the entire report has been paginated. The Chain of Custody is the final report page. This report should not be copied except in full without the approval of the laboratory.

End of project narrative.

Laurel Stodeard/Eric Baanante Laboratory Manager/Operations Manager

<u>2/26/98</u> Date

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

EXTRACTABLE PETROLEUM HYDROCARBON (EPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE06

Date Sampled: 2/13/98

Date Analyzed: 2/23/98

ESS Project ID: 980408 ESS Sample ID: 980408-06 Date Extracted: 2/18/98 Dilution Factor: 1x Analyst: JAR

EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

| Parameter | Results | MDL | Unis |
|---|---------|-----|-------|
| n-Co-n-C ₁₈ Aliphatics | ND | 29 | mg/Kg |
| n-C ₁₉ -n-C ₃₆ Aliphatics | ND | 29 | mg/Kg |
| C ₁₀ -C ₂₂ Aromatics* | ND | 29 | mg/Kg |

* Excludes Target PAHs

ND = Not Detected above Method Detection Limit (MDL)

SURROGATE RECOVERIES

| Surrogate | / % Reco | overy Advisory Range |
|-------------------------|----------|----------------------|
| Chloro-octadecane (COD) | 89 | 40 - 140% |
| Ortho-terphenyl (OTP) | 97 | 40 - 140% |

Results reported on a dry weight basis.

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

VOLATILE PETROLEUM HYDROCARBON (VPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE06 Date Sampled: 2/12/98 ESS Project ID: 980408 ESS Sample ID: 980408-06 Date Analyzed: 2/19/1998 Analyst: skh

VPH RESULTS

| Facaneter | Results | MINIMUL | Cnits-mark |
|-------------------------|---------|---------|----------------|
| n-C5-n-C, Aliphanics* | ND | 0.8 | mg/kg dry wgt. |
| n-C9-n-C12 Aliphatics** | ND | 0.8 | mg/kg dry wgt. |
| n-C9-n-C10 Aromatics | ND | 0.8 | mg/kg dry wgt |

* = Excludes Benzene, Toluene and MTBE.

** = Excludes Ethylbenzene, Xylenes (total), 1,2,4-Trimethylbenzene and Naphthalene.

Surrogate Recoveries

| Surregate | ABenner | Advimer Hanres |
|--------------------------|---------|----------------|
| FID 4-Bromoiluorobenzene | 101 | 80-120% |
| PID 1-Bromorluoropenzene | 109 | 80-120% |

TARGETED VPH ANALYTES

| Analyte | Results | MDL | Çnifa |
|------------------------|---------|-----|----------------|
| Methyl-ten-butylether | ND | 150 | ug/kg dry wgt |
| Benzene | ND | 150 | ug/kg dry wgt. |
| Toluene | ND | 150 | ug/kg dry wgt |
| Ethyibenzene | ND | 150 | ug/kg dry wgr. |
| M &P Xviene | ND | 150 | µg/kg dry wgt |
| O Xylene | ND | 150 | ug/kg dry wgt |
| Naphthalene | ND | 150 | ug/kg dry wgt |
| 1.2.4-Trimethvibenzene | ND | 150 | μg/kg drv wgt. |

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CERTIFICATE OF ANALYSIS

EXTRACTABLE PETROLEUM HYDROCARBON (EPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE07 Date Sampled: 2/13/98 Date Analyzed: 2/23/98 ESS Project ID: 980408 ESS Sampie ID: 980408-07 Date Extracted: 2/18/98 Dilution Factor: 1x Analyst: JAR

EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

| Parameter | Resuits | MDL | Units |
|---|---------|-----|-------|
| n-Co-n-C ₁₈ Aliphatics | ND | 26 | mg/Kg |
| n-C ₁₉ -n-C ₃₆ Aliphatics | ND | 26 | mg/Kg |
| C ₁₀ -C ₂₂ Aromatics [*] | ND | 26 | mg/Kg |

* Excludes Target PAHs

ND = Not Detected above Method Detection Limit (MDL)

SURROGATE RECOVERIES

| Surrogate | % Recovery | Advisory Range |
|-------------------------|------------|----------------|
| Chloro-octadecane (COD) | 84 | 40 - 140% |
| Ortho-terphenyl (OTP) | 96 | 40 - 140% |

Results reported on a dry weight basis.

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

VOLATILE PETROLEUM HYDROCARBON (VPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE07 Date Sampled: 2/12/98 ESS Project ID: 980408 ESS Sample ID: 980408-07 Date Analyzed: 2/19/1998 Analyst: skh

VPHRESULTS

| Parameter | Results | | Ualis |
|---|---------|-----|----------------|
| n-C ₅ -n-C ₅ Aliphanes* | ND | 0.9 | mg/kg dry wgt |
| n-C9-n-C12 Aliphatics** | ND | 0.9 | marka dry wat |
| n-C ₉ -n-C ₁₀ Aromatics | ND | 0.9 | mg/kg dry wgt. |

* = Excludes Benzene, Toluene and MTBE.

** = Excludes Ethylbenzene, Xylenes (total), 1,2,4-Trimethylbenzene and Naphthalene.

| Surrogate Recoveries | | |
|--------------------------|--------|---------------|
| Surrogne | Repver | Advenry Range |
| FID 4-Bromorluorobenzene | 103 | 80-120% |
| PID 4-Bromoiluorobenzene | 110 | 80-120% |

TARGETED VPH ANALYTES

| Analyte | Results | MDL | Lois |
|------------------------|---------|-----|----------------|
| Methyl-tert-burylether | ND | 180 | μg/kg άry wgt |
| Benzene | ND | 180 | µg/kg dry wgL |
| Toluene | ND | 180 | µg/kg dry wgt |
| Ethyibenzene | ND | 180 | µg/kg dry wgL |
| M &P Xyleze | ND | 180 | µg/kg dry wgt. |
| O Xviene | ND | 180 | µg/kg dry wgt |
| Naphthalene | ND | 180 | µg/kg dry wgt |
| 1,2,4-Trimethylbenzene | ND | 180 | μg/kg dry wgt. |

Tei. (401) 461-7181 Fax: 401) 461-4486

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Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

EXTRACTABLE PETROLEUM HYDROCARBON (EPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE08 Date Sampled: 2/13/98 Date Analyzed: 2/23/98 ESS Project ID: 980408 ESS Sample ID: 980408-08 Date Extracted: 2/18/98 Dilution Factor: 1x Analyst: JAR

EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

| Parameter | Results | MDL | i Umis |
|---|---------|-----|--------|
| n-C ₉ -n-C ₁₈ Aliphatics | ND | 26 | mg/Kg |
| n-C ₁₉ -n-C ₃₆ Aliphatics | ND | 26 | mg/Kg |
| C ₁₀ -C ₂₂ Aromatics* | ND | 26 | mg/Kg |

* Excludes Target PAHs

ND = Not Detected above Method Detection Limit (MDL)

SURROGATE RECOVERIES

| Surrogate | % Recovery | Advisory Range |
|-------------------------|------------|----------------|
| Chloro-octadecane (COD) | 78 | 40 - 140% |
| Ortho-terphenyl (OTP) | 94 | 40 - 140% |

Results reported on a dry weight basis.

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

VOLATILE PETROLEUM HYDROCARBON (VPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE08 Date Sampled: 2/12/98 ESS Project ID: 980408 ESS Sample ID: 980408-08 Date Analyzed: 2/19/1998 Analyst: skh

VPH RESULTS

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| Perezete | Results | ADE. | Enits |
|-------------------------|---------|------|----------------|
| n-C5-n-C5 Aliphancs* | ND | 0.5 | mg/kg dry wgt_ |
| n-Cg-n-C12 Aliphanics** | ND | 0.5 | mg/kg dry wgt_ |
| n-C9-n-C10 Aromatics | ND | 0.5 | mg/kg dry wgt. |

* = Excludes Benzene, Toluene and MTBE.

** = Excludes Ethylbenzene, Xylenes (total), 1,2,4-Trimethylbenzene and Naphthalene.

Surrogate Recoveries X Recovery Advaory Range FID 4-Bromoiluorobenzene 102 80-120% PID 4-Bromoiluorobenzene 109 80-120%

TARGETED VPH ANALYTES

| Anaiyte | Re-mite | and a second | Daita |
|------------------------|---------|--------------|----------------|
| Methyl-tert-burylether | ND | 104 | μg/kg drv wgt |
| Benzene | ND | 104 | µg/kg drv wgt |
| Toluene | ND | 104 | µg/kg dry wgL |
| Ethyibenzene | ND | 104 | µg/kg drv wgL |
| M &P Xviene | ND | 104 | µg/kg dry wgt. |
| O Xviene | ND | 104 | µg/kg dry wgt. |
| Naphthalene | ND | 104 | µg/kg dry wgt. |
| 1.2,4-Trimethylbenzene | ND | 104 | µg/kg dry wgt |

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE09 Date Sampled: 2/13/98 Date Analyzed: 2/23/98 ESS Project ID: 980408 ESS Sample ID: 980408-09 Date Extracted: 2/18/98 Dilution Factor: 1x Analyst: JAR

EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

| Parameter | Results | MDL | Units |
|---|---------|-----|-------|
| n-Co-n-C ₁₈ Aliphatics | ND | 26 | mg/Kg |
| n-C ₁₉ -n-C ₃₆ Aliphatics | ND | 26 | mg/Kg |
| C ₁₀ -C ₁₂ Aromatics* | ND | 26 | mg/Kg |

* Excludes Target PAHs

ND = Not Detected above Method Detection Limit (MDL)

SURROGATE RECOVERIES

| Surrogate de la serie de la | % Recovery | Advisory Range |
|---|------------|----------------|
| Chloro-octadecane (COD) | 87 | 40 - 140% |
| Ortho-terphenyl (OTP) | 103 | 40 - 140% |

Results reported on a dry weight basis.

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

VOLATILE PETROLEUM HYDROCARBON (VPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE09 Date Sampled: 2/12/98 ESS Project ID: 980408 ESS Sample ID: 980408-09 Date Analyzed: 2/20/1998 Analyst: skh

VPH RESULTS

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| Paremeter | Results | MILL | Unifs |
|--|---------|------|----------------|
| n-C-n-C, Aliphancs* | ND | 0.6 | mg/kg dry wgL |
| n-C ₉ -n-C ₁₂ Aliphanics** | ND | 0.6 | mg/kg dry wgt. |
| n-C ₉ -n-C ₁₀ Aromatics | ND | 0.6 | mg/kg dry wgt. |

* = Excludes Benzene, Toluene and MTBE.

** = Excludes Ethylbenzene, Xylenes (total), 1,2,4-Trimethylbenzene and Naphthalene.

| Surrogate Recoveries | | |
|--------------------------|-------------|--------------|
| Surrogate | A Recorders | ACTION RADES |
| FID 1-Bromorluorobenzene | 102 | 80-120% |
| PD 4-Bromorluorobenzene | 1 | 80-120% |

TARGETED VPH ANALYTES

| Analyte | Results | MUL | Units |
|------------------------|---------|-----|----------------|
| Methvi-tert-burviether | ND | 115 | µg/kg dry wgt |
| Benzene | ND | 115 | µg/kg dry wgt |
| Toluene | ND | 115 | µg/kg dry wgt. |
| Ethylbenzene | ND | 115 | µg/kg drv wgi |
| M &P Xviene | ND | 115 | ug/kg drv wgL |
| O Xylene | ND | 115 | µg/kg dry wgt |
| Naphthaiene | ND | 115 | µg/kg dry wgt. |
| 1.2.4-Trimethylbenzene | ND | 115 | µg/kg dry wgt |

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Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE10 Date Sampled: 2/13/98 Date Analyzed: 2/23/98 ESS Project ID: 980408 ESS Sample ID: 980408-10 Date Extracted: 2/18/98 Dilution Factor: 1x Analyst: JAR

EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

| Parameter | Results | MDL | Units |
|---|---------|-----|-------|
| n-Co-n-C ₁₈ Aliphatics | ND | 27 | mg/Kg |
| n-C ₁₉ -n-C ₃₆ Aliphatics | ND | 27 | mg/Kg |
| C ₁₀ -C ₂₂ Aromatics* | ND | 27 | mg/Kg |

* Excludes Target PAHs

ND = Not Detected above Method Detection Limit (MDL)

SURROGATE RECOVERIES

| Surrogate | % Recovery | Advisory Range |
|-------------------------|------------|----------------|
| Chloro-octadecane (COD) | 85 | 40 - 140% |
| Ortho-terpinenyl (OTP) | 101 | 40 - 140% |

Results reported on a dry weight basis.

Division of Thielsch Engineering, Inc.

CERTIFICATE OF ANALYSIS

VOLATILE PETROLEUM HYDROCARBON (VPH) ANALYSIS

Client: TMC Services Client Project ID: Camp Edwards Client Sample ID: CE10 Date Sampled: 2/12/98 ESS Project ID: 980408 ESS Sample ID: 980408-10 Date Analyzed: 2/20/1998 Analyst: skh

VPH RESULTS

| Parazetes | Resute | MDL | Units 🔷 |
|--|--------|-----|----------------|
| n-C ₅ -n-C ₅ Aliphatics* | ND | 0.6 | mg/kg dry wgt. |
| n-C9-n-C12 Aliphanics** | ND | 0.6 | mg/kg dry wgt. |
| n-C ₉ -n-C ₁₀ Aromatics | ND | 0.6 | mg/kg dry wgt. |

* = Excludes Benzene, Toluene and MTBE.

** = Excludes Ethylbenzene, Xylenes (total), 1,2,4-Trimethylbenzene and Naphthalene.

Surrogate Recoveries

| Surrogate | C. Beinger anna | Addram's Reader |
|--------------------------|-----------------|-----------------|
| FID 4-Bromofluorobenzene | 103 | 80-120% |
| PID 4-Bromoiluorobenzene | 112 | 80-120% |

TARGETED VPH ANALYTES

| Analyte | Resuits | MDL | Units |
|------------------------|---------|-----|----------------|
| Methyl-tert-burylether | ND | 125 | µg/kg dry wgL |
| Benzene | ND | 125 | µg/kg dry wgt. |
| Toluene | ND | 125 | µg/kg dry wgL |
| Ethylbenzene | ND | 125 | µg/kg dry wgt. |
| M &P Xylene | ND | 125 | µg/kg dry wgt. |
| O Xviene | ND | 125 | µg/kg dry wgt. |
| Naphthalene | ND | 125 | µg/kg dry wgL |
| 1.2.4-Trimethylbenzene | ND | 125 | ug/kg dry wgL |

1 £..... CHAIN OF CUSTODY **ESS Laboratory** Page_4_of _/___

Division of Thielsch Engineering, Inc.

| | 185 Frances Avenue, Cranston, RF 02910/2211 Tel. (401) 461-7181 - Fax (401) 461-4486 | | | | | Turn Time LStandard (2 Weel | ks) | Oth | er | | | | ESS-LAI | S PROJ | | \overline{n} | ; | |
|---------------------|---|--------------------|-------------------------------------|--------------|---|--------------------------------|----------------------|-------------------|---------|----------|-------|----------|------------------|--------|----|----------------|-------|---------|
| IVAC SEAVILLES | | | Project # Project Name CAMP FATARAS | | | | | | | | | Analy | nalysis Required | | | | | |
| Contact Person | Annie Ch | IARAQIA | Addres | ·· [. | D. Lana 4/31 | | ŗ | | T | Π | Τ | | Τ | Π | 7 | T | T | \prod |
| Liny PSELLI | | State MA | | 2 | $\mathcal{I}_{4}^{\text{Zip}}$ $\mathcal{I}_{6}^{\text{Lin}}$ | 507 · 3737 | Number of Containers | tainer | | | | | | | | | | |
| Purchase Order # | · | | - | | 1×# 502 966-4361 | | ber of (| Type of Container | IS . | 2 | | | | | | | | |
| ESS LAB Sample # | Date | Collection Time | COMP | MATRIX | Sample Identif | fication | Run N | Type | | 5 | | | | | | | | |
| Õ(| 1),1/48 | 1330 | Ý | 15 | CEOI | | 2 | 9 | X | <u> </u> | | | | | | | | |
| CIZ | 413 | 1330 | <u> </u> | 5 | CEUL | | 2 | 9 | | | | | | | | | | |
| 63 | 2/13 | 15.50 | | 15 | CEUS | | 2 | <u>G</u> | | | | | | | | | | _ |
| 04 | 2/13 | 1330 | <u> </u> | ~ | CEU4 | | \mathcal{L} | <u>G</u> | | | | | | | | | | |
| -705 | 1/13 | 1330 | X | | CEUS | | 2 | <u>G</u> | []_ | <u> </u> | | | | | | | | |
| 06 | 2/13 | 1530 | <u> </u> | | CEO6 | | 2 | <u>G</u> | _ | <u> </u> | | | | | | | | _ |
| 07 | 2/13 | /3.30 | ┠───╂╼╾ | <u>< </u> | CEOP | | 2 | <u>G</u> | _ _ | <u> </u> | | | | | | | | |
| -708 | 2/13 | 1330 | 1 | 45 | CEUR | | 2 | 9 | | | | | | | | | | |
| 09 | 2/13 | 1530 | t | < <u>5</u> | CEOQ | | 2 | G | | | | | | | | | | |
| 10 | 2/13 | 1330 | | x .5 | | | 2 | 4 | | | | | | | | | | |
| | 2/13 | 13 30 | | _ | TRIP Butnle | CEII | 1 | 9 | | <u> </u> | | | $\left \right $ | | | | | |
| | | | | | | | | | | | | | | | | | | |
| ontainer Type: | P-Puly | G-Glass | S | Sterile | : V-VOA Matri | ix: DW-Drinking | g Water | | S-Suli. | l | G₩ | Ground | Water | | WW | Waste | waler | |
| cals Intact: | Yes | No Com | inciits: | | | | | | | | | | | | | | | |
| ooler Temp: | | | | | | | | | | | | | | | | | | |
| (| ignature | Date/Time | 1 1 1 | | (Signature) Date/ 14 Jeaven 12/151 | | Signatu | re) | ſ | Date/Ti | me Re | ceived l | y: (Sign | ature) | | | Dat | e/Time |
| Refinquished by: (S | ignature) | Date/Time | Receiv | ed by: | (Signature) Date/ | Time Relinquished by: (! | Signatu | re) | |)ate/Ti | me R | ccived | oy: (Sign | ature) | | | Dit | e/Time |

| APPLICATI | |
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Hazardous Waste MANIFEST PROGRAM 79 Elm St., Hartford, CT 06106-5127

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APPENDIX D DRAINAGE STUCTURE 04CDXX1 CLOSURE REPDRT

CLOSURE REPORT

1. INTRODUCTION

1.1 DRAINAGE STRUCTURE REMOVAL PROGRAM

Massachusetts Military Reservation (MMR) is listed on the U.S. Environmental Protection Agency (USEPA) National Priority List. Remedial studies and activities are currently being conducted at MMR in accordance with guidelines and procedures of the USEPA Superfund program and the National Contingency Plan (NCP). Two types of action can be initiated under the NCP: remedial actions and removal actions. Remedial actions are long-term, permanent remediation of hazardous waste sites. Removal actions are short-term actions to abate or eliminate the release of contamination. The drainage structure removal program is a removal action program.

The National Guard Bureau (NGB) evaluated the drainage structures in a separate basewide program because of the similarity in operations and contaminants. The drainage structure removal program is based on the Phase I Sump Investigation Program conducted by ABB Environmental Services Inc. (ABB) and the Phase II Sump Investigation Program conducted by Metcalf and Eddy for the NGB. ABB characterized approximately 100 drainage structures. This characterization included sampling to identify potential contaminant sources. Also during this phase an additional '161 drainage structures were identified. Fifteen of these structures were characterized by Metcalf and Eddy as a high priority.

- A total of 185 drainage structures are included in the program. One hundred sixty-five
 of these structures are funded for removal. Currently, Jacobs Engineering Group is
 under contract through the Air Force Center for Environmental Excellence for the
 following work under the drainage structure removal program. The scope of work
 includes: The removal or abandonment in place of approximately 165 existing
 drainage structures on the Massachusetts Military Reservation and the associated
 liquid and sediment contents.
- Demolition of 12 existing vehicle maintenance slabs and foundations.
- Testing soil, liquid, sludge and sediment for contamination.
- Stockpiling and treatment or disposal of excavated drainage structure debris and surrounding soil.
- Backfilling and restoring excavations as specified.
- Preparing closure reports.

The work Jacobs is performing on the drainage structure removal program is governed by the "Final Drainage Structure Removal Program Specifications", July 1995, prepared for Massachusetts Military Reservation by Metcalf and Eddy; and the approved Jacobs' Plans (Drainage Structure Removal Program Volume I - Quality Project Plan and Volume II - Remedial Action Work Plan, January 1996, prepared for Massachusetts Military Reservation). These documents detail the procedures utilized by Jacobs during the removal of drainage structures.

1.2 CLEAN-UP LEVELS

The clean-up levels for the drainage structure removal program (DSRP) are determined by the MMR specific soil target clean-up levels (STCLs) as developed by HAZWRAP and presented in their letter, "Soil Target Clean Up Levels," dated January 30, 1996. The STCLs are risk based standards. The depth utilized for the DSRP is the 2 to 15 foot depth. Additionally, the STCLs are divided into several sub-categories: inside the flightline, outside the flightline, and clean fill. The inside and outside the flightline STCLs are based upon future use as well as risk. These standards apply to the bottom, inlet, and outlet samples and are used to determine that the drainage structure removal is a clean closure.

The clean fill STCLs apply to the paylimit sample. This data determines if the paylimit can be used for clean fill or requires treatment. The soil will be treated in the OABF if any of the contaminants tested are above the STCLs for clean fill. Additionally, if the soil fails the "twenty times" rule as defined in specification 01410, a paylimit sample will be analyzed for TCLP. If the soil passes the TCLP, it will be treated in the OABF. Any soil failing TCLP limits will be disposed off-site at a RCRA-regulated facility.

1.3 DRAINAGE STRUCTURE DESCRIPTION AND LOCATION

Drainage structure 04CDXX1 was located outside the flightline to the west of Building OMS-22, which is located in the Army National Guard Motor Pool area off of South Truck Rd. The drainage structure was a Cesspool (See Figure 1). It was not visible at the ground surface and had a depth of 10 feet. The survey coordinates are 237,242.3 North and 862,799.2 East.

The drainage structure had no cover. The structure walls were constructed of concrete barrel block with an open dirt bottom. The walls had collapsed. An 4-inch ceramic inlet pipe was observed with this structure. There was no outlet pipe present.

1.4 PREVIOUS INVESTIGATION HISTORY

1.4.1 BUILDING OMS-22 (Cesspool)

The area of Building OMS-22 is located in the Army National Guard Motor Pool. This structure was not investigated at the time of the Phase I or Phase II Investigations.

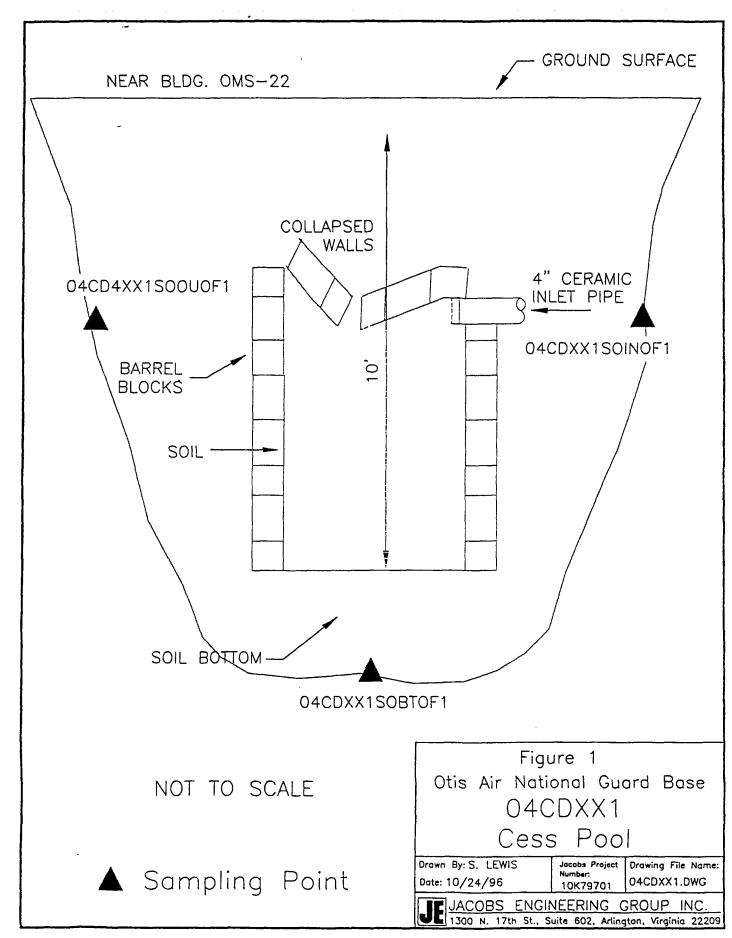
2. LIQUID REMOVAL AND DISPOSAL

No liquid was removed from this structure.

3. SEDIMENT/SLUDGE REMOVAL AND DISPOSAL

No sediment/sludge was removed from this structure. The structure was filled with soil which was removed with paylimit.

<u>Otis ANGB</u> Drainage Structure Removal Program



4. DRAINAGE STRUCTURE REMOVAL AND DISPOSAL

4.1 PROCEDURES USED TO REMOVE PAYLIMIT

4.1.1 EXCAVATION

On April 11-12, 1996, a backhoe was used to excavate and remove the drainage structure (Cesspool) and paylimit soils. The soil was removed and directly loaded onto dump trucks. Before excavation began, it was noticed that this structure was partially collapsed. Approximately 40 cubic yards of soil and debris were excavated. Final excavation size was approximately 21 feet long by 15 feet wide by 13 feet deep. Discoloration was present throughout the excavation as well as a strong fuel odor. These areas of discoloration were removed with the paylimit. Approximately 12 linear feet of 4-inch ceramic pipe were removed from the inlet during the excavation. Pipe remaining in the ground was capped with concrete. No outlet pipe was present.

Excavated soil was stored by structure in the CSA. Structure closure was observed by ABB Environmental, the Title II subcontractor to IRP. Photos are located in the photo log of the master copy of this document in the IRP office.

4.1.2 FIELD SCREENING

Field screening for volatile organic compounds and dust was conducted throughout the excavation activities. PID results were non-detect in the breathing zone. Mini-RAM results were also non-detect in the breathing zone. Headspace PID readings were also taken at the inlet, outlet, and bottom. The headspace readings were used as an initial determination of contamination. The headspace readings were 1.0 ppm for the bottom, 2.2 ppm for the inlet, and 1.2 ppm for the outlet sample locations.

4.1.3 BACKFILL

The backfill used for this excavation came from an on-base source, either the former Landfill 1 common borrow pit located on Turpentine Road or paylimit soils from drainage structures with analytical results below the clean fill STCLs. Approximately 60 cubic yards of soil were used to backfill the excavation.

4.1.4 FINAL RESTORATION

A gravel parking area was impacted by the excavation and the area was regraded in the spring.

4.2 ANALYTICAL RESULTS

The on-site and off-site laboratories analyzed the samples according to the Quality Assurance Plan and the laboratory specification (Section 01410 of the Specifications). All data was validated in accordance with HAZWRAP and DOE guidelines. The samples were analyzed for the volatile compounds utilizing EPA method 8010/8020 and EDB, the semivolatile compounds utilized EPA method 8270, pesticides and PCBs by EPA method 8080, metals by EPA method series 6000 and 7000, and TPH by EPA method 8015 modified as diesel and gasoline. The off-site laboratory has established analyte detection levels which are below the method required quantitation levels.

Otis ANGB Drainage Structure Removal Program The concentrations reported for those analytes detected below the laboratory quantitation limits are estimated and are noted as such in the following tables.

A sample was collected from the bottom, inlet, and outlet sample locations. The outlet sample was collected from the sidewall opposite the inlet sample because no outlet pipe was present. These samples were analyzed in the on-site laboratory as part of our field screening program. When the on-site laboratory results demonstrated the excavation was clean, samples were sent to the off-site laboratory for confirmation. The off-site analytical detection results for the bottom, inlet, and outlet sample locations are summarized in Tables 4-1 through 4-3, respectively. The complete on-site and off-site analytical data are in Appendix A and B, respectively. None of the compounds detected in the bottom, inlet, and outlet samples exceeded the STCLs; therefore, structure 04CDXX1 is a clean closure.

| Analyte | Result | Estimated | Exceeds Limits | STCL Outside ¹ |
|-----------------------------|------------|--------------|----------------|---------------------------|
| | <u> </u> | Quantitation | | l |
| TPH as Diesel | 61 mg/kg | | | 500 mg/kg |
| 1,4-Dichlorobenzene | 3.5 ug/kg | | | 26700 ug/kg |
| Aluminum | 2320 mg/kg | | | 54900 mg/kg |
| Arsenic | 1.7 mg/kg | | | 3.6 mg/kg |
| Barium | 11.9 mg/kg | · · | · · | 3800 mg/kg |
| Beryllium | 0.14 mg/kg | | | 1 mg/kg |
| Calcium | 354 mg/kg | | | |
| Chromium | 6.5 mg/kg | | | 274 mg/kg |
| Cobalt | 1.5 mg/kg | | | |
| Copper | 20.0 mg/kg | | | |
| Iron | 3370 mg/kg | | | |
| Lead | 19.4 mg/kg | |) | 300 mg/kg |
| Magnesium | 619 mg/kg | | | |
| Manganese | 30.4 mg/kg | | | 274 mg/kg |
| Nickel | 1.7 mg/kg | | | 1100 mg/kg |
| Potassium | 424 mg/kg | | | |
| Sodium | 80.4 mg/kg | | | |
| Vanadium | 6.5 mg/kg | | | 384 mg/kg |
| Zinc | 8.7 mg/kg | | | 10000 mg/kg |
| bis(2-Ethylhexyl) phthalate | 100 ug/kg | YES | { | 45700 ug/kg |
| Butyl benzyl phthalate | 41 ug/kg | YES | | 11000000 ug/kg |
| Fluoranthene | 46 ug/kg | YES | | 2200000 ug/kg |
| Phenanthrene | 48 ug/kg | YES | | 2200000 ug/kg |
| Pyrene | 50 ug/kg | YES | | 1650000 ug/kg |

Table 4-1 - Bottom Sample 04CDXX1SOBTOF1

¹ STCL is used when the drainage structure was located outside the flightline.

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| Analyte | Result | Estimated Quantitation | Exceeds Limits | STCL Outside ¹ |
|-----------------------------|------------|---------------------------|----------------|---------------------------|
| TPH as Diesel | 24 mg/kg | | | 500 mg/kg |
| Aluminum | 2420 mg/kg | | | 54900 mg/kg |
| Arsenic | 1.9 mg/kg | | | 3.6 mg/kg |
| Barium | 7.7 mg/kg | | | 3800 mg/kg |
| Beryllium | 0.15 mg/kg | | | 1 mg/kg |
| Calcium | 285 mg/kg | | | |
| Chromium | 4.1 mg/kg | | 1 | 274 mg/kg |
| Cobait | 1.2 mg/kg | | | _ |
| Copper | 8.8 mg/kg | ļ | | |
| Iron | 4080 mg/kg | ŀ | | |
| Lead | 6.9 mg/kg | | | 300 mg/kg |
| Magnesium | 506 mg/kg | | | |
| Manganese | 32.8 mg/kg | | | 274 mg/kg |
| Nickel | 1.6 mg/kg | | | 1100 mg/kg |
| Potassium | 334 mg/kg | | | |
| Sodium • | 85.2 mg/kg | | | |
| Vanadium | 7.8 mg/kg | | | 384 mg/kg |
| Zinc | 8.9 mg/kg | | | 10000 mg/kg |
| bis(2-Ethylhexyl) phthalate | 40 ug/kg | YES | | 45700 ug/kg |
| Butyl benzyl phthalate | 43 ug/kg | YES | | 11000000 ug/kg |

Table 4-2 - Inlet Sample 04CDXX1SOINOF1

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¹ STCL is used when the drainage structure was located outside the flightline.

Table 4-3 - Outlet Sample 04CDXX1SOOUOF1

| Analyte | Result | Estimated Quantitation | Exceeds Limits | STCL Outside |
|---------------|------------|---------------------------|----------------|--------------|
| TPH as Diesel | 55 mg/kg | | | 500 mg/kg |
| Aluminum | 3010 mg/kg | | | 54900 mg/kg |
| Arsenic | 2.5 mg/kg |] | | 3.6 mg/kg |
| Barium | 7.4 mg/kg | | | 3800 mg/kg |
| Beryllium | 0.16 mg/kg | | | 1 mg/kg |
| Calcium | 329 mg/kg | | | |
| Chromium | 3.4 mg/kg | | | 274 mg/kg |
| Cobalt | 1.1 mg/kg | | | |
| Copper | 17.4 mg/kg | | | |
| Iron | 4930 mg/kg | | | |
| Lead | 7.7 mg/kg | | | 300 mg/kg |
| Magnesium | 549 mg/kg | | | |
| Manganese | 32.8 mg/kg | | | 274 mg/kg |
| Potassium | 416 mg/kg | | | |
| Sodium | 81.7 mg/kg | | | |
| Vanadium | 8.4 mg/kg | | | 384 mg/kg |
| Zinc | 8.9 mg/kg | | | 10000 mg/kg |

¹STCL is used when the drainage structure was located outside the flightline.

<u>Otis ANGB</u> Drainage Structure Removal Program

4.3 ANALYTICAL QA

The analytical data were validated in accordance with DOE/HWP - 65/R1: "Requirements for Quality Control of Analytical Data." The validation was a Level C DQO validation and found the data to be usable. Samples were collected from the sidewalls and bottom of the excavation for analysis by an on-site laboratory. This laboratory used gas chromatography and immunoassay techniques for the organic analyses and x-ray fluorescence (XRF) for inorganic analysis. There is good agreement between the on-site and off-site analytical data, with the exception of barium results. The barium concentrations reported by the on-site laboratory are approximately 20 to 40 times the corresponding off-site data. This is caused by differences in analytical methodology. The sample digestion process in the analytical method specified for the off-site laboratory is designed to extract the contaminants from soil matrix samples, not to enable a total soil composition analysis. The XRF technique, however, is non-selective and will detect all barium in a sample, both from contamination and soil composition.

The paylimit and bottom samples were analyzed at a secondary dilution for pesticides/PCBs. There is no impact on data usability for these samples as no compound method detection limits were raised above the applicable STCLs.

Low levels of metallic analytes were detected in the preparation blanks associated with the samples from this structure. For this reason, sodium is considered an artifact in the bottom, inlet, and outlet samples and potassium is considered an artifact in the paylimit sample. As neither element has an applicable STCL, these qualifications have no impact on the overall usability of the data. All other analytes detected in blanks have no effect on the data.

4.4 SOIL FINAL DISPOSITION

The analytical results for the paylimit from this drainage structure were reviewed and are summarized in Table 4-4. The STCL was exceeded for copper. Based on these results the paylimit soils will be treated in the OABF.

| Analyte | Result | Estimated | Exceeds | STCL Clean | TCLP | 20X Rule |
|------------------------|------------|--------------|---------|-------------|----------|--------------|
| • | | Quantitation | Limits | | Limit | |
| TPH as Diesel | 160 mg/kg | | | 500 mg/kg | | |
| 1,4-Dichlorobenzene | 10 ug/kg | | | 9370 ug/kg | 7.5 mg/L | 150000 ug/kg |
| Aluminum | 1030 mg/kg | | | 26400 mg/kg | | |
| Arsenic | 0.85 mg/kg | | | 3.6 mg/kg | 5 mg/L | 100 mg/kg |
| Barium | 6.9 mg/kg | | | 3800 mg/kg | 100 mg/L | 2000 mg/kg |
| Calcium | 261 mg/kg | | | | | |
| Chromium | 1.5 mg/kg | | | 6.8 mg/kg | 5 mg/L | 100 mg/kg |
| Copper | 23.6 mg/kg | | YES | 19.3 mg/kg | | |
| Iron | 1550 mg/kg | | | | | |
| Lead | 11.0 mg/kg | 1 | | 15.8 mg/kg | 5 mg/L | 100 mg/kg |
| Magnesium | 168 mg/kg | [[| | | | |
| Manganese | 17.3 mg/kg | | | 274 mg/kg | | |
| Potassium | 195 mg/kg | | | | | |
| Selenium | 0.72 mg/kg | { | | 7.8 mg/kg | 1 mg/L | 20 mg/kg |
| Vanadium | 2.5 mg/kg | 1 | | 15.2 mg/kg | | |
| Zinc | 4.4 mg/kg | | | 16 mg/kg | | |
| Benzo(b)fluoranthene | 38 ug/kg | YES | | 5000 ug/kg | | |
| Benzo(k)fluoranthene | 15 ug/kg | YES | | 5000 ug/kg | | |
| • • • • | 220 ug/kg | YES | | 812 ug/kg | | |
| Butyl benzyl phthalate | 55 ug/kg | YES | | 99300 ug/kg | | |
| Fluoranthene | 41 ug/kg | YES | | 7810 ug/kg | | ! |
| Pyrene | 38 ug/kg | YES | | 4690 ug/kg | | |

Table 4-4 - Paylimit Sample 04CDXX1SOPLOF1

4.5 DEBRIS FINAL DISPOSITION

The debris from this structure was decontaminated. It will be sorted into suitable and non-suitable debris. The suitable debris will be used for aggregate in the OABF. The unsuitable debris will be sent to a construction debris landfill.

5. APPENDICES

A. ON-SITE ANALYTICAL DATA

B. OFF-SITE ANALYTICAL DATA

<u>Otis ANGB</u> Drainage Structure Removal Program

APPENDIX A

ON-SITE ANALYTICAL DATA

Otis ANGB Drainage Structure Removal Program

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Page No. 1

01714796

OTIS AIR NATIONAL GUARD BASE ON-SITE FIELD SCREENING SOIL SAMPLE RESULTS UNITS: PPM

Analyzed: 04/12/96

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| Analyte | 04CDXX1SOBTON1 | 04CDXX1SOINON1 | 04CDXX190INON1R | 04CDXX1800U0N1 | IN Stcl | OUT STCL |
|---------------|----------------|----------------|-----------------|----------------|------------|-------------|
| DDT | 0.200 U | 0.200 U | 0.200U | 0.200 U | , 90.000 | 6.280 |
| ран | 1.000 U | 1.000 U | 1.0000 | 1.000 U | 5.000 | 5.000 |
| PCB | 0.130 U | 0.130 U | 0.1300 | 0.130 U | 3.830 | 0.158 |
| TPH AS DIESEL | 53.000 U | 53.000 U | 53.000 U | 53.000 U | 1200.0 | 500.00 |
| TPH AS GAS | 27.000 U | 27.000 U | 27.000 U | 27.000 U | 1200.0 | 500.00 |
| (trans)-1,2-D | 0.010 U | 0.010 U | 0.0100 | 0.010 U | 10000 | 1100.0 |
| 1,1-DCE | 0.010 U | 0.010 U | 0.010U | 0.010 U | 0.010 | 0.010 |
| BENZENE | 0.010 U | 0.010 U | 0.0100 | 0.010 U | 0.010 | 0.010 |
| ETHYLBENZENE | 0.010 U | 0.010 U | 0.027 | 0.010 U | 0.700 | 0.700 |
| PCE | 0.010 U | 0.010 U | 0.0100 | . 0.010 U | 0.010 | 0.010 |
| TCE | 0.010 U | 0.010 U | 0.010U | 0.010 U | 0.010 | 0.010 |
| TOLUENE | 0.010 U | 0.010 U | 0.010U | 0.010 U | 1.000 | 1.000 |
| m+p-XYLENE | 0.010 U | 0.010 U | 0.010U | 0.010 U | 10.000 | 10.000 |
| O-XYLENE | 0.010 U | 0.010 U | 0.027 | 0.010 U | 10.000 | 10.000 |
| ARSENIC | 94.000 U | 94.000 U | 94.000 U | 94.000 U | 11.600 | 3.600 |
| BARIUM | 213.000 | 251.000 | 225.000 | 225.000 | 4070.0 | 3800.0 |
| CADMIUM | 90.000 U | 90.000 U | 90.000 U | 90.000 U | 27.300 | 26.400 |
| CHROMIUM | 130.000 U | 130.000 U | 130.000 U | 130.000 U | 1160.0 | 274.00 |
| LEAD | 50.000 U | 50.000 U | 50.000 U | 50.000 U | 1000.0 | 300.00 |
| SELENIUM | 50.000 U | 50.000 U | 50.000 U | 50.000 U | 291.00 | 274.00 |
| SILVER | 50.000 U | 50.000 U | 50.000 U | 50.000 U | 291.00 | 274.00 |

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APPENDIX B

OFF-SITE ANALYTICAL DATA

Otis ANGB Drainage Structure Removal Program

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OFF-SITE ANALYTICAL DATA BOTTOM SAMPLE

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<u>Otis ANGB</u> Drainage Structure Removal Program TEST CODE :SPH_0C1

JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

CLIENT : JG-4000 OTIS AIR FORCE BASE *SOLIDS : 94 * RESULTS IN DRY WEIGHT TEST NAME : 8010 VOA +EDB (JE) UNITS : UG/KG SAMPLE ID LAB : EE-96-42398 MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOBTOF1 DATE RECEIVED : 04/13/96 SDG # : 42395 DATE ANALYZED : 04/16/96 DILUTION FACTOR : 1.0

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMIT |
|---------------------------|---------|---|------------|
| | | - | |
| Dichlorodifluoromethane | ND. | | 5.3 |
| Chloromethane | ND | | 5.3 |
| Vinyl chloride | ND | | 1.1 |
| Bromomethane | ND | | 0.53 |
| Chloroethane | ND | | 0.85 |
| Trichlorofluoromethane ` | ND | | 0.64 |
| 1,1-Dichloroethene | ND | | 0.53 |
| Methylene chloride | ND | | 2.6 |
| trans-1,2-Dichloroethene | ND | | 0.53 |
| 1,1-Dichloroethane | ND | | 0.53 |
| cis-1,2-Dichloroethene | ND | | 0.53 |
| Chloroform | ND | | 5.3 |
| 1,1,1-Trichloroethane | ND | | 0.53 |
| Carbon tetrachloride | ND | | 0.53 |
| 1,2-Dichloroethane | ND | | 0.53 |
| Trichloroethene | ND | | 1.1 |
| 1,2-Dichloropropane | ND | | 3.2 |
| Bromodichloromethane | ND | | 2.1 |
| 2-Chloroethylvinylether | ND | | 2.1 |
| cis-1,3-Dichloropropene | ND | | 0.74 |
| trans-1,3-Dichloropropene | ' ND | | 1.6 |
| 1,1,2-Trichloroethane | ND | | 0.53 |
| Tetrachloroethene | ND | | 2.1 |
| Dibromochloromethane | ND | | 0.53 |
| Chlorobenzene | ND | | 0.85 |
| Bromoform | ND | | 0.53 |
| 1,1,2,2-Tetrachloroethane | ND | | 5.3 |
| 1,3-Dichlorobenzene | ND | | 0.85 |
| 1,4-Dichlorobenzene | ND | | 0.85 |
| 1,2-Dichlorobenzene | ND | | 0.85 |
| Ethylene dibromide | ND | | 1.1 |
| | | | |

QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT

B = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

TEST CODE :SPA_0C1

JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

| CLIENT : JG-4000 OTIS AIR FORCH | E BASE |
|----------------------------------|--------------------------|
| RESULTS IN DRY WEIGHT | SOLIDS : 94 |
| TEST NAME : 8020 VOA (JE) | UNITS : UG/KG |
| SAMPLE ID LAB : EE-96-42398 | MATRIX : SOLID |
| SAMPLE ID CLIENT: 04CDXX1SOBTOF1 | DATE RECEIVED : 04/13/96 |
| SDG # : 42395 | DATE ANALYZED : 04/16/96 |
| | DILUTION FACTOR : 1.0 |

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMIT |
|--|-----------------------------|---|--------------------------------|
| | | - | |
| MTBE | ND | | 1.6 |
| Benzene | ND | | 0.64 |
| Toluene | ND | | 11 |
| Ethylbenzene | ND | | 11 |
| Chlorobenzene | ND | | 1.5 |
| 1,3-Dichlorobenzene . | ND | | 1.5 |
| 1,4-Dichlorobenzene | 3.5 | | 1.3 |
| 1,2-Dichlorobenzene | ND | | 1.3 |
| Total Xylenes | ND | | 11 |
| Ethylbenzene Chlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene | ND ND ND 3.5 ND | | 11 1.5 1.5 1.3 1.3 |

QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT

B = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

1B SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

041BTOF1 Lab Name: E & E INC. Contract: Lab Code: EANDE Case No.: 689 SAS No.: SDG No.: 42395 Lab Sample ID: 42398 Matrix: (soil/water) SOIL Lab File ID: Sample wt/vol: 30.0 (g/mL) G I4264 Date Received: 04/13/96 Level: (low/med) LOW % Moisture: 6 decanted: (Y/N) N Date Extracted: 04/15/96 Concentrated Extract Volume: 1000 (uL) Date Analyzed: 04/16/96 Dilution Factor: Injection Volume: 2.0(uL) 1.0

GPC Cleanup: (Y/N) N pH:

CAS NO. COMPOUND

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

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|---|--|---|
| 108-95-2Phenol 111-44-4bis (2-Chloroethyl) Ether 95-57-82-Chlorophenol 541-73-11, 3-Dichlorobenzene 106-46-71, 4-Dichlorobenzene 100-51-6 | 350 350 350 350 350 350 350 350 350 350 | 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 |
| 88-74-42-Nitroaniline 131-11-3Dimethylphthalate | 1700 350 | บ บ บ |

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1C

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

041BTOF1 Lab Name: E & E INC. Contract: Lab Code: EANDE Case No.: 689 SAS No.: SDG No.: 42395 Matrix: (soil/water) SOIL Lab Sample ID: 42398 Sample wt/vol: 30.0 (g/mL) G Lab File ID: I4264 Level: (low/med) LOW Date Received: 04/13/96 % Moisture: 6 decanted: (Y/N) N Date Extracted: 04/15/96 Concentrated Extract Volume: 1000 (uL) Date Analyzed: 04/16/96 Injection Volume: 2.0(uL) Dilution Factor: 1.0 GPC Cleanup: (Y/N) N pH: CONCENTRATION UNITS:

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

| \cap | |
|--------|--|

| -64-9Dibenzofuran-14-22,4-Dinitrotoluene66-2Diethylphthalate5-72-34-Chlorophenyl-phenylether73-7Fluorene-01-64-Nitroaniline-52-14,6-Dinitro-2-methylphenol30-6N-Nitrosodiphenylamine (1)-55-34-Bromophenyl-phenylether-74-1Hexachlorobenzene86-5Phenanthrene-12-7Anthracene74-8Carbazole74-8Benzidine-00-0Pyrene68-7Butylbenzylphthalate94-1 | 1700 350 350 350 350 1700 1700 350 350 350 1700 | 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 |
|--|---|---|
| -64-9Dibenzoturan -14-22,4-Dinitrotoluene 66-2Diethylphthalate 5-72-34-Chlorophenyl-phenylether 73-7Fluorene -01-64-Nitroaniline -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1 | 350 350 350 1700 1700 350 350 350 1700 48 | 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 |
| 66-2Diethylphthalate 5-72-34-Chlorophenyl-phenylether 73-7Fluorene -01-64-Nitroaniline -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1Hexachlorophenol 01-8Pentachlorophenol 01-8Phenanthrene -12-7Anthracene 74-8Carbazole 74-2 | 350 350 1700 1700 350 350 350 1700 48 | 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 |
| 66-2Diethylphthalate 5-72-34-Chlorophenyl-phenylether 73-7Fluorene -01-64-Nitroaniline -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1Hexachlorophenol 01-8Pentachlorophenol 01-8Phenanthrene -12-7Anthracene 74-8Carbazole 74-2 | 350 350 1700 350 350 350 350 1700 48 | 0 0 0 0 0 0 0 0 0 0 0 |
| 73-7Fluorene -01-64-Nitroaniline -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Phenanthrene -12-7 | 350 1700 350 350 350 350 1700 48 | ם ס ס ס ס ט ט |
| 73-7Fluorene -01-64-Nitroaniline -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Phenanthrene -12-7 | 1700 1700 350 350 350 1700 48 | ם ם ם ם ם ם |
| -01-64-Nitroaniline -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Anthracene 74-8Carbazole 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | 1700 350 350 350 1700 48 | ם ם ם ם ם ם |
| -52-14,6-Dinitro-2-methylphenol 30-6N-Nitrosodiphenylamine (1) -55-34-Bromophenyl-phenylether -74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Anthracene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | 1700 350 350 350 1700 48 | บ บ บ บ บ |
| 74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Phenanthrene -12-7Phenanthrene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-1 | 350 350 350 1700 48 | บ บ บ |
| 74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Phenanthrene -12-7Phenanthrene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-1 | 350 350 1700 48 | ប ប |
| 74-1Hexachlorobenzene 86-5Pentachlorophenol 01-8Phenanthrene -12-7Phenanthrene -12-7Phenanthrene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-1 | 350 1700 48 | ប ប |
| 01-8Phenanthrene -12-7Phenanthrene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | 1700 48 | U |
| 01-8Phenanthrene -12-7Phenanthrene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | 48 | |
| -12-7Anthracene 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | | |
| 74-8Carbazole 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | 350 | U |
| 74-2Di-n-Butylphthalate -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13,3'-Dichlorobenzidine | 350 | U |
| -44-0Fluoranthene 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13.3'-Dichlorobenzidine | 350 | U |
| 87-5Benzidine -00-0Pyrene 68-7Butylbenzylphthalate 94-13.3'-Dichlorobenzidine | 46 | JJ |
| -00-0Pyrene 68-7Butylbenzylphthalate 94-13.3'-Dichlorobenzidine | 1700 | U |
| 68-7Butylbenzylphthalate | 50 | J |
| 94-13,3'-Dichlorobenzidine | 41 | J |
| | 700 | Ū |
| 55-3Benzo (a) Anthracene | 350 | Ū |
| -01-9Chrysene | 350 | Ū |
| -81-7bis (2-Ethylhexyl) Phthalate | 100 | Ĵ |
| -84-0Di-n-Octyl Phthalate | 350 | Ū |
| -99-2Benzo (b) Fluoranthene | 350 | Ū |
| -08-9Benzo (k) Fluoranthene | 350 | Ŭ |
| 32-8Benzo (a) Pyrene | 350 | U |
| 32-8Benzo (a) Pyrene -39-5Indeno (1, 2, 3-cd) Pyrene | 350 | Ŭ |
| 70-3Dibenz (a, h) Anthracene | | U |
| -24-2Benzo (g, h, i) Perylene | 350 | Ŭ |

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/9

EPA SAMPLE NO.

TEST CODE :SPPH 1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT\$SOLIDS : 94TEST NAME : TPH AS GASOLINEUNITS : MG/KGSAMPLE ID LAB : EE-96-42398MATRIX : SOLIDSAMPLE ID CLIENT: 04CDXX1SOBTOF1DATE RECEIVED : 04/13/96SDG #: 42395 SDG # DATE ANALYZED : 04/15/96 : 42395 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 7.0 g INJECTION VOLUME: 100 uL FINAL VOLUME: 10 mL PARAMETER RESULTS Q QNT. LIMIT ----------TPH as Gasoline ND 5.3 QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

90

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A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

TEST CODE :SCTPH 1' JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center CLIENT : JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT\$SOLIDS : 94TEST NAME : TPH AS DIESELUNITS : MG/KGSAMPLE ID LAB : EE-96-42398MATRIX : SOLIDSAMPLE ID CLIENT: 04CDXX1SOBTOF1DATE RECEIVED : 04/13/96SDG #: 42395DATE ANALYZED : 04/17/96 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 25 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 1.0 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS Q QNT. LIMIT . ----_____ TPH as Diesel 61 5.3 QUALIFIERS: C = COMMENT ND = NOT DETECTED J = ESTIMATED VALUE B = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

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TEST CODE :SP&PCB1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT SOLIDS : 94 TEST NAME : PESTICIDE-PCB UNITS : MG/KG SAMPLE ID LAB : EE-96-42398 MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOBTOF1 DATE RECEIVED : 04/13/96 SDG # : 42395 DATE ANALYZED : 04/18/96 DILUTION FACTOR : 2.0 SAMPLE VOLUME: 30 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 10 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS 0 QNT. LIMIT ----------------------Aldrin ND 0.002 alpha-BHC ND 0.002 beta-BHC ND 0.002 gamma-BHC (Lindane) ND 0.002 delta-BHC ND 0.002 Chlordane ND 0.017 4,4'-DDD ` ND 0.004 4,4'-DDE ND 0.004 4,4'-DDT ND 0.011 Dieldrin ND 0.004 Endosulfan I ND 0.004 Endosulfan II ND 0.004 Endosulfan sulfate ND 0.011 Endrin ND 0.004 Endrin aldehyde ND 0.011 Heptachlor ND 0.002 Heptachlor epoxide ND 0.002 Toxaphene ND 0.11 Methoxychlor ND 0.034 PCB-1016 ND 0.042 PCB-1221 ND 0.042 PCB-1232 ND 0.042 PCB-1242 ND 0.042 PCB-1248 ND 0.042 PCB-1254 ND 0.042 PCB-1260 ND 0.042 QUALIFIERS: C = COMMENT ND = NOT DETECTED J = ESTIMATED VALUE B = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

INORGANIC ANALYSES DATA SHEET

Lab Name: ECOLOGY_AND_ENVIRONMENT__ Contract:42398Lab Code: EANDE__ Case No.: 9600.689 SAS No.:SDG No.: 42395_Matrix (soil/water): SOIL_Lab Sample ID: 42398Level (low/med):LOW__% Solids:94.2

Concentration Units (ug/L or mg/kg dry weight): MG/KG

| | CAS No. | Analyte | Concentration | С | Q | м | |
|---------------|---|---|---|---|---|------------|---|
| | $\begin{array}{r} 7429 - 90 - 5\\ 7440 - 36 - 0\\ 7440 - 38 - 2\\ 7440 - 39 - 3\\ 7440 - 41 - 7\\ 7440 - 43 - 9\\ 7440 - 47 - 3\\ 7440 - 47 - 3\\ 7440 - 48 - 4\\ 7440 - 50 - 8\\ 7439 - 92 - 1\\ 7439 - 95 - 4\\ 7439 - 95 - 4\\ 7439 - 95 - 4\\ 7439 - 96 - 5\\ 7439 - 97 - 6\\ 7440 - 09 - 7\\ 7782 - 49 - 2\\ 7440 - 22 - 4\\ 7440 - 23 - 5\\ 7440 - 28 - 0\\ 7440 - 62 - 2\\ 7440 - 66 - 6\\ \hline \end{array}$ | Aluminum Antimony Arsenic Barium Beryllium Cadmium Calcium Chromium Chromium Cobalt Copper Iron Lead Magnesium Manganese Mercury Nickel Potassium Selenium Silver Sodium Thallium Vanadium Zinc Cyanide | $ \begin{array}{c} 2320 \\ 6.1 \\ 1.7 \\ 11.9 \\ 0.14 \\ 0.49 \\ 354 \\ 6.5 \\ 1.5 \\ 20.0 \\ 3370 \\ 19.4 \\ 619 \\ 30.4 \\ 0.11 \\ 1.7 \\ 424 \\ 0.50 \\ 0.57 \\ 80.4 \\ 0.45 \\ 6.5 \\ 8.7 \\ 8.7 \\ \end{array} $ | | | | · • • • • • • • • • • • • • • • • • • • |
| Color Before: | | Clari | ty Before: | | - | Texture: | SAND |
| Color After: | Y | Clari | ty After: C | | | Artifacts: | YES |
| Comments: | CLIENT_SAMP | LE_ID:04 | CDXX1SOBTOF1 | | | | ······ |
| | <u> </u> | F | ORM I - IN | | | | ILM03.0 |

183-

EPA SAMPLE NO.

OFF-SITE ANALYTICAL DATA INLET SAMPLE

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JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc.

TEST CODE :SPH 0C1

Analytical Services Center

CLIENT: JG-4000 OTIS AIR FORCERESULTS IN DRY WEIGHT%SOLIDS : 95TEST NAME: 8010 VOA +EDB (JE)UNITS : UG/KGSAMPLE ID LAB: EE-96-42396MATRIX : SOLIDSAMPLE ID CLIENT: 04CDXX1SOINOF1DATE RECEIVED : 04/13/96TEST: 42395DATE ANALYZED : 04/16/96DILUTTON FACTOR : 1.0 DILUTION FACTOR : 1.0

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMI |
|---------------------------|---------|---|-----------|
| Dichlorodifluoromethane | ND | - | 5.3 |
| Chloromethane | ND | | 5.3 |
| Vinyl chloride | ND | | 1.0 |
| Bromomethane | ND | | 0.53 |
| Chloroethane | ND | | 0.84 |
| Trichlorofluoromethane | ND | | 0.6 |
| 1,1-Dichloroethene | ND | | 0.5 |
| Methylene chloride | ND | | 2.6 |
| trans-1,2-Dichloroethene | ND | | 0.5 |
| 1,1-Dichloroethane | ND | | 0.5 |
| cis-1,2-Dichloroethene | ND | | 0.5 |
| Chloroform | ND | | 5.3 |
| 1,1,1-Trichloroethane | ND | | 0.5 |
| Carbon tetrachloride | ND | | 0.5 |
| 1,2-Dichloroethane | ND | | 0.5 |
| Trichloroethene | ND | | 1.0 |
| 1,2-Dichloropropane | ND | | 3.2 |
| Bromodichloromethane | ND | | 2.1 |
| 2-Chloroethylvinylether | ND | | 2.1 |
| cís-1,3-Dichloropropene | ND | • | 0.7 |
| trans-1,3-Dichloropropene | ND | | 1.6 |
| 1,1,2-Trichloroethane | ND | | 0.5 |
| Tetrachloroethene | ND | | 2.1 |
| Dibromochloromethane | ND | | 0.5 |
| Chlorobenzene | ND | | 0.8 |
| Bromoform | ND | | 0.5 |
| 1,1,2,2-Tetrachloroethane | ND | | 5.3 |
| 1,3-Dichlorobenzene | ND | | 0.8 |
| 1,4-Dichlorobenzene | ND | | 0.8 |
| 1,2-Dichlorobenzene | ND | | 0.8 |
| Ethylene dibromide | ND | | 1.0 |

J = ESTIMATED VALUE B = ALSO PRESENT IN BLANK

433

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

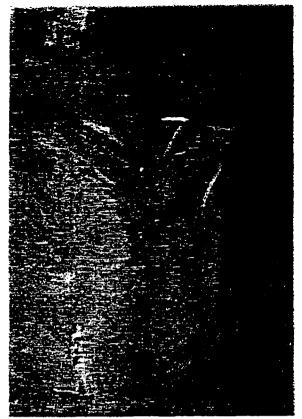
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| | Hazardous | Waste MANIFEST PI | ROGRAM | | · 1. | |
| | ase type (or print) (Form designed for use on elife (12-pitch) typewrite(.) | St., Hartford, CT 0610 | 7-3121 | | | FOR STATE USE ONL |
| Ple | | US EPA ID No. | Manuest | 2. Pac | e I Informanc | n in the shaded bloas is |
| 1 | UNIFORM HAZARDOUS | 1. (1. 2. ³ . 2. ⁷ .) : 테르 | Document No. | of | | by Federal law, out man by State law. |
| | WASTE MANIFEST | | | 1-bachel | a required a | nant Number |
| | 3. Generator's Name and Mailing Address | | | | | |
| | THE STATE WATERWATER | | | | -06 | <u>50916 2</u> |
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| | | 6. US EPA ID Nur | nper | 1.1.1 | | |
| | 5. Transporter 1 Company Name | ••• | | - | | |
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| | 7. Transporter 2 Company Name | 8. US EPA ID Nur | noer | CLS.T. | L firans Lie. Pla | 中的样子的认识 |
| | • | | | U Jrar | Phone (797 | |
| | 9. Designated Facility Name and Site Address | 10. US EPA ID Nur | noer | E-ST. | (Trans Lie Pla | R. A) Milli may war |
| 1 | | | • | F-Tean | Phone (SAF | 14715-14 - TT 121- |
| | UNITED OIL RECORDERS INC | | | | e Facility's ID (No | The second se |
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| | MERIDEN (T GELTE | 19.22.2.2.1.2.1 | | | ity s Phone 22 | |
| | | | 12. Conta | uners | - 13. Total | 14 |
| | 11. US DOT Description (Including Proper Shipping Name, Hazard C | Class, and ID Numberi | No. | Туре | | Unit Waste No. |
| | | | | | | EPA Hand |
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| E. | WOSTE COMMISTIPLE LIGHTS MUST | • | | | *** | STATE 277 |
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| | 15. Special Handling Instructions and Additional Information | 0001000 2010 000 00 000 | - CAS. 615- | 100 | | 53 - 5 |
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| | 16. GENERATOR'S CERTIFICATION: I hereby declare that the con | tents of this consignment are f | ully and accurat | ely desc | nbed above by | 207日 10日 10日 10日 10日 10日 10日 10日 10日 10日 10 |
| | i i i i i i i i i i i i i i i i i i i | abeled, and are in all respects | in proper conon | | | ay En This |
| | according to applicable international and national government (8 | | | | | of beginning of the |
| 11 | If I am a large quantity generator, I certify that I have a program i economically practicable and that I have selected the practicable | | | | | |
| | and future threat to human health and the environment; OR, if I a | im a small quantity generalor, | have made a g | ood faith | effort to minimize | my waste generation at |
| | select the best waste management method that is available to m | e and that I can alford. | | | | |
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| T A | 17. Transporter 1 Acknowledgement of Receipt of Materialis | Signature, 1 | | | 100 | Month Day |
| Å | Printed/Typed Name | Signature | | Sec. 11. | 1 Sec. 3 | |
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| Р 0 | 18. Transporter 2 Acknowledgement of Receipt of Materials | | <u> </u> | | | ••••• |
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| L' | 20. Facility Owner or Operator: Certification of receipt of hazardous m | aterials covered by this manife | st except as not | ed in ite | m 19. | |
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| | Form 8700-22 (Rev. 991) Form Approved CMB No. 2050-0039, Expires 9/30/98. | Previous edition is obsciete. | | • • | · · · · · · · · · · · · · · · · · · · | |

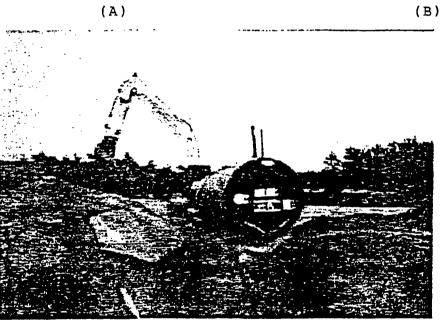
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(C)

- A. Preparation for removal of tank from Building #5218
- B. Tank grave of 12,500 gallon diesel tank OMS#22
- C. Extracted tank from OMS#22 adjacent to excavation

TEST CODE :SPA_OC1

JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

: JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT *SOLIDS : 95 * TEST NAME : 8020 VOA (JE) UNITS : UG/KG MATRIX : SOLID SAMPLE ID LAB : EE-96-42396 SAMPLE ID CLIENT: 04CDXX1SOINOF1 DATE RECEIVED : 04/13/96 SDG # : 42395 DATE ANALYZED : 04/16/96 DILUTION FACTOR : 1.0

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMIT |
|---------------------|---------|---|------------|
| | | - | |
| MTBE | ND | | 1.6 |
| Benzene | ND | | 0.63 |
| Toluene | ND | | 10 |
| Ethylbenzene | ND | | 10 |
| Chlorobenzene | ND . | | 1.5 |
| 1,3-Dichlorobenzene | . D | | 1.5 |
| 1,4-Dichlorobenzene | ND | | 1.3 |
| 1,2-Dichlorobenzene | ND | | 1.3 |
| Total Xylenes | ND | | 10 |

QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

1B SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

| Lab Name: E & E INC. | | Contract: | | 041INOF1 | |
|----------------------|-----------------|-----------|-----------------|------------|---|
| Jab Name. H a H INC. | | concrace. | ١. | | |
| Lab Code: EANDE | Case No.: 689 | SAS No.: | SDG 1 | No.: 42395 | |
| Matrix: (soil/water) | SOIL | | Lab Sample ID: | 42396 | |
| Sample wt/vol: | 30.0 (g/mL) G | | Lab File ID: | I4262 | |
| Level: (low/med) | LOW | | Date Received: | 04/13/96 | |
| % Moisture: 5 | decanted: (Y/N) | N | Date Extracted: | 04/15/96 | |
| Concentrated Extract | Volume: 1000 | (uL) | Date Analyzed: | 04/16/96 | |
| Injection Volume: | 2.0(uL) | | Dilution Factor | : 1.0 | - |
| GPC Cleanup: (Y/N) | N pH: | CONC | ENTRATION UNITS | : | |

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

| 108-95-2Phenol 111-44-4bis (2-Chloroethyl)Ether 95-57-82-Chlorophenol 541-73-11, 3-Dichlorobenzene 106-46-71, 4-Dichlorobenzene 100-51-6Benzyl Alcohol 95-50-11, 2-Dichlorobenzene 95-48-72, 2'-oxybis (1-Chloropropane) 106-44-54-Methylphenol 106-44-5Hexachloroethane 98-95-3 | 350 350 350 350 350 350 350 350 350 350 | |
|--|--|-----------------------|
| | , JJU | บ บ บ |
| 67-72-1Hexachloroethane | 350 | U U U |
| 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 65-85-0Benzoic Acid 111-91-1bis (2-Chloroethoxy)Methane | 350 350 350 1700 350 | U |
| 65-85-0Benzoic Acid 111-91-1bis (2-Chloroethoxy) Methane 120-83-22, 4-Dichlorophenol 120-82-11, 2, 4-Trichlorobenzene 91-20-3Naphthalene 106-47-84-Chloroaniline | 350 350 350 350 | U U U U |
| 120-82-11,2,4-Trichlorobenzene 91-20-3Naphthalene 106-47-8A-Chloroaniline 87-68-3 | 350 350 350 350 350 350 | U U U U U |
| 95-95-42,4,5-Trichlorophenol 91-58-72-Chloronaphthalene 88-74-42-Nitroaniline 131-11-3Dimethylphthalate | 1700 350 1700 350 | U U U |
| 91-58-72-Chloronaphthalene 88-74-42-Nitroaniline 131-11-3Dimethylphthalate 208-96-8Acenaphthylene 606-20-22,6-Dinitrotoluene 99-09-23-Nitroaniline 83-32-9Acenaphthene | 350 350 1700 350 | |

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1C SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

COMPOUND

CAS NO.

| EPA | SAMPLE | NÖ |
|-----|--------|----|
|-----|--------|----|

| | Lab Name: E & E INC. | | Contract | : | 0411NOF1 |
|---|----------------------|-----------------|----------|-----------------|------------|
| - | Lab Code: EANDE | Case No.: 689 | SAS No. | : SDG | No.: 42395 |
| | Matrix: (soil/water) | SOIL | | Lab Sample ID: | 42396 |
| | Sample wt/vol: | 30.0 (g/mL) G | | Lab File ID: | I4262 |
| | Level: (low/med) | rom | | Date Received: | 04/13/96 |
| • | % Moisture: 5 | decanted: (Y/N) | Ν | Date Extracted | : 04/15/96 |
| | Concentrated Extract | Volume: 1000 | (uL) | Date Analyzed: | 04/16/96 |
| | Injection Volume: | 2.0(uL) | | Dilution Factor | r: 1.0 |
| | GPC Cleanup: (Y/N) | .Hq | | | |

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

Q

| | I | |
|-------------------------------------|------|---|
| 51-28-52,4-Dinitrophenol | 1700 | U |
| 100-02-74-Nitrophenol | 1700 | Ū |
| 132-64-9Dibenzofuran | 350 | U |
| 121-14-22,4-Dinitrotoluene | 350 | U |
| 84-66-2Diethylphthalate | 350 | U |
| 7005-72-34-Chlorophenyl-phenylether | 350 | U |
| 86-73-7Fluorene | 350 | U |
| 100-01-64-Nitroaniline | 1700 | Ū |
| 534-52-14,6-Dinitro-2-methylphenol | 1700 | Ū |
| 86-30-6N-Nitrosodiphenylamine (1) | 350 | Ū |
| 101-55-34-Bromophenyl-phenylether | 350 | Ū |
| 118-74-1Hexachlorobenzene | 350 | Ũ |
| 87-86-5Pentachlorophenol | 1700 | Ū |
| 85-01-8Phenanthrene | 350 | Ū |
| 120-12-7Anthracene | 350 | Ū |
| 86-74-8Carbazole | 350 | Ū |
| 84-74-2Di-n-Butylphthalate | 350 | Ū |
| 206-44-0Fluoranthene | 350 | U |
| 92-87-5Benzidine | 1700 | U |
| 129-00-0Pyrene | 350 | U |
| 85-68-7Butylbenzylphthalate | 43 | J |
| 91-94-13,3'-Dichlorobenzidine | 690 | U |
| 56-55-3Benzo(a)Anthracene | 350 | U |
| 218-01-9Chrysene | 350 | U |
| 117-81-7bis(2-Ethylhexyl)Phthalate | 40 | J |
| 117-84-0Di-n-Octyl Phthalate | 350 | U |
| 205-99-2Benzo(b)Fluoranthene | 350 | U |
| 207-08-9Benzo(k)Fluoranthene | 350 | U |
| 50-32-8Benzo (a) Pyrene | 350 | U |
| 193-39-5Indeno(1,2,3-cd)Pyrene | 350 | U |
| 53-70-3Dibenz(a,h)Anthracene | | U |
| 191-24-2Benzo(g,h,i)Perylene | 350 | U |
| | . | |
| | | 1 |

(1) - Cannot be separated from Diphenylamine

3/90

TEST CODE :SPPH 1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT

 *SOLIDS : 95
 *

 INE
 UNITS : MG/KG

 396
 MATRIX : SOLID

 OINOF1
 DATE RECEIVED : 04/13/96

 DATE ANALYZED : 04/15/96

 RESULTS IN DRY WEIGHT TËST NAME : TPH AS GASOLINE SAMPLE ID LAB : EE-96-42396 SAMPLE ID CLIENT: 04CDXX1SOINOF1 SDG # : 42395 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 7.0 g FINAL VOLUME: 10 mL INJECTION VOLUME: 100 uL PARAMETER RESULTS Q QNT. LIMIT ------ ----------TPH as Gasoline ND 5.3 QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT B = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

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A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

TEST CODE :SCTPH 1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT*SOLIDS : 95TEST NAME : TPH AS DIESELUNITS : MG/KGSAMPLE ID LAB : EE-96-42396MATRIX : SOLIDSAMPLE ID CLIENT: 04CDXX1SOINOF1DATE RECEIVED : 04/13/96SDG #: 42395DATE ANALYZED : 04/17/96 : 42395 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 25 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 1.0 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS Q QNT. LIMIT _ _ _ _ _ _ _ _ _ ------TPH as Diesel 24 5.3 QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

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TEST CODE :SP&PCB1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT *SOLIDS : 95 * TEST NAME : PESTICIDE-PCB UNITS : MG/KG SAMPLE ID LAB : EE-96-42396 MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOINOF1 DATE RECEIVED : 04/13/96 SDG # : 42395 DATE ANALYZED : 04/18/96 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 30 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 10 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS Q QNT. LIMIT ------------------Aldrin 0.001 ND alpha-BHC ND 0.001 beta-BHC ND 0.001 gamma-BHC (Lindane) ND 0.001 delta-BHC ND 0.001 Chlordane ND 0.008 4,4'-DDD ' ND 0.002 4,4'-DDE ND 0.002 4,4'-DDT ND 0.005 Dieldrin ND 0.002 ND Endosulfan I 0.002 Endosulfan II 0.002 ND Endosulfan sulfate 0.005 · ND Endrin ND 0.002 Endrin aldehyde ND 0.005 Heptachlor ND 0.001 Heptachlor epoxide ND 0.001 Toxaphene ND 0.053 Methoxychlor ND 0.017 PCB-1016 ND 0.021 PCB-1221 ND 0.021 PCB-1232 ND 0.021 PCB-1242 ND 0.021 PCB-1248 ND 0.021 PCB-1254 ND 0.021 PCB-1260 ND 0.021 _____ OUALIFIERS: C = COMMENT ND = NOT DETECTED B = ALSO PRESENT IN BLANK J = ESTIMATED VALUE X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

| U. | s. | EPA | - | CLP |
|----|----|-----|---|-----|
|----|----|-----|---|-----|

Lab Name: ECOLOGY_AND_ENVIRONMENT___ Contract: _____

1 INORGANIC ANALYSES DATA SHEET EPA SAMPLE NO.

42396

| Lab Code: EAND | E_ Cas | se No.: 960 | 00.689 SAS No.: | : | SDG No.: | 42395_ |
|----------------|---|---|----------------------------|---------------------------------------|--|--------|
| Matrix (soil/w | ater): SOIL | - | | Lab Sam | ple ID: 4239 | 6 |
| Level (low/med |): LOW_ | - | | Date Re | ceived: 04/1 | 3/96 |
| % Solids: | _95.3 | 1 | | | | |
| Co | - | | /L or mg/kg dry | v weight |): MG/KG | |
| | 1 | | , | · · · · · · · · · · · · · · · · · · · | ······································ | |
| | CAS NO. | Analyte | Concentration | C Q | м | |
| | 7429-90-5 7440-36-0 7440-38-2 | Antimony Arsenic | 2420 6.0 1.9 | <u></u> | | |
| | 7440-39-3 7440-41-7 7440-43-9 7440-70-2 | Cadmium | | | | |
| | 7440-47-3 7440-48-4 7440-50-8 | Chromium Cobalt Copper | 4.1 4.1 1.2 8.8 | | | |
| | 7439-89-6 7439-92-1 7439-95-4 7439-96-5 7439-97-6 | Iron Lead Magnesium Manganese Mercury | | * | | |
| | 7440-02-0 7440-09-7 7782-49-2 7440-22-4 | Nickel Potassium Selenium Silver | 1.6 334 0.49 0.57 | B B U U N | | |
| | 7440-23-5 7440-28-0 7440-62-2 7440-66-6 | Sodium Thallium Vanadium Zinc Cyanide | 85.2 0.44 7.8 8.9 | | | |
| Color Before: | l | Clari | ty Before: | ۱ <u></u> ۱ <u></u> | _ Texture: | SAND |
| Color After: | Y | Clari | ty After: C | | Artifacts: | YES |
| Comments: | CLIENT_SAMP | LE_ID:04 | CDXX1SOINOF1 | | | |
| | <u></u> | | | | | |
| | * | F | ORM I - IN | | | ILM03 |
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Closure Report Drainage Structure 04CDXX1

OFF-SITE ANALYTICAL DATA OUTLET SAMPLE

<u>Otis ANGB</u> Drainage Structure Removal Program TEST CODE :SPH_OC1

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JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

CLIENT : JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT \$SOLIDS : 91 \$ TEST NAME : 8010 VOA +EDB (JE) UNITS : UG/KG SAMPLE ID LAB : EE-96-42397 MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOOUOF1 DATE RECEIVED : 04/13/96 SDG # DATE ANALYZED : 04/16/96 : 42395 DILUTION FACTOR : 1.0

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SAMPLE VOLUME: 5.0 g

| DichlorodifluoromethaneND5.5DichloromethaneND5.5ChloromethaneND1.1BromomethaneND0.55ChloroethaneND0.661,1-DichloroethaneND0.661,1-DichloroetheneND0.55Methylene chlorideND0.55Methylene chloroetheneND0.55istrichloroethaneND0.55istrichloroethaneND0.55cis-1,2-DichloroetheneND0.55cis-1,2-DichloroethaneND0.55ChloroformND5.51,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.55TrichloroethaneND0.55TrichloroethaneND0.55TrichloropropaneND1.11,2-DichloropropaneND2.22-ChloroethylvinyletherND2.22-ChloroethylvinyletherND0.55TetrachloroethaneND0.55TetrachloroethaneND0.55TetrachloroethaneND0.55TetrachloroethaneND0.55TetrachloroethaneND0.55ChlorobenzeneND0.55I,1,2,2-TetrachloroethaneND0.551,1,2,2-TetrachloroethaneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 <t< th=""><th>PARAMETER</th><th>RESULTS</th><th>Q</th><th>QNT. LIMIT</th></t<> | PARAMETER | RESULTS | Q | QNT. LIMIT |
|---|---------------------------|---------|---|------------|
| ChloromethaneND5.5Vinyl chlorideND1.1BromomethaneND0.55ChloroethaneND0.661.1-DichloroethaneND0.661.1-DichloroethaneND0.55Methylene chlorideND2.7trans-1.2-DichloroethaneND0.55cis-1.2-DichloroethaneND0.55cis-1.2-DichloroethaneND0.55chloroformND0.551.1-DichloroethaneND0.55cis-1.2-DichloroethaneND0.551.1.1-TrichloroethaneND0.551.2-DichloroethaneND0.551.2-DichloroethaneND0.551.2-DichloroethaneND0.551.2-DichloroethaneND0.551.2-DichloroethaneND0.551.2-DichloroethaneND0.551.2-DichloropropaneND1.11.2-DichloropropaneND2.2cis-1.3-DichloropropeneND0.77trans-1.3-DichloropropeneND0.55TetrachloroethaneND0.55ChlorobenzeneND0.551.1.2.2-TrichloroethaneND0.551.1.2.2-TrichloroethaneND0.551.1.2.2-TetrachloroethaneND0.551.1.2.2-TetrachloroethaneND0.551.1.2.2-TetrachloroethaneND0.551.1.2.2-TetrachloroethaneND0.581.3-DichlorobenzeneND0.881.4-Dichlorobenz | | | - | |
| Vinyl chlorideND1.1BromomethaneND0.55ChloroethaneND0.88TrichlorofluoromethaneND0.661,1-DichloroetheneND0.55Methylene chlorideND2.7trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55chloroformND0.55chloroformND5.51,1-TrichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.55TrichloroethaneND0.55TrichloroetheneND0.55TrichloroetheneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.55TetrachloroethaneND0.55TetrachloroethaneND0.55ChlorobenzeneND0.551,1,2-TrichloroethaneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-Dic | | ND | | 5.5 |
| BromomethaneND0.55ChloroethaneND0.88TrichlorofluoromethaneND0.661,1-DichloroetheneND0.55Methylene chlorideND2.7trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55chloroformND0.551,1-TrichloroethaneND0.551,1,1-TrichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.55TrichloroethaneND0.55TrichloroethaneND0.55TrichloroethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.55Tetras-1,3-DichloropropeneND0.55TetrachloroethaneND0.55ChlorobenzeneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.8 | | ND | | 5.5 |
| ChloroethaneND0.88TrichlorofluoromethaneND0.661,1-DichloroetheneND0.55Methylene chlorideND2.7trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55chloroformND0.551,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.2cis-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55TetrachloroetheneND0.55TetrachloroethaneND0.55TetrachloroetheneND0.551,1,2-TrichloroethaneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Vinyl chloride | ND | | 1.1 |
| TrichlorofluoromethaneND0.661,1-DichloroetheneND0.55Methylene chlorideND2.7trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55ChloroformND0.551,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND1.11,2-DichloropropaneND1.11,2-DichloropropaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.55TetrachloroethaneND0.55TetrachloroethaneND0.551,1,2-TrichloroethaneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND | Bromomethane | ND | | 0.55 |
| 1,1-DichloroetheneND0.55Methylene chlorideND2.7trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55chloroformND5.51,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.55TrichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-Dichlorobenzene </td <td>Chloroethane</td> <td>ND</td> <td></td> <td>0.88</td> | Chloroethane | ND | | 0.88 |
| Methylene chlorideND2.7trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55ChloroformND5.51,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.551,2-DichloroethaneND0.55TrichloroethaneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND0.55TetrachloroethaneND0.55ChlorobenzeneND0.551,1,2-TrichloroethaneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-Dichlorobenzene </td <td>Trichlorofluoromethane `</td> <td>ND</td> <td></td> <td>0.66</td> | Trichlorofluoromethane ` | ND | | 0.66 |
| trans-1,2-DichloroetheneND0.551,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55ChloroformND5.51,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-Dichloro | 1,1-Dichloroethene | ND | | 0.55 |
| 1,1-DichloroethaneND0.55cis-1,2-DichloroetheneND0.55ChloroformND5.51,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.551,2,2-TetrachloroethaneND0.551,3-DichloropeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneN | Methylene chloride | ND | | |
| cis-1,2-DichloroetheneND0.55ChloroformND5.51,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND0.77trans-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.551,1,2-TrichloroethaneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.881,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-Dichlorobenzene | trans-1,2-Dichloroethene | ND | | 0.55 |
| ChloroformND5.51,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.55ChlorobenzeneND0.551,3-DichloropropeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND | 1,1-Dichloroethane | ND | | 0.55 |
| 1,1,1-TrichloroethaneND0.55Carbon tetrachlorideND0.551,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.551,3-DichloropethaneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-Dichlorob | cis-1,2-Dichloroethene | ND | | 0.55 |
| Carbon tetrachlorideND0.551,2-DichloroethaneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.551,1,2,-TrichloroethaneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Chloroform | ND | | |
| 1,2-DichloroethaneND0.55TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | 1,1,1-Trichloroethane | ND | | 0.55 |
| TrichloroetheneND1.11,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.55ChlorobenzeneND0.551,1,2,2-TetrachloroethaneND0.88BromoformND0.551,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Carbon tetrachloride | ND | | 0.55 |
| 1,2-DichloropropaneND3.3BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND0.881,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | 1,2-Dichloroethane | ND | | 0.55 |
| BromodichloromethaneND2.22-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND2.2DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,2,2-TetrachloroethaneND0.881,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Trichloroethene | ND | | 1.1 |
| 2-ChloroethylvinyletherND2.2cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND2.2DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | | ND | | 3.3 |
| cis-1,3-DichloropropeneND0.77trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND2.2DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Bromodichloromethane | ND | | 2.2 |
| trans-1,3-DichloropropeneND1.61,1,2-TrichloroethaneND0.55TetrachloroetheneND2.2DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | 2-Chloroethylvinylether | ND | | 2.2 |
| 1,1,2-TrichloroethaneND0.55TetrachloroetheneND2.2DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | cis-1,3-Dichloropropene | ND | | 0.77 |
| TetrachloroetheneND2.2DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.880,880.880.88 | | ND | | 1.6 |
| DibromochloromethaneND0.55ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | | ND | | 0.55 |
| ChlorobenzeneND0.88BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | | ND | | 2.2 |
| BromoformND0.551,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Dibromochloromethane | ND | | 0.55 |
| 1,1,2,2-TetrachloroethaneND5.51,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | Chlorobenzene | ND | | 0.88 |
| 1,3-DichlorobenzeneND0.881,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | | ND | | 0.55 |
| 1,4-DichlorobenzeneND0.881,2-DichlorobenzeneND0.88 | 1,1,2,2-Tetrachloroethane | ND | | 5.5 |
| 1,2-Dichlorobenzene ND 0.88 | 1,3-Dichlorobenzene | ND | | 0.88 |
| | 1,4-Dichlorobenzene | ND | | 0.88 |
| Ethylene dibromide ND 1.1 | 1,2-Dichlorobenzene | ND | | 0.88 |
| | Ethylene dibromide | ND | | 1.1 |

QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT B = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

TEST CODE :SPA OC1

JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

CLIENT : JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT\$SOLIDS : 91TEST NAME : 8020 VOA (JE)UNITS : UG/KGSAMPLE ID LAB : EE-96-42397MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOOUOF1DATE RECEIVED : 04/13/96SDG # : 42395DATE ANALYZED : 04/16/96 DILUTION FACTOR : 1.0

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMIT |
|---------------------|---------|---|------------|
| | | - | |
| MTBE | ND | | 1.6 |
| Benzene | ND | | 0.66 |
| Toluene | ND | | 11 |
| Ethylbenzene | ND | | 11 |
| Chlorobenzene | ND | | 1.5 |
| 1,3-Dichlorobenzene | . ND | | 1.5 |
| 1,4-Dichlorobenzene | ND | | 1.3 |
| 1,2-Dichlorobenzene | ND | | 1.3 |
| Total Xylenes | ND | | 11 |

QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT

B = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

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EPA SAMPLE NO.

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

0410UOF1 Lab Name: E & E INC. Contract: Lab Code: EANDE Case No.: 689 SAS No.: SDG No.: 42395 Matrix: (soil/water) SOIL Lab Sample ID: 42397 Sample wt/vol: 30.0 (g/mL) G Lab File ID: I4263 Date Received: 04/13/96 Level: (low/med) LOW Date Extracted: 04/15/96 % Moisture: 9 decanted: (Y/N) N Concentrated Extract Volume: 1000 (uL) Date Analyzed: 04/16/96 Dilution Factor: Injection Volume: 2.0(uL) 1.0 GPC Cleanup: (Y/N) N :Hq

CAS NO. COMPOUND

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

Q

| 108-95-2Phenol 111-44-4bis (2-Chloroethyl) Ether 95-57-82-Chlorophenol 541-73-11, 3-Dichlorobenzene 106-46-71, 4-Dichlorobenzene 100-51-6Benzyl Alcohol 95-50-11, 2-Dichlorobenzene 95-48-72-Methylphenol 108-60-12, 2'-oxybis (1-Chloropropane) 106-44-54-Methylphenol 621-64-7Nitroso-Di-n-Propylamine 67-72-1Hexachloroethane 98-95-3 | 360 360 360 360 360 360 360 360 360 360 | <u>ממתמתמתמתמתמתמתמתמתמת</u> מתמתמתמתמתמת |
|--|--|--|
| 59-50-74-Chloro-3-Methylphenol 91-57-62-Methylnaphthalene 77-47-4Hexachlorocyclopentadiene 88-06-22,4,6-Trichlorophenol 95-95-42,4,5-Trichlorophenol | 360 360 360 360 1800 | บ บ บ บ บ |

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1C SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

0410UOF1 Lab Name: E & E INC. Contract: Lab Code: EANDE Case No.: 689 SAS No.: SDG No.: 42395 Matrix: (soil/water) SOIL -Lab Sample ID: 42397 Sample wt/vol: 30.0 (g/mL) G Lab File ID: I4263 Level: (low/med) LOW Date Received: 04/13/96 Date Extracted: 04/15/96 % Moisture: 9 decanted: (Y/N) N Concentrated Extract Volume: 1000 (uL) Date Analyzed: 04/16/96 Injection Volume: 2.0(uL) Dilution Factor: 1.0 GPC Cleanup: (Y/N) N pH: CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG CAS NO. COMPOUND 0

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(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90_

EPA SAMPLE NO.

TEST CODE :SPPH 1' JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center CLIENT ; JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT**%**SOLIDS : 91TEST NAME : TPH AS GASOLINEUNITS : MG/KGSAMPLE ID LAB : EE-96-42397MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOOUOF1 DATE RECEIVED : 04/13/96 DATE ANALYZED : 04/15/96. SDG # : 42395 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 7.0 g INJECTION VOLUME: 100 uL FINAL VOLUME: 10 mL PARAMETER RESULTS Q QNT. LIMIT ---------TPH as Gasoline ND 5.5 QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT. N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

89

7

TEST CODE :SCTPH 1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center CLIENT : JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT*SOLIDS : 91TEST NAME : TPH AS DIESELUNITS : MG/KGSAMPLE ID LAB : EE-96-42397MATRIX : SOLIDSAMPLE ID CLIENT: 04CDXX1SOOUOF1DATE RECEIVED : 04/13/96SDG #: 42395DATE ANALYZED : 04/17/96 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 25 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 1.0 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS Q QNT. LIMIT ---------5.5 TPH as Diesel 55 QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT. N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

TEST CODE :SP&PCB1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT SOLIDS : 91 TEST NAME : PESTICIDE-PCB UNITS : MG/KG MATRIX : SOLID SAMPLE ID LAB : EE-96-42397 SAMPLE ID CLIENT: 04CDXX1SOOUOF1 DATE RECEIVED : 04/13/96 SDG # : 42395 DATE ANALYZED : 04/18/96 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 30 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 10 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS QNT. LIMIT Q -------------------Aldrin 0.001 ND alpha-BHC ND 0.001 beta-BHC ND 0.001 gamma-BHC (Lindane) ND 0.001 delta-BHC ND 0.001 Chlordane ND 0.008 4,4'-DDD ' 0.002 ND 4,4'-DDE ND 0.002 4,4'-DDT ND 0.005 Dieldrin ND 0.002 Endosulfan I ND 0.002 Endosulfan II ND 0.002 ND Endosulfan sulfate 0.005 Endrin ND 0.002 Endrin aldehyde ND 0.005 Heptachlor ND 0.001 Heptachlor epoxide ND 0.001 Toxaphene ND 0.055 ND Methoxychlor 0.018 PCB-1016 ND 0.022 PCB-1221 ND 0.022 PCB-1232 ND 0.022 ND PCB-1242 0.022 ND PCB-1248 0.022 PCB-1254 ND 0.022 PCB-1260 ND 0.022 QUALIFIERS: C = COMMENT ND = NOT DETECTED J = ESTIMATED VALUE B = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

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| U.S. EPA - CL |
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1 INORGANIC ANALYSES DATA SHEET

| Lab Name: ECOLOGY_AN | D_ENVIRONMENT Con | tract: | 42397 | |
|----------------------|--------------------|-----------|-----------------|---|
| Lab Code: EANDE_ | Case No.: 9600.689 | SAS No.: | SDG No.: 42395_ | |
| Matrix (soil/water): | SOIL_ | Lab Sampl | e ID: 42397 | - |
| Level (low/med): | LOW | Date Rece | ived: 04/13/96 | |
| % Solids: | _91.4 | | | |

Concentration Units (ug/L or mg/kg dry weight): MG/KG

| | CAS No. | Analyte | Concentration | с | Q | м | |
|---------------|---|---|---------------|---------------|---|------------|--|
| | $\begin{array}{r} 7429 - 90 - 5\\ 7440 - 36 - 0\\ 7440 - 38 - 2\\ 7440 - 39 - 3\\ 7440 - 41 - 7\\ 7440 - 43 - 9\\ 7440 - 47 - 3\\ 7440 - 47 - 3\\ 7440 - 48 - 4\\ 7440 - 50 - 8\\ 7439 - 92 - 1\\ 7439 - 92 - 1\\ 7439 - 95 - 4\\ 7439 - 95 - 4\\ 7439 - 95 - 4\\ 7439 - 95 - 4\\ 7439 - 96 - 5\\ 7439 - 97 - 6\\ 7440 - 02 - 0\\ 7440 - 02 - 0\\ 7440 - 28 - 0\\ 7440 - 28 - 0\\ 7440 - 62 - 2\\ 7440 - 66 - 6\\ \hline \end{array}$ | Aluminum Antimony Arsenic Barium Beryllium Cadmium Calcium Chromium Cobalt Copper Iron Lead Magnesium Manganese Mercury Nickel Potassium Selenium Silver Sodium Thallium Vanadium Zinc Cyanide | | BBUB B DUBUUB | | | 14 14 14 14 14 14 14 |
| Color Before: | | Clari | ty Before: | | _ | Texture: | SAND |
| Color After: | Y | Clari | ty After: C | | _ | Artifacts: | YES |
| Comments: | CLIENT_SAMP | LE_ID:04 | CDXX1SOOUOF1 | | | | |
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FORM I - IN

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EPA SAMPLE NO.

1

Closure Report Drainage Structure 04CDXX1

OFF-SITE ANALYTICAL DATA PAYLIMIT SAMPLE

<u>Otis ANGB</u> Drainage Structure Removal Program

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TEST CODE :SPH_0C1 JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

: JG-4000 OTIS AIR FORCE BASE CLIENT SOLIDS : 92 RESULTS IN DRY WEIGHTCOLLDS : 92TEST NAME : 8010 VOA +EDB (JE)UNITS : UG/KGSAMPLE ID LAB : EE-96-42395MATRIX : SOLIDSAMPLE ID CLIENT: 04CDXX1SOPLOF1DATE RECEIVED : 04/13/96SDG # : 42395DATE ANALYZED : 04/16/96DILUTION FACTOR : 1.0 RESULTS IN DRY WEIGHT

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMIT |
|---------------------------|------------|---|------------|
| Dichlorodifluoromethane | ND | - | 5.4 |
| Chloromethane | ND | | 5.4 |
| Vinyl chloride | ND | | 1.1 |
| Bromomethane | ND | | 0.54 |
| Chloroethane | ND | | 0.87 |
| Trichlorofluoromethane | ND | | 0.65 |
| 1,1-Dichloroethene | ND | | 0.54 |
| Methylene chloride | ND | | 2.7 |
| trans-1,2-Dichloroethene | ND | | 0.54 |
| 1,1-Dichloroethane | ND | | 0.54 |
| cis-1,2-Dichloroethene | ND | | 0.54 |
| Chloroform | ND | | 5.4 |
| 1,1,1-Trichloroethane | ND | | 0.54 |
| Carbon tetrachloride | ND | | 0.54 |
| 1,2-Dichloroethane | ND | | 0.54 |
| Trichloroethene | ND | | 1.1 |
| 1,2-Dichloropropane | ND | | 3.3 |
| Bromodichloromethane | ND | | 2.2 |
| 2-Chloroethylvinylether | ND | | 2.2 |
| cis-1,3-Dichloropropene | ND | | 0.76 |
| trans-1,3-Dichloropropene | ND | | 1.6 |
| 1,1,2-Trichloroethane | ND | | 0.54 |
| Tetrachloroethene | ND | | 2.2 |
| Dibromochloromethane | ND | | 0.54 |
| Chlorobenzene | ND | | 0.87 |
| Bromoform | ND | | 0.54 |
| 1,1,2,2-Tetrachloroethane | ND | | 5.4 |
| 1,3-Dichlorobenzene | ND | | 0.87 |
| 1,4-Dichlorobenzene | ND | | 0.87 |
| 1,2-Dichlorobenzene | ND | | 0.87 |
| Ethylene dibromide | ND | | 1.1 |
| LIFIERS: C = COMMENT | ND = NOT I | | |
| DIFIERS: $C = COMMENT$ | MD = MOI T | | .120 |

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J = ESTIMATED VALUE

B = ALSO PRESENT IN BLANK

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X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

TEST CODE :SPA_OCI JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center CLIENT : JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT**%**SOLIDS : 92TEST NAME : 8020 VOA (JE)UNITS : UG/KGSAMPLE ID LAB : EE-96-42395MATRIX : SOLID SAMPLE ID LAB: EE-90-42355SAMPLE ID CLIENT: 04CDXX1SOPLOF1DATE RECEIVED : 04/13/96SDG #: 42395DATE ANALYZED : 04/16/96 DILUTION FACTOR : 1.0

SAMPLE VOLUME: 5.0 g

| PARAMETER | RESULTS | Q | QNT. LIMIT |
|---------------------|---------|---|------------|
| | | - | |
| MTBE | ND | | 1.6 |
| Benzene | ND | | 0.65 |
| Toluene | ND | | 11 |
| Ethylbenzene | ND | | 11 |
| Chlorobenzene | ND | | 1.5 |
| 1,3-Dichlorobenzene | . ND | | 1.5 |
| 1,4-Dichlorobenzene | 10 | | 1.3 |
| 1,2-Dichlorobenzene | ND | | 1.3 |
| Total Xylenes | ND | | 11 |
| | | | |

QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK

X = EXCEEDS CALIBRATION LIMIT

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

1B SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

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| Lab Name: E & E INC. | | Contract | : | 041PLOF1 | •••• |
|----------------------|-------------------|----------|-----------------|------------|------|
| Lab Code: EANDE | Case No.: 689 | SAS No.: | SDG | No.: 42395 | |
| Matrix: (soil/water) | SOIL | | Lab Sample ID: | 42395 | |
| Sample wt/vol: | 30.0 (g/mL) G | | Lab File ID: | I4265 | |
| Level: (low/med) | LOW | | Date Received: | 04/13/96 | |
| % Moisture: 8 | decanted: (Y/N) | N | Date Extracted: | 04/15/96 | - |
| Concentrated Extract | Volume: 1000 | (uL) | Date Analyzed: | 04/16/96 | |
| Injection Volume: | 2.0(uL) | | Dilution Factor | 1.0 | - |
| GPC Cleanup: · (Y/N) | N pH: | | | | |

CAS NO.

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COMPOUND

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

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| | | | : |
|---|--------|-----|---|
| 108-95-2Phenol | 360 | ט | |
| 111-44-4bis (2-Chloroethyl) Ether | 360 | υ | " |
| 95-57-82-Chlorophenol | 360 | U | |
| 541-73-11, 3-Dichlorobenzene | 360 | U | l |
| 106-46-71,4-Dichlorobenzene | 360 | U | |
| 100-51-6Benzyl Alcohol | 360 | Ŭ | - |
| 95-50-11,2-Dichlorobenzene | 360 | U | |
| | | Ŭ | |
| 95-48-72-Methylphenol 108-60-12,2'-oxybis(1-Chloropropane) 106-44-54-Methylphenol | 360 | U | 1 |
| 106-44-5 | 360 | U | |
| 106-44-54-Methylphenol 621-64-7N-Nitroso-Di-n-Propylamine | 360 | U | |
| 67-72-1Hexachloroethane | 360 | U | |
| | 360 | U | 1 |
| 98-95-3Nitrobenzene | - | | |
| 78-59-1Isophorone | 360 | U | |
| 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol | 360 | U | Į |
| 105-57-92,4-Dimethylphenol | 360 | U | |
| 65-85-0Benzoic Acid | 1700 | U | |
| 111-91-1bis(2-Chloroethoxy)Methane | 360 | U | |
| 120-83-22,4-Dichlorophenol | 360 | U | |
| 120-82-11,2,4-Trichlorobenzene | 360 | U | |
| 91-20-3Naphthalene | 360 | U | |
| 106-47-84-Chloroaniline | 360 | U | |
| 87-68-3Hexachlorobutadiene | 360 | U | |
| 59-50-74-Chloro-3-Methylphenol | 360 | U | |
| 91-57-62-Methylnaphthalene | 360 | U | 1 |
| 77-47-4Hexachlorocyclopentadiene | 360 | U | |
| 88-06-22,4,6-Trichlorophenol | 360 | U | |
| 95-95-42,4,5-Trichlorophenol | 1700 | U | } |
| 91-58-72-Chloronaphthalene | 360 | U | Ì |
| 88-74-42-Nitroaniline | 1700 | U | |
| 131-11-3Dimethylphthalate | 360 | U | 1 |
| 708-96-8 | 1 101/ | Ū | |
| 606-20-22.6-Dinitrotoluene | 360 | Ū | |
| 99-09-23-Nitroaniline | 1700 | Ū | |
| 83-32-9Acenaphthene | 360 | U | 1 |
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1C SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

041PLOF1 Lab Name: E & E INC. Contract: Lab Code: EANDE Case No.: 689 SAS No.: SDG No.: 42395 Matrix: (soil/water) SOIL Lab Sample ID: 42395 Sample wt/vol: 30.0 (g/mL) G Lab File ID: I4265 Level: (low/med) LOW Date Received: 04/13/96 Date Extracted: 04/15/96 % Moisture: 8 decanted: (Y/N) N Concentrated Extract Volume: 1000 (uL) Date Analyzed: 04/16/96 Injection Volume: 2.0(uL) Dilution Factor: 1.0 GPC Cleanup: (Y/N) N pH: CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q 51-28-5-----2,4-Dinitrophenol_____ 1700 U 100-02-7----4-Nitrophenol 1700 U 132-64-9-----Dibenzofuran 360 U 132-64-9-----Dibenzoruran 121-14-2----2,4-Dinitrotoluene_____ 360 U 84-66-2----Diethylphthalate 360 U 7005-72-3-----4-Chlorophenyl-phenylether U 360 86-73-7-----Fluorene 360 U 100-01-6-----4-Nitroaniline 1700 U 534-52-1-----4,6-Dinitro-2-methylphenol_ 1700 U 86-30-6-----N-Nitrosodiphenylamine (1) 360 U 101-55-3-----4-Bromophenyl-phenylether 360 U 118-74-1-----Hexachlorobenzene 360 U 87-86-5-----Pentachlorophenol 1700 U 85-01-8-----Phenanthrene 360 U 120-12-7-----Anthracene 360 U 86-74-8-----Carbazole 360 U 84-74-2----Di-n-ButyIphthalate U 360 206-44-0-----Fluoranthene J 41 92-87-5-----Benzidine 1700 U 129-00-0----Pyrene 38 J 85-68-7-----Butylbenzylphthalate 55 J 91-94-1-----3,3'-Dichlorobenzidine____ 720 U 56-55-3-----Benzo(a)Anthracene U 360 218-01-9----Chrysene U 360 117-81-7-----bis(2-Ethylhexyl)Phthalate 220 J

117-84-0-----Di-n-Octyl Phthalate_____

205-99-2----Benzo(b)Fluoranthene

193-39-5-----Indeno (1, 2, 3-cd) Pyrene_____

53-70-3-----Dibenz(a,h)Anthracene

(1) - Cannot be separated from Dipnenylamine

191-24-2-----Benzo(g,h,i)Perylene

207-08-9-----Benzo(k)Fluoranthene

50-32-8----Benzo(a)Pyrene

3/90

360

38

15

360

360

360

360

U

J

J

U

U

U

U

TEST CODE :SPPH 1 JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center CLIENT : JG-4000 OTIS AIR FORCE BASE \$SOLIDS : 92LINEUNITS : MG/KG2395MATRIX : SOLID RESULTS IN DRY WEIGHT TEST NAME : TPH AS GASOLINE SAMPLE ID LAB : EE-96-42395 DATE RECEIVED : 04/13/96 DATE ANALYZED : 04/15/96 SAMPLE ID CLIENT: 04CDXX1SOPLOF1 SDG # : 42395 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 7.0 g FINAL VOLUME: 10 mL INJECTION VOLUME: 100 uL PARAMETER RESULTS Q QNT. LIMIT ---------- -----TPH as Gasoline ND 5.4 _____ QUALIFIERS: C = COMMENT ND = NOT DETECTED J = ESTIMATED VALUE B = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

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TEST CODE :SCTPH 1' JOB NUMBER :9600.689 ELAP ID : 10486 Ecology and Environment, Inc. Analytical Services Center : JG-4000 OTIS AIR FORCE BASE CLIENT RESULTS IN DRY WEIGHT SOLIDS : 92 TEST NAME : TPH AS DIESEL UNITS : MG/KG SAMPLE ID LAB : EE-96-42395 MATRIX : SOLID DATE RECEIVED : 04/13/96 SAMPLE ID CLIENT: 04CDXX1SOPLOF1 DATE ANALYZED : 04/16/96 SDG # : 42395 DILUTION FACTOR : 1.0 SAMPLE VOLUME: 25 g DATE EXTRACTED: 04/15/96 FINAL VOLUME: 1.0 mL INJECTION VOLUME: 2.0 uL PARAMETER RESULTS Q QNT. LIMIT ---------- ------TPH as Diesel 160 5.4 QUALIFIERS: C = COMMENTND = NOT DETECTEDJ = ESTIMATED VALUEB = ALSO PRESENT IN BLANK X = EXCEEDS CALIBRATION LIMIT.

N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE

JOB NUMBER :9600.689 ELAP ID : 10486

Ecology and Environment, Inc. Analytical Services Center

TEST CODE :SP&PCB1

CLIENT : JG-4000 OTIS AIR FORCE BASE RESULTS IN DRY WEIGHT SOLIDS : 92 TEST NAME : PESTICIDE-PCB : MG/KG UNITS SAMPLE ID LAB : EE-96-42395 MATRIX : SOLID SAMPLE ID CLIENT: 04CDXX1SOPLOF1 DATE RECEIVED : 04/13/96 SDG # DATE ANALYZED : 04/18/96 : 42395 DILUTION FACTOR : 5.0 SAMPLE VOLUME: 30 q DATE EXTRACTED: 04/15/96 FINAL VOLUME: 10 mL INJECTION VOLUME: 2.0 uL PARAMETER 0 QNT. LIMIT RESULTS _ _ _ _ _ _ _ _ _ _ _ _ ----------Aldrin ND 0.005 alpha-BHC ND 0.005 beta-BHC ND 0.005 0.005 gamma-BHC (Lindane) ND delta-BHC ND 0.005 Chlordane ND 0.043 4,4'-DDD ' ND 0.011 4,4'-DDE ND 0.011 4,4'-DDT ND 0.027 Dieldrin ND 0.011 Endosulfan I ND 0.011 Endosulfan II 0.011 ND Endosulfan sulfate ND 0.027 0.011 Endrin ND Endrin aldehyde ND 0.027 Heptachlor ND 0.005 Heptachlor epoxide ND 0.005 Toxaphene ND 0.27 ND 0.087 Methoxychlor PCB-1016 ND 0.11 0.11 PCB-1221 ND PCB-1232 ND 0.11 0.11 PCB-1242 ND PCB-1248 ND 0.11 PCB-1254 ND 0.11 ND 0.11 PCB-1260 _____ ND = NOT DETECTED QUALIFIERS: C = COMMENT B = ALSO PRESENT IN BLANK J = ESTIMATED VALUE X = EXCEEDS CALIBRATION LIMIT N = ANALYTE WAS NOT CONFIRMED BY ALTERNATE PROCEDURE A = PHENOMENON OF METHODOLOGY WITH ACID PRESERVATION

| U.S. | . EPA | - | CLP |
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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

42395

- Lab Name: ECOLOGY_AND_ENVIRONMENT___Contract: ________ Lab Code: EANDE____Case No.: 9600.689 SAS No.: _____SDG No.: 42395____ Matrix (soil/water): SOIL____Lab Sample ID: 42395 Level (low/med): LOW_____Date Received: 04/13/96
- ***** % Solids: 91.8

Concentration Units (ug/L or mg/kg dry weight): MG/KG

Concentration | C CAS No. Analyte Q М P 7429-90-5 1030 Aluminum Ū **P**_ 7440-36-0 6.2 Antimony_ p⁻ 0.85 Arsenic___ B 7440-38-2 P⁻ 7440-39-3 Barium 6.9 B 7440-41-7 Beryllium 0.07 U **P**⁻ 7440-43-9 Cadmium 0.50|0 P 261 B P 7440-70-2 Calcium -1.5 P 7440-47-3 Chromium <u>0</u>.90|0 P 7440-48-4 Cobalt 23.6 P 7440-50-8 Copper 1550 P 7439-89-6 Iron P 7439-92-1 11.0 Lead 168 B P^{-} 7439-95-4 Magnesium 7439-96-5 Manganese 17.3 P^{-} Mercury___ CV 7439-97-6 0.11 0 Ρ 7440-02-0 Nickel 1.5 U _____195|B 7440-09-7 P Potassium 0.72 P 7782-49-2 Selenium ០.59 ប៊ 7440-22-4 P Silver 70.9 \mathbf{P}^{-} Sodium 7440-23-5 U 7440-28-0 Ь_ Thallium 0.46 U \mathbf{P}^{-} 7440-62-2 Vanadium⁻ 2.5 в P_ 7440-66-6 Zinc 4.4 NR Cyanide Clarity Before: ____ Color Before: Texture: SAND Clarity After: C____ Color After: Artifacts: YES Υ____ Comments: CLIENT_SAMPLE_ID: 04CDXX1SOPLOF1 FORM I - IN ILM03.0