

**Waste Management Division
PO Box 95, 29 Hazen Drive
Concord, NH 03302**

Type of Submittal (Check One-Most Applicable)

<input type="checkbox"/> Work Scope <input type="checkbox"/> Reimbursement Request	<input type="checkbox"/> Remedial Action <ul style="list-style-type: none"> • Remedial Action Plan • Bid Plans and Specifications • Remedial Action Implementation Report
<input type="checkbox"/> UST Facility Report <input type="checkbox"/> AST Facility Report	<input type="checkbox"/> Treatment System and POE O&M <input type="checkbox"/> Activity and Use Restriction
<input type="checkbox"/> Emergency/Initial Response Action <input type="checkbox"/> Groundwater Quality Assessment	<input type="checkbox"/> Temporary Surface Water Discharge Permit
<input type="checkbox"/> Initial Site Characterization <input type="checkbox"/> Site Investigation <ul style="list-style-type: none"> • Site Investigation Report • Supplemental Site Investigation Report • GMZ Delineation • Source Area Investigation • Data Submittal • Annual Summary Report <input checked="" type="checkbox"/> Unsolicited Brownfields Submittal <input type="checkbox"/> Closure Documentation	<input type="checkbox"/> Groundwater Management Permit <ul style="list-style-type: none"> • Permit Application • Renewal Application • Deed Recordation Documentation • Abutter Notification Documentation • Release of Recordation <input type="checkbox"/> Data Submittal <input type="checkbox"/> Annual Summary Report

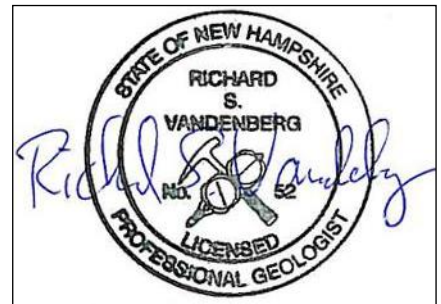
PHASE II ENVIRONMENTAL SITE ASSESSMENT REPORT

Former Polyclad Laminates Site
45 Tannery Street
Franklin, New Hampshire
NHDES Site #199902062

Prepared For:
Lakes Region Planning Commission
103 Main Street – Suite #3
Meredith, NH 03253
Phone: (603) 279-8171
Contact: Mr. Kimon Koulet

Prepared By:
CREDERE ASSOCIATES, LLC
776 Main Street
Westbrook, Maine 04092
Phone: (207) 828-1272 ext. 20
Contact: Jonathan O'Donnell

June 27, 2012



Recommended Risk Category (check one)

<input type="checkbox"/> 1. Immediate Human Health Risk (Impacted water supply well, etc.)	<input type="checkbox"/> 4. Surface Water Impact	<input checked="" type="checkbox"/> 7. Alternate Water Available/Low Level Groundwater Contamination (<1,000 X AGQS)
<input type="checkbox"/> 2. Potential Human Health Risk (Water supply well within 1,000' or Site within SWPA)	<input type="checkbox"/> 5. No Alternate Water Available/No Existing Wells in Area	<input type="checkbox"/> 8. No AGQS Violation/No Source Remaining
<input type="checkbox"/> 3. Free Product or Source Hazard	<input type="checkbox"/> 6. Alternate Water Available/High Level Groundwater Contamination (>1,000 X AGQS)	<input type="checkbox"/> Closure Recommended



CREDERE ASSOCIATES, LLC

776 Main Street
Westbrook, Maine 04092
Phone: 207-828-1272
Fax: 207-887-1051

June 27, 2012

Mr. Kimon Koulet
Executive Director
Lakes Region Planning Commission
103 Main Street, Suite #3
Meredith, New Hampshire 03253

**Subject: Phase II Environmental Site Assessment
Former Polyclad Laminates Site
45 Tannery Street, Franklin, NH
NHDES Site # 199902062**

Dear Mr. Koulet:

Attached is the Phase II Environmental Site Assessment for the Former Polyclad Laminates Site located at 45 Tannery Street in Franklin, New Hampshire. **Sections 11** and **12** of the attached report present our conclusions and recommendations regarding the Site. Copies of this report have been forwarded to the New Hampshire Department of Environmental Services (NHDES), the United States Environmental Protection Agency (U.S. EPA), and the City of Franklin.

Please do not hesitate to contact me at (207) 828-1272 ext. 20 if you have any questions, comments, or require additional information regarding this investigation.

Very truly,

CREDERE ASSOCIATES, LLC

Jonathan O'Donnell
Geo-Environmental Engineer

cc: Elizabeth Dragon, City of Franklin
Richard Lewis, City of Franklin
John Liptak, NHDES
Jerry Minor-Gordon, U.S. EPA





Phase II Environmental Site Assessment Report

Former Polyclad Laminates Site

45 Tannery Street

Franklin, New Hampshire

NHDES Site #199902062



Prepared for:

Lakes Region Planning Commission

103 Main Street, Suite #3

Meredith, New Hampshire 03253

Attn: Kimon Koulet

Executive Director

June 27, 2012



In Reference to:

Crede Project No. 10001086

Submitted by:

Crede Associates, LLC

776 Main Street

Westbrook, ME 04092

TABLE OF CONTENTS

Section	Title	Page No.
EXECUTIVE SUMMARY		1
1. INTRODUCTION.....		1-1
2. PROJECT BACKGROUND.....		2-1
2.1	Site Description.....	2-1
2.2	Summary of Previous Work.....	2-1
2.2.1	<i>Recognized Environmental Conditions.....</i>	<i>2-1</i>
2.2.2	<i>Other Potential Environmental Conditions</i>	<i>2-3</i>
2.2.3	<i>History of Underground Storage Tanks at the Site.....</i>	<i>2-4</i>
2.2.4	<i>Other Documented Previous Environmental Work Cited by Nobis</i>	<i>2-4</i>
2.3	Potential Future Site Use	2-5
3. PHASE II SCOPE OF WORK		3-1
4. PHASE II FIELD ACTIVITIES		4-1
4.1	Ground Penetrating Radar Survey	4-1
4.2	Surficial Soil Investigation	4-1
4.3	Test Pit Investigation	4-2
4.4	Sub-Slab Soil Gas Investigation	4-5
4.5	Soil Boring and Monitoring Well Installation	4-5
4.6	Groundwater Sampling	4-8
4.7	Building Material Debris Sampling	4-9
4.7.1	<i>Potential Asbestos Containing Materials</i>	<i>4-9</i>
4.7.2	<i>Potential PCB-Containing Materials</i>	<i>4-9</i>
5. SUMMARY OF REGULATORY STANDARDS		5-1
5.1	Soil.....	5-1
5.2	Groundwater	5-1
5.3	Soil Gas.....	5-1
5.4	Potential Asbestos-Containing Materials.....	5-1
5.5	Potential PCB-Containing Materials.....	5-1
6. PHASE II RESULTS.....		6-1
6.1	Site Geology	6-1
6.1.1	<i>Surficial Geology.....</i>	<i>6-1</i>
6.1.2	<i>Bedrock Geology</i>	<i>6-1</i>
6.1.3	<i>Site Hydrogeology.....</i>	<i>6-1</i>
6.2	Ground Penetrating Radar Results.....	6-1
6.3	Surficial Soil Sample Results.....	6-2
6.4	Test Pit Sample Results	6-2
6.5	Sub-Slab Soil Gas Sample Results	6-3
6.6	Soil Boring Sample Results	6-3



6.7	Groundwater Sample Results.....	6-3
6.8	Building Material/Debris Results	6-4
6.8.1	<i>Potential Asbestos Containing Material.....</i>	<i>6-4</i>
6.8.2	<i>Potential PCB-Containing Material.....</i>	<i>6-4</i>
7.	QUALITY ANALYSIS/QUALITY CONTROL.....	7-1
7.1	Precision.....	7-1
7.2	Bias	7-2
7.3	Accuracy	7-3
7.4	Representativeness.....	7-3
7.5	Comparability	7-3
7.6	Completeness	7-4
8.	CONCEPTUAL SITE MODEL	8-1
8.1	Contaminants of Concern	8-1
8.2	Site Geology and Hydrogeology.....	8-1
8.3	Definitions of Exposure Pathways and Potential Receptors	8-1
9.	DEVIATIONS.....	9-1
10.	DATA GAPS.....	10-1
11.	CONCLUSIONS	11-1
11.1	Recognized Environmental Conditions	11-1
11.2	Additional Environmental Concerns.....	11-2
12.	RECOMMENDATIONS.....	12-1
13.	SIGNATURES OF ENVIRONMENTAL PROFESSIONALS	13-1
14.	LIMITATIONS.....	14-1



LIST OF TABLES

Table 1	Summary of Exploration Locations and Sampling Methods
Table 2	Summary of Photoionization Detector (PID) Field Screening Results
Table 3	Summary of X-Ray Fluorescence (XRF) Meter Field Screening Results
Table 4	Summary of Groundwater Elevation Data
Table 5	Summary of Analytical Results for Soil Samples
Table 6	Summary of Analytical Results for Building Material Samples
Table 7	Summary of Analytical Results for Soil Gas Samples
Table 8	Summary of Analytical Results for Groundwater Samples
Table 9	Summary of Duplicate Sample Analyses

LIST OF FIGURES

Figure 1	Site Location Plan
Figure 2	Phase II Environmental Site Assessment Sample Location Plan Showing Recognized Environmental Conditions
Figure 3	UST-6 Area Enlargement Plan
Figure 4	Soil and Building Material Exceedance Location Plan
Figure 5	Summary of Groundwater Results Plan
Figure 6	July 5, 2011 Groundwater Elevation Contour Plan
Figure 7	Buried Tannery Waste Location Plan
Figure 8	Conceptual Site Model

LIST OF APPENDICES

Appendix A	Site-Specific Quality Assurance Project Plan Addendum
Appendix B	Site Photographs
Appendix C	Test Pit Logs
Appendix D	Soil Boring Logs
Appendix E	Groundwater Sampling Logs
Appendix F	Laboratory Analytical Reports



EXECUTIVE SUMMARY

Crede Associates, LLC (Crede) has conducted a Phase II Environmental Site Assessment (ESA) at the former Polyclad Laminates Site (the Site) located at 45 Tannery Street in Franklin, New Hampshire, in general conformance with the American Society for Testing Materials (ASTM) Standard Guide for Environmental Site Assessments: *Phase II Environmental Site Assessment Process* E 1903-97 (reapproved 2002). Phase II ESA activities were performed specifically to confirm or dismiss *recognized environmental conditions* (RECs) and other ASTM *non-scope considerations* (NCs) identified during a Phase I ESA that was completed for the Site by Nobis Engineering, Inc. (Nobis) in April 2011 and additional environmental concerns identified by Crede during the development of a site-specific quality assurance project plan (SSQAPP) addendum for the Site.

Crede's Phase II ESA work, which is outlined in the New Hampshire Department of Environmental Services (NHDES) and the United States Environmental Protection Agency (U.S. EPA) approved SSQAPP addendum, included the following: a ground penetrating radar (GPR) survey of the Site to identify subsurface features; advancement of five (5) soil borings; excavation of thirteen (13) test pits; installation of three (3) monitoring wells; and collection and laboratory analysis of two (2) sub-slab soil gas samples, four (4) surficial soil samples, eighteen (18) subsurface soil samples, seven (7) building material/debris samples, and groundwater samples from an existing monitoring well and the three (3) newly installed monitoring wells.

Based on the findings of this work, Crede's conclusions regarding the identified RECs and additional environmental concerns are as follows:

Recognized Environmental Conditions

- REC-1, which was associated the presence of a 12,000-gallon fuel oil underground storage tank (UST) (identified as UST-6) that was closed-in-place in the vicinity of the loading dock, is dismissed because no evidence of contamination was identified in soil samples collected from the sides and adjacent to the bottom of the closed-in-place UST in test pits CA-TP-7 and CA-TP-7A. In addition, observations made by Crede indicated that this UST was filled with clean sand at the time of closure-in-place.
- REC-2, which was associated with the potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-8) shown on a historical Sanborn Map, is dismissed because no UST was identified in this location during the GPR survey and no evidence of contamination exceeding applicable standards was identified in test pits CA-TP-8, CA-TP-10, or CA-TP-11.
- REC-3, which was associated with the potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-9) shown on a historical Sanborn Map, is dismissed because no UST was identified in this location during the GPR survey, and no evidence of



contamination was identified in test pit CA-TP-9 or in soil gas sample locations CA-SG-1 through CA-SG-5, which were completed to evaluate this REC.

- REC-4, which was associated with a former fuel dispenser and a number of filling caps formerly located in the vicinity of the closed-in-place UST-6, is dismissed because no evidence of contamination exceeding applicable standards was identified in soil samples collected from test pit CA-TP-12, which was advanced in the suspected location a former fuel dispenser.
- REC-5, which was associated with the presence of a sheen of degraded gasoline previously noted in 2008 in existing monitoring well MW-2, is confirmed, because groundwater from a newly installed monitoring well (CA-2) located 30 feet southwest of MW-2 contained 1,2,4-trimethylbenzene at a concentration exceeding the applicable NHDES Ambient Groundwater Quality Standard (AGQS), suggesting that a plume of degraded gasoline is present beneath the Site.
- REC-6, which was associated with a subsurface structure identified during a previous GPR survey east of the former building and believed to be a former concrete septic tank, is dismissed because this structure was determined to be a part of the municipal sewer system and not a septic tank, and therefore does not represent a pathway to the subsurface environment.
- REC-7, which was associated with the presence of buried waste related to the previous usage of the Site as a tannery, is dismissed. While evidence of buried waste was observed in the form of leather hides in test pits CA-TP-9, CA-TP-13, and CA-TP-14, no evidence of contamination at concentrations exceeding applicable standards was observed in association with the buried waste. However, this condition represents a solid waste issue that should be properly managed.
- REC-8, which was associated with pipes that were previously noted emitting unknown solvent odors and elevated total volatile organic readings in the southwestern portion of the building, is dismissed, because no evidence of contamination exceeding applicable standards was identified in soil gas (soil gas sample locations CA-SG-6 through CA-SG-11), soil (boring CA-3), or in groundwater (monitoring well CA-3) samples collected from this area.
- REC-9, which was associated with former press pits located within the manufacturing slab that are stained with petroleum, is dismissed because no evidence of contamination was identified in soil samples collected from adjacent to the bottoms of the press pits (borings CA-4 and CA-5).
- REC-10, which was associated with a pile of demolition debris located on the southwestern portion of the Site, is dismissed because no potential lead-based paint or PCB-containing materials were observed in the pile, and no asbestos was identified in a floor tile sample collected from the pile.



Additional Environmental Concerns

The following additional environmental concerns were identified by Credere during development of the SSQAPP and assessed during this Phase II ESA:

- Potential impact associated with the former use of the Site for railroad activities is confirmed, because benzo[a]pyrene was detected above the applicable NHDES Soil Remediation Standards (SRS) in surficial soil sample CA-SS-3. Benzo[a]pyrene is a polycyclic aromatic hydrocarbon (PAH) which is known to be associated with railroad activities.
- The potential impact of a former electrical transformer located on the west side of the building is dismissed, because no PCBs were identified in soil collected from an electrical pit in this area of the Site. Additionally, two samples of stained concrete collected from this area of the Site contained no PCBs above laboratory practical quantification limits.
- A soil sample collected from test pit CA-TP-15 at 5 to 7 feet below ground surface (bgs) contained the PAHs benzo[a]anthracene, benzo[b]fluoranthene, benzo[k]fluoranthene, benzo[a]pyrene, indeno(1,2,3-cd)pyrene, and dibenzo(a,h)anthracene at concentrations which exceeded their applicable NHDES SRS. No staining was observed in this test pit nor was this sample associated with any observed buried wastes. Additionally, no PAHs were identified in soil in contact with buried waste in test pits CA-TP-13 or CA-TP-14. Although this test pit was not in the location of or downgradient of any identified Site USTs, it is Credere's opinion that this contamination may be associated with a past petroleum release from the former UST-10, which was located approximately 40 feet away or from another unknown on-site source.
- Building material sample results indicate that blue 12-inch square floor tile remnants remaining attached to the concrete building slab are asbestos-containing material (ACM). It is not clear if any of these tiles are present in the demolition pile remaining on-site.

Based on observations and results of this Phase II ESA for the Site, Credere makes the following recommendations:

- Credere recommends that delineation of all buried solid waste at the Site be conducted, including tannery hides, other tannery wastes, and building demolition debris, per NHDES Solid Waste Rules Env-Sw 309 in support of a Pre-1981 landfill registration.
- Subsequent to the delineation of all buried solid waste at the Site, Credere recommends that the Site be registered as a Pre-1981 landfill according to NHDES Solid Waste Rules Env-Sw 309.



- Credere recommends that delineation of the identified PAH contaminated surficial soil in the areas of CA-SS-3 and CA-TP-15 be conducted. Credere recommends that a Remedial Action Plan (RAP) be developed to address the PAH contaminated soil.
- Credere recommends additional investigation to assess the full extent and magnitude of degraded gasoline contamination identified at monitoring well CA-2. This additional assessment should supplement the development of a groundwater flow model of the Site in order to define the nature, extent, and/or potential fate and transport characteristics of the identified impacted media. Credere also recommends that this petroleum contamination issue be referred to the NHDES Petroleum Remediation Program as the observed contamination is likely associated with one of the former USTs at the Site. If accepted into this program, future investigation work will be reimbursed by the program.
- Credere recommends that the identified ACM floor tile remaining on the former building slab be properly handled according to all applicable NHDES Asbestos and Solid Waste Rules prior to the development of the Site.



1. INTRODUCTION

This report presents the results of a Phase II Environmental Site Assessment (ESA) conducted by Credere Associates, LLC (Credere) at the former Polyclad Laminates site (the Site) located at 45 Tannery Street in Franklin, New Hampshire as part of the Lakes Region Planning Commission's (LRPC) Brownfields Assessment Program. **Figure 1** shows the general location of the Site in Franklin.

The Phase II ESA was completed in general conformance with the American Society for Testing Materials (ASTM) Standard Guide for Environmental Site Assessments: *Phase II Environmental Site Assessment Process* E 1903-97 (reapproved 2002).

The field program used during this Phase II ESA was completed in accordance with the U.S. Environmental Protection Agency (EPA)-approved June 28, 2011, Site-Specific Quality Assurance Project Plan (SSQAPP) Addendum. The SSQAPP is an addendum to the previously approved New Hampshire Generic QAPP RFA #08166 and #09036, which was prepared for all of Credere's EPA work in New Hampshire. The SSQAPP addendum is included in **Appendix A**. Photographs taken during the completion of this Phase II ESA are included in **Appendix B**.



2. PROJECT BACKGROUND

2.1 SITE DESCRIPTION

The Site is composed of a single vacant 1.97-acre parcel of land located at 45 Tannery Street in Franklin, New Hampshire, that is situated along the Pemigewasset River. The Site consists of a vacant lot with a concrete slab-on-grade foundation and a pile of debris remaining from the previous demolition of the Site building. **Figure 2** shows the former Site building footprint and other pertinent Site features.

According to Nobis Engineering, Inc. (Nobis)'s April 11, 2011, Phase I ESA report for the Site, which was reviewed as a part of the development of the SSQAPP, the Site was first developed in 1892. By approximately 1930, the first building was constructed on the Site. The most recent building was demolished in 2008 after it collapsed due to heavy winter snow cover.

According to the Nobis Phase I ESA report, the Site appears to have been used for railroad purposes when the Boston & Maine and Bristol Railroad transected the Site beginning in 1892. It is not clear when railroad activities ceased on the Site. However, records do show that the Hingston Leather Company, Inc. and Louis Verza Leather Company were using the Site for tannery activities beginning in the 1950s until the mid to late 1950s. Nobis indicated that there was very little environmental information available for the Site as it pertains to the former tannery operations.

Between 1979 and 1980, Polyclad Laminates, Inc. began operations on the Site. Polyclad Laminates, Inc. manufactured component materials used in multi-layered circuit boards called "pre-preg." The process involved dipping fiberglass cloth in a dip tank filled with epoxy resin. Polyclad Laminates used the Site for the manufacture of pre-preg from between 1979 and 1980 until approximately 2006.

2.2 SUMMARY OF PREVIOUS WORK

The following is a summary of the findings and conclusions of Nobis as reported in their April 2011 Phase I ESA for the Site:

2.2.1 Recognized Environmental Conditions

1. The presence of a 12,000-gallon fuel oil underground storage tank (UST) (identified as UST-6) that is located in the vicinity of the loading dock and was closed-in-place (date uncertain) represented a REC.
2. The potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-8) shown on a Sanborn Map represented a REC because the presence of this tank has not



been confirmed and no documentation regarding closure of the tank or information on the integrity of the tank was identified during the Phase I ESA.

3. The potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-9) shown on a Sanborn Map represented a REC because the presence of this tank has not been confirmed and documentation regarding closure of the tank or information on the integrity of the tank was not identified during the Phase I ESA.
4. Due to the potential for undocumented past releases to the environment, Nobis identified a REC associated with presence of a former gasoline dispenser and a number of filling caps formerly located in the vicinity of former UST-6 because these items may have been associated with additional undocumented USTs in this area, and documentation regarding the removal of this dispenser or the disposition of related piping with respect to potential releases has not been confirmed.
5. The presence of a sheen of degraded gasoline previously noted in 2008 in monitoring well MW-2 for which a source was not identified, represents a REC due to the potential for an unidentified source of petroleum contamination to be present at the Site. [Note: According to a Phase II ESA Report prepared by Delta Consultants on April 10, 2008, previous sampling data collected in 2007 indicates that no volatile organic compounds (VOCs), metals, or polychlorinated biphenyls (PCBs) were detected in excess of New Hampshire Soil Remediation Standards (SRS) in a soil sample collected at 30 to 32 feet bgs during drilling of this well and no VOCs or semi-volatile organic compounds (SVOCs) were detected in excess of Ambient Groundwater Quality Standards (AGQS) in the groundwater sample collected from this well. However, a petroleum sheen was noted during sampling of the groundwater and the tentatively identified compounds (TICs) noted during laboratory analysis revealed that the sheen was likely gasoline.]
6. The potential that a subsurface structure identified during a ground penetrating radar (GPR) survey east of the former building, which was suspected to be a former concrete septic tank, represents a REC because of the potential for previous undocumented subsurface storage or disposal of hazardous substances and/or petroleum products in and around this structure.
7. The presence of buried waste related to the previous usage of the Site as a tannery represents a REC because the understanding of the extent and disposition of the waste (leather waste or unsuitable materials) is limited and may have led to improper disposal of hazardous substances.
8. Nobis concluded that, because of the potential for undocumented releases of hazardous substances in the past or in the future, pipes that were previously noted emitting unknown solvent odors and elevated total volatile organic compounds detected in the former chemical mixing room in the southwestern portion of the building, represented a REC



because the source of these odors (which may be a formerly closed-in-place acetone tank) is not known.

9. Two former press pits located within the manufacturing slab that are stained with petroleum represent a REC and they have not been fully assessed.
10. A pile of unknown debris located on the southwestern portion of the Site that is consistent with demolition debris represents a REC for its potential to contain waste ACM, lead-based paint [and other contaminants].

2.2.2 Other Potential Environmental Conditions

Based on review of the April 2011 Nobis Phase I ESA, Credere offers the following list of other potential environmental conditions:

1. Nobis documented that the Site was previously used for railroad activities beginning as early as 1892. Nobis concluded that this past use of the Site represented a *de minimis condition* because any contaminants from this past use would likely be deemed a background condition. Credere agrees with this assertion, but it is our opinion that non-point source background type releases of polycyclic aromatic hydrocarbons (PAHs), which would be considered the primary contaminants of concern, still represent a threat to human health if they exceed applicable New Hampshire Department of Environmental Services (NHDES) soil standards and should be addressed as a part of future redevelopment of the Site. Therefore, it is our opinion that this past use should be assessed as a part of this Phase II ESA.
2. Nobis indicated in the Phase I ESA that a data gap in the form of snow cover was noted during the Site Reconnaissance portion of their work. Given that the ground could not be observed, confirmation of exterior surface conditions is warranted prior commencing Phase II ESA field work because additional RECs may be identified that require assessment. As a result, a supplemental Site Reconnaissance is recommended as the first task of this Phase II ESA.
3. A former electrical transformer was located on the west side of the building. No information was noted in the previous reports regarding the potential presence of PCBs in the dielectric fluid associated with this transformer. However, considering the timeframe when this transformer would have been in operation, it may have contained PCBs, and a potential release of dielectric fluid from this former transformer may have impacted environmental conditions at the Site. For this reason, the potential presence of PCBs is a concern that should be considered.



2.2.3 History of Underground Storage Tanks at the Site

The operational history of USTs is an important element in the environmental history of the Site. As result, the information below, as reported by Nobis in the April 2011 Phase I ESA, provides an inventory of the USTs which have been identified at the Site. Additional details concerning UST history can be reviewed in the previously submitted Nobis Phase I ESA.

UST ID	Capacity (gallons)	Material Held	Current Status
UST -1	6,000	Methyl Cellosolve	Removed, 1998
UST-2	4,500	CS 350 Acetone	Closed in-place, 1998
UST-3	4,500	Epoxy Resin	Closed in-place, 1999
UST-4	6,000	CS 350 Acetone	Closed in-place, 1999
UST-5	4,000	Gasoline	Removed, date uncertain.
UST-6	12,000	Fuel Oil	Closed in-place, date uncertain
UST-7	13,500	Epoxy Resin	Removed 2008
UST-8	4,000	Fuel Oil	Status unknown, approximate location from 1964 Sanborn Map
UST -9	4,000	Fuel Oil	Status unknown, approximate location from 1929 Sanborn Map
UST-10	8,000	Waste Petroleum Distillates	Removed 2008

2.2.4 Other Documented Previous Environmental Work Cited by Nobis

The Nobis Phase I ESA details the completion of several other investigations, which were completed outside the Brownfields due diligence process, likely as part of a previous effort to sell the Site. Documented previous works include the following:

1. A January 12, 1999 Underground Storage Tank (UST) Closure Report UST-1 prepared by Les A. Cartier and Associates, Inc.
2. A July 15, 1999 UST Closure Report UST-2, UST-3 and UST-4 prepared by Environmental Science & Engineering (ESE).
3. A 2005 Phase I Environmental Site Assessment and Limited Compliance Review prepared by Delta Consultants.
4. An April 10, 2008 Phase II Environmental Site Assessment prepared by Delta Consultants.
5. An October 15, 2008 UST Closure Report UST-7 and UST-10 prepared Delta Consultants.



The conditions identified by Nobis in **Section 2.2** cumulatively considered all the above referenced previous investigations. Credere also reviewed and considered this cumulative work as a part of the development of the SSQAPP for the Site.

2.3 POTENTIAL FUTURE SITE USE

The City of Franklin has partnered with LRPC to assess the Site so that the City can redevelop the Site into a new Water Department Building.



3. PHASE II SCOPE OF WORK

Crede performed this Phase II ESA to assess the Site considering the anticipated re-use scenario and the identified environmental conditions noted in **Section 2**. An SSQAPP Addendum was developed that outlined the work to be completed, methodologies to be used, and data quality objectives for the project (see **Appendix A**). The Phase II ESA tasks completed included the following:

1. Ground Penetrating Radar (GPR) survey of the Site to identify subsurface features
2. Four (4) surficial soil samples were collected to and submitted for off-site laboratory analysis.
3. Thirteen (13) test pits were excavated at the Site and subsurface soil samples from twelve (12) test pits were collected and submitted for off-site laboratory analysis.
4. Five (5) soil borings were advanced at the Site and subsurface soil samples were collected and submitted for off-site laboratory analysis.
5. Three of the soil borings were completed as monitoring wells (CA-1, CA-2, and CA-3), and groundwater was collected from the new monitoring wells and existing monitoring well MW-2, and submitted for off-site laboratory analysis.
6. Building material/debris samples were collected from the former building foundation slab and from a demolition debris pile at the Site and submitted for off-site laboratory analysis.

Deviations from the scope of work described in the SSQAPP Addendum are summarized in **Section 9** of this report.



4. PHASE II FIELD ACTIVITIES

This sampling program was developed to confirm or dismiss the RECs identified during the Phase I ESA (see SSQAPP Addendum in **Appendix A**) and other potential environmental conditions identified by Credere. All soil, groundwater, and potential PCB-containing building material samples collected by Credere for laboratory analysis were submitted to Absolute Resource Associates of Portsmouth, New Hampshire for analysis. Soil gas samples collected by Credere for laboratory analysis were submitted to Alpha Analytical of Mansfield, Massachusetts. Potential asbestos-containing building materials were submitted to EMSL Analytical, Inc. of Woburn, Massachusetts. Requirements relative to Chain of Custody, Data Management and Documentation, Data Validation, and Data Usability Assessments contained in the SSQAPP were followed. **Figure 2** shows the approximate locations of the building, pertinent Site features, and sample locations.

4.1 GROUND PENETRATING RADAR SURVEY

On June 28, 2011, Credere oversaw a GPR survey of the Site by DigSmart of Maine (DigSmart) to identify subsurface features at the Site potentially associated with environmental conditions at the Site. DigSmart scanned the subsurface with a GPR by walking in continuous and consecutive north-south transects over the areas of the Site associated with RECs (see **Appendix A**). The GPR technician located the lateral extent of subsurface anomalies in areas of concern at the Site. The location of subsurface anomalies was marked on the surface in pink spray paint.

4.2 SURFICIAL SOIL INVESTIGATION

On June 28 and July 5, 2011, Credere collected three (3) surficial soil samples (CA-SS-1, CA-SS-2, and CA-SS-3) to assess possible impacts to the surface from former railroad activities at the Site. The locations of these samples at the Site are depicted on **Figure 2**. The collected samples were submitted for off-site laboratory analysis of PAHs.

All surficial soil samples were collected from 0 to 2 feet below ground surface (bgs) in accordance with standard operating procedures (SOPs) HWRB-11, HWRB-12, HWRB-15, DR#024, DR#025, and Credere-004. Any visible organic debris and/or grass or degraded asphalt was removed from samples prior to placement in laboratory glassware.

Each collected soil sample was logged and visual and/or olfactory evidence of contamination was noted. Samples were then field screened for VOCs with a 10.6 electron volt (eV) photoionization detector (PID) calibrated to a 100 parts per million by volume (ppm_v) isobutylene standard with the instrument response factor set to 1.0. Surficial soil samples were also screened for RCRA 8 metals with an Innov-X X-ray fluorescence (XRF) meter.

Exploration locations and methodologies used are summarized in **Table 1**, PID field screening results are included in **Table 2**, and XRF field screening results are included in **Table 3**.



4.3 TEST PIT INVESTIGATION

On June 28, 2011, Credere oversaw test pit investigation activities which were performed by J.A. Eaton Inspection Services, Inc. of Windham, Maine to assess RECs identified in the previous Phase I ESA and to investigate subsurface anomalies identified during the GPR survey. Please note that six (6) test pits were previously excavated part of a 2008 investigation by Delta Consultants (TP-1 to TP-6). Test pits TP-1 and TP -2 were located off of the Site in the adjacent parking area to the south of the Site. Test pits TP-3 to TP-6 were located on the Site, and are depicted on **Figure 2**. Thirteen (13) test pits were excavated by Credere during this Phase II ESA work, and are identified as CA-TP-7, CA-TP-7A, and CA-TP-8 through CA-TP-18. Soil samples were continuously collected from each test pit. Each collected soil sample was logged and visual and/or olfactory screened for evidence of contamination and was then screened for VOCs with PID and for metals using an XRF as described in **Section 4.2**.

The rationale for each test pit and soil samples submitted for laboratory analysis are provided below. The locations of each test pit are described in **Table 1** and are also depicted on **Figure 2**. Copies of test pit logs are included as **Appendix C**.

CA-TP-7

Test pit CA-TP-7 was excavated along the western side of a closed-in-place UST, believed to be UST-6, which is a 12,000-gallon former fuel oil UST. The top of the UST was located at 4 feet bgs. A ductile iron pipe was located parallel to the UST at 8 feet bgs, which prevented the advancement of the test pit past this depth. No soil sample was submitted for laboratory analysis from this test pit, and a second test pit was excavated to assess soil near the bottom of the UST.

CA-TP-7A

Test pit CA-TP-7A was advanced along the southern side of the closed-in-place UST believed to be UST-6, which is a 12,000-gallon former fuel-oil UST. The top of the UST was located at 4 feet bgs and the bottom of the UST was located at approximately 11 feet bgs. Samples from deeper than 12 feet bgs were collected using a split-spoon sampler driven from a drill rig because this depth was beyond the capacity of the excavator used at the Site (see **Section 4.5**). The 10 to 12 foot bgs soil sample was submitted for laboratory analysis of NHDES *Petroleum and Hazardous Waste Full List of Analytes for Volatile Organics* (NHDES Full List VOCs), total petroleum hydrocarbons (TPH), lead, and PAHs because this sample was just below and adjacent to the bottom of UST-6.

CA-TP-8

Test pit CA-TP-8 was excavated west of the former building in the area where a reported 4,000-gallon fuel oil UST (identified as UST-8) was shown on a historical Sanborn Map. No evidence of a UST was identified in this test pit. No evidence of contamination was observed in this test pit, although the upper 4.5 feet of material was comprised of

building demolition debris, consisting of mostly crushed concrete block, and an older asphalt paving layer was encountered at 4.5 feet bgs. The material beneath the asphalt layer was composed of sand with some gravel. The 6 to 8 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, TPH, and PAHs because this was perceived to be the intersection of fill materials and native material.

CA-TP-9

Test pit CA-TP-9 was excavated beneath the northern end of the building slab where a reported 4,000-gallon fuel oil UST (identified as UST-9) is shown on a historical Sanborn Map. No evidence of a UST was identified during the GPR survey or in this test pit. The test pit location was between a patch in the concrete slab and an area of GPR anomalies beneath the slab. Tannery waste in the form of hides was observed from 2 to 4 feet bgs. Asphalt pieces were encountered at approximately 8 feet bgs. The 6 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, TPH, and PAHs because this was immediately beneath the buried tannery waste.

CA-TP-10

Test pit CA-TP-10 was excavated west of the former building in the area where a reported 4,000-gallon fuel oil UST (identified as UST-8) was shown on a historical Sanborn Map. No evidence of a UST was identified in this test pit. No evidence of contamination was observed in this test pit, although the upper 4.5 feet of material was comprised of building demolition debris, consisting of mostly crushed concrete block. An older asphalt paving layer was present at 4.5 feet bgs. The material beneath the asphalt layer was composed of sand with some gravel. The 0 to 2 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs and total Pb because of elevated metals detected during XRF field screening.

CA-TP-11

Test pit CA-TP-11 was excavated west of the former building in the area where a reported 4,000-gallon fuel oil UST (identified as UST-8) was shown on a historical Sanborn Map. No evidence of a UST was identified in this test pit. No evidence of contamination was observed in this test pit, although the upper 3 feet of material was comprised of building demolition debris, consisting of mostly crushed concrete block, and an older asphalt paving layer was present at 3 feet bgs. The 6.5 to 9 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs and total Pb because of elevated metals detected during XRF field screening.

CA-TP-12

Test pit CA-TP-12 was advanced west of the former building in the suspected location of a former fuel dispenser. A concrete sump extended to 2 feet bgs. A soil sample collected from 3 foot bgs was submitted for laboratory analysis of NHDES Full List VOCs and total Pb because petroleum staining was observed at this depth.



CA-TP-13

Test pit CA-TP-13 was excavated east of the former building in a location where the GPR survey indicated possible waste fill materials. Tannery waste in the form of hides was observed from 4 to 5 feet bgs. However, the 2 to 4 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, RCRA 8 metals, PAHs, and PCBs because this sample was in contact with the buried tannery waste and elevated chromium (Cr) was noted in this sample during XRF field screening.

CA-TP-14

Test pit CA-TP-14 was excavated east of the former building foundation in a location where the GPR survey indicated possible waste fill materials. Tannery waste in the form of hides was observed from 4 to 8 feet bgs. The 6 to 8 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, RCRA 8 metals, and PAHs because this sample was in contact with the buried tannery waste and elevated chromium was noted during XRF field screening. This sample was also analyzed for hexavalent Cr (Cr VI) based on the results of total Cr analysis.

CA-TP-15

Test pit CA-TP-15 was excavated east of the former building foundation to assess possible buried tannery waste. No tannery waste was present in this test pit. The 5 to 7 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, RCRA 8 metals, and PAHs because of the presence of wastes at this depth in test pit CA-TP-14.

CA-TP-16

Test pit CA-TP-16 was excavated west of the former building foundation to assess possible buried tannery waste. No tannery waste was present in this test pit. However, sand and gravel fill was observed in the 0 to 2 foot bgs interval, and native material was observed below this interval. Therefore, the 0 to 2 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, RCRA 8 metals, and PAHs to characterize fill material that did not appear to be associated with tannery waste at the Site.

CA-TP-17

Test pit CA-TP-17 was advanced north of the former building foundation to assess possible buried tannery waste. No tannery waste was present in this test pit. However, the 4 to 7 foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, RCRA 8 metals, and PAHs because of the presence of wastes at this depth in test pits CA-TP-13 and CA-TP-14.

CA-TP-18

Test pit CA-TP-18 was excavated north of the former building foundation to assess possible buried tannery waste. No tannery waste was present in this test pit. The 3 to 5



foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, RCRA 8 metals, and PAHs because of the presence of wastes at this depth in test pits CA-TP-13 and CA-TP-14.

4.4 SUB-SLAB SOIL GAS INVESTIGATION

On June 28, 2011, Credere installed eleven (11) soil gas collection points beneath the concrete slab of the former Site building. Five (5) soil gas collection points (CA-SG-1 to CA-SG-5) were installed beneath the northern end of the slab to assess a suspected former UST (REC-3), and six (6) soil gas collection points (CA-SG-6 to CA-SG-11) were installed beneath the southern end of the slab to assess previously observed organic vapor odors (REC-8) in this area.

A 0.5-inch diameter hole was drilled through the slab and underlying sub-base material in each sampling location. A steel soil gas/pore-water probe was then installed in the drilled hole to a depth of 6 inches beneath the slab bottom. The annulus around the sampler was then sealed with activated bentonite clay and the probe was allowed to equilibrate for one hour. At this time, the internal rod was removed from the soil gas/pore-water sampler and the soil gas was field screened with a ppbRAE Plus PID, which is a low level PID capable of measuring VOCs in air below 1 ppm_v. The PID was connected to the sampler using low-density polyethylene (LDPE) tubing and calibrated to a 10 parts per million by volume (ppm_v) isobutylene standard with the instrument response factor set to 1.0.

One location from each area (CA-SG-1 to CA-SG-5 and CA-SG-6 to CA-SG-11) with the highest field screening VOC concentration was then sampled for laboratory analysis for VOCs using a vacuum sealed Summa canister connected to the sampler using LDPE tubing with an elapsed sampling time of one hour. The Summa canisters were then submitted for laboratory analysis of Toxic Organic Compounds in Ambient Air by EPA Method TO-15.

4.5 SOIL BORING AND MONITORING WELL INSTALLATION

Five (5) soil borings (CA-1 through CA-5) were advanced at the Site on June 29, 2011. Soil boring locations are depicted on **Figure 2**. T&K Well Drilling, a licensed NH well driller, from Troy, NH was retained by Credere and performed the drilling activities. Credere was onsite during the work to provide oversight of the drilling and well installation and to collect soil samples for field screening and laboratory analysis.

Soil borings were advanced using a truck-mounted hollow stem auger drill rig. Soil samples were continuously collected and logged using split-spoon soil sampling. Borings CA-4 and CA-5 were advanced to 10 feet bgs before sampling to assess the soils adjacent to the bottom of the former press pits. Each collected soil sample was field logged and then screened for VOCs with a PID and for metals using an XRF as described in **Section 4.2**.



Monitoring wells were installed in soil boring locations CA-1, CA-2, and CA-3. Each monitoring well was completed with two-inch diameter PVC pipe and 0.010-inch wide slotted PVC screen. The annulus was packed with size No. 2 washed sand to approximately 1-foot above the screen depth and a bentonite seal was placed on top of the annulus and at the ground surface to prevent surface water infiltration. Following installation, each monitoring well was developed and allowed to equilibrate for five days prior to sampling. **Table 1** is a summary of the exploration methods used at each location including the media that was sampled. Field screening results are included in **Table 2** and **Table 3**. Copies of soil boring logs are included as **Appendix D**.

The rationale for the location of each soil boring and the soil sample submitted for laboratory analyses is provided below:

CA-1

Soil boring CA-1 was advanced to a depth of 35 feet bgs in the location of soil gas sampling point CA-SG-2, which had the highest VOC concentration observed during field screening of sub-slab soil gas sampling from locations CA-SG-1 to CA-SG-5. These soil gas sampling points were positioned in an area where a former UST (UST-9) was believed to have been located. The 26 to 28-foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs, TPH, and PAHs because the water table was present within this depth interval, and although an aromatic odor was observed throughout the entire boring depth during advancement of the boring, no PID field screening evidence of contamination was observed during drilling at this location.

Soil boring CA-1 was completed as a monitoring well, screened from 25 to 35 feet bgs to intersect the water table.

CA-2

Soil boring CA-2 was advanced to a depth of 35 feet bgs in an area east of the former Site building where buried tannery waste was observed during test pitting and near existing well MW-2 (the location where a sheen was observed on groundwater in 2008). CA-2 was originally planned to be located downgradient of the suspected septic tank, but the boring/well was moved to generally assess the groundwater flow from the rest of the site when it was discovered that the suspect septic tank was actually part of the Franklin sewer system. The 6 to 8-foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs and RCRA 8 metals plus Cu, Ni, and Zn because VOCs and elevated Cr indicative of tannery waste were detected in this interval during field screening.

Soil boring CA-2 was completed as a monitoring well, screened from 25 to 35 feet bgs to intersect the water table.

CA-3

Soil boring CA-3 was advanced to a depth of 35 feet bgs in the location of soil gas sampling point CA-SG-10, which had the highest VOC concentration observed during field screening of sub-slab soil gas sampling from locations CA-SG-6 to CA-SG-11. These soil gas sampling points were positioned in an area where organic odors had previously been observed emanating from a pipe during the Phase I ESA (REC-8). The 2 to 4-foot bgs soil sample was submitted for laboratory analysis of NHDES Full List VOCs and RCRA 8 metals plus Cu, Ni, and Zn. This sample interval was selected because Cr, Pb, and Hg were identified in the sample during XRF field screening at concentrations which may exceed the NHDES SRS.

Soil boring CA-3 was completed as a monitoring well, and was screened from 25 to 35 feet bgs to intersect the water table.

CA-4 and CA-5

Soil borings CA-4 and CA-5 were advanced immediately adjacent to the former press pits located within the former building footprint. One attempt was made to advance CA-5 within one of the former press pits, which are now filled with demolition debris. However, the bottom of the press pit, located at 10 feet bgs, is composed of reinforced concrete and prevented full penetration of the augers (i.e. refusal). The locations of CA-4 and CA-5 were then moved to be immediately adjacent to the press pits on the downgradient (east) side, and borings were advanced to 10 feet bgs before beginning sampling.

CA-4 and CA-5 were advanced to 10 feet bgs, and then samples were collected continuously from 10 to 14 feet bgs. An aromatic odor was observed in both borings that was similar to the odor observed in boring CA-1; however, no evidence of contamination was identified with PID field screening. The 10 to 12 foot bgs soil sample from CA-4 and the 12 to 14 foot bgs soil sample from CA-5 were submitted for laboratory analysis of NHDES Full List VOCs, TPH, RCRA 8 metals plus Cu, Ni, and Zn, SVOCs, and PCBs to assess soil adjacent to the bottom of the press pits.

CA-TP-7A (Soil Boring Portion)

Samples from 10 to 16 feet bgs from test Pit CA-TP-7A were collected using a split-spoon sampler driven from a drill rig because this depth was beyond the capacity of the excavator used at the Site. The 10 to 12 foot bgs soil sample was submitted for laboratory analysis of NHDES *Petroleum and Hazardous Waste Full List of Analytes for Volatile Organics* (NHDES Full List VOCs), total petroleum hydrocarbons (TPH), total lead (Pb), and polycyclic aromatic hydrocarbons (PAHs) because this sample was adjacent to the end of the tank bottom.



4.6 GROUNDWATER SAMPLING

Credeire returned to the Site on July 5, 2011, to sample the newly installed monitoring wells and previously installed monitoring well MW-2.

Each of the new monitoring wells was surveyed for location and relative groundwater elevations were measured for the new monitoring wells and existing monitoring well MW-2. Based on the results of the survey and measured groundwater elevations, a model of groundwater flow was interpolated.

Depth to groundwater at the Site at the date of sampling was approximately 27 feet bgs, which is at the practical suction limits of Credeire's low-flow peristaltic pump. Credeire therefore opted to deviate from the EPA-approved SSQAPP and collect groundwater by purging the wells with a submersible pump. Samples were then collected with a single use bailer in accordance with NHDES Standard Operating Procedures (HWRB-2, HWRB-4 and HWRB-6) included in Credeire's NHDES/EPA approved generic NH QAPP.

A minimum of three well volumes of groundwater were pumped from each well prior to using the bailer for collecting samples. Field measurements were collected for temperature, pH, conductivity, oxidation-reduction potential, and turbidity at approximate 5-minute intervals.

Because bailers were used to sample the wells, Credeire did not have the proper equipment to field filter the samples for metals from CA-2 and CA-3. Credeire therefore opted to collect unfiltered samples in the HNO₃-preserved laboratory HDPE bottles. These samples were therefore *total* metals, which was a deviation from the specification for *dissolved* metals that was included in the EPA approved SSQAPP. Results revealed that chromium Cr and Pb were present in the CA-2 sample in excess of their Ambient Groundwater Quality Standards (AGQS). As a result, Credeire then returned to the Site on July 14, 2011 and re-sampled well CA-2. The collected sample was filtered in the field through a 0.2 micron filter, and the collected sample was submitted for laboratory analysis of *dissolved* RCRA 8 metals. Groundwater sampling logs are included in **Appendix E**, and groundwater and monitoring well elevation field data are included as **Table 4**.

All collected groundwater samples were submitted to ARA for laboratory analysis. Groundwater sample from location CA-1 was analyzed for NHDES Full List VOCs. The initial sample from location CA-2 was analyzed for NHDES Full List VOCs, SVOCs, and *total* RCRA 8 metals plus Cu, Ni, and Zn. A second sample from CA-2 was analyzed for dissolved RCRA 8 metals. The groundwater sample collected from CA-3 was analyzed for NHDES Full List VOCs and total RCRA 8 metals. The groundwater sample collected from MW-2 was analyzed for NHDES Full List VOCs and total Pb.



4.7 BUILDING MATERIAL DEBRIS SAMPLING

4.7.1 Potential Asbestos Containing Materials

On June 28, 2011, NH-certified Asbestos Inspector Judd Newcomb (Certification # AI 000383) of Credere inventoried suspect asbestos-containing material (ACM) bulk products in building demolition debris and on the former building foundation at the Site. Credere identified four (4) potential ACMs and collected three (3) samples of each material for laboratory analysis for asbestos in accordance with NHDES Env-A 1800. Samples were submitted to EMSL of Woburn, Massachusetts for analysis. Copies of generated laboratory reports are included as **Appendix F**.

The following is a description of the collected potential ACM-containing bulk product samples:

- Potential ACM bulk product sample D-1 was collected from light blue 12 inch-square floor tiles remaining on the former building slab.
- Potential ACM bulk product sample D-2 was collected from white speckled floor tile remaining on the former building slab.
- Potential ACM bulk product sample D-3 was collected from gray floor tile remaining on the former building slab.
- Potential ACM bulk product sample D-4 was collected from off-white floor tile pieces in the debris pile located on the southern end of the former building slab.

4.7.2 Potential PCB-Containing Materials

On June 28, 2011, Credere inventoried suspect PCB-containing bulk products in building demolition debris and on the building foundation at the Site. Examples of suspect products typically include, but are not limited to, paint, caulking, sealants, grout, mastic, glazing, insulation, transformers, capacitors, electrical equipment, used motor/hydraulic oil, fluorescent light ballasts, cable insulation, thermal insulation, adhesives and tapes, plastics, carbonless copy paper, floor finishes, gaskets, ceiling tile coatings, flooring sealants, roofing materials, and siding materials. Consistent with this inventory and the results of previous investigations at similar sites, Credere identified and sampled one (1) potentially PCB-containing bulk product. The collected material was a blue floor coating (sample D-Floor Covering) that was noted coating a portion of the former building foundation. This sample was submitted to ARA for analysis of PCBs.

Additionally, Credere collected a soil sample for an electrical pit (identified as Transformer Pit) and two concrete samples (CA-CC-1 and CA-CC-2) from oil-stained concrete in an area where a transformer may have been located which may have been impacted by possible PCB-containing dielectric fluid released from a former transformer.



5. SUMMARY OF REGULATORY STANDARDS

As a part of this Phase II ESA, Credere collected soil, groundwater, soil gas and building material samples to confirm or dismiss the presence of contaminants associated with the RECs and other potential environmental concerns identified at the Site, and to assess the potential for future risk which may result during anticipated redevelopment. Sample results were compared to the applicable state and federal standards and guidelines described below.

5.1 SOIL

Concentrations in soil samples were compared to New Hampshire's Soil Remediation Standards (SRS) detailed in NHDES Env-Or 600 Contaminated Site Management.

5.2 GROUNDWATER

Concentrations in groundwater samples were compared to New Hampshire's Ambient Groundwater Quality Standards (AGQS) detailed in NHDES Env-Or 600 Contaminated Site Management.

5.3 SOIL GAS

Concentrations of VOCs in sub-slab soil gas were compared to the NHDES Vapor Intrusion Screening Levels for Commercial Soil Gas and the Vapor Intrusion Mitigation Decision Criteria, both updated July 2011. Analytes in soil gas without screening levels were considered to be unregulated.

5.4 POTENTIAL ASBESTOS-CONTAINING MATERIALS

Concentrations of asbestos in building material and debris samples were compared to the limit of 1% by weight detailed in NHDES Env-A 1800 Asbestos Management and Control.

5.5 POTENTIAL PCB-CONTAINING MATERIALS

Bulk products that contain concentrations of total PCBs equal to or in excess of 50 mg/kg are defined as PCB bulk product wastes in accordance with 40 CFR 761.3. These materials are regulated for disposal under 40 CFR 761.62. Bulk products that have been analyzed to contain total PCBs at a concentration of equal to or greater than 1 mg/kg but less than 50 mg/kg (and not as a result of dilution) are not regulated for disposal as long as they remain in use. However, if these materials are removed from use, they must be disposed of at a facility that is licensed to accept these materials in accordance with the applicable state regulations. Bulk products which have been analyzed to contain total PCBs at a concentration of less than 1 mg/kg are unrestricted for future use and/or disposal.



Bulk materials which have been analyzed to contain total PCB concentrations equal to or in excess of 1 mg/kg as a result of contact with a PCB bulk product waste are defined as a PCB remediation waste in accordance with 40 CFR 761.3. These materials must be disposed of in accordance with 40 CFR 761.61 or under a special approval which meets the requirements presented at 40 CFR 761.79(h). Bulk materials which have been analyzed to contain total PCBs at concentrations of less than 1 mg/kg are unrestricted for future use and/or disposal.



6. PHASE II RESULTS

The following subsections present the results of the data collected during the field work portion of this Phase II ESA.

6.1 SITE GEOLOGY

6.1.1 Surficial Geology

During the advancement of soil borings for the current and previous Phase II ESA work, the native surficial geology of the Site was observed to consist of sand deposits with some gravel. Areas of the northern and eastern portions of the Site were observed to contain up to 8 feet or more of fill material, some of which included thin layers of tannery waste in the form of scrap hides (see **Figure 7**).

6.1.2 Bedrock Geology

No bedrock was encountered at the Site during test pit or soil boring activities. According to the Generalized Bedrock Geologic Map of New Hampshire compiled by the United States Geological Survey, the Site is underlain primarily by aluminous schist, calc-silicate granofels, and bimodal metavolcanic rocks of Silurian age.

6.1.3 Site Hydrogeology

Groundwater elevations are based on water levels recorded on July 5, 2011, and are depicted relative to the elevation of the top of monitoring well MW-2 of 334.89 feet based on the elevation of this well used in previous investigations of the Site.

The data show that the groundwater gradient at the time of sampling was almost flat across the Site, with a slight hydraulic gradient of approximately 0.0008 ft/ft, flowing to the northeast towards the Pemigwasset River (**Figure 6**). Please note that the gradient and flow direction were calculated from a limited number of wells. Additional wells may be needed to confirm the predominant groundwater flow direction.

6.2 GROUND PENETRATING RADAR RESULTS

DigSmart identified a potential UST west of the former building foundation. DigSmart also identified subsurface anomalies on the northern and eastern sides of the former building and beneath the northern portion of the foundation slab. These locations were investigated during test pitting activities at the Site.



6.3 SURFICIAL SOIL SAMPLE RESULTS

No visual and/or olfactory evidence of contamination was observed in the field in any of the surficial soil samples collected at the Site. PID readings were non-detect for all surficial soil samples collected. XRF field screening detected Cr in sample CA-SS-2 at a concentration that exceeded the applicable NHDES SRS. No concentrations of any other RCRA 8 metals were detected in surficial soil samples exceeding the applicable NHDES SRS. Surficial soil sample field screening results are summarized in **Table 2** and **Table 3**.

Laboratory analyzed surficial soil sample CA-SS-3 contained the PAH benzo[a]pyrene at a concentration (0.71 mg/kg) which exceeded the applicable NHDES SRS of 0.70 mg/kg. Results for surficial soil samples CA-SS-1 and CA-SS-2 indicated that these samples contained no PAHs above practical quantitation limits (PQLs). Surficial soil sample laboratory results are summarized in **Table 5**. Locations where NHDES SRS were exceeded are depicted on **Figure 4**.

6.4 TEST PIT SAMPLE RESULTS

Thirteen (13) test pits were excavated by Crede during this Phase II ESA work (CA-TP-7, CA-TP-7A, and CA-TP-8 to CA-TP-18). Test Pits TP-1 to TP-6 were previously completed by others. Test pits TP-1 and TP-2 were reportedly located off of the Site in the adjacent paved parking lot on Tannery Street.

Tannery waste in the form of thin layers of hides were observed in test pits CA-TP-9 (2 to 6 feet bgs), CA-TP-13 (4 to 5 feet bgs), and CA-TP-14 (4 to 8 feet bgs). No other visual and/or olfactory evidence of contamination was observed in the field in any of the test pit soil samples collected at the Site. PID readings ranged from non-detect to 2.4 ppm_v for all test pit soil samples collected.

XRF field screening indicated that Cr was present at concentrations exceeding the applicable NHDES SRS of 130 mg/kg for Cr VI in test pits CA-TP-8 (6 to 8 feet bgs), CA-TP-9 (0 to 4 feet bgs), CA-TP-13 (2 to 4 feet bgs), and CA-TP-14 (6 to 8 feet bgs). XRF field screening also indicated that arsenic (As) was present at concentrations exceeding the applicable NHDES SRS of 11 mg/kg in test pits CA-TP-8 (0 to 4 feet bgs), CA-TP-10 (0 to 4 feet bgs), and CA-TP-11 (0 to 2 and 4.5 to 6.5 feet bgs). Test Pit soil sample field screening results are summarized in **Table 2** and **Table 3**.

The laboratory analytical sample collected from test pit CA-TP-14 at 6 to 8 feet bgs contained total Cr at a concentration (260 mg/kg) which exceeded the NHDES SRS of 130 mg/kg for Cr VI. This sample was then analyzed for Cr VI, and was found to contain a concentration of Cr VI (4.5 mg/kg) that was below the applicable NHDES SRS. This indicates that only 1.7% of the Cr in this sample was in the Cr VI state. Therefore, no exceedance of the applicable NHDES SRS was identified.



The laboratory analytical sample collected from test pit CA-TP-15 at 5 to 7 feet bgs contained the PAHs benzo[a]anthracene, benzo[b]fluoranthene, benzo[k]fluoranthene, benzo[a]pyrene, indeno(1,2,3-cd)pyrene, and dibenzo(a,h)anthracene at concentrations which exceeded their applicable NHDES SRS.

No other soil samples from the Site test pits exceeded applicable NHDES SRS for any of the laboratory analyzed parameters. Test pit soil sample laboratory analytical results are summarized in **Table 5**. Locations where NHDES SRS were exceeded are depicted on **Figure 4**.

6.5 SUB-SLAB SOIL GAS SAMPLE RESULTS

No VOCs in soil gas exceeded the applicable vapor intrusion guidelines for commercial soil gas in samples collected from the Site during this Phase II ESA. Soil gas laboratory results are summarized on **Table 7**.

6.6 SOIL BORING SAMPLE RESULTS

As indicated previously, an aromatic odor was observed in all soil samples collected from soil boring CA-1, CA-4, and CA-5. However, no other visual and/or olfactory evidence of contamination was observed in the field in any of the soil boring soil samples collected at the Site. PID readings ranged from non-detect to 10.2 ppm_v for all boring soil samples collected.

XRF field screening indicated that Cr was present at concentrations exceeding the applicable NHDES SRS of 130 mg/kg for Cr VI in soil borings CA-1 (2 to 4 feet bgs), CA-2 (0 to 2 and 6 to 8 feet bgs), and CA-3 (0 to 18, 20 to 28, and 30 to 32 feet bgs). XRF field screening also indicated that Hg was present at concentrations exceeding the applicable NHDES SRS of 6 mg/kg in soil boring CA-3 (0 to 20 and 22 to 32 feet bgs). XRF field screening also indicated that Pb was present at concentrations exceeding the applicable NHDES SRS of 400 mg/kg in soil boring CA-3 (2 to 4 feet bgs). Soil boring sample field screening results are summarized in **Table 2** and **Table 3**.

In contrast to the field screening results, laboratory results for soil samples collected from locations CA-1, CA-2, CA-3, CA-4, and CA-5 contained no contaminants exceeding the applicable NHDES SRS. Soil boring sample laboratory results are summarized on **Table 5**.

6.7 GROUNDWATER SAMPLE RESULTS

Groundwater from well CA-2 was observed to have a petroleum odor during the two times it was sampled (July 5 and 14, 2011). 1,2,4-trimethylbenzene was quantified at 710 µg/L in the groundwater sample collected from monitoring well CA-2 on July 5, 2011 exceeding the applicable NHDES AGQS of 330 µg/L.



Concentrations of *total* Cr and Pb exceeding their respective NHDES AGQS were quantified in the groundwater sample collected from monitoring well CA-2 on July 5, 2011; however concentrations of *dissolved* Cr and Pb collected on July 14, 2011 were below the applicable NHDES AGQS. Therefore, no exceedances of the applicable NHDES AGQS for Cr and Pb were identified.

No other analytes exceeded their applicable NHDES AGQS in groundwater samples collected from the Site during this Phase II ESA. Groundwater sample laboratory results are summarized on **Table 8**.

6.8 BUILDING MATERIAL/DEBRIS RESULTS

6.8.1 Potential Asbestos Containing Material

Light blue 12-inch square floor tiles (Bulk product sample D-1) remaining on the former building slab were found to be ACM (containing 5% chrysotile). These floor tiles are present in one area of the former building floor slab (see **Figure 4**). These tiles are in a highly degraded state. No other samples were analyzed to be ACM. Building material laboratory results are summarized on **Table 6**. The laboratory analytical report is included in **Appendix F**.

6.8.2 Potential PCB-Containing Material

The blue floor coating (sample D-Floor Covering) that covers much of the northern portion of the former building foundation was found to contain no PCBs above the laboratory practical quantitation limit (PQL).

Soil from the electrical pit (“Transformer Pit”) and two concrete samples (CA-CC-1 and CA-CC-2) from oil-stained concrete in an area where a transformer may have been located were found to contain no PCBs above the laboratory PQL.

Building material laboratory results are summarized on **Table 6**. The laboratory analytical report is included in **Appendix F**.



7. QUALITY ANALYSIS/QUALITY CONTROL

The contracted laboratory, Absolute Resource Associates (ARA) of Portsmouth, New Hampshire, provided Level II analytical data according to EPA protocols, EPA laboratory data validation guidance, and the SSQAPP. The laboratories provided the following information in analytical reports:

- Data results sheets
- Method blank results
- Surrogate recoveries and acceptance limits
- Duplicate results/acceptance limits
- Spike/duplicate results/acceptance limits
- Laboratory control sample results
- Description of analytical methods and results
- Other pertinent results/limits as deemed appropriate

As outlined in the SSQAPP, at the completion of the field tasks and upon receipt of the analytical results, a data usability analysis was conducted to document the precision, bias, accuracy, representativeness, comparability, and completeness of the results. The following sections present this analysis.

7.1 PRECISION

Precision measures the reproducibility of measurements. The precision measurement is established using the relative percent difference (RPD) between the sample and duplicate results. Relative percent differences were calculated for soil samples where both sample and duplicate values were greater than five times (5X) the PQL of the analyte. The acceptable upper limit for RPD is 35%. The RPD is calculated as follows:

$$\text{RPD (\%)} = \frac{(\text{Sample Result} - \text{Duplicate Result}) \times 100}{\text{Mean of the Two Results}}$$

The following duplicate samples were collected during this Phase II ESA:

- CA-TP-DUP-1 (duplicate soil sample collected at test pit location CA-TP-14 from 6 to 8 feet bgs) and analyzed for VOCs, PAHs, and RCRA 8 metals plus Cu, Ni, and Zn.
- CA-TP-DUP-2 (duplicate soil sample collected at test pit location CA-TP-9 from 6 feet bgs) and analyzed for TPH.

Table 9 summarizes the duplicate sample results and RPDs.

The duplicate concentrations of barium in samples CA-TP-14/CA-TP-DUP-1 had a RPD of 64.7%, which exceeded the limit of 35%. However, both barium concentrations (47 and 92



mg/kg were below the NHDES SRS of 1,000 mg/kg. Therefore, Credere does not believe that this lack of precision affects the conclusions of this Phase II ESA.

7.2 BIAS

Bias is the systematic or persistent distortion of a measurement process that causes errors in one direction. Bias assessments are made using personnel, equipment, and spiking materials or reference materials as independent as possible from those used in the calibration of the measurement system. Bias assessments were based on the analysis of spiked samples so that the effect of the matrix on recovery is incorporated into the assessment. A documented spiking protocol and consistency in following that protocol are important in obtaining meaningful data quality estimates.

Matrix spike and matrix spike duplicate samples (MS/MSD) were used to assess bias as prescribed in the specified methods. Acceptable recovery values were within the recoveries specified by each of the analysis methods. Control samples for assessing bias were analyzed at a rate as specified in the analytical SOPs and specified analytical methods.

The laboratory provides quality control non-conformance reports that indicate if Laboratory Control Samples/Laboratory Control Sample Duplicates (LCS/LCSD) and/or MS/MSD had low, failing, or high recoveries, and if the sample result was affected. Likewise, the laboratory reports any compounds that had failing RPDs in the LCS/LCSD pair or the MS/MSD pair. This indicates the percent difference between the laboratory sample and its duplicate or the spike and its duplicate. According to the laboratory, unless noted in the non-conformance summary, all of the quality control criteria for these analyses were within acceptable limits. Specific comments from the laboratory included:

VOCs

The MLCS/D4295 did not meet the acceptance criteria for chloromethane. This compound showed high recovery. There is no impact to the data as this analyte was not detected in the associated samples. The MLCS/D4295 did not meet the acceptance criteria for dichlorodifluoromethane. The MLCSD4295 did not meet the acceptance criteria for bromomethane. The recovery was acceptable in the LCS. These compounds are known to be problematic in the method.

The MLCS/D4307 did not meet the acceptance criteria for dichlorodifluoromethane and vinyl chloride. These compounds are known to be problematic in the method.

PCBs

The relative percent difference between the LCS4305 and LCSD4305 was outside the acceptance criteria for PCB-1016 and PCB-1260. The percent recovery for these analytes in each QC parameter was within the acceptance criteria. No impact to the data suspected.



PAHs/SVOCs

The LCS/D4327 did not meet the acceptance criteria for phenol and 4-nitrophenol. These compounds are known to be problematic in the method.

7.3 ACCURACY

Accuracy is a statistical measurement of correctness and includes components of random error (variability due to imprecision) and systemic error. It, therefore, reflects the total error associated with a measurement. A measurement is accurate when the value reported does not differ from the true value or known concentration of the spike or standard. For VOCs and PAHs, surrogate compound recoveries are also used to assess accuracy and method performance for each sample analyzed. Analysis of performance evaluation samples are also used to provide additional information for assessing the accuracy of the analytical data being produced. Both accuracy and precision are calculated for each analytical batch, and the associated sample results are interpreted by considering these specific measurements.

The laboratory provides a non-conformance summary that reports if all of the quality control criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for analysis were within acceptable limits. According to the laboratory, unless noted in the non-conformance summary, all of the quality control criteria for these analyses were within acceptable limits.

7.4 REPRESENTATIVENESS

Objectives for representativeness are defined for each sampling and analysis task and are a function of the investigative objectives. Representativeness was accomplished during this project through use of standard field, sampling, and analytical procedures.

All objectives for sampling and analytical representativeness for samples that were analyzed, as specified in the SSQAPP Addendum, were met.

7.5 COMPARABILITY

Comparability is the confidence with which one data set can be compared to another data set. The objective for this QA/QC program is to produce data with the greatest possible degree of comparability. Comparability was achieved by using standard methods for sampling and analysis, reporting data in standard units, normalizing results to standard conditions and using standard and comprehensive reporting formats. Complete field documentation was used, including standardized data collection forms to support the assessment of comparability. Historical comparability shall be achieved through consistent use of methods and documentation procedures throughout the project.



7.6 COMPLETENESS

Completeness is calculated by comparing the number of samples successfully analyzed to the number of samples collected. The goal for completeness is 95 percent. The completeness for this project was 100 percent, as there were no samples that could not be analyzed due to holding time violations, samples spilled or broken, or any other reason.



8. CONCEPTUAL SITE MODEL

This Phase II ESA was designed to provide further understanding of the contaminants at the Site and to aid in changing the Site use from industrial to recreational. The following section is a description of the Conceptual Site Model (CSM), which incorporates information from this investigation.

8.1 CONTAMINANTS OF CONCERN

Based on the results of this Phase II ESA, the following COCs were identified at the Site:

- 1,2,4-trimethylbenzene in groundwater
- Benzo[a]pyrene in surficial soil
- The PAHs benzo[a]anthracene, benzo[b]fluoranthene, benzo[k]fluoranthene, benzo[a]pyrene, indeno(1,2,3-cd)pyrene, and dibenzo(a,h)anthracene in subsurface soil
- Asbestos in floor tiles on the former building foundation slab

8.2 SITE GEOLOGY AND HYDROGEOLOGY

The Site is located near the west bank of the Pemigewasset River in a developed area of Franklin, New Hampshire. The eastern adjoining property between the Site and river is currently undeveloped. Area topography generally slopes east towards the river. Storm water at the Site likely follows surficial topography and is directed into the river.

Soil boring and test pit data obtained from this Phase II ESA and from reports that were reviewed as a part of the Phase I ESA indicates that soil beneath the Site consists of a mixture of sand and gravel with some areas of fill material that includes tannery waste (hides) on the east side of the Site. The water table was identified at approximately 28 feet bgs.

8.3 DEFINITIONS OF EXPOSURE PATHWAYS AND POTENTIAL RECEPTORS

To aid in a thorough understanding of the environmental concerns present at the Site, a graphical presentation of the identified COCs and potential migration pathways to receptors is included as **Figure 8**. Exposure Pathways and Potential Receptors depicted on the CSM are defined as follows.

Exposure Pathways describe how a human or environmental receptor comes into contact with contaminants which may be present at the Site. Exposure pathways presented in the CSM include the following:



- **Inhalation:** This pathway is primarily associated with groundwater contamination within 30 feet of an occupied structure when groundwater elevation is less than 15 feet below surface grade, or when depth to groundwater is unknown. This pathway is applicable when receptors may inhale impacted media in the form of vapor.
- **Dermal Absorption:** Exposure via dermal absorption occurs when receptors are exposed to chemical concentrations present in soil, groundwater, or surface water through direct contact with the skin.
- **Active Ingestion:** The Active Ingestion pathway represents exposure which may occur through the active ingestion of contaminant concentrations via a drinking water supply well or through agricultural products.
- **Incidental Uptake:** This pathway is applicable when receptors may incidentally ingest or inhale impacted media in the form of dust or airborne particulates.

Potential Receptors are categorized by duration of exposure and intensity of use at the Site. The receptor categories for this project described in the CSM include the following:

- **Commercial:** Commercial receptors are those which are present at the Site for long durations but with low intensity exposure such as indoor office workers.
- **Site Worker:** Site workers are present at the Site for short durations though intensity of use is high, such as during non-routine activities including construction or utility work. Examples include outdoor commercial workers and construction workers.
- **Visitor:** Visitors are characterized by low duration, i.e. less than two hours per day, and low intensity usage such as that which would occur during activities such as walking, shopping, and bird watching.
- **Terrestrial and Aquatic Biota:** These receptors include flora and fauna which may be exposed to contaminants in their respective land-based or aquatic environments.



Please note that the Site is not anticipated to be used for residential purposes. Therefore, residential receptors are not evaluated as part of this CSM.

Based on known history of the Site, the identified COCs may have been released to the environment through surficial and subsurface releases associated with previous industrial usage, filling with tannery waste, the degradation of hazardous building materials, and previous bulk petroleum storage. Primary impacted media at the Site include surficial soil, subsurface soil, and groundwater.

The PAH benzo[a]pyrene in surficial soil and asbestos in exposed floor tiles may migrate through aeolian dispersion and impact on-site or off-site receptors. Both asbestos and PAHs have low propensity to leach and are unlikely to affect subsurface conditions through infiltration and leaching. Potential receptors identified for the Site include future commercial workers, site workers (i.e. construction workers, etc.), and visitors, as well as terrestrial biota. Potential exposure pathways for these human receptors to PAHs or asbestos include incidental uptake (including inhalation). Receptors may also be exposed to PAHs via dermal absorption.

The PAHs in subsurface soil are unlikely to migrate unless disturbed during future Site work. Human receptors identified for the Site include future site workers (i.e. construction workers, etc.) who could be exposed via dermal absorption or incidental ingestion (including inhalation).

The 1,2,4-trimethylbenzene identified in groundwater at the Site has likely leached and infiltrated from an upgradient on-site subsurface soil source area. This potential source area was not defined by this Phase II ESA. Human receptors identified for the Site include future site workers (i.e. construction workers, etc.) who could be exposed via dermal absorption or incidental ingestion (including inhalation).



9. DEVIATIONS

The following deviations were made from the SSQAPP Addendum (see **Appendix A**) during the course of the investigation:

1. Soil boring CA-2 and test pit CA-TP-13 were moved from their proposed locations in the SSQAPP. These sampling locations were proposed to assess a possible septic tank which was suspected of discharging to the adjacent Pemigwasset River. Based on field observations and information obtained from City of Franklin public works employee, it was determined that this tank was connected to the City sanitary sewer system. Based on this information, Credere determined that this tank did not require assessment, and soil boring CA-2 and test pit CA-TP-13 were moved approximately 30 to 40 feet to the southwest to assess possible buried tannery waste.
2. Soil borings CA-4 and CA-5 were not installed through the bottoms of the former press pits, because the drill rig augers could not penetrate the reinforced concrete at the bottom. Instead, these soil borings were moved several feet to the east and were advanced immediately adjacent to the sidewalls of the press pits.
3. Due to the depth to groundwater at the Site, Credere did not have the proper equipment to field filter metals samples from CA-2 and CA-3 when they were first sampled on July 5, 2011. Credere therefore opted to collect unfiltered samples in the HNO₃-preserved laboratory HDPE bottles. These samples therefore represented *total* metals samples which deviated from the specification in Credere's NHDES and EPA approved SSQAPP. The SSQAPP called for the collection of samples for *dissolved* metals. To meet the SSQAPP requirement, Credere returned to the Site on July 14, 2011 and re-sampled monitoring well CA-2, and the sample was field filtered prior to placing it in appropriate glassware. It is worth noting that monitoring well CA-3 was not re-sampled because no contaminants were detected above AGQS in this well.



10. DATA GAPS

The source of gasoline contamination of groundwater detected at monitoring well CA-2 has not been identified, and the extent of groundwater contamination has not been fully delineated. No other data gaps have been identified for the Site.



11. CONCLUSIONS

Our conclusions in relation to the identified RECs and other environmental concerns, and the investigation results are presented below:

11.1 RECOGNIZED ENVIRONMENTAL CONDITIONS

- REC-1, which was associated the presence of a 12,000-gallon fuel oil underground storage tank (UST) (identified as UST-6) that was closed-in-place in the vicinity of the loading dock, is dismissed because no evidence of contamination was identified in soil samples collected from the sides and adjacent to the bottom of the closed-in-place UST in test pits CA-TP-7 and CA-TP-7A. In addition, observations made by Credere indicated that this UST was filled with clean sand at the time of closure-in-place.
- REC-2, which was associated with the potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-8) shown on a historical Sanborn Map, is dismissed because no UST was identified in this location during the GPR survey and no evidence of contamination exceeding applicable standards was identified in test pits CA-TP-8, CA-TP-10, or CA-TP-11.
- REC-3, which was associated with the potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-9) shown on a historical Sanborn Map, is dismissed because no UST was identified in this location during the GPR survey, and no evidence of contamination was identified in test pit CA-TP-9 or in soil gas sample locations CA-SG-1 through CA-SG-5, which were completed to evaluate this REC.
- REC-4, which was associated with a former fuel dispenser and a number of filling caps formerly located in the vicinity of the closed-in-place UST-6, is dismissed because no evidence of contamination exceeding applicable standards was identified in soil samples collected from test pit CA-TP-12, which was advanced in the suspected location a former fuel dispenser.
- REC-5, which was associated with the presence of a sheen of degraded gasoline previously noted in 2008 in existing monitoring well MW-2, is confirmed, because groundwater from a newly installed monitoring well (CA-2) located 30 feet southwest of MW-2 contained 1,2,4-trimethylbenzene at a concentration exceeding the applicable NHDES Ambient Groundwater Quality Standard (AGQS), suggesting that a plume of degraded gasoline is present beneath the Site.
- REC-6, which was associated with a subsurface structure identified during a previous GPR survey east of the former building and believed to be a former concrete septic tank, is dismissed because this structure was determined to be a part of the municipal sewer system and not a septic tank, and therefore does not represent a pathway to the subsurface environment.



- REC-7, which was associated with the presence of buried waste related to the previous usage of the Site as a tannery, is dismissed. While evidence of buried waste was observed in the form of leather hides in test pits CA-TP-9, CA-TP-13, and CA-TP-14, no evidence of contamination at concentrations exceeding applicable standards was observed in association with the buried waste. However, this condition represents a solid waste issue that should be properly managed.
- REC-8, which was associated with pipes that were previously noted emitting unknown solvent odors and elevated total volatile organic readings in the southwestern portion of the building, is dismissed, because no evidence of contamination exceeding applicable standards was identified in soil gas (soil gas sample locations CA-SG-6 through CA-SG-11), soil (boring CA-3), or in groundwater (monitoring well CA-3) samples collected from this area.
- REC-9, which was associated with former press pits located within the manufacturing slab that are stained with petroleum, is dismissed because no evidence of contamination was identified in soil samples collected from adjacent to the bottoms of the press pits (borings CA-4 and CA-5).
- REC-10, which was associated with a pile of demolition debris located on the southwestern portion of the Site, is dismissed because no potential lead-based paint or PCB-containing materials were observed in the pile, and no asbestos was identified in a floor tile sample collected from the pile.

11.2 ADDITIONAL ENVIRONMENTAL CONCERNS

The following additional environmental concerns were identified by Credere during development of the SSQAPP and assessed during this Phase II ESA:

- Potential impact associated with the former use of the Site for railroad activities is confirmed, because benzo[a]pyrene was detected above the applicable NHDES SRS in surficial soil sample CA-SS-3. Benzo[a]pyrene is a PAH which is known to be associated with railroad activities.
- The potential impact of a former electrical transformer located on the west side of the building is dismissed, because no PCBs were identified in soil collected from an electrical pit in this area of the Site. Additionally, two samples of stained concrete collected from this area of the Site contained no PCBs above laboratory practical quantification limits.
- A soil sample collected from test pit CA-TP-15 at 5 to 7 feet below ground surface (bgs) contained the PAHs benzo[a]anthracene, benzo[b]fluoranthene, benzo[k]fluoranthene, benzo[a]pyrene, indeno(1,2,3-cd)pyrene, and dibenzo(a,h)anthracene at concentrations which exceeded their applicable NHDES SRS. No staining was observed in this test pit nor was this sample associated with any observed buried wastes. Additionally, no PAHs



were identified in soil in contact with buried waste in test pits CA-TP-13 or CA-TP-14. Although this test pit was not in the location of or downgradient of any identified Site USTs, it is Credere's opinion that this contamination may be associated with a past petroleum release from the former UST#10, which was located approximately 40 feet away or from another unknown on-site source.

- Building material sample results indicate that blue 12-inch square floor tile remnants remaining attached to the concrete building slab are asbestos-containing material (ACM). It is not clear if any of these tiles are present in the demolition pile remaining on-site.



12. RECOMMENDATIONS

Based on observations and results of this Phase II ESA for the Site, Credere makes the following recommendations:

- Credere recommends that delineation of all buried solid waste at the Site be conducted, including tannery hides, other tannery wastes, and building demolition debris, per NHDES Solid Waste Rules Env-Sw 309 in support of a Pre-1981 landfill registration.
- Subsequent to the delineation of all buried solid waste at the Site, Credere recommends that the Site be registered as a Pre-1981 landfill according to NHDES Solid Waste Rules Env-Sw 309.
- Credere recommends that delineation of the identified PAH contaminated surficial soil in the areas of CA-SS-3 and CA-TP-15 be conducted. Credere recommends that a Remedial Action Plan (RAP) be developed to address the PAH contaminated soil.
- Credere recommends additional investigation to assess the full extent and magnitude of degraded gasoline contamination identified at monitoring well CA-2. This additional assessment should supplement the development of a groundwater flow model of the Site in order to define the nature, extent, and/or potential fate and transport characteristics of the identified impacted media. Credere also recommends that this petroleum contamination issue be referred to the NHDES Petroleum Remediation Program as the observed contamination is likely associated with one of the former USTs at the Site. If accepted into this program, future investigation work will be reimbursed by the program.
- Credere recommends that the identified ACM floor tile remaining on the former building slab be properly handled according to all applicable NHDES Asbestos and Solid Waste Rules prior to the development of the Site.

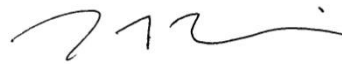


13. SIGNATURES OF ENVIRONMENTAL PROFESSIONALS

The following Environmental Professionals performed this Phase II ESA in conformance with ASTM Standard Guide E 1903-97 (reapproved 2002). The following individual(s) meet the qualifications for individuals completing or overseeing all appropriate inquiries, and possess sufficient specific education, training, and experience necessary to exercise professional judgment to develop opinions and conclusions regarding the existence of environmental conditions on the Site. Any work completed on this Phase II ESA by an individual who is not considered an environmental professional was completed under the supervision or responsible charge of the environmental professional.



Jonathan O'Donnell
Geo-Environmental Engineer



Jedd Steinglass
Senior Geologist



Richard S. Vandenberg, PG
Senior Geologist



14. LIMITATIONS

This report has been prepared as part of a contract agreement between Credere Associates, LLC and LRPC for their Brownfields program. This agreement was established in order to provide LRPC with information upon which it can rely concerning the existence or likely existence of various environmental contaminants on or adjacent to the Site.

This report does not reflect:

1. Conditions in untested areas.
2. Variations in chemical concentrations that can occur between sample locations.
3. The total understanding of potential influences of off-site areas or historical uses that may have contributed or currently contribute to Site contamination, particularly relating to groundwater and subsurface soil conditions. The limited evaluation of off-site contamination sources was based on available data and records.
4. The potential presence of compound sources was based on available data and records.
5. The potential presence of analytes that were not analyzed for or that may be present below minimum Practical Quantification Limits for the methods tested.
6. The conditions of groundwater and/or surface water beyond available data.
7. Variation in the Site conditions that occurred at a time other than when the Site inspection was completed.

In the event that any conditions different from those described herein are encountered at a later time, Credere Associates, LLC requests an opportunity to review such differences and modify the assessment and conclusions of this report. This report was prepared expressly for the purpose described. The information in this report may not be suitable for any other use without adaptation for the specific purpose intended. Any such reuse of this report, without adaptation, shall be at the sole risk and liability of the party undertaking the reuse.



TABLES



**TABLE 1
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF EXPLORATION LOCATIONS AND SAMPLING METHODS**

Location Name	Media Sampled	Type of Exploration	Location	Sampling Method			
CA-SG-1	Soil Gas		Beneath northern end of foundation slab.	Summa sampling canister via 1/4 inch stainless steel probe through floor slab, sealed in place with bentonite			
CA-SG-2							
CA-SG-3							
CA-SG-4							
CA-SG-5							
CA-SG-6							
CA-SG-7							
CA-SG-8							
CA-SG-9							
CA-SG-10							
CA-SG-11							
CA-TP-7	Subsurface Soil	Test Pit	West side of closed-in-place UST-6	Excavator			
CA-TP-7A		Test Pit/Soil Boring	South side of closed-in-place UST-6	Excavator, split-spoon sampler			
CA-TP-8		Test Pit	West of slab loading dock area in suspected location of UST-8.	Excavator			
CA-TP-9			Under northern portion of foundation near suspected location of UST-9.				
CA-TP-10			West of slab loading dock area in suspected location of UST-8.				
CA-TP-11			West of slab loading dock area in suspected location of UST-8.				
CA-TP-12			West of slab, beneath fuel dispenser sump.				
CA-TP-13			East of slab, subsurface anomaly detected by Ground Penetrating Radar.				
CA-TP-14			East of slab, subsurface anomaly detected by Ground Penetrating Radar.				
CA-TP-15			East of slab.				
CA-TP-16			West of slab.				
CA-TP-17			North of slab.				
CA-TP-18			North of slab.				
Transformer Pit			Surficial Soil		Surficial Soil Sample	West of slab, loading dock area, grab sample from soil at pit bottom.	Pre-cleaned shovel
CA-SS-1						Co-located with CA-TP-17.	
CA-SS-2						Co-located with CA-TP-14.	
CA-SS-3						South of slab,	
CA-1			Subsurface Soil and Groundwater		Soil Boring & Monitoring Well	Co-located with CA-SG-2.	Split-spoon sampler. Purge groundwater sampling using submersible pump.
CA-2	East of slab, buried tannery waste area.						
CA-3	Co-located with CA-SG-10.						
CA-4	Subsurface Soil	Soil Boring	Immediately east of press pits.				
CA-5			Immediately east of press pits.				
MW-2	Groundwater	Monitoring Well	East of slab, buried tannery waste area.	Purge groundwater sampling using submersible pump.			
CC-1	Building Material/Debris		Stained concrete in area of possible transformer	Single-use drill bit and scupula			
CC-2			On northern end of slab.	Hand.			
D-Floor Covering				On central portion of slab.	Hand		
D-1					Hand		
D-2					Hand		
D-3			From debris pile, southern end of slab.	Hand			
D-4	Hand						

TABLE 2
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF PHOTOIONIZATION DETECTOR (PID) FIELD SCREENING RESULTS

Location	Sample Depth (feet bgs)	Sample Date	PID Results (ppm.)	Evidence of Contaminated or Petroleum Impacted Soils
Test Pit Sample				
CA-TP-7	0-2	06/28/11	ND	No visual evidence observed.
	3-5		ND	
	5-7		ND	
	7-8		ND	
CA-TP-7A	4	06/30/11	0.5	No visual evidence observed.
	6		1.4	
	8		1.3	
	9		1.4	
	9.5		1.3	
	10-12		1.9	
	12-14		1.1	
	14-16		0.4	
CA-TP-8	0-2	06/28/11	1.2	No visual evidence observed.
	2-4		ND	
	4.5-6		1.1	
	6-8		ND	
	8-9		ND	
CA-TP-9	0	06/30/11	0.7	No visual evidence observed.
	2		1.3	
	4		1.8	
	6		2.4	
	8		2.0	
CA-TP-10	0-2	06/28/11	ND	No visual evidence observed.
	2-4		ND	
	4.5-6		ND	
	6-8		ND	
	8-9		ND	
CA-TP-11	0-2	06/28/11	ND	No visual evidence observed.
	3-4.5		ND	
	4.5-6.5		ND	
	6.5-9		ND	
CA-TP-12	2 (In sump)	06/30/11	ND	In sump. Wet black organic material. No odor.
	2 (below concrete)		0.7	
	3		1.2	
	4		0.5	
CA-TP-13	0-2	06/28/11	ND	No visual evidence observed.
	2-4		ND	
	4-5		ND	
	6-8		ND	
CA-TP-14	2-4	06/28/11	ND	No visual evidence observed.
	4-6		ND	
	6-8		ND	
CA-TP-15	0-2	06/28/11	ND	No visual evidence observed.
	2-4		ND	
	5-7		ND	
CA-TP-16	0-2	06/28/11	ND	No visual evidence observed.
	2-4		ND	
	5-7		ND	
CA-TP-17	2-4	06/28/11	ND	No visual evidence observed.
	4-7		ND	
CA-TP-18	0-1.5	06/28/11	ND	No visual evidence observed.
	1.5-3		ND	
	3-5		ND	

TABLE 2
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF PHOTOIONIZATION DETECTOR (PID) FIELD SCREENING RESULTS

Location	Sample Depth (feet bgs)	Sample Date	PID Results (ppm _v)	Evidence of Contaminated or Petroleum Impacted Soils
Surficial Soil Samples				
CA-SS-1	0-2	06/28/11	ND	No visual evidence observed.
CA-SS-2	0-2	06/28/11	ND	No visual evidence observed.
CA-SS-3	0-2	07/05/11	ND	No visual evidence observed.
Transformer Pit	Pit Bottom	06/28/11	ND	No visual evidence observed.
Subsurface Soil Samples				
CA-1	0-2	06/29/11	ND	No visual evidence observed. Slight aromatic odor observed.
	2-4		ND	
	4-6		ND	
	6-8		ND	
	8-10		ND	
	10-12		ND	
	12-14		ND	
	14-16		ND	
	16-18		ND	
	18-20		ND	
	20-22		ND	
	22-24		ND	
	24-26		ND	
CA-2	0-2	06/29/11	ND	No visual evidence observed.
	2-4		ND	
	4-6		ND	
	6-8		10.2	
	8-10		7.9	
	10-12		1.2	
	12-14		ND	
	14-16		ND	
	16-18		ND	
	18-20		ND	
	20-22		ND	
	22-24		ND	
	24-26		ND	
CA-3	0-2	06/29/11	ND	No visual evidence observed.
	2-4		ND	
	4-6		ND	
	6-8		1.2	
	8-10		ND	
	10-12		ND	
	12-14		ND	
	14-16		ND	
	16-18		ND	
	18-20		ND	
	20-22		ND	
	22-24		ND	
	24-26		ND	
CA-4	10-12	06/29/11	NS	No visual evidence observed. Slight aromatic odor observed.
CA-5	12-14	06/29/11	ND	No visual evidence observed. Slight aromatic odor observed.
	10-12		ND	
	12-14		ND	

Notes:

Samples were field screened using a Thermo OVM 580B PID; the PID was calibrated using 100 ppm isobutylene and a response factor of 1.0.

ND - VOCs not detected with PID

NS - no sample from this depth

ppm_v - parts per million by volume

bgs - below ground surface

TABLE 3
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF X-RAY FLUORESCENCE (XRF) METER FIELD SCREENING RESULTS

Location	Sample Depth (feet bgs)	Sample Date	NHDES Soil Remediation Standard and Metal Concentration (mg/kg)											
			Cr	Ni	Cu	Zn	As	Se	Ag	Cd	Ba	Hg	Pb	
			130*	400	NE	1,000	11	180	89	33	1,000	6	400	
Surficial Soil Samples														
CA-SS-1	0-2	6/28/2011	<LOD	<LOD	<LOD	23	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	21	
CA-SS-2	0-2	6/28/2011	260	<LOD	<LOD	51	<LOD	<LOD	<LOD	<LOD	483	<LOD	24	
CA-SS-3	0-2	7/5/2011	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	48	
Subsurface Soil Samples														
CA-1	0-2	6/29/2011	112	<LOD	<LOD	19	<LOD	4	<LOD	<LOD	<LOD	<LOD	30	
	2-4		136	<LOD	<LOD	20	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	17	
	4-6		<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	21	
	6-8		<LOD	<LOD	24	13	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	20	
	8-10		<LOD	<LOD	<LOD	19	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	16	
	10-12		<LOD	40	<LOD	14	<LOD	<LOD	<LOD	<LOD	321	<LOD	21	
	12-14		<LOD	<LOD	<LOD	12	8	<LOD	<LOD	<LOD	<LOD	<LOD	14	
	14-16		<LOD	<LOD	<LOD	11	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	19	
	16-18		<LOD	<LOD	<LOD	12	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	23	
	18-20		<LOD	<LOD	<LOD	17	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	21	
	20-22		<LOD	<LOD	<LOD	13	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	18	
22-24	<LOD	<LOD	<LOD	15	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	20			
24-26	<LOD	<LOD	<LOD	14	<LOD	<LOD	<LOD	<LOD	<LOD	365	<LOD	18		
26-28	<LOD	<LOD	<LOD	20	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	20		
CA-2	0-2	6/29/2011	148	<LOD	<LOD	37	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	28	
	2-4		<LOD	<LOD	<LOD	29	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	25	
	4-6		<LOD	<LOD	<LOD	35	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	20	
	6-8		151	<LOD	<LOD	38	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	20	
	8-10		<LOD	<LOD	<LOD	24	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	20	
	10-12		<LOD	<LOD	<LOD	19	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	17	
	12-14		<LOD	<LOD	<LOD	19	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	27	
	14-16		<LOD	<LOD	<LOD	24	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	19	
	16-18		<LOD	<LOD	<LOD	31	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	17	
	18-20		<LOD	<LOD	<LOD	19	<LOD	<LOD	<LOD	<LOD	<LOD	380	<LOD	9
	20-22		<LOD	<LOD	<LOD	35	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	19	
22-24	<LOD	<LOD	<LOD	26	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	16			
24-26	114	<LOD	30	38	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	23			
26-28	<LOD	<LOD	<LOD	25	<LOD	<LOD	<LOD	<LOD	<LOD	421	<LOD	16		
28-30	<LOD	<LOD	<LOD	22	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	15		
CA-3	0-2	6/29/2011	472	<LOD	<LOD	10	<LOD	5	<LOD	<LOD	<LOD	18	185	
	2-4		423	<LOD	<LOD	22	<LOD	5	<LOD	<LOD	<LOD	25	401	
	4-6		164	<LOD	<LOD	13	<LOD	<LOD	<LOD	<LOD	<LOD	19	199	
	6-8		248	<LOD	<LOD	18	<LOD	<LOD	<LOD	<LOD	<LOD	16	215	
	8-10		461	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	14	123	
	10-12		518	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	14	146	
	12-14		647	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	17	127	
	14-16		674	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	15	108	
	16-18		187	<LOD	<LOD	13	<LOD	<LOD	<LOD	<LOD	<LOD	10	119	
	18-20		105	<LOD	<LOD	9	<LOD	<LOD	<LOD	<LOD	<LOD	15	106	
	20-22		173	<LOD	<LOD	10	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	101	
	22-24		434	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	13	106	
	24-26		330	<LOD	<LOD	10	<LOD	<LOD	<LOD	<LOD	<LOD	14	105	
26-28	280	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	9	88			
28-30	89	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	14	85			
30-32	150	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	8	85			
CA-4	10-12	6/29/2011	<LOD	<LOD	23	41	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	19	
	12-14		<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	25	
CA-5	10-12	6/29/2011	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	22	
	12-14		<LOD	<LOD	<LOD	34	<LOD	<LOD	<LOD	<LOD	<LOD	<LOD	27	

<LOD - Concentration less than instrument level of detection

Highlighted cells have concentrations that exceed NHDES Soil Remediation Standards

* = The regulatory threshold for Cr VI was used because it is the most stringent standard for chromium

XRF - X-Ray Fluorescence Detector

bgs - below ground surface

**TABLE 4
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF GROUNDWATER ELEVATION DATA**

MONITORING WELL ID		CA-1	CA-2	CA-3	MW-2
Pertinent Well Construction Details	WELL DEPTH ⁽¹⁾ (FEET)	34.42	34.41	34.57	33.96
	LENGTH OF SCREEN (FEET)	10	10	10	10
	DEPTH TO TOP OF SCREEN ⁽¹⁾ (FEET)	25	25	25	25
	TOP OF WELL ELEVATION ⁽²⁾ (FEET)	335.26	334.82	335.25	334.89
GROUNDWATER DEPTH ⁽¹⁾ (FEET) July 5, 2011		28.02	27.57	27.75	27.64
GROUNDWATER ELEVATION ⁽³⁾ (FEET) July 5, 2011		307.24	307.25	307.50	307.25

Notes:

- (1) All depth levels were gauged from top of well PVC riser. Well depth measurements were taken at the time of groundwater sampling, and may reflect loss of well depth due to sedimentation.
- (2) Top of Well Elevations were related, via stadia survey, to elevation of existing well MW-2 (334.89 ft) from Delta Phase II ESA, April 10, 2008. Elevations of wells CA-1 through CA-3 were surveyed by Credere on July 5, 2011.
- (3) Groundwater elevations have been calculated by subtracting the depth to groundwater from the well elevations.

TABLE 6
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF ANALYTICAL RESULTS FOR BUILDING MATERIAL SAMPLES

Parameter	Regulatory Standard	Building Material Sample Location, Date, and Depth (feet)						
	NHDES SRS ⁽¹⁾ (mg/kg) or Standard for ACM ⁽²⁾ (% by volume)	CA-CC-1	CA-CC-2	D-1	D-2	D-3	D-4	D-Floor Covering
		6/30/2011	6/30/2011	6/28/2011	6/28/2011	6/28/2011	6/28/2011	6/28/2011
		0	0	0	0	0	0	0
Asbestos Analysis of Bulk Materials (% by Volume) via EPA 600/R-93/116 Method using Polarized Light Microscopy								
Asbestos	1	NS	NS	5	ND < 1	ND < 1	ND < 1	NS
Polychlorinated Biphenyls (mg/kg) EPA Method 8082								
Total PCBs	1	ND < 0.6	ND < 0.6	NS	NS	NS	NS	ND < 0.03

NOTES:

⁽¹⁾ New Hampshire Soil Remediation Standards from the Risk Characterization Management Policy Env-Or 606.19, Soil Remediation Criteria.

⁽²⁾ NHDES Env-A 1800

ND < 0.1 = Not detected above quantitation limit (i.e. 0.1 mg/kg)

NS = Not sampled for the analysis.

Bold Exceeds laboratory quantitation limit

Exceeds NH DES Soil Remediation Standards.

**TABLE 7
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF ANALYTICAL RESULTS FOR SOIL GAS
SAMPLES**

Parameter	Regulatory Standard	Sample Location, Date, and Depth (feet)	
	Commercial Soil Gas, Vapor Intrusion Screening Levels ⁽²⁾ ($\mu\text{g}/\text{m}^3$)	CA-SG-2	CA-SG-10
		6/28/2011	6/28/2011
		1	1
⁽¹⁾Volatile Organics in Air ($\mu\text{g}/\text{m}^3$), Method TO-15			
Dichlorodifluoromethane	NE	235	16
Ethanol	NE	7050	5760
Acetone	NE	651	211
Trichlorofluoromethane	NE	23.6	106
Isopropanol	NE	381	295
Methylene chloride	NE	57.3	ND < 34.7
2-Butanone	NE	30.1	32.7
Tetrahydrofuran	NE	14.2	16.4
n-Hexane	NE	30.9	25.3
Benzene	95	28.4	30
Cyclohexane	NE	7.64	7.71
Heptane	NE	15.8	19.1
Toluene	73,000	154	177
Ethylbenzene	250	8.69	17.2
Total Xylene	1,500	83.1	69.4
1,2,4-Trimethylbenzene	220	12.7	ND < 9.83

NOTES:

⁽¹⁾ Only analytes identified above detection limit are summarized .

⁽²⁾ NHDES Vapor Intrusion Guidance Document, July 5, 2011

NE = No regulatory guideline established

ND < 0.2 = Not detected above quantitation limit (i.e. 0.2 mg/kg)

NS = Not sampled for the analysis.

Bold Exceeds laboratory quantitation limit

TABLE 8
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF ANALYTICAL RESULTS FOR GROUNDWATER SAMPLES

Parameter	Regulatory Standard	Sample Location, Date, and Concentration (µg/L)				
	NH AGQS ⁽²⁾ (µg/L)	CA-1	CA-2		CA-3	MW-2
		7/5/2011	7/5/2011	7/14/2011*	7/5/2011	7/5/2011
(1) Volatile Organic Compounds (ug/L) EPA Method 8260B						
Propylbenzene, n-	260	ND < 2	29	NS	ND < 2	ND < 2
Trimethylbenzene, 1,3,5-	330	4	320	NS	ND < 2	ND < 2
Trimethylbenzene, 1,2,4-	330	9	710	NS	ND < 2	ND < 2
Butylbenzene, sec-	260	ND < 2	43	NS	ND < 2	ND < 2
Isopropyltoluene, p-	260	ND < 2	100	NS	ND < 2	ND < 2
(1) Semi-Volatile Organic Compounds(µg/L) EPA Method 8270C						
All Compounds	Various	NS	All ND	NS	NS	NS
Metals SW3051A (ug/L)						
Arsenic	10	NS	21	ND < 8*	ND < 8	NS
Barium	2,000	NS	130	ND < 50*	ND < 50	NS
Cadmium	5	NS	ND < 4	ND < 4*	ND < 4	NS
Chromium	100	NS	130	ND < 50*	ND < 50	NS
Copper	1,300	NS	230	NS	NS	NS
Lead	15	NS	45	ND < 8*	ND < 8	8
Mercury	2	NS	ND < 2	ND < 2*	ND < 2	NS
Nickel	100	NS	ND < 50	NS	NS	NS
Selenium	50	NS	ND < 50	ND < 50*	ND < 50	NS
Silver	100	NS	ND < 7	ND < 7*	ND < 7	NS
Zinc	NE	NS	120	NS	NS	NS

NOTES:

⁽¹⁾ Only analytes above detection level are summarized.

⁽²⁾ New Hampshire Code of Administrative Rules Ambient Groundwater Quality Standards (AGQS), effective July 23, 2008.

* - Sample collected on July 14, 2011 was field filtered and is for dissolved metals. All samples collected on July 5, 2011 were not field filtered and are for total metals.

NE = No regulatory guideline established.

ND < 0.2 = Not detected above quantitation limit (i.e. 0.2 ug/L).

NS = Not Sampled.

Bold Exceeds laboratory quantitation limit.

Exceeds NHDES AGQS

Exceeds NHDES AGQS in unfiltered total metals sample but NOT in filtered dissolved metals sample.

**TABLE 9
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET - FRANKLIN NEW HAMPSHIRE
NHDES SITE NO. 199902062
SUMMARY OF DUPLICATE SAMPLE ANALYSES**

Parameter	NHDES Threshold ⁽¹⁾	Quantitation Limit (ug/g)	5x Quantitation Limit	Sample ⁽²⁾	Duplicate	Relative Percent Difference
VOCs						
QA/QC (CA-TP-DUP-1, 06/28/11); duplicate of CA-TP-14 6-8'						
All parameters non-detect.						
PAHs						
QA/QC (CA-TP-DUP-1, 06/28/11); duplicate of CA-TP-14 6-8'						
All parameters non-detect.						
Metals						
QA/QC (CA-TP-DUP-1, 06/28/11); duplicate of CA-TP-14 6-8'						
Barium	1,000	3.0	15.0	47	92	-64.7%
Chromium	130	3.0	15.0	260	300	-14.3%
Copper	NE	3.0	15.0	14	16	-13.3%
Lead	400	0.7	3.5	6.2	7	-12.1%
Zinc	1,000	3.0	15.0	17	20	-16.2%
TPH						
QA/QC (CA-TP-DUP-2, 06/30/11); duplicate of CA-TP-9 6'						
All parameters non-detect.						

NOTES:

⁽¹⁾ New Hampshire Soil Remediation Standards from the Risk Characterization Management Policy Env-Or 606.19, Soil Remediation Criteria and Env-Or 603.3 Ambient Groundwater Quality Standards.

⁽²⁾ Only analytes above detection level and five times the quantitation limit are summarized herein.

NA - Not applicable

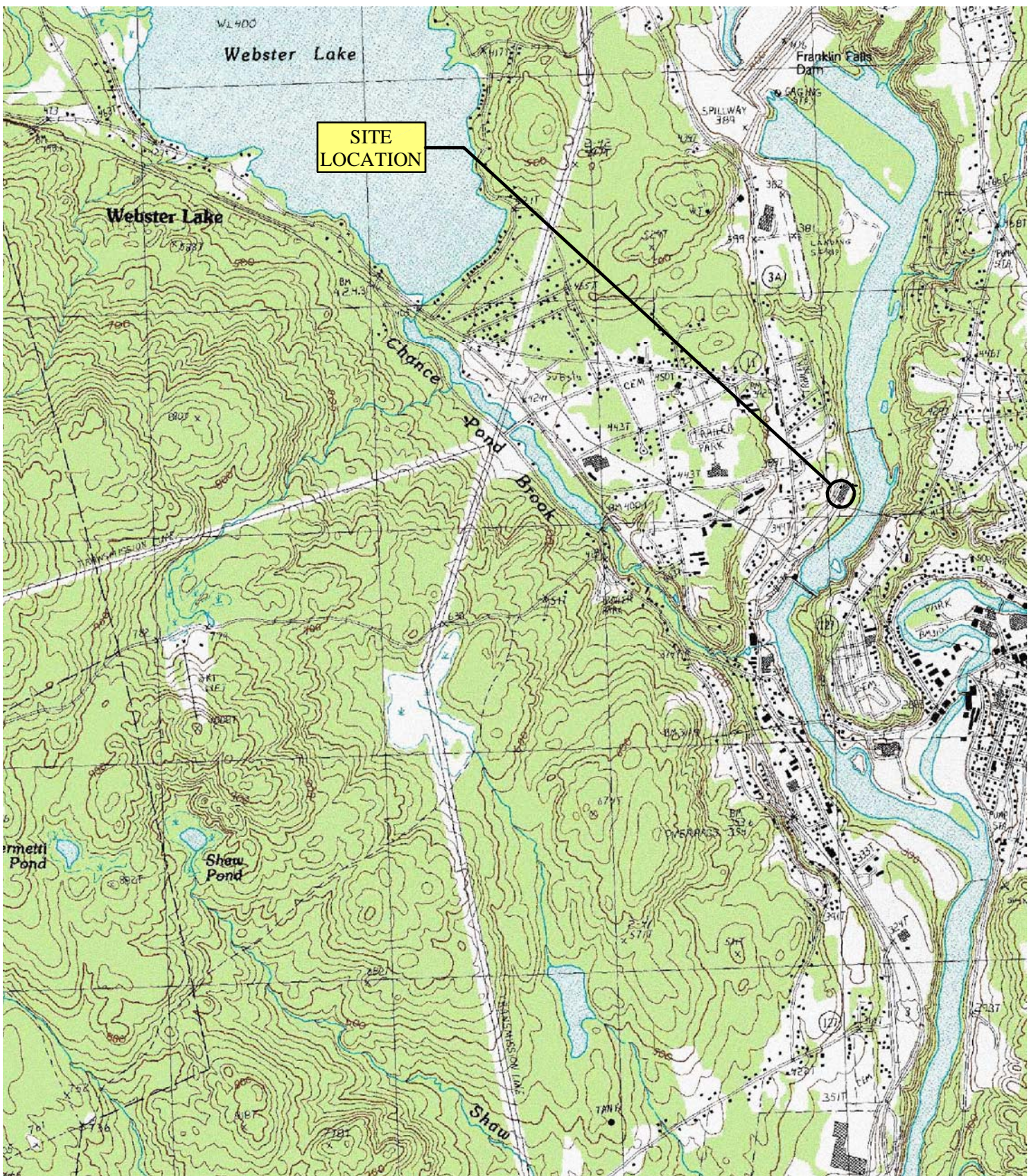
NE - Not established

ND - All analyte concentrations were below the analytical method practical quantitation limit

Exceeds Relative Percent Difference quality control limit of 35% for solid samples as specified in the Project QAPP

FIGURES





USGS 7.5 MINUTE FRANKLIN, NH QUADRANGLE (1987)

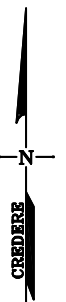
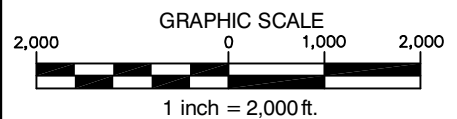
DRAWN BY: WTE	DATE: 2/24/12
CHECKED BY: RSV/JSS	PROJECT: 10001086

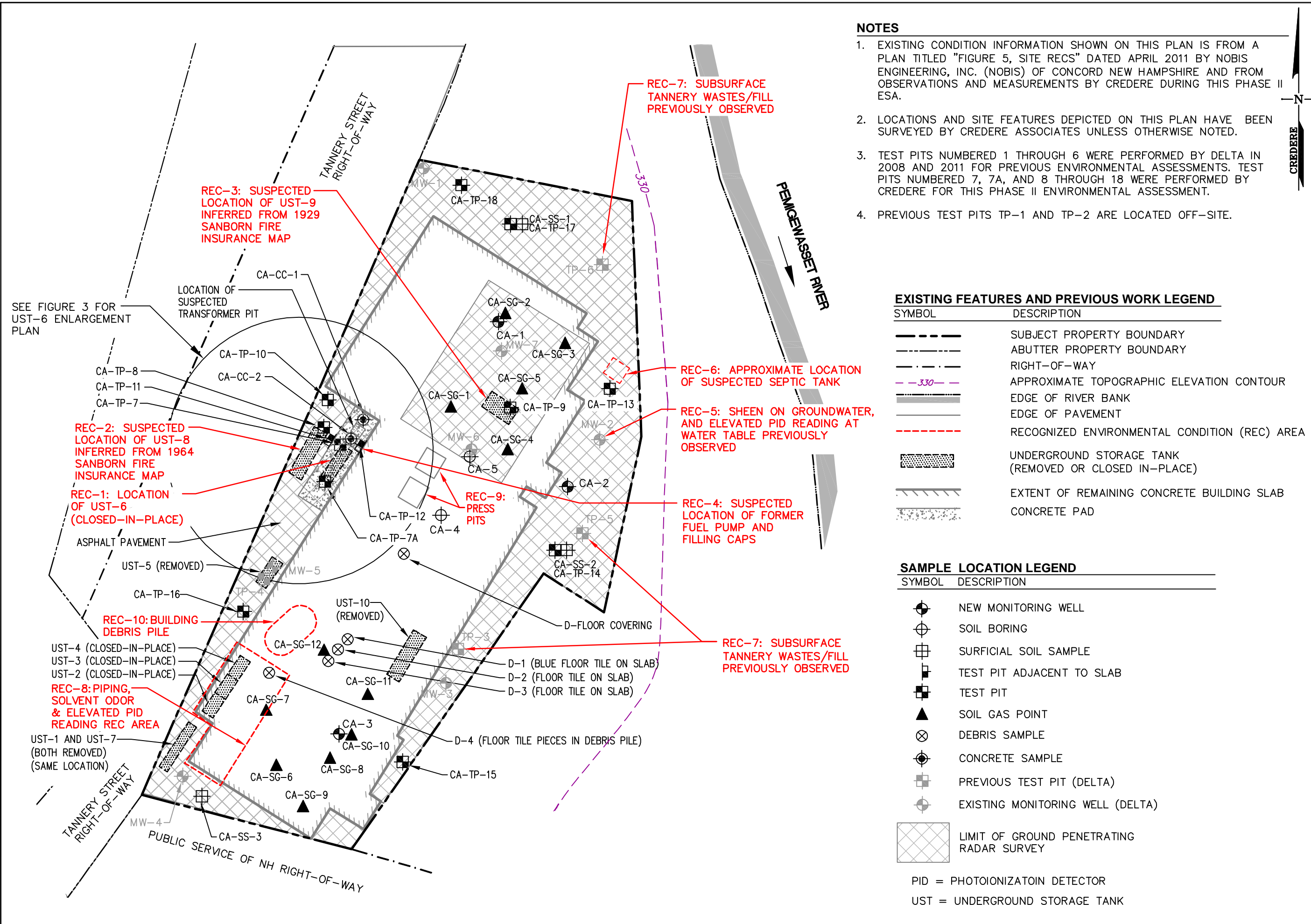
FIGURE 1 - SITE LOCATION PLAN



CREDERE ASSOCIATES, LLC
 776 MAIN STREET
 WESTBROOK, MAINE 04092
 TEL: 207.828.1272
 FAX: 207.887.1051
 WWW.CREDERELLC.COM

FORMER POLYCLAD LAMINATES SITE
 45 TANNERY STREET
 FRANKLIN, NH
 NHDES #199902062





NOTES

1. EXISTING CONDITION INFORMATION SHOWN ON THIS PLAN IS FROM A PLAN TITLED "FIGURE 5, SITE RECS" DATED APRIL 2011 BY NOBIS ENGINEERING, INC. (NOBIS) OF CONCORD NEW HAMPSHIRE AND FROM OBSERVATIONS AND MEASUREMENTS BY CREDERE DURING THIS PHASE II ESA.
2. LOCATIONS AND SITE FEATURES DEPICTED ON THIS PLAN HAVE BEEN SURVEYED BY CREDERE ASSOCIATES UNLESS OTHERWISE NOTED.
3. TEST PITS NUMBERED 1 THROUGH 6 WERE PERFORMED BY DELTA IN 2008 AND 2011 FOR PREVIOUS ENVIRONMENTAL ASSESSMENTS. TEST PITS NUMBERED 7, 7A, AND 8 THROUGH 18 WERE PERFORMED BY CREDERE FOR THIS PHASE II ENVIRONMENTAL ASSESSMENT.
4. PREVIOUS TEST PITS TP-1 AND TP-2 ARE LOCATED OFF-SITE.

EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
---	SUBJECT PROPERTY BOUNDARY
- - - -	ABUTTER PROPERTY BOUNDARY
- · - · -	RIGHT-OF-WAY
- - 330 - -	APPROXIMATE TOPOGRAPHIC ELEVATION CONTOUR
▬▬▬▬▬▬	EDGE OF RIVER BANK
▬▬▬▬▬▬▬▬	EDGE OF PAVEMENT
- - - - -	RECOGNIZED ENVIRONMENTAL CONDITION (REC) AREA
▨▨▨▨▨▨	UNDERGROUND STORAGE TANK (REMOVED OR CLOSED IN-PLACE)
▨▨▨▨▨▨▨▨	EXTENT OF REMAINING CONCRETE BUILDING SLAB
▨▨▨▨▨▨▨▨▨▨	CONCRETE PAD

SAMPLE LOCATION LEGEND

SYMBOL	DESCRIPTION
⊕	NEW MONITORING WELL
⊙	SOIL BORING
⊠	SURFICIAL SOIL SAMPLE
⊞	TEST PIT ADJACENT TO SLAB
⊞	TEST PIT
▲	SOIL GAS POINT
⊗	DEBRIS SAMPLE
⊙	CONCRETE SAMPLE
⊞	PREVIOUS TEST PIT (DELTA)
⊙	EXISTING MONITORING WELL (DELTA)
▨▨▨▨▨▨	LIMIT OF GROUND PENETRATING RADAR SURVEY

PID = PHOTOIONIZATOIN DETECTOR
 UST = UNDERGROUND STORAGE TANK

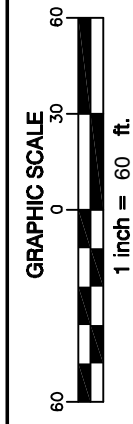
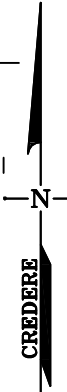


FIGURE 2

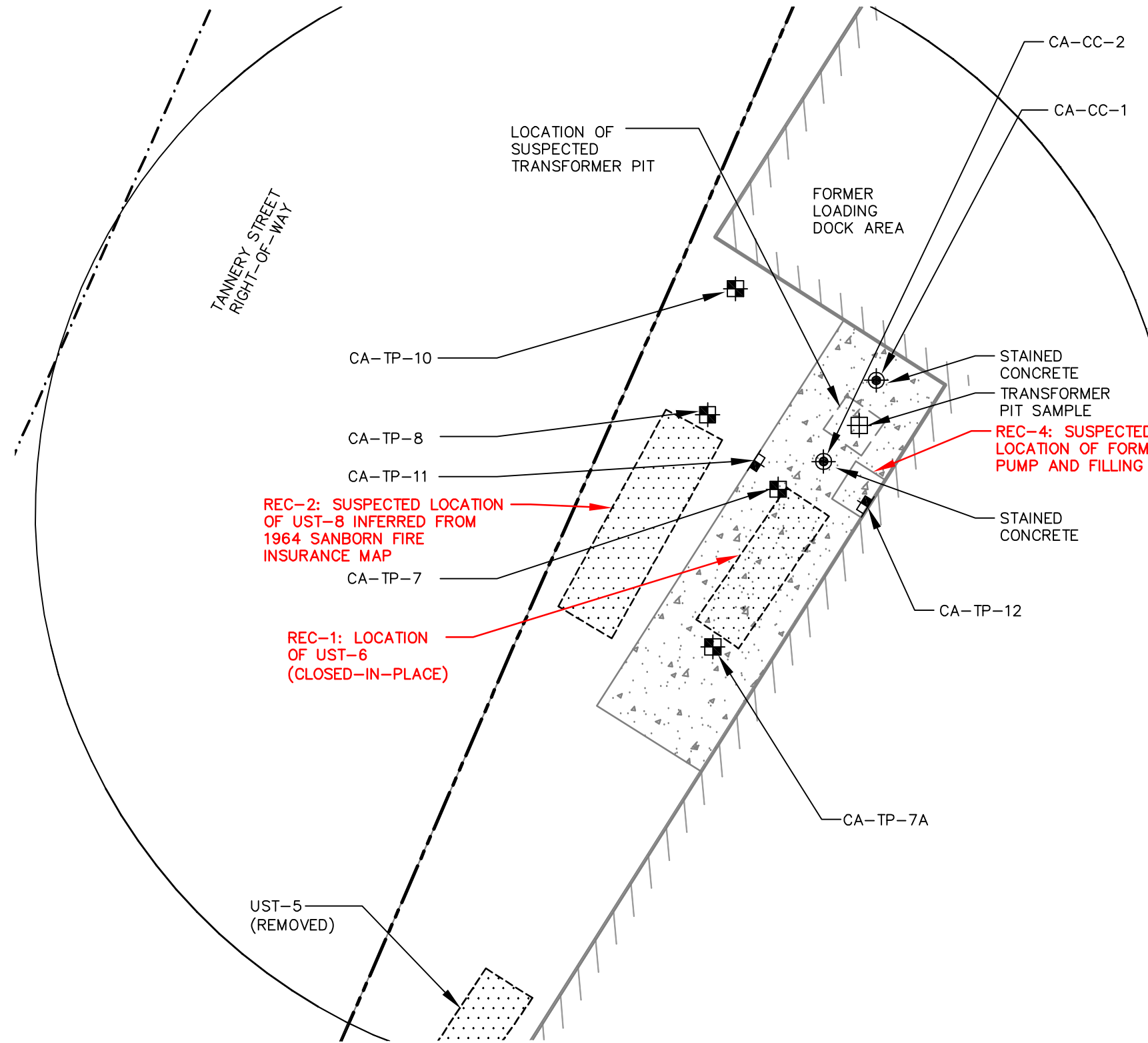
PHASE II ENVIRONMENTAL SITE ASSESSMENT SAMPLE LOCATION PLAN SHOWING RECOGNIZED ENVIRONMENTAL CONDITIONS

FORMER POLYCLAD LAMINATES SITE
 45 TANNERY STREET
 FRANKLIN, NH
 NHDES #199902062

DRAWN BY: WTE DATE: 2/24/12
 CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
 776 MAIN STREET
 WESTBROOK, MAINE 04092
 FAX: 207.887.1051
 TEL: 207.828.1272
 WWW.CREDERELLC.COM



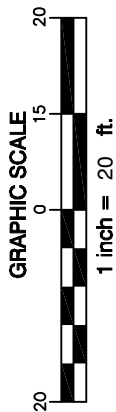
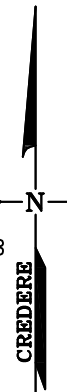


EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
	SUBJECT PROPERTY BOUNDARY
	ABUTTER PROPERTY BOUNDARY
	RIGHT-OF-WAY
	EDGE OF PAVEMENT
	EXTENT OF REMAINING CONCRETE BUILDING SLAB
	CONCRETE PAD
	UNDERGROUND STORAGE TANK (REMOVED OR CLOSED IN-PLACE)

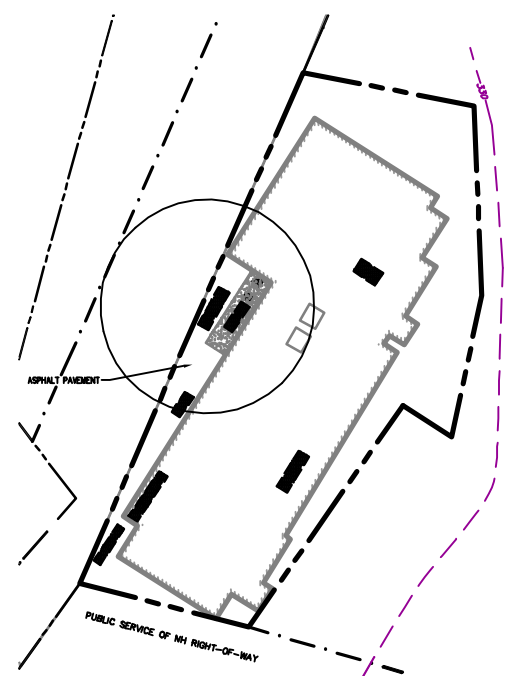
SAMPLE LOCATION LEGEND

SYMBOL	DESCRIPTION
	TEST PIT
	TEST PIT ADJACENT TO SLAB
	CONCRETE SAMPLE
	SURFICIAL SOIL SAMPLE



**FIGURE 3
UST-6 AREA ENLARGEMENT PLAN**

FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET
FRANKLIN, NH
NHDES #199902062

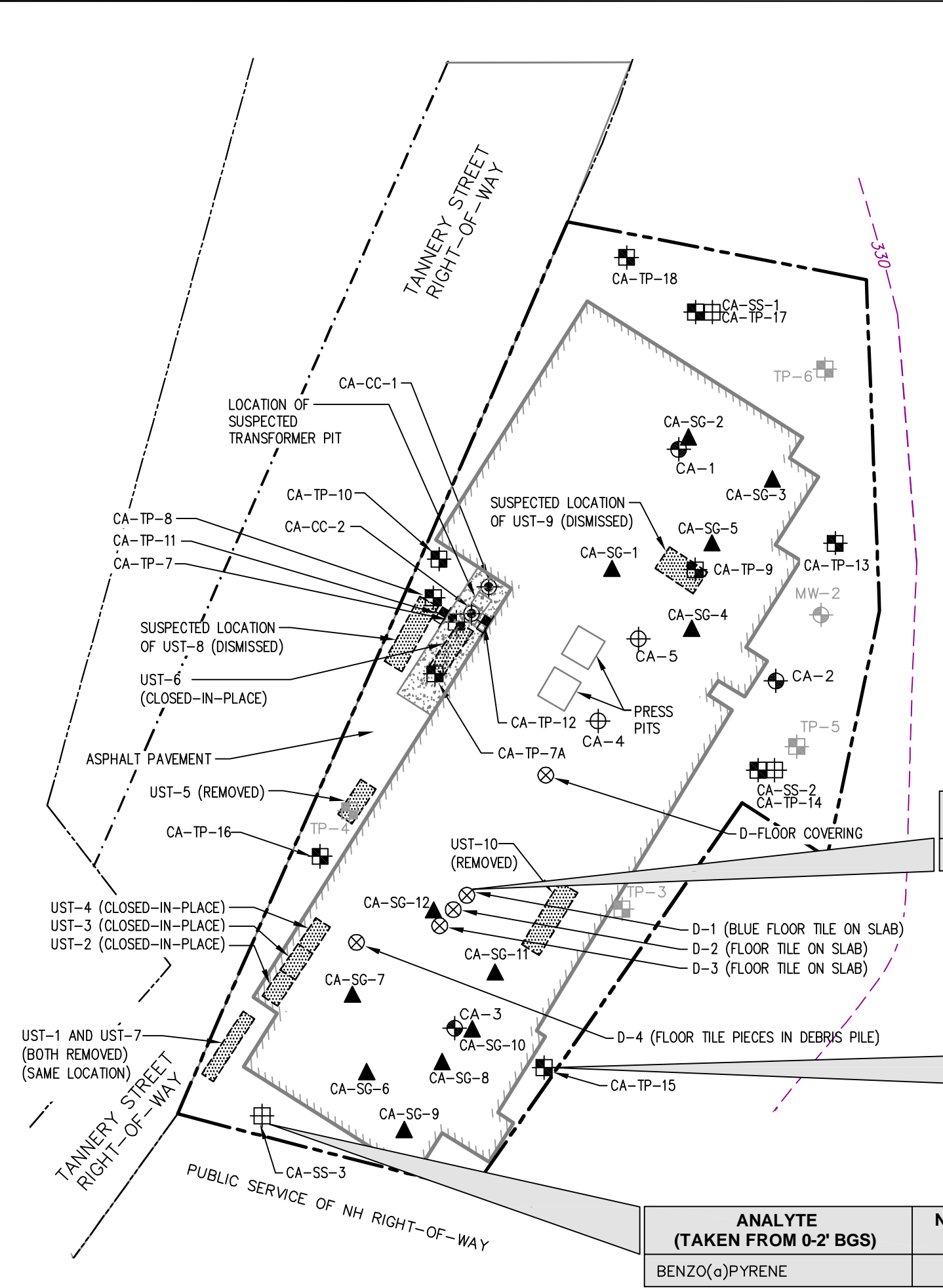


**SUBJECT PROPERTY OVERVIEW
NTS**

DRAWN BY: WTE DATE: 2/24/12
CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
776 MAIN STREET
WESTBROOK, MAINE 04092
FAX: 207.887.1051
TEL: 207.828.1272
WWW.CREDERELLC.COM





NOTES

- EXISTING CONDITION INFORMATION SHOWN ON THIS PLAN IS FROM A PLAN TITLED "FIGURE 5, SITE RECS" DATED APRIL 2011 BY NOBIS ENGINEERING, INC. (NOBIS) OF CONCORD NEW HAMPSHIRE AND FROM OBSERVATIONS AND MEASUREMENTS BY CREDERE DURING THIS PHASE II ESA.
- LOCATIONS AND SITE FEATURES DEPICTED ON THIS PLAN HAVE BEEN SURVEYED BY CREDERE ASSOCIATES.
- ONLY ANALYTES DETECTED EXCEEDING THEIR APPLICABLE STANDARDS ARE DEPICTED ON THIS FIGURE. DATA FOR ALL ANALYTES AND LOCATIONS ARE SHOWN IN TABLES 5 AND 6.
- ANALYTE CONCENTRATIONS SHOWN IN BOLD REPRESENT EXCEEDENCES OF THE APPLICABLE STANDARD.

EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
(Dashed line)	SUBJECT PROPERTY BOUNDARY
(Dotted line)	ABUTTER PROPERTY BOUNDARY
(Dash-dot line)	RIGHT-OF-WAY
(Dashed purple line)	APPROXIMATE ELEVATION CONTOUR
(Solid line)	EDGE OF RIVER
(Line with hatching)	EDGE OF PAVEMENT
(Line with diagonal hatching)	EXTENT OF REMAINING CONCRETE BUILDING SLAB
(Stippled area)	CONCRETE PAD
(Patterned box)	UNDERGROUND STORAGE TANK (REMOVED OR CLOSED IN-PLACE)

SAMPLE LOCATION LEGEND

SYMBOL	DESCRIPTION
(Circle with crosshair)	NEW MONITORING WELL
(Circle with center dot)	EXISTING MONITORING WELL (DELTA)
(Circle with crosshair)	SOIL BORING
(Square with crosshair)	SURFICIAL SOIL SAMPLE
(Square with center dot)	TEST PIT ADJACENT TO SLAB
(Square with crosshair)	TEST PIT
(Square with center dot)	PREVIOUS TEST PIT (DELTA)
(Triangle)	SOIL GAS POINT
(Circle with crosshair)	DEBRIS PILE SAMPLE
(Circle with center dot)	CONCRETE SAMPLE

SRS = SOIL REMEDIATION STANDARDS

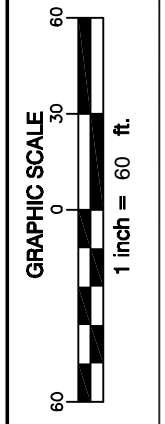
ANALYTE	NHDES SRS (%)	CONC. (%)
ASBESTOS	1	5

ANALYTE (TAKEN FROM 0-2' BGS)	NHDES SRS (mg/kg)	CONC. (mg/kg)
BENZO(a)PYRENE	0.70	0.71

ANALYTE (TAKEN FROM 5-7' BGS)	NHDES SRS (mg/kg)	CONC. (mg/kg)
BENZO(a)ANTHRACENE	1.0	80
BENZO(b)FLUORANTHENE	1.0	46
BENZO(k)FLUORANTHENE	4.0	80
BENZO(a)PYRENE	0.70	66
INDENO(1,2,3-cd)PYRENE	1.0	18
DIBENZO(a,h)ANTHRACENE	0.70	9.6

**FIGURE 4
SOIL AND BUILDING MATERIAL EXCEEDANCE LOCATION PLAN**

FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET
FRANKLIN, NH
NHDES #199902062



DRAWN BY: WTE DATE: 2/24/12
CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
776 MAIN STREET
WESTBROOK, MAINE 04092
FAX: 207.887.1051
TEL: 207.828.1272
WWW.CREDERELLC.COM

ANALYTE	AGQS (ug/L)	CONC. (ug/L)
VOCs	-	ALL < AGQS
METALS	-	NOT SAMPLED
SVOCs	-	NOT SAMPLED

NOTES

- EXISTING CONDITION INFORMATION SHOWN ON THIS PLAN IS FROM A PLAN TITLED "FIGURE 5, SITE RECS" DATED APRIL 2011 BY NOBIS ENGINEERING, INC. (NOBIS) OF CONCORD NEW HAMPSHIRE AND FROM OBSERVATIONS AND MEASUREMENTS BY CREDERE DURING THIS PHASE II ESA.
- LOCATIONS AND SITE FEATURES DEPICTED ON THIS PLAN HAVE BEEN SURVEYED BY CREDERE ASSOCIATES.
- ANALYTE CONCENTRATIONS SHOWN IN BOLD REPRESENT EXCEEDENCES OF THE APPLICABLE STANDARD.

EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
	SUBJECT PROPERTY BOUNDARY
	ABUTTER PROPERTY BOUNDARY
	RIGHT-OF-WAY
	APPROXIMATE ELEVATION CONTOUR
	EDGE OF RIVER
	EDGE OF PAVEMENT
	EXTENT OF REMAINING CONCRETE BUILDING SLAB
	CONCRETE PAD
	UNDERGROUND STORAGE TANK (REMOVED OR CLOSED IN-PLACE)

SAMPLE LOCATION LEGEND

SYMBOL	DESCRIPTION
	MONITORING WELL
	EXISTING MONITORING WELL (DELTA)

ANALYTICAL RESULTS TABLE LEGEND

- SRS = SOIL REMEDIATION STANDARDS
- AGQS = AMBIENT GROUNDWATER QUALITY STANDARDS
- PQL = LABORATORY PRACTICAL QUANTITATION LIMIT
- VOCs = VOLATILE ORGANIC COMPOUNDS
- SVOCs = SEMI-VOLATILE ORGANIC COMPOUNDS

ANALYTE	AGQS (ug/L)	CONC. (ug/L)
VOCs	-	ALL < PQL
LEAD	15	8
SVOCs	-	NOT SAMPLED

ANALYTE	AGQS (ug/L)	CONC. (ug/L)
VOCs		
TRIMETHYLBENZENE, 1, 2, 4-	330	710
ALL OTHER VOCs	-	ALL < AGQS
SVOCs	-	ALL < AGQS
METALS	-	ALL < AGQS

ANALYTE	AGQS (ug/L)	CONC. (ug/L)
VOCs	-	ALL < PQL
METALS	-	ALL < PQL
SVOCs	-	NOT SAMPLED

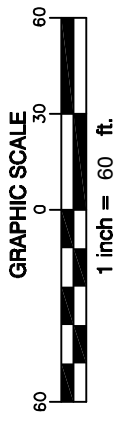
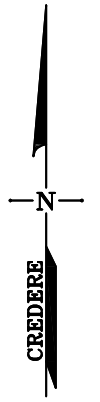
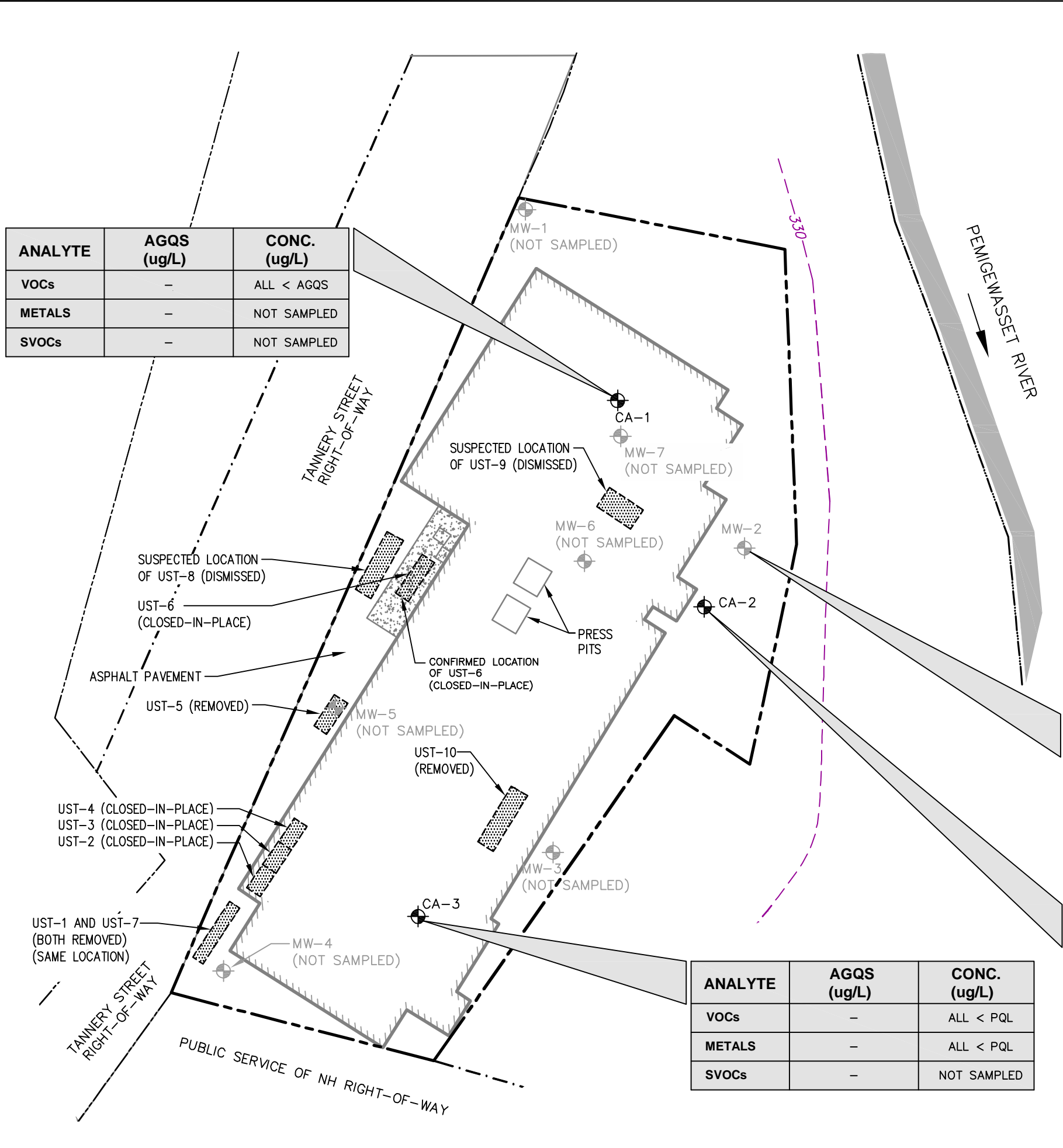


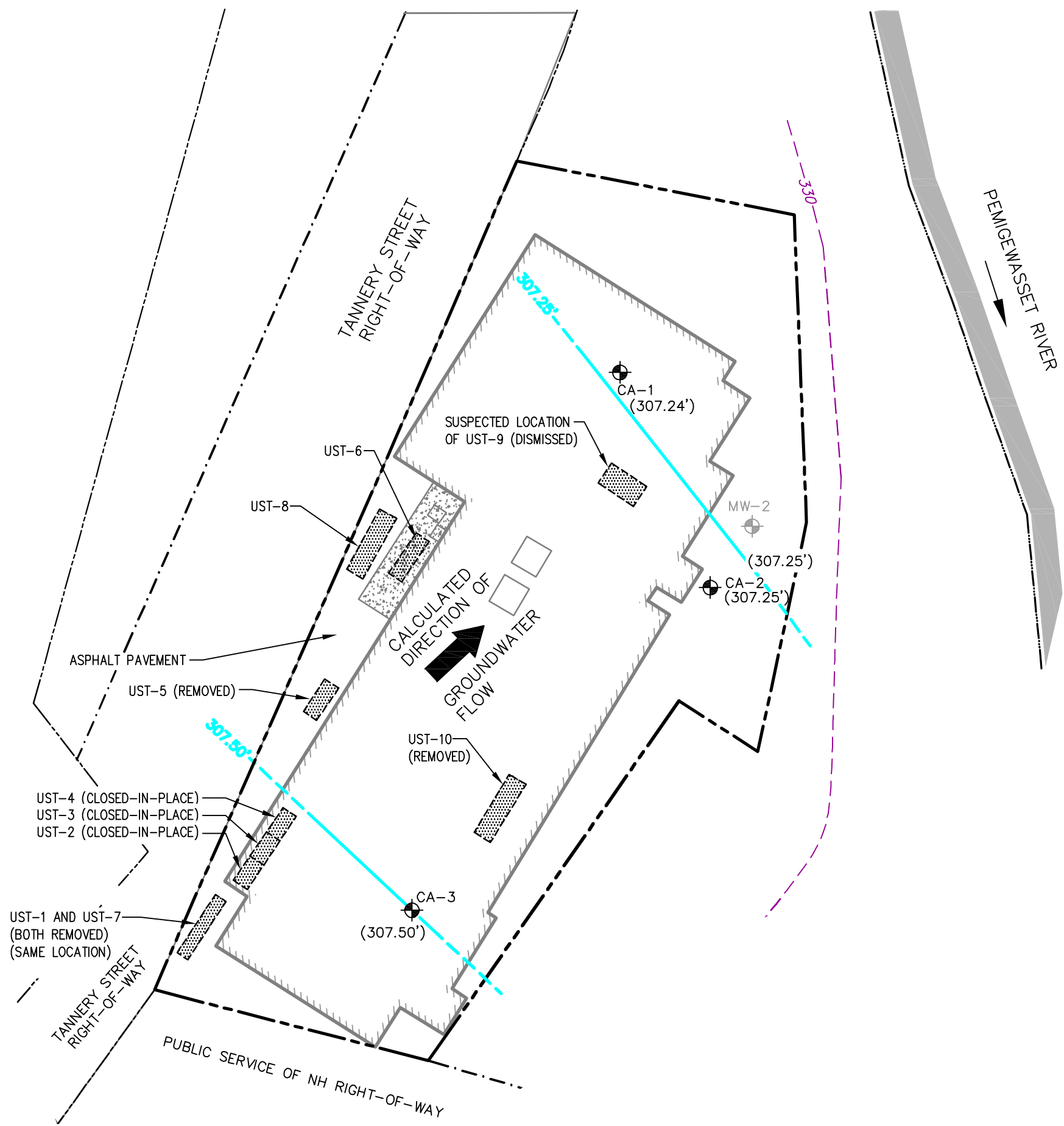
FIGURE 5
SUMMARY OF GROUNDWATER RESULTS PLAN

FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET
FRANKLIN, NH
NHDES #199902062

DRAWN BY: WTE DATE: 2/24/12
CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
776 MAIN STREET
WESTBROOK, MAINE 04092
FAX: 207.887.1051
TEL: 207.828.1272
WWW.CREDERELLC.COM





NOTES

1. EXISTING CONDITION INFORMATION SHOWN ON THIS PLAN IS FROM A PLAN TITLED "FIGURE 5, SITE RECS" DATED APRIL 2011 BY NOBIS ENGINEERING, INC. (NOBIS) OF CONCORD NEW HAMPSHIRE AND FROM OBSERVATIONS AND MEASUREMENTS BY CREDERE DURING THIS PHASE II ESA.
2. SAMPLING LOCATIONS DEPICTED ON THIS PLAN HAVE BEEN SURVEYED BY CREDERE ASSOCIATES.
3. GROUNDWATER CONTOURS ARE BASED ON ELEVATION MEASUREMENTS TAKEN ON JULY 5, 2011. ELEVATIONS ARE RELATIVE TO BENCHMARK MW-2 (334.89') DEPICTED IN DELTA'S 2008 PHASE II ESA FOR THE SITE.
4. HYDRAULIC GRADIENT IS EQUAL TO APPROXIMATELY 0.001 FT/FT.
5. GROUNDWATER MONITORING WELL PLACEMENT WAS BASED ON THE INITIAL INFERRED DIRECTION OF GROUNDWATER FLOW AND/OR TO INVESTIGATE RECS AND OTHER CONDITIONS OBSERVED DURING PHASE II PERFORMANCE.

EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
	SUBJECT PROPERTY BOUNDARY
	ABUTTER PROPERTY BOUNDARY
	RIGHT-OF-WAY
	APPROXIMATE TOPOGRAPHIC ELEVATION CONTOUR
	CALCULATED GROUNDWATER CONTOUR
	ASSUMED GROUNDWATER CONTOUR
	EDGE OF RIVER BANK
	EDGE OF PAVEMENT
	EXTENT OF REMAINING CONCRETE BUILDING SLAB
	CONCRETE PAD
	UNDERGROUND STORAGE TANK (REMOVED OR CLOSED IN-PLACE)

SAMPLE LOCATION LEGEND

SYMBOL	DESCRIPTION
	NEW MONITORING WELL
	EXISTING MONITORING WELL (DELTA)
	(307.25') GROUNDWATER ELEVATION CALCULATED FROM JULY 5, 2011 DEPTH TO WATER

FIGURE 6
JULY 5, 2011 GROUNDWATER ELEVATION CONTOUR PLAN

FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET
FRANKLIN, NH
NHDES #199902062

DRAWN BY: WTE DATE: 2/21/12
CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
776 MAIN STREET
WESTBROOK, MAINE 04092
FAX: 207.887.1051
TEL: 207.828.1272
WWW.CREDERELLC.COM



NOTES

- EXISTING CONDITION INFORMATION SHOWN ON THIS PLAN IS FROM A PLAN TITLED "FIGURE 5, SITE RECS" DATED APRIL 2011 BY NOBIS ENGINEERING, INC. (NOBIS) OF CONCORD NEW HAMPSHIRE AND FROM OBSERVATIONS AND MEASUREMENTS BY CREDERE DURING THIS PHASE II ESA.
- LOCATIONS AND SITE FEATURES DEPICTED ON THIS PLAN HAVE BEEN SURVEYED BY CREDERE ASSOCIATES.

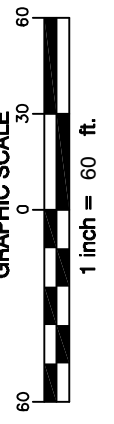
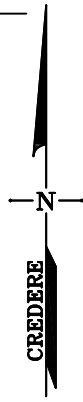
EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
	SUBJECT PROPERTY BOUNDARY
	ABUTTER PROPERTY BOUNDARY
	RIGHT-OF-WAY
	APPROXIMATE ELEVATION CONTOUR
	EDGE OF RIVER
	EDGE OF PAVEMENT
	UNDERGROUND STORAGE TANK (REMOVED OR CLOSED IN-PLACE)
	EXTENT OF REMAINING CONCRETE BUILDING SLAB
	CONCRETE PAD
	AREA WITH OBSERVED BURIED TANNERY WASTE*

* EXTENT IS ESTIMATED BASED ON OBSERVED CONDITIONS AND MAY NOT BE REPRESENTATIVE OF AREAS NOT DIRECTLY ASSESSED DURING PHASE II ESA ACTIVITIES.

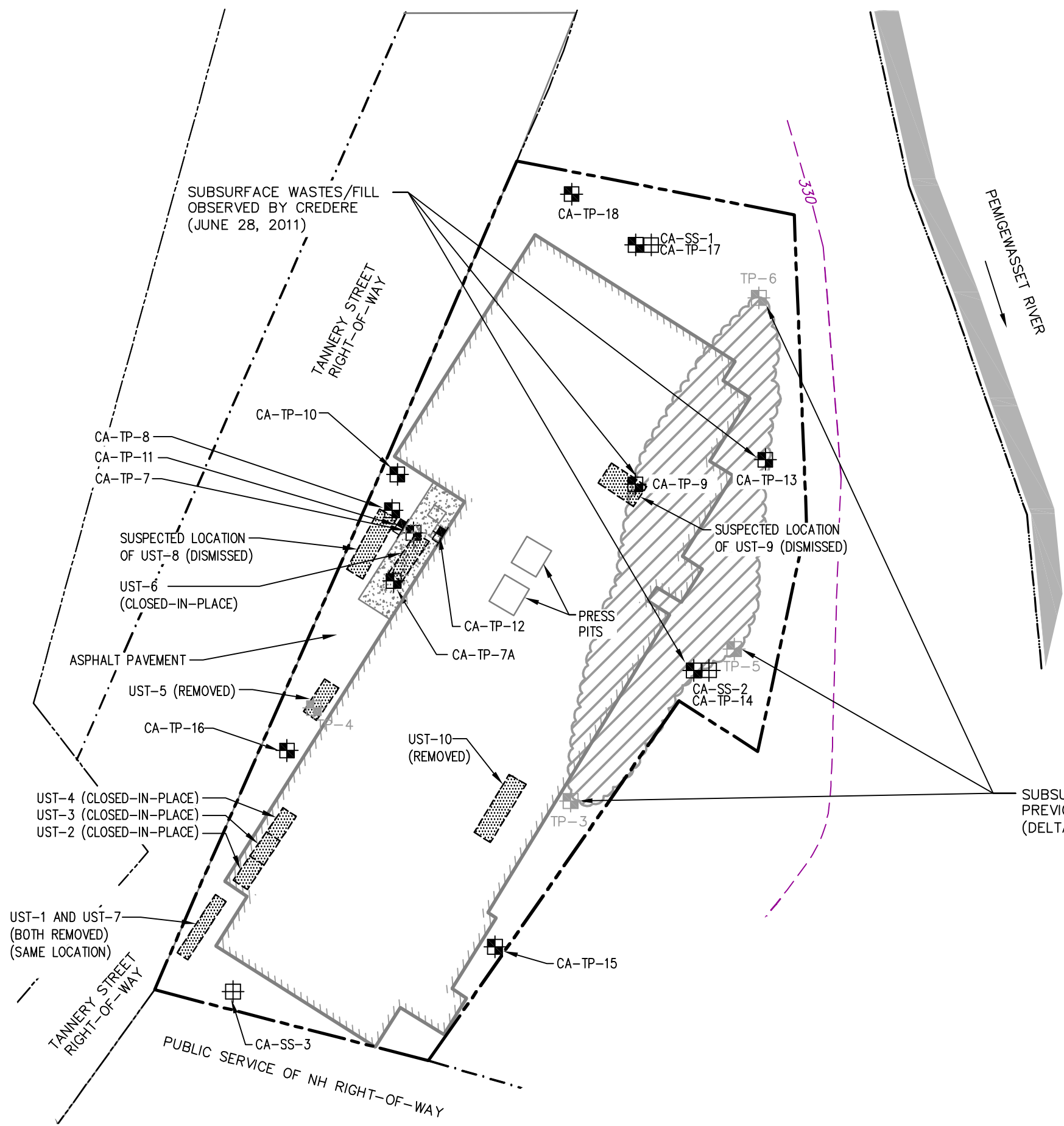
SAMPLE LOCATION LEGEND

SYMBOL	DESCRIPTION
	SURFICIAL SOIL SAMPLE
	TEST PIT
	PREVIOUS TEST PIT (DELTA)
	TEST PIT ADJACENT TO SLAB



**FIGURE 7
BURIED TANNERY WASTE LOCATION PLAN**

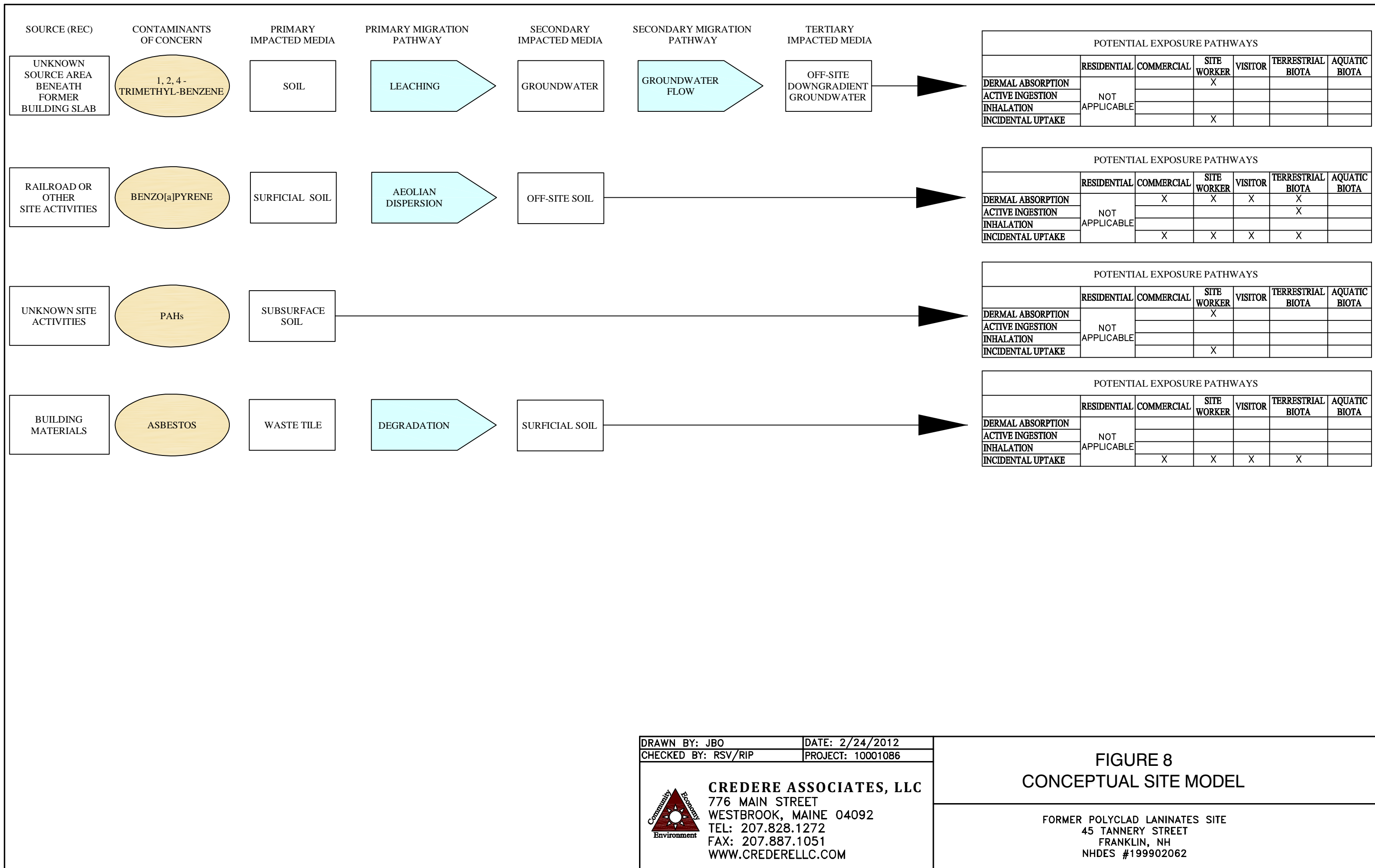
FORMER POLYCLAD LAMINATES SITE
45 TANNERY STREET
FRANKLIN, NH
NHDES #199902062



DRAWN BY: WTE DATE: 2/24/12
CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
776 MAIN STREET
WESTBROOK, MAINE 04092
FAX: 207.887.1051
TEL: 207.828.1272
WWW.CREDERELLC.COM





DRAWN BY: JBO	DATE: 2/24/2012
CHECKED BY: RSV/RIP	PROJECT: 10001086



CREDERE ASSOCIATES, LLC
 776 MAIN STREET
 WESTBROOK, MAINE 04092
 TEL: 207.828.1272
 FAX: 207.887.1051
 WWW.CREDERELLC.COM

FIGURE 8
CONCEPTUAL SITE MODEL

FORMER POLYCLAD LANINATES SITE
 45 TANNERY STREET
 FRANKLIN, NH
 NHDES #199902062

APPENDIX A

SITE-SPECIFIC QUALITY ASSURANCE PROJECT PLAN ADDENDUM



**Waste Management Division
PO Box 95, 29 Hazen Drive
Concord, NH 03302**

Type of Submittal (Check One-Most Applicable)

<input type="checkbox"/> Work Scope <input type="checkbox"/> Reimbursement Request	<input type="checkbox"/> Remedial Action <ul style="list-style-type: none"> • Remedial Action Plan • Bid Plans and Specifications • Remedial Action Implementation Report
<input type="checkbox"/> UST Facility Report <input type="checkbox"/> AST Facility Report	<input type="checkbox"/> Treatment System and POE O&M <input type="checkbox"/> Activity and Use Restriction
<input type="checkbox"/> Emergency/Initial Response Action <input type="checkbox"/> Groundwater Quality Assessment	<input type="checkbox"/> Temporary Surface Water Discharge Permit
<input type="checkbox"/> Initial Site Characterization <input type="checkbox"/> Site Investigation <ul style="list-style-type: none"> • Site Investigation Report • Supplemental Site Investigation Report • GMZ Delineation • Source Area Investigation • Data Submittal • Annual Summary Report <input checked="" type="checkbox"/> Unsolicited Site-Specific Quality Assurance Project Plan Addendum <input type="checkbox"/> Closure Documentation	<input type="checkbox"/> Groundwater Management Permit <ul style="list-style-type: none"> • Permit Application • Renewal Application • Deed Recordation Documentation • Abutter Notification Documentation • Release of Recordation <input type="checkbox"/> Data Submittal <input type="checkbox"/> Annual Summary Report

**SITE SPECIFIC QUALITY ASSURANCE PROJECT
PLAN ADDENDUM**

Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES#199902062
Brownfields Grant #BF-96111801

Prepared For:
Lakes Region Planning Commission
103 Main Street, Suite #3
Meredith, NH 03253
Phone: (603) 279-8171
Contact: Mr. Kimon Koulet

Prepared By:
CREDERE ASSOCIATES, LLC
776 Main Street
Westbrook, ME 04902
Phone: (207) 828-1272 ext. 35
Contact: Richard Vandenberg, CG, PG

June 27, 2011

Recommended Risk Category (check one)

<input type="checkbox"/> 1. Immediate Human Health Risk (Impacted water supply well, etc.)	<input type="checkbox"/> 4. Surface Water Impact	<input type="checkbox"/> 7. Alternate Water Available/Low Level Groundwater Contamination (<1,000 X AGQS)
<input type="checkbox"/> 2. Potential Human Health Risk (Water supply well within 1,000' or Site within SWPA)	<input type="checkbox"/> 5. No Alternate Water Available/No Existing Wells in Area	<input type="checkbox"/> 8. No AGQS Violation/No Source Remaining
<input type="checkbox"/> 3. Free Product or Source Hazard	<input type="checkbox"/> 6. Alternate Water Available/High Level Groundwater Contamination (>1,000 X AGQS)	<input type="checkbox"/> Closure Recommended

1. TITLE AND APPROVAL PAGE

SITE-SPECIFIC QUALITY ASSURANCE PROJECT PLAN (SSQAPP) ADDENDUM TO GENERIC QAPP RFA #08166 AND #09036 Revision 0

Site Information:

Former Polyclad Laminates Property
45 Tannery Street, New Hampshire
New Hampshire Department of Environmental Services (NHDES) Site No. 199902062

Funding Source:

Lakes Region Planning Commission's (LRPC's) Brownfields Assessment Program
United States Environmental Protection Agency (EPA) Brownfields Grant # BF96111801


Prepared By:

Richard S. Vandenberg, CG, PG
Crede Associates, LLC
776 Main Street, Westbrook, Maine 04092
(207) 828-1272

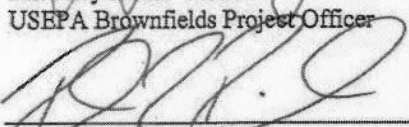
Date Prepared:

June 27, 2011

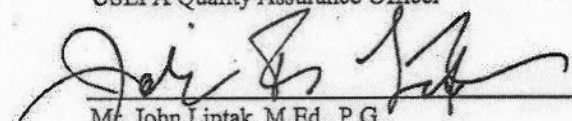
Below is a listing of the names, titles, signatures, and signature dates of officials approving this SSQAPP:




Ms. Jerry Minor-Gordon
USEPA Brownfields Project Officer
Date: 6/28/11




Mr. Robert Reinhart
USEPA Quality Assurance Officer
Date: 6-27-11



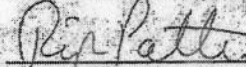
Mr. John Liptak, M.Ed., P.G.
Brownfields Project Manager
Date: 6-27-11



for Mr. Vincent R. Perelli
New Hampshire DES QA Manager
Date: 6/27/11



Mr. Richard S. Vandenberg, CG, PG
Crede Associates, LLC Project QA Manager
Date: June 27, 2011



Mr. Robert I Patten, PE, LBED-AP, LSP
Crede Associates, LLC Project Manager
Date: June 27, 2011

TABLE OF CONTENTS

1. TITLE AND APPROVAL PAGE	1
2. INTRODUCTION.....	4
3. FINDINGS OF THE PHASE I ESA	5
3.1 Site Description and History	5
3.2 Other documented previous environmental work Cited by Nobis	5
3.3 history of tanks at the property	6
3.4 Identified Recognized Environmental Conditions	7
3.5 other potential Environmental Conditions	8
4. POTENTIAL REDEVELOPMENT SCENARIO.....	10
5. CONCEPTUAL SITE MODEL	11
5.1 Contaminants of Concern	11
5.2 Geology, Storm water Flow, and Groundwater Flow	12
5.3 Definitions of Exposure Pathways and Potential Receptors	12
6. SAMPLING DESIGN.....	15
7. FIELD ACTIVITY METHODOLOGY	21
7.1 GPR Survey and Electronic Tracing.....	21
7.2 Soil Borings, Test Pitting, Surficial Soil Sampling, and Field Screening.....	21
7.3 Monitoring Well Installation and Survey.....	22
7.4 Groundwater Sampling	23
7.5 soil gas Sampling	23
7.6 Building material debris Surveys.....	23
8. REGULATORY STANDARDS	24
8.1 Soil.....	24
8.2 Groundwater	24
8.3 Soil gas.....	24
8.4 ACM	24
8.5 Lead-Based Paint	25
8.6 PCB Bulk Products	25
9. PROPOSED PROJECT SCHEDULE	26



FIGURES

- Figure 1** Site Location Map
Figure 2 Credere Organization and Responsibility Chart
Figure 3 Proposed Phase II Sample Location Plan
Figure 4 Conceptual Site Model

TABLES

- Table 1** Potential Contaminants of Concern (imbedded in text)
Table 2 Sample Reference Table



2. INTRODUCTION

The Lakes Region Planning Commission (LRPC) has received a United States Environmental Protection Agency (USEPA) Brownfields Hazardous Substance Assessment Grant to conduct environmental investigations at sites within the 30 member communities of the Lake Winnepesaukee Region. The investigations provide the basis for reuse planning specific to each site's community needs. The assessment of each site will include the completion of Phase I and Phase II Environmental Site Assessments (ESAs) and may also potentially include the development of cleanup and reuse options for selected sites.

On behalf of LRPC's Brownfields Assessment Program, this document is a Site-Specific Quality Assurance Project Plan (SSQAPP) Addendum for the Former Polyclad Laminates Property located at 45 Tannery Street, Franklin, New Hampshire (the Site). **Figure 1** shows the general location of the Site in Franklin and **Figure 2** is Credere's Organization and Responsibility Chart for the project and **Figure 3** is a plan showing the locations of proposed sampling work.

This SSQAPP presents the following information:

1. A summary of the pertinent findings of the Phase I ESA
2. The potential redevelopment scenario for the Site
3. A conceptual site model
4. Credere's proposed sampling design including recommended sample locations, and analytical methods for proposed Phase II ESA activities
5. Credere's proposed field activity methodology
6. Regulatory standards applicable to the Site
7. A proposed project schedule

This SSQAPP was prepared to be used in concert with Credere Associates, LLC (Credere) Generic Quality Assurance Project Plan (QAPP) Rev. 2 (USEPA RFA #08166 and #09036) which was prepared for all of Credere's USEPA work in New Hampshire. The quality assurance and quality control (QA/QC) procedures outlined in Credere's Generic QAPP will be followed for this investigation program including sample collection, handling, and analysis, chain of custody, data management and documentation, data validation, and data usability assessments.



3. FINDINGS OF THE PHASE I ESA

A Phase I ESA was completed by Nobis Engineering, Inc. (Nobis) for the Site in April 2011. The Phase I ESA was conducted using EPA Brownfields funds under Brownfields Coalition Assessment Grant #RP-96128501 at the direction of the New Hampshire Department of Environmental Services (NHDES).

The following is a summary of the pertinent findings from the Phase I ESA:

3.1 SITE DESCRIPTION AND HISTORY

The Site is composed of a single vacant 1.97-acre parcel of land located at 45 Tannery Street in Franklin, New Hampshire, that is situated along the Pemigewasset River. The Site consists of a vacant lot with only the former building's concrete slab-on-grade foundation and a small pile of debris. **Figure 3** shows the former site building footprint and other pertinent Site features.

According to Nobis's Phase I ESA report, which was reviewed as a part of the development of this SSQAPP, the Site was first developed in 1892. By approximately 1930, the first building was constructed on the Site. The most recent building was demolished in 2008 after it collapsed due to heavy winter snow cover.

According to the Nobis Phase I ESA report, the Site appears to be have used for railroad purposes when the Boston & Maine and Bristol Rail Road transected the Site beginning in 1892. It is not clear when railroad activities ceased on the Site. However, records do show that the Hingston Leather Company, Inc. and Louis Verza Leather Company were using the Site for tannery activities beginning in the 1950's until the mid to late 1970's. Nobis indicated that there was very little environmental information available for the Site as it pertains to the former tannery operations.

Between 1979 and 1980, Polyclad Laminates, Inc. began operations on the Site. Polyclad Laminates, Inc. manufactured component materials used in multi-layered circuit boards called "pre-preg". The process involved dipping fiberglass cloth in a dip tank filled with epoxy resin. Polyclad Laminates used the Site for the manufacture of pre-preg from between 1979 and 1980 until around 2006.

3.2 OTHER DOCUMENTED PREVIOUS ENVIRONMENTAL WORK CITED BY NOBIS

The Nobis Phase I details the completion of several other investigations, which were completed outside the Brownfields due diligence process, likely as part of a previous effort to sell the Site. Documented previous works include the following:



1. A January 12, 1999 Underground Storage Tank (UST) Closure Report UST-1 prepared by Les A. Cartier and Associates, Inc.
2. A July 15, 1999 UST Closure Report UST-2, UST-3 and UST-4 prepared by Environmental Science & Engineering (ESE).
3. A 2005 Phase I Environmental Site Assessment and Limited Compliance Review prepared by Delta Consultants.
4. An April 10, 2008 Phase II Environmental Site Assessment prepared by Delta Consultants.
5. An October 15, 2008 UST Closure Report UST-7 and UST-10 prepared Delta Consultants.

The conditions identified by Nobis in **Section 3.4** below cumulatively considered all the above referenced previous investigations. Credere also reviewed and considered this cumulative work as a part of the development of this SSQAPP.

3.3 HISTORY OF TANKS AT THE SITE

The operational history of USTs is an important element in the environmental history of the Site. As result, the information below, as reported by Nobis in the April 2011 Phase I ESA, provides an inventory of the USTs which have been identified at the Site. Additional details concerning UST history can be reviewed in the previously submitted Nobis Phase I ESA.

UST ID	Capacity (gallons)	Material Held	Current Status
UST -1	6,000	Methyl Cellosolve	Removed, 1998
UST-2	4,500	CS 350 Acetone	Closed in-place, 1998
UST-3	4,500	Epoxy Resin	Closed in-place, 1999
UST-4	6,000	CS 350 Acetone	Closed in-place, 1999
UST-5	4,000	Gasoline	Removed, date uncertain.
UST-6	12,000	Fuel Oil	Closed in-place, date uncertain
UST-7	13,500	Epoxy Resin	Removed 2008
UST-8	4,000	Fuel Oil	Approximate location from 1964 Sanborn Map
UST -9	4,000	Fuel Oil	Approximate location under slab from 1929 Sanborn Map
UST-10	8,000	Waste Petroleum Distillates	Removed 2008



3.4 IDENTIFIED RECOGNIZED ENVIRONMENTAL CONDITIONS

The following is a summary of the *recognized environmental conditions* (RECs) that were identified by Nobis and reported in their April 2011 Phase I ESA for the Site:

1. The presence of a 12,000-gallon fuel oil UST (identified as UST-6) that is located in the vicinity of the loading dock and was closed in-place (date uncertain) represented a REC.
2. The potential presence of a reported 4,000-gallon fuel oil UST (identified as UST-8) shown on a Sanborn Map, represented a REC because the presence of this tank has not been confirmed and no documentation regarding closure of the tank or information on the integrity of the tank was identified during the Phase I ESA.
3. The potential presence of a reported 4,000-gallon fuel oil (identified as UST-9) shown on a Sanborn Map, represented a REC because the presence of this tank has not been confirmed and no documentation regarding closure of the tank or information on the integrity of the tank was identified during the Phase I ESA.
4. Due to the potential for undocumented past releases to the environment, Nobis identified a REC associated with presence of a former gasoline dispenser and a number of filling caps formerly located in the vicinity of former UST-6 because these items may have been associated with additional undocumented USTs in this area, and documentation regarding the removal of this dispenser or the disposition of related piping with respect to potential releases has not been confirmed.
5. The presence of a sheen of degraded gasoline previously noted in 2008 in monitoring well MW-2 for which a source was not identified, represents a REC due to the potential for an unidentified source of petroleum contamination to be present at the Site. [Note: According to a Phase II Environmental Site Assessment Report prepared by Delta Consultants on April 10, 2008, previous sampling data collected in 2007 indicates that no VOCs, metals or polychlorinated biphenyls (PCBs) were detected in excess of New Hampshire Soil Remediation Standards (SRS) in a soil sample collected (30 to 32 feet) during drilling of this well and no volatile organic compounds (VOCs) or semi-volatile organic compounds (SVOCs) were detected in excess of Ambient Groundwater Quality Standards (AGQS) in the groundwater sample collected from this well. However, a petroleum sheen was noted during sampling of the groundwater and the tentatively identified compounds (TICs) noted during laboratory analysis revealed that the sheen was likely gasoline.]
6. The potential that a subsurface structure identified during a GPR survey east of the former building, which was concluded to be a former concrete septic tank, represents a REC because of the potential for previous undocumented subsurface storage or disposal of hazardous substances and/or petroleum products in and around this structure.



7. The presence of buried waste related to the previous usage of the Site as a tannery represents a REC because the understanding of the extent and disposition of the waste (leather waste or unsuitable materials) is limited and may have lead to improper disposal of hazardous substances.
8. Nobis concluded that because of the potential for undocumented releases of hazardous substances in the past or in the future, pipes that were previously noted emitting unknown solvent odors and elevated total volatile organic readings in the former chemical mixing room in the southwestern portion of the building, represented a REC because the source of these odors (which may be a formerly closed in-place acetone tank) is not known.
9. Two former press pits located within the manufacturing slab that are stained with petroleum represent a REC and they have not been fully assessed.
10. A pile of unknown debris located on the southwestern portion of the Site that is consistent with demolition debris represents a REC for its potential to contain waste ACM, lead-based paint [and other contaminants].

3.5 OTHER POTENTIAL ENVIRONMENTAL CONDITIONS

Based on review of the April 2011 Nobis Phase I ESA, Credere offers the following list of other potential environmental conditions:

1. Nobis documented that the Site was previously used for railroad activities beginning as early as 1892. Nobis concluded that this past use of the Site represented a *de minimus condition* because any contaminants from this past use would likely be deemed a background condition. Credere agrees with this assertion, but it is our opinion that non-point source background type releases of polycyclic aromatic hydrocarbons (PAHs), which would be considered the primary contaminants of concern, still represent a threat to human health if they exceed applicable NHDES soil standards and should be addressed as a part of future redevelopment of the Site. Therefore, it is our opinion that this past use should be assessed as a part of this Phase II ESA.
2. Nobis indicated in the Phase I ESA that a data gap in the form of snow cover was noted during the Site Reconnaissance portion of their work. Given that the ground could not be observed, confirmation of exterior surface conditions is warranted prior commencing Phase II ESA field work because additional RECs may be identified that require assessment. As a result, a supplemental Site Reconnaissance is recommended as the first task of this Phase II ESA.
3. A former electrical transformer was located on the west side of the building. No information was noted in the previous reports regarding the potential presence of polychlorinated biphenyl's (PCBs) in the transformer. However, considering the timeframe when this transformer would have been in operation, it may have contained



PCBs. For this reason, the potential presence of PCBs is a concern that should be considered.



4. POTENTIAL REDEVELOPMENT SCENARIO

The City of Franklin has partnered with LRPC to assess the Site so that the City can redevelop the Site into a new Water Department Building or other related municipal use.



5. CONCEPTUAL SITE MODEL

The Conceptual Site Model (CSM) includes a description of source areas and/or RECs, the nature and extent of the identified or suspected releases, potential contaminants of concern (COCs), impacted media, transport mechanisms, and potential human and environmental receptors.

5.1 CONTAMINANTS OF CONCERN

Based on the findings of the Phase I ESA, the following potential COCs were identified for each of the RECs identified at the Site:

Table 1 - Potential Contaminants of Concern (COCs)			
REC	REC Nickname	COCs	Potential Source
REC-1	UST-6	Volatile Organic Compounds (VOCs) PAHs Total Petroleum Hydrocarbons (TPH)	Fuel oil releases from the closed in-place 12,000 gallon UST (UST-6)
REC-2	UST-8	VOCs PAHs TPH	Fuel oil releases from the potential presence of a 4,000 gallon UST (UST-8) that was documented on a 1964 Sanborn Map.
REC-3	UST-9	VOCs PAHs TPH	Fuel oil releases from the potential presence of a 4,000 gallon UST (UST-9) that was documented on a 1929 Sanborn Map.
REC-4	Filling Caps and Dispenser near UST-6	VOCs Lead (Pb)	Gasoline from past releases related to the use of filling caps, a dispenser, and undocumented USTs.
REC-5	Degraded Gasoline in MW-2	VOCs Pb	Gasoline from an undocumented source.
REC-6	Concrete Septic Tank	VOCs Semi-volatile Organic Compounds (SVOCs) PCBs RCRA 8 Metals (Ag, As, Ba, Cd, Cr, Hg, Pb, and Se) Plus Zn, Cu and Ni	Tannery and multi-layered board manufacturing operational waste disposed of in the septic system and released to the environment.
REC-7	Buried Tannery Waste	VOCs PAHs RCRA 8 Metals (Ag, As, Ba, Cd, Cr, Hg, Pb, and Se) Plus Zn, Cu, and Ni	Leaching of contaminants from tannery wastes disposed outside around the building.
REC-8	Solvent Odor Emitting Pipe in Chemical Mixing Room	VOCs	Acetone releases from a pipe that may lead to a previously closed in-place tank.



Table 1 - Potential Contaminants of Concern (COCs)			
REC	REC Nickname	COCs	Potential Source
REC-9	Press Pits	VOCs SVOCs TPH PCBs RCRA 8 Metals (Ag, As, Ba, Cd, Cr, Hg, Pb, and Se) Plus Zn, Cu, and Ni	Releases of tannery and/or multi-layered board manufacturing waste through the press pits.
REC-10	Building Waste Debris Pile	Lead Asbestos Containing Materials (ACMs) PCBs	Building materials containing hazardous substances.
Other	Rail road Use of Site	PAHs	Non-point sources releases associated with past rail road activities on the Site.
Other	Transformer	PCBs	Former Electric Transformer on the west side of the building.

5.2 GEOLOGY, STORM WATER FLOW, AND GROUNDWATER FLOW

The Site is located near the west bank of the Pemigewasset River in a developed area of Franklin, New Hampshire. The eastern adjoining property between the Site and river is currently undeveloped. Area topography generally slopes east towards the river. Storm water at the Site likely follows surficial topography and is directed into the river.

Soil boring data contained in reports that were reviewed as a part of the Phase I ESA indicate that soil beneath the Site consists of a mixture of sand and gravel fill material with some areas of tannery waste (hides) noted on the east side of the Site. A fine to medium light brown to dark brown sand horizon is present below the fill horizon. The water table was identified between 28 and 30 feet below the surface.

5.3 DEFINITIONS OF EXPOSURE PATHWAYS AND POTENTIAL RECEPTORS

To aid in a thorough understanding of the environmental concerns present at the Site, a graphical presentation of the identified COCs and potential migration pathways to receptors is included as **Figure 4**. Exposure Pathways and Potential Receptors depicted on the CSM are defined as follows.

Exposure Pathways describe how a human or environmental receptor comes into contact with contaminants which may be present at the Site. Exposure pathways presented in the CSM include the following:



- **Inhalation:** This pathway is primarily associated with groundwater where petroleum contaminated groundwater is within 30 feet (horizontally or vertically), or non-petroleum contaminated groundwater is within 100 feet (horizontally or vertically) of an occupied structure, or when depth to groundwater is unknown. In addition, this pathway is applicable when receptors may incidentally inhale impacted media in the form of dust, vapor, or airborne particulates.
- **Dermal Absorption:** Exposure via dermal absorption occurs when receptors are exposed to chemical concentrations present in soil, groundwater, or surface water through direct contact with the skin.
- **Active Ingestion:** The active ingestion pathway represents exposure which may occur through the active ingestion of contaminant concentrations via a drinking water supply well or through agricultural products.
- **Incidental Ingestion** This pathway is applicable when receptors may incidentally ingest impacted media in the form of dust or airborne particulates.

Potential Receptors are categorized by duration of exposure and intensity of use at the Site. The receptor categories for this project described in the CSM include the following:

- **Commercial:** Commercial receptors are those which are present at the Site for long durations but with low intensity exposure such as indoor office workers.
- **Site Worker:** Site workers are present at the Site for short durations though intensity of use is high, such as during non-routine activities including construction or utility work. Examples include outdoor commercial workers and construction workers.
- **Visitor:** Visitors are characterized by low duration, i.e. less than two hours per day, and low intensity usage such as that which would occur during activities such as walking, shopping, and bird watching.
- **Terrestrial and Aquatic Biota:** These receptors include flora and fauna which may be exposed to contaminants in their respective land-based or aquatic environments.



Based on known history of the Site, the identified contaminants of concern may have been released to the environment through surficial and subsurface releases associated with previous industrial usage, filling with tannery waste, the degradation of potentially hazardous building materials, and previous bulk petroleum storage. Primary impacted media at the Site include surficial and subsurface soil, and groundwater.

Contaminants present in surficial soil may migrate through aeolian dispersion and impact off-site receptors, or they may affect subsurface conditions through infiltration and leaching. Similarly, releases which have impacted subsurface soil conditions have the potential to infiltrate and leach to the overburden groundwater aquifer. Following the redevelopment of the Site, impacted groundwater presents a potential risk to indoor air conditions through contaminant migration via vaporization and diffusion.

Human receptors identified for the Site include future commercial workers at the Site, site workers (i.e. construction workers, etc.), and visitors. Also, current and future potential environmental receptors include the Pemigewasset River and terrestrial and aquatic biota.

Exposure pathways to commercial workers at the site and visitors potentially include dermal absorption and incidental ingestion during routine activities. An inhalation exposure pathway also may exist for commercial workers through vapor intrusion. Site construction workers may be exposed through dermal absorption, inhalation, and incidental ingestion during excavation activities. Terrestrial and aquatic biota are potentially exposed through active ingestion of impacted groundwater; however, the active ingestion pathway does not currently nor is expected to pose a potential future pathway for human receptors, as the Site and adjacent properties in the vicinity are served by municipal water.



6. SAMPLING DESIGN

The following section describes the proposed work to be conducted to confirm or dismiss each REC and includes the number of samples that will be collected during the Phase II ESA sampling program along with the proposed analysis. Proposed sample locations are depicted on **Figure 3** and sampling methodologies are described in **Section 7**. Detailed descriptions of laboratory methods are included on attached **Table 2**.

- REC-1: This REC will be assessed by first conducting a ground penetrating radar (GPR) survey of the area thought to contain the closed in-place 12,000-gallon fuel oil UST (UST-6) to determine its exact location. One test pit (CA-TP-7) will be dug immediately adjacent to the tank (**Figure 3**). During test pit activities, removed soil will be field screened with a properly calibrated photoionization detector (PID) at 1 to 2 foot depth intervals. Multiple test pits may be needed next to the tank to thoroughly evaluate the condition of the soil around it. One soil sample will be collected from the test pit (and interval) exhibiting the highest PID measurement. If PID measurements indicate that little to no residual VOCs are present, then a sample will be collected from adjacent to the tank, but at the same level as the bottom of the tank. For this and all in-place tanks that are being evaluate via test pit work, please note that Credere will consider the position of the tank when digging test pits. If the tank appears to be tipped (i.e. not horizontal), then Credere will direct the digging of the test pit along the deepest side or end. The collected soil sample will be sent to Absolute Resource Associates, LLC (Absolute Resource) of Portsmouth, NH to be analyzed for VOCs, PAHs, and TPH.
- REC-2: This REC will be assessed by first conducting a GPR survey of the area thought to contain the 4,000-gallon fuel oil UST (UST-8) to determine its location. If the tank has not been removed, one test pit (CA-TP-8) will be dug immediately adjacent to the tank to assess the condition of the soil adjacent to the tank (**Figure 3**). Similar to above, multiple test pits may be needed next to the tank to thoroughly evaluate the condition of the soil around it. If the tank appears to have been removed and the tank grave can be discerned, Credere will install the test pit directly in the center of the perceived tank grave. During test pitting, removed soil will be field screened with a properly calibrated PID. One soil sample will be collected from the test pit (and interval) exhibiting the highest PID measurement. If PID measurements indicate that little to no residual VOCs are present, then a sample will be collected from adjacent to the tank, but at the level of the bottom of the tank. The collected soil sample will be sent to Absolute Resource to be analyzed for VOCs, PAHs, and TPH.
- REC-3: This REC will be assessed by first conducting a visual inspection for evidence (e. g. fill, vents, concrete scars) of UST-9, a 4,000-gallon fuel oil UST. If the visual survey proves unsuccessful, a GPR survey will be conducted to locate it. If the tank has not been removed, portions of the concrete slab will be removed and one test pit (CA-TP-9) will be dug immediately adjacent to the tank to assess the condition of the soil



adjacent to the tank (**Figure 3**). Multiple test pits may be needed next to the tank to thoroughly evaluate the condition of the soil around it. If the tank appears to have been removed and the tank grave can be discerned, Credere will install the test pit directly in the center of the perceived tank grave. During test pitting, removed soil will be field screened with a properly calibrated PID at 1 to 2 foot depth intervals. One soil sample will be collected from the test pit (and interval) exhibiting the highest PID measurement. If PID measurements indicate that little to no residual VOCs are present, then a sample will be collected from adjacent to the tank, but at the level of the bottom of the tank. The collected soil sample will be sent to Absolute Resource to be analyzed for VOCs, PAHs, and TPH.

If a tank or evidence of tank is not identified during this work, five shallow soil probes (CA-SG-1 to CA-SG-5) will be installed through the concrete slab to collect soil gas to determine if the vadose zone under the slab has been impacted by contamination from the tank (i.e. looking for light component of fuel oil - naphthalene). The collected soil gas will be analyzed with a properly calibrated PID equipped with an 11.7 electron volt lamp. If this work reveals that volatile organics are present, one soil gas sample will be sent to Columbia Analytical for VOC analysis to identify the number and type of VOCs present. In this scenario, Credere will then install a boring (CA-1) adjacent to the soil gas probe exhibiting the highest result. During drilling, if field screening results indicate that soil contamination appears to extend down to the water table, then a groundwater monitoring well would be installed at this location. A groundwater sample will be collected to assess the groundwater for VOCs. In the absence of detectable concentrations of VOCs and other direct evidence of a tank, Credere will rely on groundwater quality data collected from MW-2 (which will be sampled as a part of assessing REC-5) to assess the potential for a source area of contamination to be present as part of this REC.

REC-4: The REC associated with the ‘fill caps and dispenser’ will be assessed by first conducting a GPR survey and/or electronic line tracing to determine if any tanks remain associated with this equipment. If this proves unsuccessful, then the filling caps will be dug up and manually traced. If a tank or tanks are identified associated with the fill caps, two test pits (CA-TP-10 and CA-TP-11) will then be dug adjacent to the tanks in a fashion similar to that described for REC-1, REC-2, and REC-3. For the purposes of this investigation, Credere has assumed that two tanks will be found and one test pit will be dug next to each. However, Credere also anticipates that one additional test pit will be dug in the location of the former dispenser (CA-TP-12). During test pitting, removed soil will be field screened with a properly calibrated PID at 1 to 2 foot intervals. Only the soil sample collected from each test pit (and interval) exhibiting the highest PID measurement will be sent to the laboratory. If PID measurements indicate that little to no residual VOCs are present, then one sample will be collected from the fill/native soil interface below the dispenser or the perceived level of bottom of any tank or tank grave(s). All collected soil samples will be sent to

Absolute Resource and analyzed for VOCs and lead. If no evidence of tanks are identified in this area, then only the one test pit (CA-TP-12) will be dug.

REC-5: The REC associated with the degraded gasoline identified in existing monitoring well MW-2 will be assessed in the following fashion:

- 1) Collect a groundwater sample from existing monitoring well MW-2 using low flow sampling techniques and analyze it for VOCs and dissolved lead.

Depending on the results of work conducted to assess REC-3, one of the following two paths will be taken to further evaluate REC-5 if the presence of degraded gasoline is confirmed in monitoring well MW-2.

- 2) Credere will evaluate field data collected from soil probes (CA-SG-1 to CA-SG-5) and monitoring well CA-1 to help locate the potential gasoline source for the MW-2 contamination. The results of this evaluation will be included in the Phase II ESA along with a specific recommendation for future work, if needed.
- 3) If the soil probes and upgradient monitoring well work was not completed as a part of the assessment of REC-3, then Credere will install five soil gas probes (CA-SG-1 to CA-SG-5) to locate the gasoline source under the building slab upgradient from MW-2, if it is present. Monitoring well CA-1 will be installed in the perceived area of highest VOC contamination based on field screening. CA-1 will be installed upgradient of MW-2 even if the soil gas results do not show the presence of a soil gas source under the building. As indicated as a part of the investigation of REC-3, one soil gas sample collected from the soil probes and will be sent to the laboratory for VOCs analysis. In addition, one soil sample will be collected from CA-1 and submitted to the laboratory for analysis of TPH, VOCs, and lead by EPA 200.7/6010 and one groundwater sample will be collected from CA-1 and submitted to the laboratory for VOCs and dissolved lead by EPA 200.7/6010.

REC-6: The REC associated with the suspected concrete septic tank will be assessed by first conducting a GPR survey to locate the tank. Once the tank is located, a test pit (CA-TP-13) will be dug immediately adjacent to the tank. Multiple test pits may be needed next to the tank to thoroughly evaluate the condition of the soil around it. During test pitting, removed soil will be field screened with a properly calibrated PID and X-Ray Fluorescence Meter (XRF) at 1 to 2 foot depth intervals. However, only one soil sample will be collected from the test pit (and interval) exhibiting the greatest evidence of contamination [visually, olfactory, PID and/or XRF measurement]. In the absence of evidence of contamination, the sample will be collected from adjacent to the tank, but at the level of the bottom of the tank. The collected soil sample will be sent to Absolute Resource to be analyzed for VOCs, SVOCs, PCBs and RCRA 8 Metals plus Zn, Cu, and Ni. In addition, if present, one (1) sample of sludge will be collected from the tank and analyzed for VOCs, SVOCs, PCBs, and RCRA 8 Metals plus Zn, Cu, and Ni. This sample will be field screened as described above.

The test pit work will help assess if there are pipes directed towards an old leachfield or discharge area. Additional geophysical work may be conducted to trace out the observed discharge line. At the terminus of the discharge, a boring (CA-2) will be drilled by T&K Drilling. During drilling, soils will be screened in the field via a properly calibrated PID and XRF. These samples will be visually inspected for evidence of contamination. In addition, the soil sample exhibiting the highest PID measurement or greatest visual extent of contamination, during drilling will be sent to the laboratory for analysis of VOCs, SVOCs, PCBs, and RCRA 8 metals. The boring will be finished in the water table and a monitoring well will be constructed to allow for the collection of a groundwater sample at this location. Groundwater will be collected from this location and will be analyzed for VOC, SVOCs, and dissolved RCRA 8 Metals plus Zn, Cu, and Ni. However, no PCB samples will be collected from the groundwater because we do not expect this parameter to dissolve in groundwater.

- REC-7: The REC related to the presence of buried tannery waste (hides) will be assessed by first conducting a GPR survey to define, if possible, the horizontal extent of waste on the Site. Following the completion of GPR, up to five test pits (CA-TP-14 to CA-TP-18) will be dug to confirm the GPR results and to define the vertical limits of waste observed. Two of these test pits will be dug near the northern property boundary to determine if these wastes extend off-site. During test pitting, soil will be visually screened and evaluated with a properly calibrated PID and XRF. One representative soil sample from each test pit will be collected for laboratory analysis. Soil in direct contact with tannery waste will be selected to represent a worst case scenario. Samples will be sent to the Absolute Resource for VOCs, PAHs, and RCRA 8 Metals plus Zn, Cu and Ni.
- REC-8: The REC associated with the pipes in the former chemical mixing room that was previously emitting solvent odors will be assessed by installing six (6) sub-slab soil gas probes (CA-SG-6 to CA-SG-11) to determine if past or present solvents being emitted are impacting the vadose zone beneath the concrete slab with VOCs. Locations CA-SG-6 to CA-SG-11 will be installed to evaluate this REC. Representative soil gas will be removed from each soil gas point and screened in the field with a properly calibrated PID. The sample with the highest PID measurement will sent to the laboratory for VOC analysis to determine the number and type of VOCs present. In addition, any piping present on-Site noted near closed in-place UST#2 would be field screened with a PID for evidence of solvent emission. If this data shows that the source of the solvent odors is likely a previously closed in-place tank, then a recommendation to remove the tank will be made in the Phase II ESA Report. A soil boring and monitoring well (CA-3) will be installed to assess the impact to Site soil and groundwater adjacent to the soil gas probe location that exhibited the highest field screening results. CA-3 will be installed at the location shown on **Figure 3** if no vapor source is identified during the soil probe work. This work will include appropriate



field screening as described above and the collection of one (1) soil sample and one (1) groundwater sample for laboratory VOC analysis.

REC-9: The REC associated with the press pits will be assessed by advancing two (2) soil shallow borings (CA-4 and CA-5) through the concrete bottom of the pits. Soil samples will be collected directly under the concrete slab to determine if substances added to the pit have released to the soil beneath the pits. Collected samples will be field screened with a PID and grab samples will be collected from both borings and laboratory analyzed for TPH, VOCs, RCRA 8 Metals plus Zn, Cu, and Ni, SVOCs, TPH, and PCBs.

REC-10: The REC associated with the debris pile identified on the concrete slab by Nobis will be assessed by collecting representative building material samples (Debris-1, etc.) and analyzing them for lead, asbestos, and PCBs. To accomplish this work, Credere will retain the services of Absolute Resources because the pile likely contains asbestos and sampling of the waste must be accomplished in accordance with NHDES requirements.

OTHER: Due to snow cover present during Nobis's Phase I Site Reconnaissance, the first task of this Phase II ESA will be to perform a supplemental Site Reconnaissance so that the ground can be viewed for evidence of RECs. If additional RECs are identified, the work plan will be revised and reappraised.

The former use of the Site for rail road activities will be assessed in two parts:

- 1) Credere will review available topographic maps and other historic sources to determine the former location of the rail road activities on-Site, and
- 2) Collect three (3) surficial samples (CA-SS-1, CA-SS-2, and CA-SS-3) from this area and analyzing them for PAHs. The locations shown on **Figure 3** are approximate and once research is completed the location of these surficial soil samples will be moved to the former rail-road right-of-way. If it cannot be location of the former right-of-way determined, the locations depicted on **Figure 3** will be sampled.

The former use of the of a potential PCB containing transformer on the west side of the Site building will be assessed by collecting two concrete samples of the transformer pad from immediately adjacent to the former transformer's location and analyzing them for PCBs.

Credere considered the collection of a background soil sample from the Site for metals and PAH analysis, but in order to be valid this would have to be from an area likely not impacted by previous site activities. Because all areas of the Site may have been impacted by previous historical site uses, no background sample is specified for this Phase II ESA. A background study with nearby off-site sampling may be recommended if the soil appear to be impacted by metals and/or PAHs within expected background ranges. However, during soil sampling work



the on-site geologist will be evaluating all collected soil samples for evidence of anthropogenic and native minerals that may present in samples because they could affect the laboratory results. These findings of this work will be discussed in our Phase II Report. Additional detail on what the field geologist will be looking for in soil samples is discussed in **Section 7**.

Table 2 includes the number and type of samples that are proposed be collected, cross-referenced with the appropriate standard operating procedure (SOP) that will be used from Credere’s Generic QAPP. The data collected from these activities will serve as the basis for evaluating the Site conditions and will determine if any additional subsurface investigation and/or remedial actions are necessary. **Figure 3** shows the proposed locations where media will be collected. Requirements relative to Chain of Custody, Data Management and Documentation, Data Validation, and Data Usability Assessments contained in the Generic QAPP will be followed.



7. FIELD ACTIVITY METHODOLOGY

Field activity methodologies for assessing the RECs and other potential environmental concerns are summarized in the following subsections. Field activities will be conducted in accordance with the SOPs included in Credere's Generic QAPP Rev. 2 (USEPA RFA #08166 and #09036) and the rationale discussed in **Section 6**. Where field observations and/or field screening results indicate the presence of additional source areas or potentially impacted media, additional exploration locations or samples may be added to determine the horizontal and/or vertical extent of contamination. The number and locations of these additional soil samples or exploration locations will be dependent on field data, site constraints and professional judgment. All decisions regarding delineation will be recorded in the field logbook, and all locations will be documented. All soil samples obtained for the purposes of contamination delineation will be collected and field-analyzed in accordance with Credere's standard SOPs outlined on **Table 2**. If Credere determines these additional sample locations should be tested for analytes not described in **Table 2**, the USEPA QA Manager will be contacted, and pending the outcome of the communication, an email update will be provided to the USEPA describing the additional sample analysis, methods and SOPs.

7.1 GPR SURVEY AND ELECTRONIC TRACING

DigSmart of Maine, Inc. will conduct the GPR survey and will perform the work in accordance with the CREDERE SOP #007 provided in the generic QAPP.

7.2 SOIL BORINGS, TEST PITTING, SURFICIAL SOIL SAMPLING, AND FIELD SCREENING

Soil samples will be collected from soil borings and test pits in accordance with NHDES SOP HWRB-11 for soil sampling. Soil borings will be completed using hollow stem auger methodologies. During soil boring advancement, soil samples will be continuously collected in 2-foot split spoon samplers. During test pitting, soil samples will be collected from each approximate 2-foot depth interval. Each soil sample will be individually logged, and evidence of contamination will be noted. If changes in strata occur (e.g. a transition from sand to clay), the sample may be appropriately divided or additional samples may be collected and individually logged. To assist with the data evaluation, the field geologist will note the presence of materials that may affect laboratory soil sample results on the logs. These materials include:

- Presence or absence of mineral soils (i.e. pyrite or other likely arsenic bearing minerals as well as abundant micas)
- Presence or absence of asphalt
- Presence or absence of coal clinkers, coal ash or wood ash

Each collected soil sample will then be field screened for total VOCs using a photoionization detector (PID) in accordance with NHDES SOP HWRB-12. The PID will be calibrated with a 100 part per million by volume (ppm_v) isobutylene gas and a response factor of 1.0. Soil



samples collected for laboratory analysis of VOCs will be collected and preserved in accordance with NHDES SOP VOCs/Soil-2000.

Each soil sample will also be field screened for RCRA-8 metals with an Innov-X[®] Alpha 4000 portable X-ray fluorescence meter (XRF) in accordance with Maine DEP SOP DR#025. Because this is a screening level event to determine which, if any, metals are present in Site soil, samples will be screened directly in the collected polybags. Prior to XRF screening, sample bags will be kneaded to homogenize the soil.

Soil samples will be collected in accordance with NHDES SOP HWRB-11. Surficial soil samples from soil borings will be collected from the 0 to 2 foot below ground surface (bgs) interval. Surficial soil samples collected from hand dug locations will be collected from the 0 to 1 foot bgs interval. In each case, asphalt and base materials, landscaping materials, and other organic detritus will be removed prior to sampling.

Subsurface soil samples will be selected based on the sampling rationale discussed in **Section 6**. In general, subsurface soil samples will be selected based on the highest PID field screening result for VOCs, where visual or olfactory observations of contamination are observed, or from the depth of the water table interface if no other indicators of contamination are noted.

7.3 MONITORING WELL INSTALLATION AND SURVEY

7.3.1 Monitoring Well Installation and Development

As discussed in **Section 6**, selected soil borings will be completed as groundwater monitoring wells per EPA SOP EPASOP#2048. Unless Site conditions (e.g. shallow bedrock or confining layers) warrant modified well construction, each monitoring well will be installed using 2.0-inch PVC well materials including a well point or cap installed to span the water table, 10 feet of 0.010-inch slotted pipe, and solid pipe to grade. Washed No. 1 size silica sand will be placed around the slotted pipe to an elevation approximately two (2) feet above the slotted pipe to establish a well annulus. At least one (1) foot of bentonite chips will be placed on top of each well annulus to prevent stormwater infiltration into the wells. Each well will then be completed to grade with a flush mounted road box protected by a minimum of a 1-foot by 1-foot concrete pad. Each monitoring well will be developed by over pumping and agitation, and then allowed to equilibrate for at least two weeks prior to sampling.

7.3.2 Rod and Level Survey

Following installation, each monitoring well will be surveyed with a rod and level to determine the top of well and ground elevation at each location in accordance with ASTM E 1364-95. The monitoring wells will be tied to the existing monitoring well network at the Site and the west adjoining property.



7.4 GROUNDWATER SAMPLING

Prior to sampling, the depth to groundwater will be measured (and the presence of free floating product will be noted) in each groundwater monitoring well in accordance with NHDES SOP HWRB-1 to allow for the calculation of groundwater elevations and the determination of groundwater flow direction and gradients. Each monitoring well will then be sampled using low-flow sampling techniques and dedicated equipment in accordance with NHDES SOP HWRB-9. Each groundwater sample will be collected directly in laboratory glassware and will be submitted for off-site laboratory analysis.

7.5 SOIL GAS SAMPLING

NHDES's July 2006 Vapor Intrusion Guidance will be consulted for the installation of soil gas points and collection of representative soil gas samples. However, because this work is for evaluating sources of contaminants, not evaluating for vapor intrusion, we plan to deviate from the guidance and install points just beneath the concrete slab which is closer to the suspected source elevation. Probes will be situated 1 to 2 feet beneath the slab. Probes will be installed with a properly decontaminated hand held hammer drill. Once in place, probes will be sealed in place with modeling clay. To ensure that valid soil gas samples are collected as part of assessment, the use of a tracer compound may be used to assess for surface/annular seal leaks around the top of the soil gas probe. Before representative samples are collected between 1 and 5 purge volumes of soil gas will be removed from each probe. Samples will be evaluated in the field with a properly calibrated PID. Laboratory samples will be send to Columbia Analytical Laboratories for TO-15 Analysis for VOCs.

7.6 BUILDING MATERIAL DEBRIS SURVEYS

A NHDES certified asbestos inspector will be contracted to perform an asbestos survey of the debris pile at the Site. During the asbestos survey, samples will be collected from each suspect media (i.e. piping insulation, flooring, ceiling tiles) in accordance with NHDES Certified Administrative Rule ENV-1800: Asbestos Management Control. Asbestos samples will be submitted to a State approved laboratory for asbestos analysis by polarized light microscopy (PLM) by USEPA 600/R-93/116.

Based on the age of the building, there is the potential for lead-based paint and PCB building materials to have been used in the former building and may be present in the remaining debris pile. As a result, the building debris pile will be sampled for lead and PCBs using the laboratory analysis listed on **Table 2**.



8. REGULATORY STANDARDS

Sample results will be compared to the applicable state and/or federal standards/guidelines described below.

8.1 SOIL

Concentrations in soil samples will be compared to New Hampshire's Soil Remediation Standards (SRS) detailed in NHDES Env-Or 600 Contaminated Site Management. Where guidelines are not available, soil concentrations will be compared to other appropriate regulatory standards and guidelines, e.g., USEPA Regional Screening Levels (RSL) for Chemical Contaminants at Superfund Sites, RSL Table Updated April 2009 for soil. If standards or guidelines do not exist, action levels will be triggered if the sample analytical results exceed available published background levels. Please note that concentrations of total chromium identified at the Site will be compared to the SRS for hexavalent chromium as this elemental oxidation state is the applicable toxicological driver.

8.2 GROUNDWATER

Groundwater sample results will be compared to the New Hampshire Ambient Groundwater Quality Standards (AGQS) detailed in NHDES Env-Or 600 Contaminated Site Management and USEPA Maximum Contaminant Levels (MCLs) for Drinking Water. In the event that no AGQS or MCLs exist for a particular contaminant, results will be compared to the USEPA Region 9 RSLs.

8.3 SOIL GAS

Soil gas sample results will not be compared to any NHDES or EPA standard because the results are being used to determine if a source or sources of VOCs are present beneath the slab. Moreover, no building currently exists at the Site or within 100 feet of the slab so data collected will only be used to determine the type of VOCs present and to locate borings and monitoring wells.

8.4 ACM

Asbestos sampling will be completed in accordance with NHDES Certified Administrative Rule ENV-1800: Asbestos Management Control. Levels of asbestos greater than or equal to 1% by volume as determined by weight, visual evaluation, and/or point count analysis within representative bulk samples obtained from the Site will result in the definition of such bulk materials as ACM.



8.5 LEAD-BASED PAINT

Concentrations of lead in paint as determined through the use of a XRF analyzer will be compared to a limit of 1.0 mg/cm² or 0.5% by weight. All construction work involving exposure or potential exposure to lead is covered by the Occupation Safety and Health Administration's Lead in Construction Standard 29 CFR 1926.62.

8.6 PCB BULK PRODUCTS

Building materials that have been analyzed to contain concentrations of total PCBs equal to or in excess of 50 ppm are defined as a PCB bulk product waste in accordance with 40 CFR 761.3. These materials are regulated for disposal under 40 CFR 761.62. Materials that have been analyzed to contain total PCBs at a concentration of equal to or greater than 1 ppm but less than 50 ppm are not regulated by the Toxic Substance Control Act (TSCA) for disposal as long as they remain in use. However, if this material is removed from use, such as through demolition, they must be disposed of at a facility that is licensed to accept this waste. Building materials which have been analyzed to contain total PCBs at a concentration of less than 1 ppm are unrestricted for future use and/or disposal.



9. PROPOSED PROJECT SCHEDULE

The following schedule is proposed for the Phase II ESA activities at the Site:

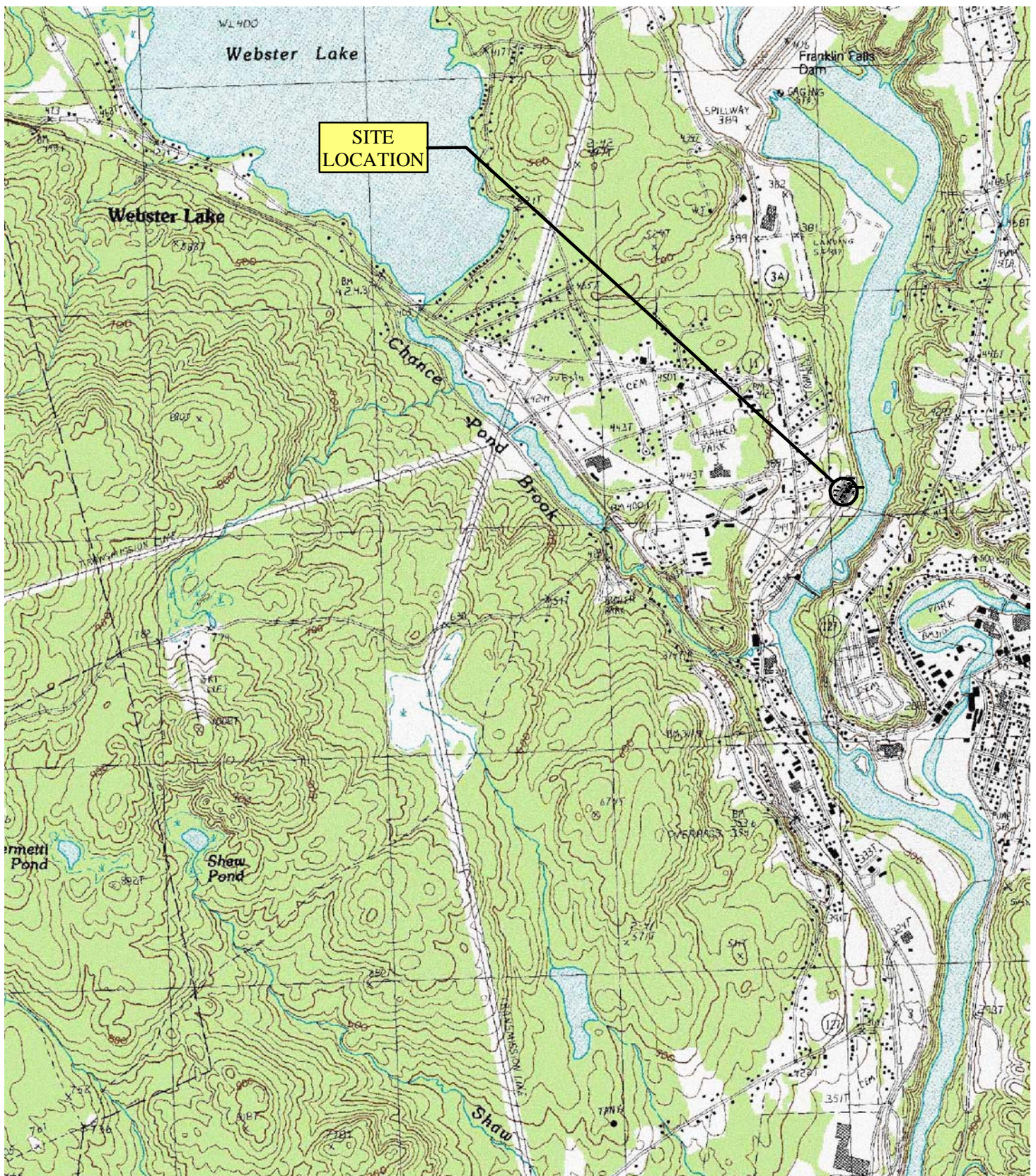
DATE	ACTION
Week of June 27, 2011	Finalize SSQAPP
Week of June 28, 2011	Implement Field Program
Week of July 7, 2011	Receive Laboratory Analytical Data
Week of July 30, 2011	Submit Draft Phase II ESA Report
Week of August 14, 2011	Submit Final Phase II ESA Report



FIGURES

Figure 1Site Location Map
Figure 2 Credere Organization and Responsibility Chart
Figure 3 Proposed Phase II Sample Location Plan
Figure 4Conceptual Site Model





USGS 7.5 MINUTE FRANKLIN, NH QUADRANGLE (1987)

DRAWN BY: SWC/JRN DATE: 12/14/10
 CHECKED BY: RSV/JSS PROJECT: 10001086

FIGURE 1 - SITE LOCATION PLAN



CREDERE ASSOCIATES, LLC
 776 MAIN STREET
 WESTBROOK, MAINE 04092
 TEL: 207.828.1272
 FAX: 207.887.1051
 WWW.CREDERELLC.COM

**FORMER POLYCLAD LANINATES, INC
 PROPERTY**
 45 TANNERY STREET
 FRANKLIN, NH
 NHDES #199902062

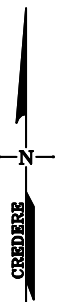
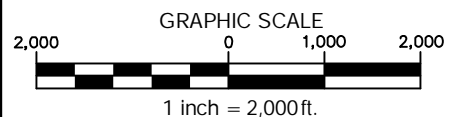
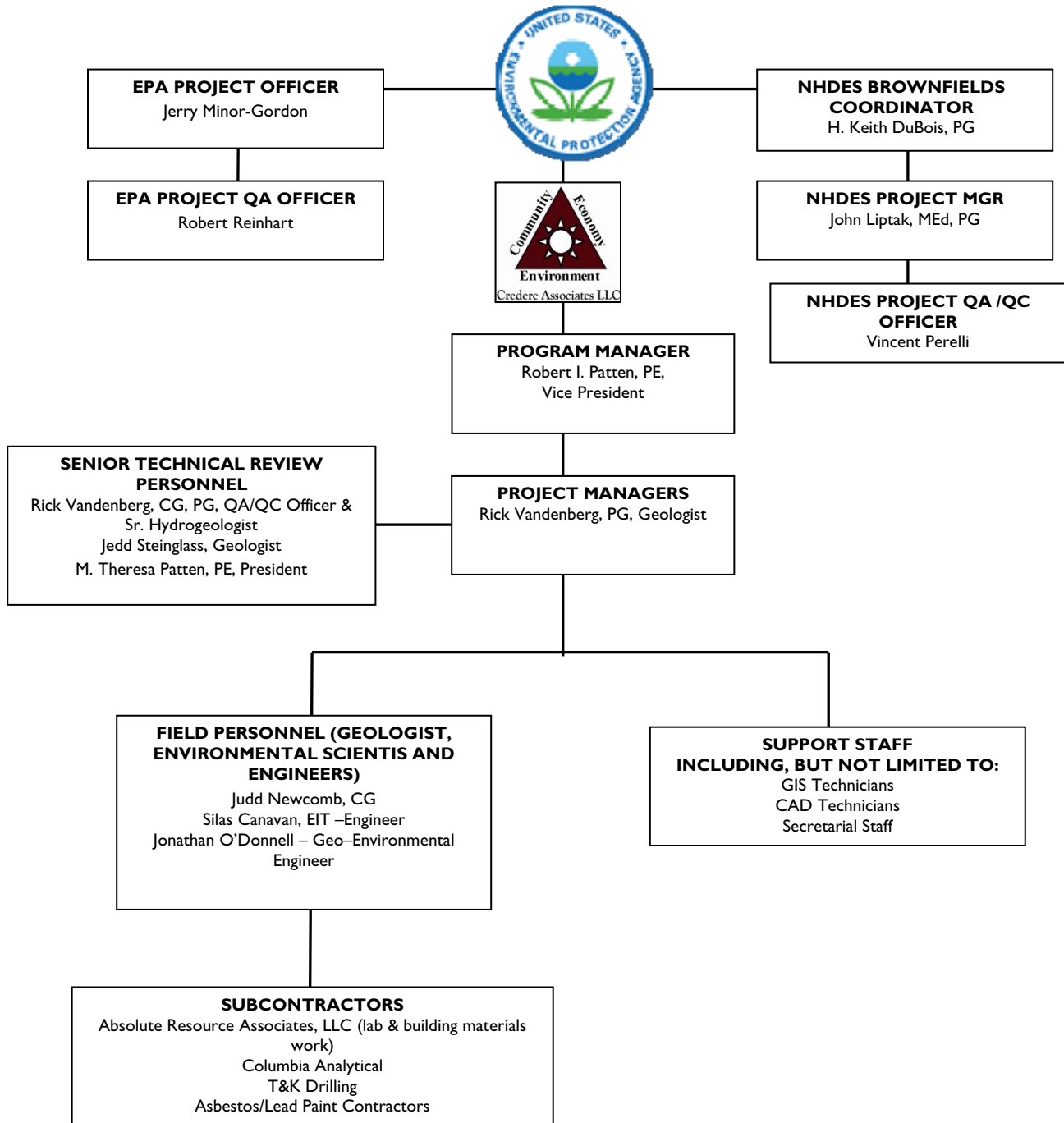
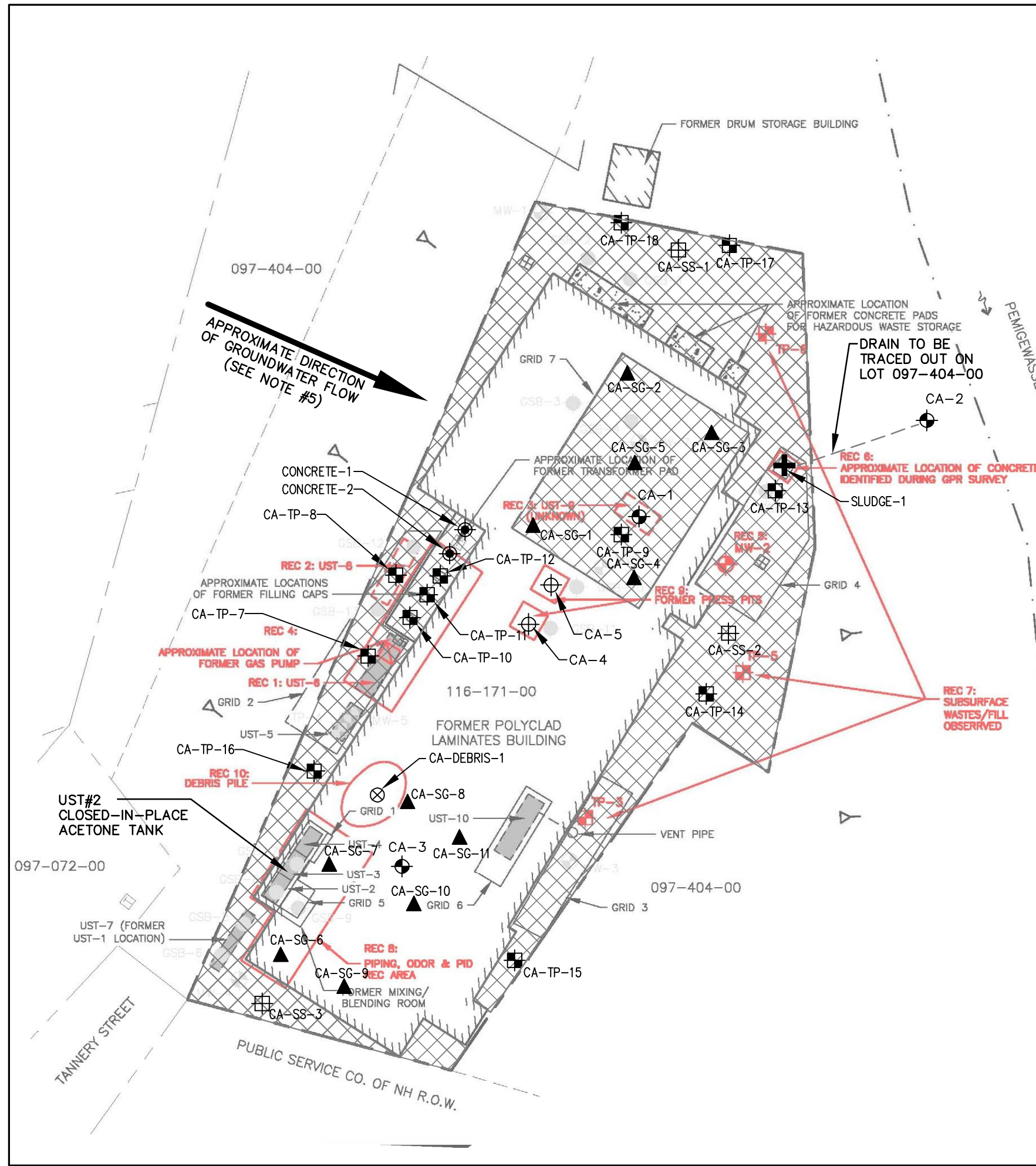


Figure 2: Credere Organization and Responsibility Chart





- NOTES:
- EXISTING CONDITION INFORMATION SHOWN ON THIS PLAN IS FROM A PLAN TITLED "FIGURE 5, SITE RECS" DATED APRIL 2011 BY NOBIS ENGINEERING, INC. (NOBIS) OF CONCORD NEW HAMPSHIRE.
 - LOCATIONS AND SITE FEATURES DEPICTED ON THIS PLAN ARE APPROXIMATE AND HAVE NOT BEEN FIELD-VARIFIED BY CREDERE ASSOCIATES.
 - ITEMS SHOWN IN RED ARE RECS DISCUSSED IN THE PHASE I DOCUMENT PREPARED BY NOBIS ON APRIL 11, 2011.
 - SAMPLE LOCATIONS MAY BE ADJUSTED. SEE SSQAPP TEXT FOR EXPLANATION.
 - GROUNDWATER FLOW DIRECTION WAS TAKEN FROM THE APRIL 12, 2008 DELTA CONSULTANTS PHASE II ENVIRONMENTAL SITE ASSESSMENT REPORT.

EXISTING FEATURES AND PREVIOUS WORK LEGEND

SYMBOL	DESCRIPTION
—	PROPERTY BOUNDARY
- - -	OTHER PROPERTY LINE
- · - · -	EASEMENT
—	STREAM/RIVER
116-171-00	MAP AND LOT NUMBER
→	RIVER FLOW DIRECTION
⊠	FORMER TRANSFORMER PAD
⊞	CATCH BASIN
∇	SLOPE
⊕	MONITORING WELL (2008)
⊞	TEST PIT (2008)
●	SOIL BORING (2008)
⊠	APPROXIMATE GPR SURVEY GRID LOCATION
■	FORMER UST LOCATION (CONFIRMED)
□	FORMER UST LOCATION (UNCONFIRMED)

PROPOSED SAMPLE WORK LOCATION LEGEND

SYMBOL	DESCRIPTION
⊕	MONITORING WELL
⊞	SOIL BORING
⊞	SURFICIAL SOIL SAMPLE
⊞	TEST PIT
▲	SOIL GAS POINT
⊗	DEBRIS PILE SAMPLE
⊕	CONCRETE SAMPLE
+	SLUDGE SAMPLE FROM TANK
⊞	LIMIT OF GROUND PENETRATING RADAR SURVEY

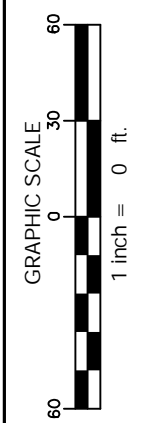
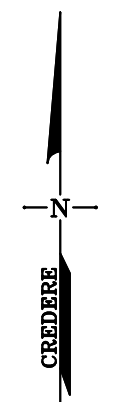
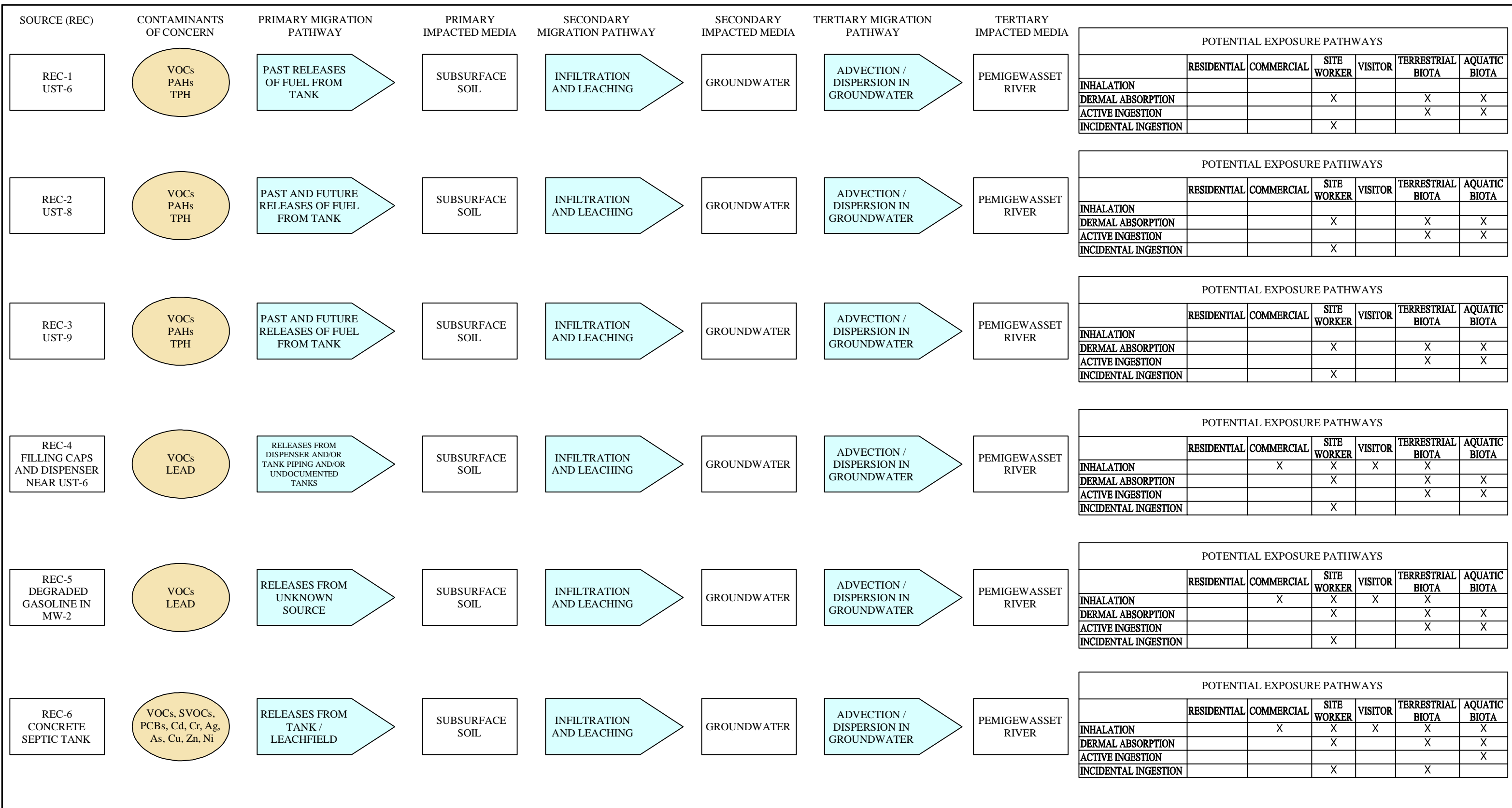


FIGURE 3
PROPOSED PHASE II SAMPLE LOCATION PLAN

FORMER POLYCLAD LAMINATES, INC PROPERTY
45 TANNERY STREET
FRANKLIN, NH
NHDES #199902062

DRAWN BY: SWC DATE: 5/5/11
CHECKED BY: RSV PROJECT: 10001086

CREDERE ASSOCIATES, LLC
776 MAIN STREET
WESTBROOK, MAINE 04092
FAX: 207.887.1051
TEL: 207.828.1272
WWW.CREDERELLC.COM



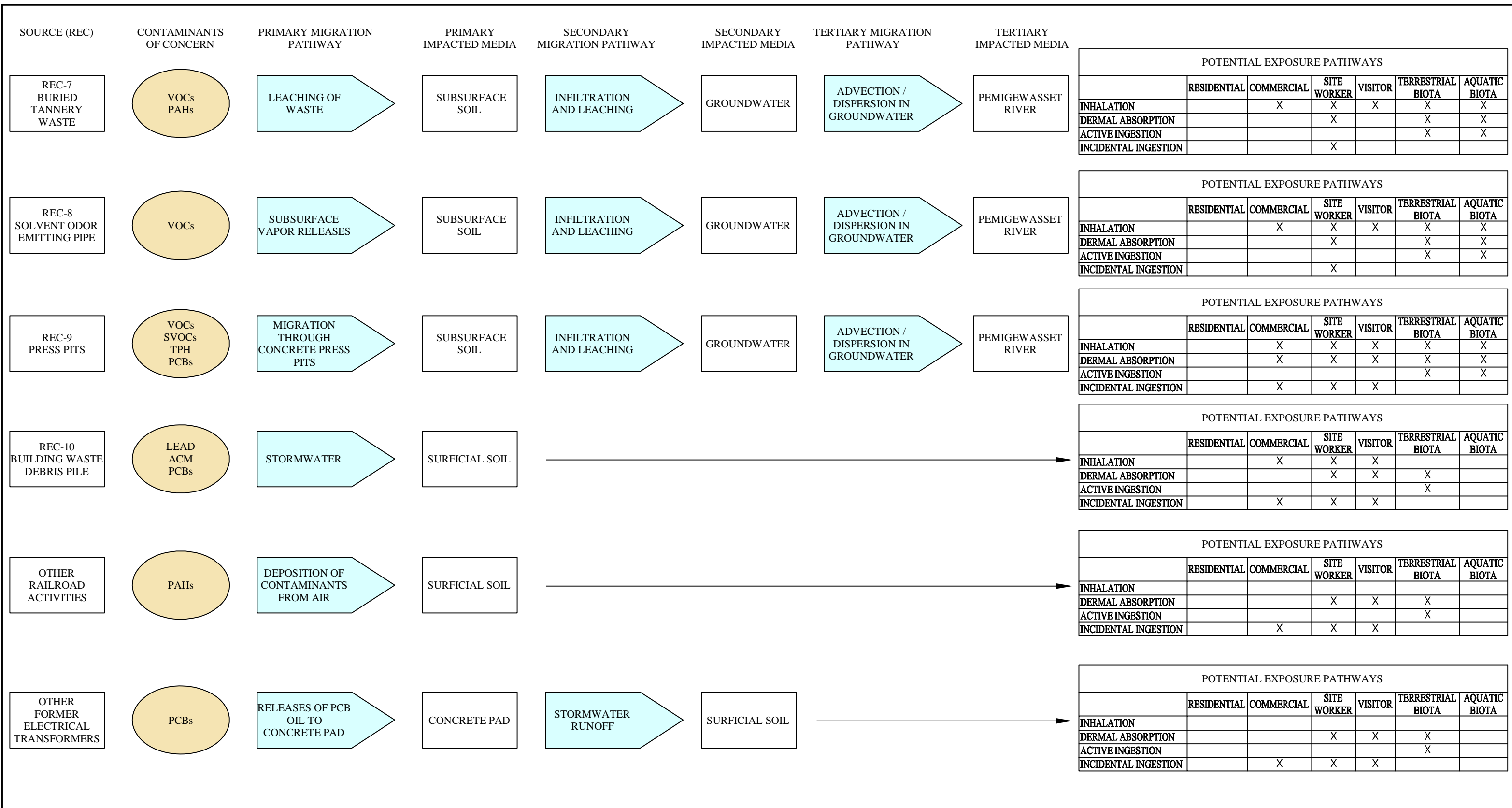
DRAWN BY: RSV	DATE: 5/7/11
CHECKED BY: RIP/JSS	PROJECT: 10001086



CREDERE ASSOCIATES, LLC
 776 MAIN STREET
 WESTBROOK, MAINE 04092
 TEL: 207.828.1272
 FAX: 207.887.1051
 WWW.CREDERELLC.COM

**FIGURE
 CONCEPTUAL SITE MODEL
 SHEET 1 OF**

FORMER POLYCLAD LANINATES, INC PROPERTY
 45 TANNERY STREET
 FRANKLIN, NH
 NHDES #199902062



DRAWN BY: RSV DATE: 5/7/11
 CHECKED BY: RIP/JSS PROJECT: 10001086



CREDERE ASSOCIATES, LLC
 776 MAIN STREET
 WESTBROOK, MAINE 04092
 TEL: 207.828.1272
 FAX: 207.887.1051
 WWW.CREDERELLC.COM

FIGURE
CONCEPTUAL SITE MODEL
SHEET OF

FORMER POLYCLAD LANINATES, INC PROPERTY
 45 TANNERY STREET
 FRANKLIN, NH
 NHDES #199902062

TABLES

Table 1..... Potential Contaminants of Concern
Table 2..... Sample Reference Table



**Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062**

Media to be Collected	Proposed Sample IDs	Associated RECs	Sample Design	Sample Depth (ft bgs)	Field SOPs to be Used	Field Analysis/ Observations	No. of Samples for Analysis	No. of Field Dups	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Surficial Soil	CA-SS-1 CA-SS-2 CA-SS-3	OTHER	Three surficial soil samples will be collected to determine if former rail road operations have impacted the Site.	0-1	Credere-004 HWRB-11 HWRB-12 HWRB-15 HWRB-17 DR#012 DR#024 DR#025	Visual & Olfactory PID Headspace XRF Screening	3	Soil samples will be duplicated at a rate of 5% per the generic QAPP for a total of one (1) based on the proposed total number of samples indicated in this table.	- PAHs by EPA Method 8270	PAHs - 4 oz. amber glass with Teflon-lined cap	RL-13	Absolute Resource Associates (formerly Resource Laboratories)
	Subsurface Soil (Soil Borings & Test Pits)	CA-TP-7*	REC-1	One subsurface sample will be collected during test pit work to determine if the closed in-place 12,000 gallon tank (UST-6) has impacted the subsurface soil beneath the Site.	Soil field screened every 2-foot interval during drilling of borings and at every 1 to 2 feet during test pitting. Laboratory samples collected at highest field screening detection, visual/olfactory evidence of contamination.	Credere-002 Credere-004 HWRB-11 HWRB-12 HWRB-15 HWRB-17 VOCs/SOIL-2000 DR#012 DR#024 DR#025	Visual & Olfactory PID Headspace XRF Screening	1	Soil samples will be duplicated at a rate of 5% per the generic QAPP for a total of one (1) based on the proposed total number of samples indicated in this table.	- TPH by EPA Method 8015 - VOCs by EPA Method 8260 - PAHs by EPA Method 8270	TPH - 4 oz. amber glass VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids PAHs - 4 oz. amber glass with Teflon-lined cap	
CA-TP-8*		REC-2	One subsurface sample will be collected during test pit work to determine if the 4,000 gallon tank (UST-8) remains in place and if any releases have impacted the subsurface soil beneath the Site.	1								
CA-TP-9* or CA-1		REC-3	One subsurface sample will be collected during test pit work to determine if the 4,000 gallon tank (UST-9) remains in place and if any releases have impacted the subsurface soil beneath the Site. If evidence of a tank is not identified during this work, five soil gas probes will be installed and field screen to determine if any VOC sources are present beneath the building. All soil gas samples will be field screened, but one soil gas sample will be collected and sent to the laboratory for analysis to confirm the volatile organics that were measured in the field.	1								
CA-TP-10* CA-TP-11* CA-TP-12*		REC-4	Up to 3 subsurface samples will be collected during test pit work to assess any tank present connected to the "cap and fill" and gasoline dispenser.	3				- VOCs by EPA Method 8260 - Lead by EPA 200.7/6010				VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids -Lead - 4 oz. glass with Teflon-lined cap

**Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062**

Media to be Collected	Proposed Sample IDs	Associated RECs	Sample Design	Sample Depth (ft bgs)	Field SOPs to be Used	Field Analysis/ Observations	No. of Samples for Analysis	No. of Field Dups	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Subsurface Soil (Continued from page 1) (Soil Borings & Test Pits)	TP-13* CA-2	REC-6	Two subsurface samples are anticipated as a part of test pit and boring work that will be conducted and aimed at assessing the suspect concrete septic tank and the terminus of the discharge from the tank.			Visual & Olfactory PID Headspace XRF Screening	2	Soil samples will be duplicated at a rate of 5% per the generic QAPP for a total of one (1) based on the proposed total number of samples indicated in this table.	- VOCs by EPA Method 8260 - RCRA 8 Metal plus Zn, Cu, and Ni by EPA Method 6010 - PAHs by EPA Method 8270 - PCBs by EPA Method 8082	VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids Metals - 4 oz. glass with Teflon-lined cap PAHs - 4 oz. amber glass with Teflon-lined cap PCBs - 4 oz glass with Teflon lined cap	RL-9 RL-5 RL-7 RL-13 RL-4	
	CA-TP-14* CA-TP-15* CA-TP-16* CA-TP-17* CA-TP-18*	REC-7	Up to five (5) subsurface soil samples are anticipated during test pit work aimed at delineating the extent of buried tannery waste at the Site.				Up to 5		- VOCs by EPA Method 8260 - RCRA 8 Metal plus Zn, Cu, and Ni by EPA Method 6010 - PAHs by EPA Method 8270	VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids Metals - 4 oz. glass with Teflon-lined cap PAHs - 4 oz. amber glass with Teflon-lined cap	RL-9 RL-5 RL-13	
	CA-3	REC-8	Six (6) soil gas probes are anticipated to evaluate this REC (see soil gas sampling below) and if the result indicate a source of VOCs under the building slab, then a boring will be installed through the slab at the perceived center of the vapor source. A groundwater well will be installed and sampled appropriately.				1		- VOCs by EPA Method 8260	VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids	RL-9	
	CA-4 CA-5	REC-9	Two shallow borings will be advanced through the former press pits to assess the condition of the soil directly beneath the concrete slab.				2		- TPH by EPA Method 8015 - VOCs by EPA Method 8260 - RCRA 8 Metal plus Zn, Cu, and Ni by EPA Method 6010 - SVOCs by EPA Method 8270 - PCBs by EPA Method 8082	TPH- - 4 oz. glass with Teflon-lined cap VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids Metals - 4 oz. glass with Teflon-lined cap SVOCs - 4 oz. amber glass with Teflon-lined cap PCBs - 4 oz glass with Teflon lined cap	RL-7 RL-9 RL-5 RL-13 RL-4	

**Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062**

Media to be Collected	Proposed Sample IDs	Associated RECs	Sample Design	Sample Depth (ft bgs)	Field SOPs to be Used	Field Analysis/Observations	No. of Samples for Analysis	No. of Field Dups	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Soil Gas	CA-SG-1 CA-SG-2 CA-SG-3 CA-SG-4 CA-SG-5	REC-3	Five(5) soil gas probes (CA-SG-1 to CA-SG-5) are anticipated to evaluate this REC and if the results indicate a source of VOCs under the building slab other than the former closed in place acetone tank near the chemical mixing building, then a boring will be installed through the slab at the perceived center of the vapor source. A groundwater well will be installed and sampled appropriately. One soil gas sample will be sent to the laboratory to confirm the field screening results and determine the volatile organic compounds that were measured in the field. The sample that will be sent to the laboratory will be determined in the field.	1 to 2 feet below the concrete slab	WMD-06-01 Credere-004 DR#027 DR#012 HWRB-15	PID	1	1	- VOCs by EPA Method 8260	VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids	VOA-TO15	Columbia Analytical
	CA-SG-6 CA-SG-7 CA-SG-8 CA-SG-9 CA-SG-10 CA-SG-11	REC-8	Six (6) soil gas probes (CA-SG-6 to CA-SG-11) are anticipated to evaluate this REC and if the result did a source of VOCs under the building slab other than the former closed in place acetone tank near the chemical mixing building, then a boring will be installed through the slab at the perceived center of the vapor source. A groundwater well will be installed and sampled appropriately. One soil gas sample will be sent to the laboratory to confirm the field screening results and determine the volatile organic compounds that were measured in the field. The sample that will be sent to the laboratory will be determined in the field.				1	1	- VOCs by EPA Method 8260	VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids	VOA-TO15	

*Test pit number have been carried on from the previous work (i.e. Credere's numbers start at 7)

**Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062**

Media to be Collected	Proposed Sample IDs	Associated RECs	Sample Design	Field SOPs to be Used	Field Analysis/ Observations	No. of Samples for Analysis	No. of Field Dups	No. of Trip Blanks	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Groundwater	CA-1	REC-3	As a part of assessing REC-3, if contamination extends into groundwater as noted by field screening a monitoring well will be installed and sampled for VOCs.	Credere-004 HWRB-1 HWRB-3 HWRB-9 HWRB-15 HWRB-17 DR#012 EPA SOP EPASOP#2048	Visual & Olfactory Field Parameters: Temperature, PH, Dissolved Oxygen, Turbidity, Conductivity, Oxidation-Reduction Potential	1	Groundwater samples will be duplicated at a rate of 5% per the generic QAPP for a total of one (1) based on the proposed total number of samples indicated in this table.	1	- VOCs by EPA Method 8260	VOCs - (2) 40 ml VOA with HCL	RL-9	Absolute Resource Associates (formerly Resource Laboratories)
	MW-2	REC-5	As a part of assessing REC-5 existing monitoring well MW-2 will be assessed for free floating product and will be sampled to determine if it still contains degraded gasoline.			1			- VOCs by EPA Method 8260 -Lead by EPA 200.7/6010	VOCs - (2) 40 ml VOA with HCL Lead - 1 liter platic bottle, nitric perserved	RL-9 RL-5	
	CA-2	REC-6	As a part of assessing REC-6, if contamination at the septic tank discharge location extends into groundwater as noted by field screening, a monitoring well will be installed and sampled for VOCs.			1			- VOCs by EPA Method 8260 - SVOCs by EPA Method 8270 -Dissolved RCRA 8 Metal plus Zn, Cu, and Ni by EPA Method 6010	VOCs - (2) 40 ml VOA with HCL PAHs - 1 Liter amber bottle, unpreserved Metals - 1 liter platic bottle, nitric perserved	RL-9 RL-13 RL-5	
	CA-3	REC-8	A monitoring well may be installed to assess the condition of the groundwater in the general area of this REC, or in the area of high soil gas conditions. The installed monitoring well will be sampled for VOCs.			1			- VOCs by EPA Method 8260	VOCs - (2) 40 ml VOA with HCL	RL-9	

**Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062**

Media to be Collected	Proposed Sample IDs	Associated RECs/NCs	Sample Design	Field SOPs to be Used	Field Analysis/ Observations	No. of Samples for Analysis	No. of Field Dups	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Building Materials	Debris-	REC-10	An asbestos survey will be conducted to evaluate the "ASTM Non-scope" consideration related to potential ACMs in the debris pile at the Site.	According to NH DES Env-A-1800	Visual Inspection and Bulk Sampling	According to NH DES Env-A-1800	According to NH DES Env-A-1800	Polarized Light Microscopy byEPA 600/R-93/116	Plastic bags, labeled (no preservation)	Included in SSQAPP Addendum No. 4	ProScience Laboratories (subcontractor to Absolute Resource Associates)
	Debris-		Painted debris will be screened with the XRF for lead-based paint.	DR#024 DR#025 Credere-004	XRF Screening	up to 10	NA	EPA Method 6200	NA	NA	NA
	Debris-		One (1) representative bulk sample will be collected from the building debris pile.	EPA SOP No. 2011 for Chip, Wipe, and Sweep Sampling; EIASOP_POROUSSAMPLING1; Credere-004	Visual Inspection and Bulk Sampling	1	1	PCBs via EPA Method 8082 using extraction Method 3540	PBCs - 4 oz. glass with Teflon-lined cap, no preservative	RL-4	Absolute Resource Associates (formerly Resource Laboratories)
	Debris-		An asbestos survey will be conducted to evaluate the "ASTM Non-scope" consideration related to potential ACMs in the debris pile at the Site.	According to NH DES Env-A-1800	Visual Inspection and Bulk Sampling	According to NH DES Env-A-1800	According to NH DES Env-A-1800	Polarized Light Microscopy byEPA 600/R-93/116	Plastic bags, labeled (no preservation)	Included in SSQAPP Addendum No. 4	ProScience Laboratories (subcontractor to Absolute Resource Associates)

**Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062**

Media to be Collected	Proposed Sample IDs	Associated RECs/NCs	Sample Design	Field SOPs to be Used	Field Analysis/ Observations	No. of Samples for Analysis	No. of Field Dups	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Building Materials	Debris-	REC-10	Painted debris will be screened with the XRF for lead-based paint.	DR#024 DR#025 Crede-re-004	XRF Screening	up to 10	NA	EPA Method 6200	NA	NA	NA
	Concrete -1 Concrete -2	OTHER	Two (2) concrete samples will be collected from the from the concrete around the former transformer location	EPA SOP No. 2011 for Chip, Wipe, and Sweep Sampling; EIASOP_POROUSSAMPLING1; Crede-re-004	Visual Inspection and Bulk Sampling	2	1	PCBs via EPA Method 8082 using extraction Method 3540	PBCs - 4 oz. glass with Teflon-lined cap, no preservative	RL-4	Absolute Resource Associates (formerly Resource Laboratories)

Table 2: Sample Reference
Former Polyclad Laminates Property
45 Tannery Street
Franklin, New Hampshire
NHDES #199902062

Media to be Collected	Proposed Sample IDs	Associated RECs	Sample Design	Sample Depth (ft bgs)	Field SOPs to be Used	Field Analysis/ Observations	No. of Samples for Analysis	No. of Field Dups	Analytical Method	Sample Container information & Preservative (per location)	Lab SOPs	Laboratory To be Used
Sludge	SLUDGE-1	REC-6	One sludge sample will be collected from the septic tank to assess the contents for hazardous substances.	Bottom of Tank	Credera-004 HWRB-11 HWRB-12 HWRB-15 HWRB-17 VOCs/SOIL-2000 DR#012 DR#022	Visual & Olfactory PID Headspace XRF Screening	1	No duplicate will be collected because only one sludge sample is proposed.	- VOCs by EPA Method 8260 - RCRA 8 Metal plus Zn, Cu, and Ni by EPA Method 6010 - PAHs by EPA Method 8270 - PCBs by EPA Method 8082	VOCs - (1) 40 ml VOA w/5 ml methanol, (1) 40 ml VOA for % solids Metals - 4 oz. glass with Teflon-lined cap PAHs - 4 oz. amber glass with Teflon-lined cap PCBs - 4 oz glass with Teflon lined cap	RL-9 RL-5 RL-7 RL-13 RL-4	Absolute Resource Associates, LLC

APPENDIX B
SITE PHOTOGRAPHS





Picture 1

View of western portion of the Site, facing northeast, where potential UST was identified during GPR survey, presumed to be closed-in-place UST-6.



Picture 2

View of location of former fuel dispenser on western portion of the Site, facing northeast.



Picture 3

View of fill building demolition debris in Test Pit CA-TP-8 on western portion of the Site.



Picture 4

View of sub-slab soil gas sampling at sampling location SG-5, located within northern portion of former Site building footprint.



Picture 5

View of location of asbestos-containing floor tiles on former building slab, indicated by white arrows, facing southeast.



Picture 6

View of tannery hides visible in Test Pit CA-TP-13, located on northeastern portion of the Site.



Picture 7
View of tannery
hides visible in Test
Pit CA-TP-13,
located on
northeastern
portion of the Site.



Picture 8
View of tannery
waste in soil from
test pit CA-TP-14,
located on
northeastern
portion of the Site.

APPENDIX C
TEST PIT LOGS



**CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272**

TEST PIT DATA:

PROJECT NAME:	<u>Former Polyclad Laminates Site</u>	DATE:	<u>6/28/2011</u>
PROJECT NUMBER:	<u>10001086</u>	ACTIVITY:	
SAMPLE LOCATION ID:	<u>CA-TP-7</u>	START:	<u>1300</u>
CREDERE REPRESENTATIVE:	<u>Jonathan O'Donnell</u>	END:	<u>1430</u>
CONTRACTOR/FOREMAN:	<u>J.A. Eaton Inspections/Jeff Eaton</u>		

NOTES:

Test pit on western side of closed-in-place UST (presumed to be UST-6). Ductile iron pipe parallel and adjacent to UST prevented advancement of test pit beyond 8 feet below ground surface (bgs). No laboratory sample collected from this test pit.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES	
0	0-2	S-1	Slightly moist	0.0	Red-brown, medium to coarse SAND and Gravel. No volatile organic odor. Side of UST observed. Pipe adjacent to UST Side.	
1						
2	No sample from 2 to 3 feet bgs.					
3	3-5	S-2	Slightly moist	0.0		
4						
5	5-7	S-3	Slightly moist	0.0		
6						
7	7-8	S-4	Slightly moist	0.0		
8	End of Exploration at 8' bgs.					
PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis						

CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/30/2011

PROJECT NUMBER: 10001086

SAMPLE LOCATION ID: CA-TP-7A ACTIVITY: START: 1535
END: 1720








CREDERE REPRESENTATIVE: Judd Newcomb

CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton; T & K Drilling/Sean McGarry

NOTES:

Test pit on southern side of closed-in-place UST (presumed to be UST-6) because test pit TP-7 was stopped due to a pipe near the UST western side. Test pit TP-7A advanced to 10 feet bgs using the excavator. Samples from 10 to 16 feet bgs collected by split-spoon sampler via a drill rig because the limit of the excavator's reach was approximately 10 feet.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES						
0	4	S-1	Dry	0.5	<p style="text-align: center;">Light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p> <p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>						
1											
2											
3											
4	6	S-2	Dry	1.4		<p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>					
5											
6	8	S-3	Dry	1.3			<p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>				
7											
8	9	S-4	Dry	1.4				<p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>			
9	9.5	S-5	Dry	1.3							
10	10-12*	S-6	Dry	1.9					<p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>		
11											
12	12-14	S-7	Dry	1.1						<p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>	
13											
14	14-16	S-8	Dry	0.4							<p style="text-align: center;">Top of UST at 4 feet bgs. Observed 3" fill pipe (broken off during excavation) and UST full of sand. No volatile organic odor. Side of UST # 6 visible (4 to 10' bgs).</p> <div style="text-align: center;">  </div> <p style="text-align: center;">Medium dense, light-brown, fine to medium SAND and Gravel. No volatile organic odor.</p>
15											
16					<p style="text-align: center;">End of Exploration at 16' bgs.</p>						
	<p>PID - Photo-Ionization Detector ppm_v - parts per million by volume * - Submitted for laboratory analysis</p>										

**CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG**
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/28/2011
 PROJECT NUMBER: 10001086
 SAMPLE LOCATION ID: CA-TP-8 ACTIVITY: START: 0900
 CREDERE REPRESENTATIVE: Jonathan O'Donnell END: 0955
 CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton

NOTES:

Test pits CA-TP-8, CA-TP-10, and CA-TP-11 were advanced in the area reported to contain UST # 8. This area appears to be an former loading dock that had an asphalt paved area that slopes down to the north, but is now filled with demolition debris.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-2	S-1	Dry	1.2	Building demolition debris. Crushed concrete masonry unit, metal scrap, and sand. No volatile organic odor.
1					
2	2-4	S-2	Dry	ND	
3					
4	Asphalt paving layer at 4.5 feet bgs.				
5	4-6	S-3	Dry	1.1	Red-brown, medium to coarse SAND and Gravel starting at 4.5' bgs.No volatile organic odor.
6					
7	6-8*	S-4	Dry	ND	
8					
9	End of Exploration at 9' bgs.				
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis				

**CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272**

TEST PIT DATA:

PROJECT NAME:	<u>Former Polyclad Laminates Site</u>	DATE:	<u>6/30/2011</u>
PROJECT NUMBER:	<u>10001086</u>	ACTIVITY:	
SAMPLE LOCATION ID:	<u>CA-TP-9</u>	START:	<u>1345</u>
CREDERE REPRESENTATIVE:	<u>Judd Newcomb</u>	END:	<u>1520</u>
CONTRACTOR/FOREMAN:	<u>J.A. Eaton Inspections/Jeff Eaton</u>		

NOTES:

Located beneath northern end of former building slab, in between patch in concrete and area of metal GPR anomaly identified by DigSmart. Test pit advanced to investigate possible former UST.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0					<p>Medium dense, light-brown, medium to coarse SAND and Gravel. No volatile organic odor.</p> <p>Thin layers of tannery hides present from 2 to 6 feet bgs.</p> <p style="text-align: center;">↓</p> <p>Asphalt pieces at approximately 8 feet bgs.</p>
1	2	S-1	Dry	1.3	
2					
3	4	S-2	Dry	1.8	
4					
5	6*	S-3	Dry	2.4	
6					
7	8	S-4	Dry	2.0	
8	End of Exploration at 8' bgs.				

PID - Photo-Ionization Detector
ppm_v - parts per million by volume
* - Submitted for laboratory analysis

CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/28/2011

PROJECT NUMBER: 10001086

SAMPLE LOCATION ID: CA-TP-10 ACTIVITY: START: 1000
END: 1000

CREDERE REPRESENTATIVE: Jonathan O'Donnell

CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton

NOTES:

Test pits CA-TP-8, CA-TP-10, and CA-TP-11 were advanced in the area reported to contain UST # 8. This area appears to be an former loading dock that had an asphalt paved area that slopes down to the north, but is now filled with demolition debris.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-2*	S-1	Dry	ND	Building demolition debris. Crushed CMU, metal scrap, and sand. No volatile organic odor.
1					
2	2-4	S-2	Dry	ND	
3					
4	4.5-6	S-3	Dry	ND	Asphalt paving layer at 4.5 feet bgs.
5					
6	6-8	S-4	Dry	ND	Red-brown, medium to coarse SAND and Gravel from 4.5 to 9 feet bgs. No volatile organic odor.
7					
8	8-9	S-5	Dry	ND	
9	End of Exploration at 9' bgs.				
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis CMU = concrete masonry unit				

CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/28/2011
PROJECT NUMBER: 10001086
SAMPLE LOCATION ID: CA-TP-11 ACTIVITY: START: 1100
CREDERE REPRESENTATIVE: Jonathan O'Donnell END: 1215
CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton

NOTES:

Test pits CA-TP-8, CA-TP-10, and CA-TP-11 were advanced in the area reported to contain UST # 8. This area appears to be an former loading dock that had an asphalt paved area that slopes down to the north, but is now filled with demolition debris.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-2	S-1	Dry	ND	Building demolition debris. Crushed CMU, metal scrap, and sand. No volatile organic odor.
1					
2	No sample from 2 to 3 feet bgs.				Asphalt paving layer at 3 feet bgs.
3	3-4.5	S-2	Dry	ND	Red-brown, medium to coarse SAND and Gravel. No volatile organic odor.
4	4.5-6.5	S-3	Dry	ND	
5					
6					
7	6.5-9*	S-4	Dry	ND	
8					
9	End of Exploration at 9' bgs.				
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis CMU = concrete masonry unit				

**CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272**

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/30/2011
 PROJECT NUMBER: 10001086
 SAMPLE LOCATION ID: CA-TP-12 ACTIVITY:
 START: 0935
 END: 1010
 CREDERE REPRESENTATIVE: Judd Newcomb
 CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton

NOTES:

Test pit advanced through bottom of concrete sump surrounded by stained concrete, believed to be the former location of a fuel dispenser. Bottom of the sump was broken with excavator, and the test pit advanced further by hand using a clean shovel.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	2	S-1	Wet	ND	Wet, black, organic detritus (material from interior of 2' deep concrete sump). No volatile organic odor. Concrete bottom of sump at 2 feet bgs.
1					
2	3*	S-2	Wet	1.2	Red-brown, medium to coarse SAND and Gravel. Wet. Staining beneath concrete at 2 feet bgs. No volatile organic odor.
3	4	S-3	Wet	0.5	
4	End of Exploration at 4' bgs.				
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis				

CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME:	<u>Former Polyclad Laminates Site</u>	DATE:	<u>6/28/2011</u>
PROJECT NUMBER:	<u>10001086</u>	ACTIVITY:	
SAMPLE LOCATION ID:	<u>CA-TP-14/CA-SS-2</u>	START:	<u>1510</u>
CREDERE REPRESENTATIVE:	<u>Jonathan O'Donnell</u>	END:	<u>1550</u>
CONTRACTOR/FOREMAN:	<u>J.A. Eaton Inspections/Jeff Eaton</u>		

NOTES:

Test pit on northeastern portion of the Site, in area where the GPR survey indicated that buried wastes may be present.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0					Red-brown, medium to coarse SAND and Gravel. No volatile organic odor.
1	0-2**	S-1	Dry	ND	
2					
3	2-4	S-2	Dry	ND	
4					Red-brown, medium to coarse SAND and Gravel mixed with thin layers of tannery hides present from 4 to 8 feet bgs. No volatile organic odor.
5	4-6	S-3	Dry	ND	
6					
7	6-8*	S-4	Dry	ND	
8	End of Exploration at 8' bgs.				
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis ** - Collected as surficial soil sample CA-SS-2				

**CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272**

TEST PIT DATA:

PROJECT NAME:	<u>Former Polyclad Laminates Site</u>	DATE:	<u>6/28/2011</u>
PROJECT NUMBER:	<u>10001086</u>	ACTIVITY:	
SAMPLE LOCATION ID:	<u>CA-TP-15</u>	START:	<u>1555</u>
CREDERE REPRESENTATIVE:	<u>Jonathan O'Donnell</u>	END:	<u>1625</u>
CONTRACTOR/FOREMAN:	<u>J.A. Eaton Inspections/Jeff Eaton</u>		

NOTES:

Test pit on southeastern portion of the Site to assess possible buried tannery waste.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-2	S-1	Dry	ND	Red-brown, medium to coarse SAND and Gravel. No volatile organic odor.
1					
2					
3	2-4	S-2	Dry	ND	
4					
4	No sample from 4 to 5 feet bgs.				
5	5-7*	S-3	Dry	ND	Red-brown, medium to coarse SAND and Gravel. Appears native. No volatile organic odor.
6					
7	End of Exploration at 7' bgs.				

PID - Photo-Ionization Detector
ppm, - parts per million by volume
* - Submitted for laboratory analysis

CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/28/2011

PROJECT NUMBER: 10001086

SAMPLE LOCATION ID: CA-TP-16 ACTIVITY: START: 1635
END: 1705

CREDERE REPRESENTATIVE: Jonathan O'Donnell

CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton

NOTES:

Test pit on southwestern portion of the Site to assess possible buried tannery waste.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-2*	S-1	Dry	ND	Red-brown, medium to coarse SAND and Gravel. No volatile organic odor.
1					
2	2-4	S-2	Dry	ND	
3					
4	No sample from 4 to 5 feet bgs.				
5	5-7	S-3	Dry	ND	Red-brown, medium to coarse SAND and Gravel. Appears native. No volatile organic odor.
6					
7	End of Exploration at 7' bgs.				
8					
9					
10					
12					
13					
14					
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis				

**CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272**

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 6/28/2011
PROJECT NUMBER: 10001086
SAMPLE LOCATION ID: CA-TP-18 ACTIVITY:
START: 1440
END: 1450
CREDERE REPRESENTATIVE: Jonathan O'Donnell
CONTRACTOR/FOREMAN: J.A. Eaton Inspections/Jeff Eaton

NOTES:

Test pit north of building foundation.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-1.5	S-1	Dry	ND	Asphalt upper three inches (kept separate from soil samples).
1					Red-brown, medium to coarse SAND and Gravel. No volatile organic odor.
2	1.5-3	S-2	Dry	ND	
3	3-5*	S-3	Dry	ND	
4					
5	End of Exploration at 5' bgs.				

PID - Photo-Ionization Detector
ppm_v - parts per million by volume
* - Submitted for laboratory analysis

CREDERE ASSOCIATES, LLC
TEST PIT SAMPLING LOG
776 Main Street, Westbrook, Maine 04092 - 207-828-1272

TEST PIT DATA:

PROJECT NAME: Former Polyclad Laminates Site DATE: 7/5/2011

PROJECT NUMBER: 10001086

SAMPLE LOCATION ID: CA-SS-3 ACTIVITY:

CREDERE REPRESENTATIVE: Jonathan O'Donnell START: 1155

CONTRACTOR/FOREMAN: None END: 1205

NOTES:



Surficial soil sample south of building foundation.

FIELD ANALYSIS DATA:

DEPTH (FT)	SAMPLE DEPTH (FT)	SAMPLE NUMBER	MOISTURE	PID (ppm)	SOIL DESCRIPTION / NOTES
0	0-2	S-1*	Dry	ND	Grass and Organic in upper two inches. Organic material seperated from sample. Red-brown, medium to coarse SAND. No volatile organic odor.
1					
2	End of Exploration at 5' bgs.				
3					
4					
5					
	PID - Photo-Ionization Detector ppm _v - parts per million by volume * - Submitted for laboratory analysis				

APPENDIX D
SOIL BORING LOGS



Geologic Log													
SITE INFORMATION				WELL SPECIFICATIONS									
 Credere Associates, LLC 776 Main Street Westbrook, Maine 04092				Project Number/Client: 10001086/Former Polyclad Laminates Site				Well Depth (feet) from TOW: 35					
				Site Location: 45 Tannery Street, Franklin, NH				Screen Length (feet): 10					
				NHDES Site#: 199902062				Date Start/Finish: 06-29-11					
				TOW Elev. (feet): 335.26				Ground Elevation: NM					
				Credere, LLC Representative: Jonathan O'Donnell				Well Material: 2" PVC - ND10-inch PVC Slotted Pipe and Riser					
				CONTRACTOR				DRILLING EQUIPMENT					
Drilling Contractor: T&K Drilling				Equipment: 4 1/4" ID Hollow Stem Auger									
Foreman: Sean McGarry				Casing Diameter: NA									
Drilling Method: Auger				Casing Material: NA									
Sample Information						Soil Description and Classification	Strata	USCS Code	Equipment Installed		Depth		
Depth	Sample No.	Pen/Rec (In.)	Depth (FL)	Blows (/0.5')	PID (ppm _v) (RF=1.0)				Depth	Depth			
0						Dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.	Poorly Graded Sand and Gravel	SP	Flush mount well box	ND			
1	S-1	24/12	0-2	10-19-16-10	ND						2-inch expansion plug	1	
2						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						2	
3	S-2	24/19	2-4	10-7-7-8	ND							3	
4						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.					2-inch PVC riser	4	
5	S-3	24/17	4-6	4-4-6-6	ND						Backfill: Auger Cuttings	5	
6						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						6	
7	S-4	24/15	6-8	4-5-7-7	ND							7	
8						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						8	
9	S-5	24/18	8-10	7-7-7-9	ND							9	
10						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						10	
11	S-6	24/13	10-12	4-7-8-9	ND							11	
12						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						12	
13	S-7	24/18	12-14	8-8-7-9	ND							13	
14						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						14	
15	S-8	24/17	14-16	6-7-8-13	ND							15	
16						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.						16	
17	S-9	24/17	16-18	11-12-13-15	ND							17	
18	Boring Continued on Page 2 of 2										18		

Remarks:


Soils described using "Modified Burmister Method"

TOW Elevation - Elevation of top of well PVC riser, based on TOW elevation of MW-2 at 334.89 feet

Well Depth shown is actual installed depth. Depth measured during groundwater sampling may reflect a loss of depth due to sedimentation.

NM - Not Measured

*Sample submitted for laboratory analysis.



 Graphically shows depth of the water table inferred from field observations during soil boring activities.

Boring No:

CA-1

Page 1 of 2

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Geologic Log											
SITE INFORMATION						WELL SPECIFICATIONS					
 Credere Associates, LLC 776 Main Street Westbrook, Maine 04092						Project Number/Client: 10001086/Former Polyclad Laminates Site			Well Depth (feet) from TOW: 35		
						Site Location: 45 Tannery Street, Franklin, NH			Screen Length (feet): 10		
						NHDES Site#: 199902062			Date Start/Finish: 06-29-11		
						TOW Elev. (feet): 335.26			Ground Elevation: NM		
						Credere, LLC Representative: Jonathan O'Donnell			Well Material: 2" PVC - ND10-inch PVC Slotted Pipe and Riser		
						CONTRACTOR			DRILLING EQUIPMENT		
						Drilling Contractor: T&K Drilling			Equipment: 4 1/4" ID Hollow Stem Auger		
						Foreman: Sean McGarry			Casing Diameter: NA		
						Drilling Method: Auger			Casing Material: NA		
Depth	Sample Information					Soil Description and Classification	Strata	USCS Code	Equipment Installed		Depth
	Sample No.	Pen/Rec (In.)	Depth (Ft.)	Blows (/0.5')	PID (ppm _v) (RF=L0)						
18						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.			2-inch PVC riser	Backfill: Auger Cuttings	18
19	S-10	24/17	18-20	6-8-10-13	ND						19
20						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.			Bentonite		20
21	S-11	24/17	20-22	6-8-9-9	ND						21
22						Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.					22
23	S-12	24/16	22-24	9-9-10-10	ND						23
24						Dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.					24
25	S-13	24/9	24-26	6-14-23-27	ND						25
26						Dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.					26
27	S-14*	24/16	26-28	19-19-17-17	ND						27
28	End of sampling at 28'. Boring advanced to 35 feet for installation of well.										28
29						29					
30						Clean Sand Filter pack					30
31						2-inch PVC 0.010 slotted pipe					31
32											32
33											33
34											34
35	Well Set at 35'										35

Remarks:


Soils described using "Modified Burmister Method"

TOW Elevation - Elevation of top of well PVC riser, based on TOW elevation of MW-2 at 334.89 feet

Well Depth shown is actual installed depth. Depth measured during groundwater sampling may reflect a loss of depth due to sedimentation.

NM - Not Measured

*Sample submitted for laboratory analysis.



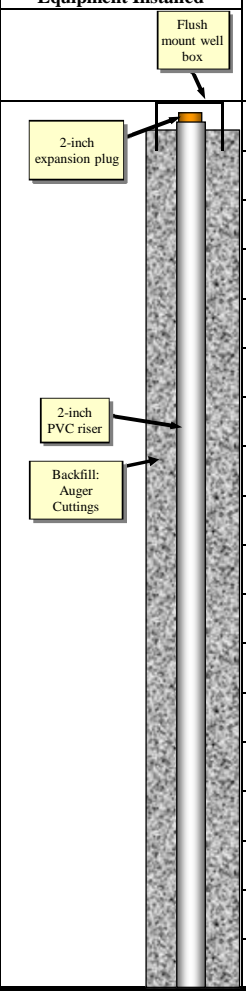
 Graphically shows depth of the water table inferred from field observations during soil boring activities.

Boring No:


CA-1

Page 2 of 2



Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Geologic Log																	
SITE INFORMATION						WELL SPECIFICATIONS											
 Credere Associates, LLC 776 Main Street Westbrook, Maine 04092						Project Number/Client: 10001086/Former Polyclad Laminates Site			Well Depth (feet) from TOW: 35 (field measurement)								
						Site Location: 45 Tannery Street, Franklin, NH			Screen Length (feet): 10								
						NHDES Site#: 199902062			Date Start/Finish: 06-29-11			TOW Elev. (feet): 334.82 Ground Elevation: NM					
						Credere, LLC Representative: Jonathan O'Donnell						Well Material: 2" PVC - 0.010-inch PVC Slotted Pipe and Riser					
						CONTRACTOR						DRILLING EQUIPMENT					
						Drilling Contractor: T&K Drilling			Equipment: 4 1/4" ID Hollow Stem Auger			Casing Diameter: NA			Casing Material: NA		
Foreman: Sean McGarry						Drilling Method: Auger											
Sample Information						Soil Description and Classification	Strata	USCS Code	Equipment Installed			Depth					
Depth	Sample No.	Pen/Rec (In.)	Depth (FL)	Blows (/0.5')	PID (ppm _v) (RF=1.0)				Depth	Depth	Depth						
0						Medium dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.	Poorly Graded Sand and Gravel	SP				0					
1	S-1	24/10	0-2	5-6-5-4	ND							1					
2						Loose, coarse SAND and Gravel. No petroleum or volatile odor.							2				
3	S-2	24/14	2-4	4-2-2-2	ND								3				
4						Loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							4				
5	S-3	24/8	4-6	3-2-2-3	ND								5				
6						Loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							6				
7	S-4*	24/11	6-8	2-2-3-2	10.2								7				
8						Very loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							8				
9	S-5	24/13	8-10	1-1-1-1	7.9								9				
10						Very loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							10				
11	S-6	24/12	10-12	1-1-1-1	1.2								11				
12						Very loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							12				
13	S-7	24/15	12-14	1-1-1-2	ND								13				
14						Loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							14				
15	S-8	24/9	14-16	1-1-4-5	ND								15				
16						Loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.							16				
17	S-9	24/13	16-18	7-3-4-3	ND								17				
18	Boring Continued on Page 2 of 2										18						

Remarks:

Soils described using "Modified Burmister Method"
 TOW Elevation - Elevation of top of well PVC riser, based on TOW elevation of MW-2 at 334.89 feet
 Well Depth shown is actual installed depth. Depth measured during groundwater sampling may reflect a loss of depth due to sedimentation.
 REF - Refusal, 100 blows and less than 1 " of depth gained.
 NM - Not Measured
 *Sample submitted for laboratory analysis.
 Graphically shows depth of the water table inferred from field observations during soil boring activities.

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Geologic Log											
SITE INFORMATION			WELL SPECIFICATIONS								
 Credere Associates, LLC 776 Main Street Westbrook, Maine 04092			Project Number/Client: 10001086/Former Polyclad Laminates Site			Well Depth (feet) from TOW: 35 (field measurement)					
			Site Location: 45 Tannery Street, Franklin, NH			Screen Length (feet): 10					
			NHDES Site#: 199902062			Date Start/Finish: 06-29-11		TOW Elev. (feet): 334.82		Ground Elevation: NM	
			Credere, LLC Representative: Jonathan O'Donnell			Well Material: 2" PVC - 0.010-inch PVC Slotted Pipe and Riser					
			CONTRACTOR			DRILLING EQUIPMENT					
			Drilling Contractor: T&K Drilling			Equipment: 4 1/4" ID Hollow Stem Auger					
			Foreman: Sean McGarry			Casing Diameter: NA					
			Drilling Method: Auger			Casing Material: NA					
Depth	Sample Information					Soil Description and Classification	Strata	USCS Code	Equipment Installed		
	Sample No.	Pen/Rec (In.)	Depth (FL)	Blows (/0.5')	PID (ppm _v) (RF=1.0)				Depth	Depth	
18						Very loose, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.			2-inch PVC riser	18	
19	S-10	24/9	18-20	3-1-2-1	ND					19	
20						Medium dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.			Backfill: Auger Cuttings	20	
21	S-11	24/14	20-22	5-8-24-24	ND					21	
22						Dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.				22	
23	S-12	24/11	22-24	25-26-23-REF	ND					23	
24						Very dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.			Bentonite	24	
25	S-13	24/16	24-26	16-25-33-24	ND					25	
26						Dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.				26	
27	S-14	24/12	26-28	23-19-11-6	ND					27	
28						Medium dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.				28	
29	S-15	24/18	28-30	5-11-7-7	ND					29	
30						Medium dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.			Clean Sand Filter pack	30	
31	S-16	24/7	30-32	4-7-5-4	ND					31	
32	End of sampling at 32'. Boring advanced to 35 feet for installation of well.									32	
33	NOTE: Strong aromatic odor in water drained from drilling auger during withdrawal of auger.									33	
34										34	
35	Well Set at 35'									35	

Remarks:

Soils described using "Modified Burmister Method"


TOW Elevation - Elevation of top of well PVC riser, based on TOW elevation of MW-2 at 334.89 feet

Well Depth shown is actual installed depth. Depth measured during groundwater sampling may reflect a loss of depth due to sedimentation.

REF - Refusal, 100 blows and less than 1" of depth gained.

NM - Not Measured

*Sample submitted for laboratory analysis.


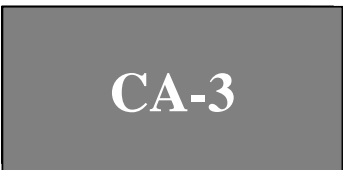
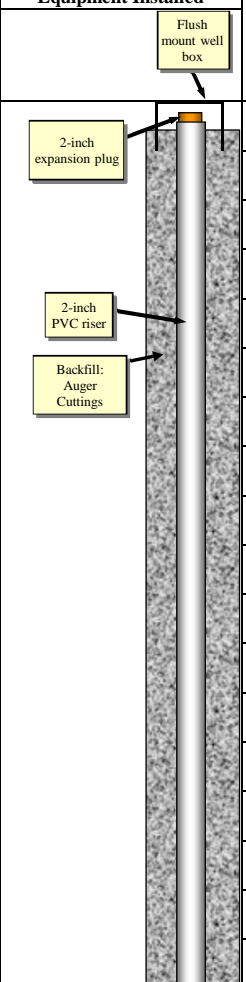
 Graphically shows depth of the water table inferred from field observations during soil boring activities.

Boring No:

CA-2

Page 2 of 2

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Geologic Log											
SITE INFORMATION				WELL SPECIFICATIONS							
 Credere Associates, LLC 776 Main Street Westbrook, Maine 04092				Project Number/Client: 10001086/Former Polyclad Laminates Site				Well Depth (feet) from TOW: 35 (field measurement)			
				Site Location: 45 Tannery Street, Franklin, NH				Screen Length (feet): 10			
				NHDES Site#: 199902062				Date Start/Finish: 06-29-11			
				Credere, LLC Representative: Jonathan O'Donnell				Well Material: 2" PVC - 0.010-inch PVC Slotted Pipe and Riser			
				CONTRACTOR				DRILLING EQUIPMENT			
				Drilling Contractor: T&K Drilling				Equipment: 4 1/4" ID Hollow Stem Auger			
Foreman: Sean McGarry				Casing Diameter: NA							
Drilling Method: Auger				Casing Material: NA							
Sample Information						Soil Description and Classification	Strata	USCS Code	Equipment Installed		Depth
Depth	Sample No.	Pen/Rec (In.)	Depth (Ft.)	Blows (/0.5')	PID (ppm _v) (RF=1.0)				Depth	Depth	
0						Dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.	Poorly Graded Sand and Gravel	SP		0	
1	S-1	24/11	0-2	4-14-17-18	ND					1	
2						Dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.				2	
3	S-2*	24/11	2-4	23-12-26-21	ND					3	
4						Dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.				4	
5	S-3	24/12	4-6	12-21-24-20	ND					5	
6						Very dense, brown, medium to coarse SAND and Gravel. No petroleum or volatile odor.				6	
7	S-4	24/9	6-8	39-24-35-14	1.2					7	
8						Loose, coarse SAND and Gravel. No petroleum or volatile odor.				8	
9	S-5	24/13	8-10	3-4-5-7	ND		9				
10						Medium dense, light brown fine SAND. No petroleum or volatile odor.	Well Graded Sand	SW	10		
11	S-6	24/13	10-12	8-6-11-9	ND				11		
12						Medium dense, light brown fine SAND. No petroleum or volatile odor.			12		
13	S-7	24/14	12-14	7-10-12-12	ND				13		
14						Medium dense, light brown fine SAND. No petroleum or volatile odor.			14		
15	S-8	24/13	14-16	10-11-13-15	ND				15		
16						Medium dense, light brown fine SAND. No petroleum or volatile odor.			16		
17	S-9	24/16	16-18	10-12-15-14	ND		17				
18	Boring Continued on Page 2 of 2										18

Remarks:


Soils described using "Modified Burmister Method"

TOW Elevation - Elevation of top of well PVC riser, based on TOW elevation of MW-2 at 334.89 feet

Well Depth shown is actual installed depth. Depth measured during groundwater sampling may reflect a loss of depth due to sedimentation.

NM - Not Measured

*Sample submitted for laboratory analysis.


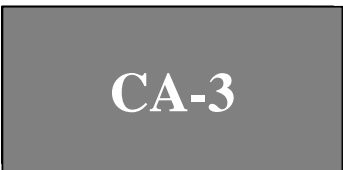
 Graphically shows depth of the water table inferred from field observations during soil boring activities.

Boring No:

CA-3

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Page 1 of 2

Geologic Log																	
SITE INFORMATION						WELL SPECIFICATIONS											
 Crede Associates, LLC 776 Main Street Westbrook, Maine 04092						Project Number/Client: 10001086/Former Polyclad Laminates Site			Well Depth (feet) from TOW: 35 (field measurement)								
						Site Location: 45 Tannery Street, Franklin, NH			Screen Length (feet): 10								
 CA-3						NHDES Site#: 199902062			Date Start/Finish: 06-29-11			TOW Elev. (feet): 335.25			Ground Elevation: NM		
						Crede, LLC Representative: Jonathan O'Donnell						Well Material: 2" PVC - 0.010-inch PVC Slotted Pipe and Riser					
						CONTRACTOR						DRILLING EQUIPMENT					
						Drilling Contractor: T&K Drilling						Equipment: 4 1/4" ID Hollow Stem Auger					
Foreman: Sean McGarry						Casing Diameter: NA											
Drilling Method: Auger						Casing Material: NA											
Sample Information						Soil Description and Classification	Strata	USCS Code	Equipment Installed		Depth						
Depth	Sample No.	Pen/Rec (In.)	Depth (FL)	Blows (/0.5')	PID (ppm _v) (RF=1.0)				Depth								
18						Dense, light brown fine SAND. No petroleum or volatile odor.			2-inch PVC riser	Backfill: Auger Cuttings	18						
19	S-10	24/15	18-20	14-15-17-17	ND						19						
20						Medium dense, light brown fine SAND. No petroleum or volatile odor.			Bentonite		20						
21	S-11	24/16	20-22	9-10-15-16	ND						21						
22						Dense, light brown fine SAND. No petroleum or volatile odor.					22						
23	S-12	24/15	22-24	15-15-17-19	ND						23						
24						Medium dense, light brown fine SAND. No petroleum or volatile odor.					24						
25	S-13	24/17	24-26	5-10-12-15	ND						25						
26						Medium dense, light brown fine SAND. No petroleum or volatile odor.					26						
27	S-14	24/15	26-28	14-15-14-15	ND						27						
28						Medium dense, light brown fine SAND. No petroleum or volatile odor.					28						
29	S-15	24/16	28-30	3-6-8-10	ND						29						
30						Loose, light brown fine SAND. No petroleum or volatile odor.					30						
31	S-16	24/15	30-32	3-3-4-5	ND						31						
32	End of sampling at 32'. Boring advanced to 35 feet for installation of well.										32						
33											33						
34											34						
35	Well Set at 35'										35						

Remarks:


Soils described using "Modified Burmister Method"

TOW Elevation - Elevation of top of well PVC riser, based on TOW elevation of MW-2 at 334.89 feet

Well Depth shown is actual installed depth. Depth measured during groundwater sampling may reflect a loss of depth due to sedimentation.

NM - Not Measured

*Sample submitted for laboratory analysis.

 Graphically shows depth of the water table inferred from field observations during soil boring activities.

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Geologic Log											
SITE INFORMATION						WELL SPECIFICATIONS					
<p>Crede Associates, LLC 776 Main Street Westbrook, Maine 04092</p>						Project Number/Client: 10001086/Former Polyclad Laminates Site		Well Depth (feet) from TOW: N/A			
						Site Location: 45 Tannery Street, Franklin, NH		Screen Length (feet): N/A			
NHDES Site#: 199902062				Date Start/Finish: 06-29-11		TOW Elev. (feet): N/A		Ground Elevation: NM			
						Crede, LLC Representative: Jonathan O'Donnell		Well Material: N/A			
						CONTRACTOR			DRILLING EQUIPMENT		
						Drilling Contractor: T&K Drilling			Equipment: 4 1/4" ID Hollow Stem Auger		
						Foreman: Sean McGarry			Casing Diameter: NA		
Drilling Method: Auger			Casing Material: NA								
Depth	Sample Information					Soil Description and Classification	Strata	USCS Code	Equipment Installed		
	Sample No.	Pen/Rec (In.)	Depth (Ft.)	Blows (/0.5')	PID (ppm _v) (RF=1.0)					Depth	
0	No Sampling from 0 to 10 feet below ground surface. Purpose of boring was to assess soil beneath the press pits, beginning at 10 feet below ground surface.							No Monitoring Well Installed		0	
1										0	
2										0	
3										0	
4										0	
5										0	
6										0	
7										0	
8										0	
9										0	
10	S-1*	24/12	10-12	1-2-2-3	ND	Loose, brown, medium to coarse SAND and Gravel. Slight aromatic odor.	Poorly Graded Sand	SP		10	
11										11	
12	S-2	24/16	12-14	3-4-2-4	ND	Loose, brown, medium to coarse SAND and Gravel. Slight aromatic odor.	Poorly Graded Sand	SP		12	
13										13	
14	End of sampling at 14'									14	

Remarks:



Soils described using "Modified Burmister Method"

NM - Not Measured

*Sample submitted for laboratory analysis.

Boring No: CA-4

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

Geologic Log													
SITE INFORMATION						WELL SPECIFICATIONS							
 <p>Crede Associates, LLC 776 Main Street Westbrook, Maine 04092</p>						Project Number/Client: 10001086/Former Polyclad Laminates Site			Well Depth (feet) from TOW: N/A				
						Site Location: 45 Tannery Street, Franklin, NH			Screen Length (feet): N/A				
						NHDES Site#: 199902062		Date Start/Finish: 06-29-11		TOW Elev. (feet): N/A		Ground Elevation: NM	
						Crede, LLC Representative: Jonathan O'Donnell						Well Material: N/A	
CONTRACTOR						DRILLING EQUIPMENT							
Drilling Contractor: T&K Drilling						Equipment: 4 1/4" ID Hollow Stem Auger							
Foreman: Sean McGarry						Casing Diameter: NA							
Drilling Method: Auger						Casing Material: NA							
Sample Information						Soil Description and Classification	Strata	USCS Code	Equipment Installed		Depth		
Depth	Sample No.	Pen/Rec (In.)	Depth (Ft.)	Blows (/0.5')	PID (ppm _v) (RF=1.0)								
0						No Sampling from 0 to 10 feet below ground surface. Purpose of boring was to assess soil beneath the press pits, beginning at 10 feet below ground surface.			No Monitoring Well Installed	0			
1										1			
2										2			
3										3			
4										4			
5										5			
6										6			
7										7			
8										8			
9										9			
10	S-1	24/12	10-12	1-5-6-4	ND	Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.	Poorly Graded Sand	SP		10			
11										11			
12	S-2*	24/13	12-14	4-6-6-7	ND	Medium dense, brown, medium to coarse SAND and Gravel. Slight aromatic odor.				12			
13										13			
14	End of sampling at 14'									14			

Remarks:

Soils described using "Modified Burmister Method"

NM - Not Measured

*Sample submitted for laboratory analysis.

Boring No: CA-5

Stratification lines represent approximate boundaries between soil types, transitions may be gradual. Water level readings have been made at times and under conditions stated. Fluctuations of groundwater may occur due to other factors than those present at the time measurements were made.

APPENDIX E
GROUNDWATER SAMPLING LOGS



**LOW FLOW SAMPLING LOG
CREDERE ASSOCIATES**



Crederre Associates LLC

PROJECT NAME: Polyclad

DATE: 7 15 2011

PROJECT NUMBER: 10001086

LOCATION ACTIVITY

SAMPLE LOCATION ID: CA-1

START: 1440
END: 1520

WELL DATA:

WELL DEPTH (ft): 34.42 MEASURED HISTORICAL TOP OF WELL TOP OF CASING WATER LEVEL EQUIPMENT USED:
 FROM GRADE ELECT. COND. PROBE

WATER DEPTH (ft): 28.02 MEASURED HISTORICAL FROM GRADE FLOAT ACTIVATED PROBE
 PRESSURE TRANSDUCER

WELL MATERIAL: WELL PROTECTIVE CASING CONCRETE COLLAR AMBIENT AIR VOC: _____ PPM
 PVC LOCKED: SECURE: INTACT: YES NO
 SS YES YES YES WEL MOUTH VOC: _____ PPM
 _____ NO NO NO

EQUIPMENT DATA:

PURGING SAMPLING METER ID DECONTAMINATION FLUIDS USED:

<input type="checkbox"/>	<input type="checkbox"/>	PERISTALTIC PUMP	<input type="checkbox"/>	pH	_____	<input type="checkbox"/>	DISTILLED WATER
<input checked="" type="checkbox"/>	<input type="checkbox"/>	SUBMERSIBLE	<input type="checkbox"/>	Specific Conductivity	_____	<input type="checkbox"/>	DEIONIZED WATER
<input type="checkbox"/>	<input type="checkbox"/>	BLADDER PUMP	<input type="checkbox"/>	Dissolved Oxygen	_____	<input type="checkbox"/>	POTABLE WATER
<input type="checkbox"/>	<input type="checkbox"/>	HAND PUMP	<input type="checkbox"/>	ORP	_____	<input type="checkbox"/>	TSP SOLUTION
<input type="checkbox"/>	<input type="checkbox"/>	DEDICATED HDPE	<input type="checkbox"/>	Turbidity	_____	<input type="checkbox"/>	ALCONOX SOLUTION
<input type="checkbox"/>	<input type="checkbox"/>	NEW HDPE				<input type="checkbox"/>	NONE
<input type="checkbox"/>	<input type="checkbox"/>	DEDICATED LDPE				<input type="checkbox"/>	_____
<input type="checkbox"/>	<input type="checkbox"/>	NEW LDPE				<input type="checkbox"/>	_____
<input type="checkbox"/>	<input type="checkbox"/>	FILTER				<input type="checkbox"/>	_____
<input type="checkbox"/>	<input checked="" type="checkbox"/>	<u>SANITIZED</u>					

FIELD ANALYSIS DATA:

PUMP ON: _____ PUMP OFF: _____ STABLE FLOW RATE (ml/min): _____ MEASURED ESTIMATED

TIME	TEMP (°C)	pH	COND. (mS)	ORP (mV)	D.O. (mg/l)	TURBID. (ntu)	COMMENTS
1440	15.13	5.86	.190	116.7	6.44		v BROWN TURBID/NO ODOOR (3 GALLONS)
1444	13.71	5.96	.181	130.6	7.11		(5 GALLONS) TURBID
1450	13.37	6.26	.175	126.5	13.50		(10 GALLONS) TURBID
1456	14.07	6.23	.172	125.8	10.74	331	(14 GALLONS) SLIGHTLY LESS TURBID

SAMPLE DATA:

SAMPLE BOTTLE ID	PRESERVATION METHOD	SAMPLE CONTAINER #	LABORATORY ANALYSIS
TIME LOCATION	METHOD	TYPE	ANALYSIS
CA-2 1515			
(Vials only) ↓			

NOTES:

SAMPLER _____

**LOW FLOW SAMPLING LOG
CREDERE ASSOCIATES**



PROJECT NAME: PLYMOUTH
 PROJECT NUMBER: 10001086
 SAMPLE LOCATION ID: CA-2

DATE: 7/5/11

LOCATION ACTIVITY
 START: 1350
 END: _____

WELL DATA:

WELL DEPTH (ft): 34.41 MEASURED HISTORICAL TOP OF WELL TOP OF CASING
 WATER DEPTH (ft): 127.57 MEASURED HISTORICAL FROM GRADE

WATER LEVEL EQUIPMENT USED:
 ELECT. COND. PROBE FLOAT ACTIVATED PROBE
 PRESSURE TRANSDUCER _____

WELL MATERIAL: PVC SS _____
 WELL LOCKED: YES NO
 PROTECTIVE CASING SECURE: YES NO
 CONCRETE COLLAR INTACT: YES NO
 AMBIENT AIR VOC: _____ PPM
 WELL MOUTH VOC: _____ PPM

EQUIPMENT DATA:

PURGING SAMPLING PERISTALTIC PUMP SUBMERSIBLE BLADDER PUMP HAND PUMP DEDICATED HDPE NEW HDPE DEDICATED LDPE NEW LDPE FILTER BAILER

ANALYSIS: pH Specific Conductivity Dissolved Oxygen ORP Turbidity

METER ID: 451
Peristaltic 2090

DECONTAMINATION FLUIDS USED:
 DISTILLED WATER
 DEIONIZED WATER
 POTABLE WATER
 TSP SOLUTION
 ALCONOX SOLUTION
 NONE

FIELD ANALYSIS DATA:

PUMP ON: _____ PUMP OFF: _____ STABLE FLOW RATE (ml/min): _____ [] MEASURED [] ESTIMATED

TIME	TEMP (°C)	pH	COND. (mS)	ORP (mV)	D.O. (mg/l)	TURBID. (ntu)	COMMENTS
1355	3.93	6.16	.203	6.40	10.85		v. hot out of gate/draw turbid (3 gal)
1400	14.67	5.97	.204	-13.6	1.39		(6 gallons) more same
1405	13.27	6.11	.199	-60.9	0.04	99.7	(8 gallons) more still
1410	19.12	6.43	.224	-65.8	0.52		(10 gallons)
1415	14.06	6.19	.195	-81.9	-3.28		(15 gallons) clearer
1420	14.10	6.54	0.196	< 71.2	0.71	29.7	(18 gal) clearer

SAMPLE DATA:

SAMPLE BOTTLE ID	PRESERVATION	SAMPLE CONTAINER	LABORATORY ANALYSIS	
TIME	METHOD	#	TYPE	
1420 CA-2	HNO3	1	HDPE	Metal
	HCl	2	10A	VOCs
	none	1	amber	Silica

NOTES:

BLACK, OPAQUE TURBID, SMELLS LIKE PETROLEUM

_____ SAMPLER

LOW FLOW SAMPLING LOG
 CREDERE ASSOCIATES



PROJECT NAME: POLYGLAD

DATE: 7 15 2011

PROJECT NUMBER: 10001086

LOCATION ACTIVITY

SAMPLE LOCATION ID: CA-3

START: 15 15
 END: 15 45

WELL DATA:

WELL DEPTH (ft): 34.57 MEASURED TOP OF WELL WATER LEVEL EQUIPMENT USED:
 HISTORICAL TOP OF CASING ELECT. COND. PROBE
 FROM GRADE FLOAT ACTIVATED PROBE
 WATER DEPTH (ft): 27.75 MEASURED _____ PRESSURE TRANSDUCER
 HISTORICAL _____

WELL MATERIAL: WELL LOCKED: PROTECTIVE CASING SECURE: CONCRETE COLLAR INTACT: AMBIENT AIR VOC: _____ PPM
 PVC YES YES YES YES
 SS NO NO NO NO
 _____ NO NO NO WELL MOUTH VOC: _____ PPM

EQUIPMENT DATA:

PURGING SAMPLING			METER ID	DECONTAMINATION
<input checked="" type="checkbox"/>	<input type="checkbox"/>	PERISTALTIC PUMP	<input type="checkbox"/> pH	FLUIDS USED:
<input checked="" type="checkbox"/>	<input type="checkbox"/>	SUBMERSIBLE	<input type="checkbox"/> Specific Conductivity	<input type="checkbox"/> DISTILLED WATER
<input type="checkbox"/>	<input type="checkbox"/>	BLADDER PUMP	<input type="checkbox"/> Dissolved Oxygen	<input type="checkbox"/> DEIONIZED WATER
<input type="checkbox"/>	<input type="checkbox"/>	HAND PUMP	<input type="checkbox"/> ORP	<input type="checkbox"/> POTABLE WATER
<input type="checkbox"/>	<input type="checkbox"/>	DEDICATED HDPE	<input type="checkbox"/> Turbidity	<input type="checkbox"/> TSP SOLUTION
<input type="checkbox"/>	<input type="checkbox"/>	NEW HDPE	_____	<input type="checkbox"/> ALCONOX SOLUTION
<input type="checkbox"/>	<input type="checkbox"/>	DEDICATED LDPE	_____	<input type="checkbox"/> NONE
<input type="checkbox"/>	<input type="checkbox"/>	NEW LDPE	_____	_____
<input type="checkbox"/>	<input checked="" type="checkbox"/>	FILTER	_____	_____
<input type="checkbox"/>	<input checked="" type="checkbox"/>	<u>MIRAC</u>	_____	_____

FIELD ANALYSIS DATA:

PUMP ON: _____ PUMP OFF: _____ STABLE FLOW RATE (ml/min): _____ MEASURED ESTIMATED

TIME	TEMP (°C)	pH	COND. (mS)	ORP (mV)	D.O. (mg/l)	TURBID. (ntu)	COMMENTS
1518	14.15	6.91	.302	145.3	10.45		Ignored Turbid WATER No ODOUR
1524	17.03	6.07	.319	157.6	15.24	36.4	(5 Gallons)
1536	14.74	5.97	.304	166.5	15.24		(10 Gallons) No ODOUR - 50697 Turbid

SAMPLE DATA:

SAMPLE BOTTLE ID	PRESERVATION METHOD	SAMPLE CONTAINER #	LABORATORY ANALYSIS
TIME LOCATION		TYPE	
1530 CA-3	_____	_____	_____
_____	_____	_____	_____
_____	_____	_____	_____
_____	_____	_____	_____

NOTES:

SAMPLER _____

**LOW FLOW SAMPLING LOG
CREDERE ASSOCIATES**



PROJECT NAME: POLYCLAD
 PROJECT NUMBER: 10001086
 SAMPLE LOCATION ID: MW-2

DATE: 7.5.11

LOCATION ACTIVITY
 START: 1250
 END: 1330

WELL DATA:

WELL DEPTH (ft): 33.96 [] MEASURED [] HISTORICAL [] TOP OF WELL [] TOP OF CASING [] FROM GRADE [] _____
 WATER DEPTH (ft): 27.64 [] MEASURED [] HISTORICAL [] _____

WATER LEVEL EQUIPMENT USED:
 ELECT. COND. PROBE
 FLOAT ACTIVATED PROBE
 PRESSURE TRANSDUCER

WELL MATERIAL: [] PVC [] SS [] _____
 WELL LOCKED: [] YES [] NO
 PROTECTIVE CASING SECURE: [] YES [] NO
 CONCRETE COLLAR INTACT: [] YES [] NO
 AMBIENT AIR VOC: _____ PPM
 WELL MOUTH VOC: _____ PPM

EQUIPMENT DATA:

PURGING SAMPLING
 PERISTALTIC PUMP [] pH
 SUBMERSIBLE [] Specific Conductivity
 BLADDER PUMP [] Dissolved Oxygen
 HAND PUMP [] ORP
 DEDICATED HDPE [] Turbidity
 NEW HDPE
 DEDICATED LDPE
 NEW LDPE
 FILTER
 BALER

METER ID
451
LAUREL 2020

DECONTAMINATION FLUIDS USED:
 DISTILLED WATER
 DEIONIZED WATER
 POTABLE WATER
 TSP SOLUTION
 ALCONOX SOLUTION
 NONE

FIELD ANALYSIS DATA:

PUMP ON: _____ PUMP OFF: _____ STABLE FLOW RATE (ml/min): _____ [] MEASURED [] ESTIMATED

TIME	TEMP (°C)	pH	COND. (mS)	ORP (mV)	D.O. (mg/l)	TURBID. (ntu)	COMMENTS
1301	16.04	6.28	0.258	-0.4	2.14	32.2	2 gal slight odor
1306	12.95	6.06	0.234	7.5	1.84	-	5 gal
1309	12.82	6.29	0.255	12.7	4.07		6 gal slight odor
1321	12.04	6.18	0.228	18.4	9.3		8 gal

SAMPLE DATA:

SAMPLE BOTTLE ID	PRESERVATION	SAMPLE CONTAINER	LABORATORY ANALYSIS
TIME LOCATION	METHOD	# TYPE	
1320 MW-2	NO3	1 LDPE	NO3
	NO2	2 NO2	PH

NOTES: Purge sampling, well volume ≈ 0.8 ft³

SR

SAMPLER

**PURGE + BAILER
LOW-FLOW SAMPLING LOG
CREDERE ASSOCIATES**



PROJECT NAME: POLYCLAD
 PROJECT NUMBER: 10001086
 SAMPLE LOCATION ID: CA-2

DATE: 7, 14, 11
 LOCATION ACTIVITY
 START: 1435
 END: 1545

WELL DATA:

WELL DEPTH (ft): 34.40 MEASURED TOP OF WELL WATER LEVEL EQUIPMENT USED:
 HISTORICAL TOP OF CASING ELECT. COND. PROBE
 HISTORICAL FROM GRADE FLOAT ACTIVATED PROBE
 WATER DEPTH (ft): 28.10 MEASURED _____ PRESSURE TRANSDUCER
 HISTORICAL _____ _____
 WELL MATERIAL: WELL PROTECTIVE CASING CONCRETE COLLAR
 PVC LOCKED: SECURE: INTACT: AMBIENT AIR VOC: N/A PPM
 SS YES YES YES WELL MOUTH VOC: N/A PPM
 _____ NO NO NO

EQUIPMENT DATA:

PURGING SAMPLING PERISTALTIC PUMP pH METER ID DECONTAMINATION
 SUBMERSIBLE Specific Conductivity YSI FLUIDS USED:
 BLADDER PUMP Dissolved Oxygen _____ DISTILLED WATER
 HAND PUMP ORP _____ DEIONIZED WATER
 DEDICATED HDPE Turbidity Lamella 2020 POTABLE WATER
 NEW HDPE _____ _____ TSP SOLUTION
 DEDICATED LDPE _____ _____ ALCONOX SOLUTION
 NEW LDPE _____ _____ NONE
 FILTER _____ _____ _____
 BAILER

FIELD ANALYSIS DATA:

PUMP ON: _____ PUMP OFF: _____ STABLE FLOW RATE (ml/min): _____ MEASURED ESTIMATED

TIME	TEMP (°C)	pH	COND. (mS)	DO ORP (mV)	D.O. (mg/l)	TURBID. (ntu)	COMMENTS
1515	16.21	5.99	0.184	3.25	135.1	485	
1525	14.19	6.13	0.191	10.11	140.7		
1534	13.42	6.21	0.193	8.35	142.7	95	
1539	13.49	6.25	0.198	10.31	142.7	70	

SAMPLE DATA:

SAMPLE BOTTLE ID	PRESERVATION METHOD	SAMPLE CONTAINER #	LABORATORY ANALYSIS
TIME LOCATION		TYPE	
<u>1539 CA-2</u>	<u>ANO₃</u>	<u>1 LDPE</u>	<u>RCRA 8 Metals</u>

NOTES: PURGED 15 gallons

SBO
SAMPLER

APPENDIX F

LABORATORY ANALYTICAL REPORTS





Absolute Resource *associates*

124 Heritage Avenue #10 Portsmouth, NH 03801

Jonathan O'Donnell
CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: None
Job ID: 21886
Date Received: 6/29/11

Project: POLYCLAD 10001086

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

Sue Sylvester
Principal, General Manager

Date of Approval: 7/11/2011
Total number of pages: 58

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SS-1	Solid	6/28/2011 15:12	21886-001	PAHs in solid by 8270 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-2	Solid	6/28/2011 16:28	21886-002	PAHs in solid by 8270 Percent Dry Matter for Sample Calc by SM2540B,G
CA-TP-8	Solid	6/28/2011 10:48	21886-003	TPH in solids by 8100 PAHs in solid by 8270 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-10	Solid	6/28/2011 11:50	21886-004	Solid Digestion for ICP Analysis Lead in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-11	Solid	6/28/2011 12:31	21886-005	Solid Digestion for ICP Analysis Lead in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-13	Solid	6/28/2011 15:00	21886-006	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-14	Solid	6/28/2011 15:50	21886-007	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-TP-14	Solid	6/28/2011 15:50	21886-007	Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-15	Solid	6/28/2011 16:25	21886-008	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-16	Solid	6/28/2011 17:00	21886-009	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-17	Solid	6/28/2011 14:40	21886-010	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-18	Solid	6/28/2011 14:48	21886-011	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-TP-18	Solid	6/28/2011 14:48	21886-011	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-DUP-1	Solid	6/28/2011	21886-012	PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
D-Floor Covering	Solid	6/28/2011 15:35	21886-013	PCBs in soil by 8082
Transformer Pit	Solid	6/28/2011 12:55	21886-014	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-3	Solid	6/29/2011 10:47	21886-015	Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
Trip Blank	Solid	6/28/2011	21886-016	VOCs in solid by 8260 Petro & Haz Waste

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-003

Sample ID: CA-TP-8

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/28/11 10:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-003

Sample ID: CA-TP-8

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/28/11 10:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	90	78-114	%	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
toluene-D8 SUR	92	88-110	%	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
4-bromofluorobenzene SUR	97	86-115	%	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B
a,a,a-trifluorotoluene SUR	112	70-130	%	1	LMM	6/30/11	4295	6/30/11	17:35	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-004

Sample ID: CA-TP-10

Matrix: Solid

Percent Dry: 88.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 11:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
trichlorofluoromethane	0.3	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
acetone	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
t-butanol (TBA)	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
2-butanone (MEK)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
tetrahydrofuran (THF)	< 0.7	0.7	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,4-dioxane	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
2-hexanone	< 0.7	0.7	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-004

Sample ID: CA-TP-10

Matrix: Solid

Percent Dry: 88.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 11:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
m&p-xylenes	0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	88	78-114	%	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
toluene-D8 SUR	94	88-110	%	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
4-bromofluorobenzene SUR	96	86-115	%	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B
a,a,a-trifluorotoluene SUR	106	70-130	%	1	LMM	6/30/11	4295	6/30/11	18:06	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-005

Sample ID: CA-TP-11

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 12:31

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-005

Sample ID: CA-TP-11

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 12:31

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	87	78-114	%	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
toluene-D8 SUR	94	88-110	%	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
4-bromofluorobenzene SUR	98	86-115	%	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B
a,a,a-trifluorotoluene SUR	98	70-130	%	1	LMM	6/30/11	4295	6/30/11	18:37	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-006

Sample ID: CA-TP-13

Matrix: Solid

Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
tetrahydrofuran (THF)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
2-hexanone	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-006

Sample ID: CA-TP-13

Matrix: Solid

Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	84	78-114	%	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
toluene-D8 SUR	91	88-110	%	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
4-bromofluorobenzene SUR	95	86-115	%	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B
a,a,a-trifluorotoluene SUR	101	70-130	%	1	LMM	6/30/11	4295	6/30/11	19:08	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-007

Sample ID: CA-TP-14

Matrix: Solid

Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
tetrahydrofuran (THF)	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
2-hexanone	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-007

Sample ID: CA-TP-14

Matrix: Solid

Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	85	78-114	%	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
toluene-D8 SUR	91	88-110	%	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
4-bromofluorobenzene SUR	93	86-115	%	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B
a,a,a-trifluorotoluene SUR	104	70-130	%	1	LMM	6/30/11	4295	6/30/11	19:39	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-008

Sample ID: CA-TP-15

Matrix: Solid

Percent Dry: 75.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 16:25

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
acetone	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
t-butanol (TBA)	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
2-butanone (MEK)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
tetrahydrofuran (THF)	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,4-dioxane	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
2-hexanone	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-008

Sample ID: CA-TP-15

Matrix: Solid

Percent Dry: 75.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 16:25

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	81	78-114	%	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
toluene-D8 SUR	90	88-110	%	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
4-bromofluorobenzene SUR	93	86-115	%	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B
a,a,a-trifluorotoluene SUR	124	70-130	%	1	LMM	6/30/11	4295	6/30/11	20:10	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-009

Sample ID: CA-TP-16

Matrix: Solid

Percent Dry: 98.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 17:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-009

Sample ID: CA-TP-16

Matrix: Solid

Percent Dry: 98.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 17:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	82	78-114	%	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
toluene-D8 SUR	89	88-110	%	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
4-bromofluorobenzene SUR	97	86-115	%	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B
a,a,a-trifluorotoluene SUR	96	70-130	%	1	LMM	6/30/11	4295	6/30/11	20:42	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-010

Sample ID: CA-TP-17

Matrix: Solid

Percent Dry: 81.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
acetone	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
t-butanol (TBA)	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
tetrahydrofuran (THF)	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,4-dioxane	< 3	3	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
2-hexanone	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-010

Sample ID: CA-TP-17

Matrix: Solid

Percent Dry: 81.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	82	78-114	%	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
toluene-D8 SUR	90	88-110	%	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
4-bromofluorobenzene SUR	95	86-115	%	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B
a,a,a-trifluorotoluene SUR	118	70-130	%	1	LMM	6/30/11	4295	6/30/11	21:13	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-011

Sample ID: CA-TP-18

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
acetone	< 3	3	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
t-butanol (TBA)	< 3	3	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
tetrahydrofuran (THF)	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,4-dioxane	< 3	3	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
2-hexanone	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-011

Sample ID: CA-TP-18

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	88	78-114	%	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
toluene-D8 SUR	94	88-110	%	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
4-bromofluorobenzene SUR	95	86-115	%	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B
a,a,a-trifluorotoluene SUR	120	70-130	%	1	LMM	6/30/11	4295	7/5/11	15:54	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-012

Sample ID: CA-TP-DUP-1

Matrix: Solid

Percent Dry: 93.7% Results expressed on a dry weight basis.

Sampled: 6/28/11

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-012

Sample ID: CA-TP-DUP-1

Matrix: Solid

Percent Dry: 93.7% Results expressed on a dry weight basis.

Sampled: 6/28/11

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	96	78-114	%	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
toluene-D8 SUR	98	88-110	%	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
4-bromofluorobenzene SUR	97	86-115	%	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B
a,a,a-trifluorotoluene SUR	101	70-130	%	1	LMM	6/30/11	4295	7/1/11	16:06	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-015

Sample ID: CA-3

Matrix: Solid

Percent Dry: 95.5% Results expressed on a dry weight basis.

Sampled: 6/29/11 10:47

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
acetone	< 3	3	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
t-butanol (TBA)	< 3	3	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
2-butanone (MEK)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
tetrahydrofuran (THF)	< 0.7	0.7	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,4-dioxane	< 3	3	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.6	0.6	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
2-hexanone	< 0.7	0.7	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-015

Sample ID: CA-3

Matrix: Solid

Percent Dry: 95.5% Results expressed on a dry weight basis.

Sampled: 6/29/11 10:47

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	96	78-114	%	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
toluene-D8 SUR	101	88-110	%	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
4-bromofluorobenzene SUR	96	86-115	%	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B
a,a,a-trifluorotoluene SUR	101	70-130	%	1	LMM	6/30/11	4295	7/1/11	16:39	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-016

Sample ID: Trip Blank

Matrix: Solid

Sampled: 6/28/11

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
tetrahydrofuran (THF)	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.4	0.4	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
2-hexanone	< 0.5	0.5	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-016

Sample ID: Trip Blank

Matrix: Solid

Sampled: 6/28/11

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	92	78-114	%	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
toluene-D8 SUR	90	88-110	%	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
4-bromofluorobenzene SUR	93	86-115	%	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B
a,a,a-trifluorotoluene SUR	98	70-130	%	1	LMM	6/30/11	4295	6/30/11	17:05	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-001

Sample ID: CA-SS-1

Matrix: Solid

Percent Dry: 82.6% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:12

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
phenanthrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
benzo(a)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
chrysene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
benzo(b)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
benzo(k)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
benzo(a)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
benzo(g,h,i)perylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	85	43-116	%	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D
o-terphenyl SUR	86	33-141	%	1	AJD	7/1/11	4298	7/5/11	16:50	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-002

Sample ID: CA-SS-2

Matrix: Solid

Percent Dry: 82.9% Results expressed on a dry weight basis.

Sampled: 6/28/11 16:28

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
phenanthrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
benzo(a)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
chrysene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
benzo(b)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
benzo(k)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
benzo(a)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
benzo(g,h,i)perylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	82	43-116	%	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D
o-terphenyl SUR	84	33-141	%	1	AJD	7/1/11	4298	7/5/11	21:16	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-003

Sample ID: CA-TP-8

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/28/11 10:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	84	43-116	%	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D
o-terphenyl SUR	88	33-141	%	1	AJD	7/1/11	4298	7/5/11	21:54	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-006

Sample ID: CA-TP-13

Matrix: Solid

Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	82	43-116	%	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D
o-terphenyl SUR	89	33-141	%	1	AJD	7/1/11	4298	7/5/11	19:22	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-007

Sample ID: CA-TP-14

Matrix: Solid

Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	78	43-116	%	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D
o-terphenyl SUR	84	33-141	%	1	AJD	7/1/11	4298	7/5/11	20:00	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-008

Sample ID: CA-TP-15

Matrix: Solid

Percent Dry: 75.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 16:25

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 3.3	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
2-methylnaphthalene	< 3.3	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
acenaphthylene	< 3.3	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
acenaphthene	5.3	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
dibenzofuran	3.6	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
fluorene	7.7	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
phenanthrene	100	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
anthracene	33	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
fluoranthene	160	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
pyrene	160	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
benzo(a)anthracene	80	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
chrysene	74	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
benzo(b)fluoranthene	46	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
benzo(k)fluoranthene	80	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
benzo(a)pyrene	66	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
indeno(1,2,3-cd)pyrene	18	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
dibenzo(a,h)anthracene	9.6	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
benzo(g,h,i)perylene	16	3.3	ug/g	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	81	43-116	%	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D
o-terphenyl SUR	82	33-141	%	5	AJD	7/1/11	4298	7/7/11	11:29	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-009

Sample ID: CA-TP-16

Matrix: Solid

Percent Dry: 98.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 17:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	81	43-116	%	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D
o-terphenyl SUR	86	33-141	%	1	AJD	7/1/11	4298	7/5/11	17:28	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-010

Sample ID: CA-TP-17

Matrix: Solid

Percent Dry: 81.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
phenanthrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
benzo(a)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
chrysene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
benzo(b)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
benzo(k)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
benzo(a)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
benzo(g,h,i)perylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	83	43-116	%	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D
o-terphenyl SUR	90	33-141	%	1	AJD	7/1/11	4298	7/5/11	18:06	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-011

Sample ID: CA-TP-18

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
phenanthrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
benzo(a)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
chrysene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
benzo(b)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
benzo(k)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
benzo(a)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
benzo(g,h,i)perylene	< 0.6	0.6	ug/g	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	83	43-116	%	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D
o-terphenyl SUR	88	33-141	%	1	AJD	7/1/11	4298	7/5/11	18:44	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-012

Sample ID: CA-TP-DUP-1

Matrix: Solid

Percent Dry: 93.7% Results expressed on a dry weight basis.

Sampled: 6/28/11

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	86	43-116	%	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D
o-terphenyl SUR	91	33-141	%	1	AJD	7/1/11	4298	7/5/11	20:38	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-006

Sample ID: CA-TP-13

Matrix: Solid Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	63	30-150	%	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A
decachlorobiphenyl SUR	72	30-150	%	1	JLZ	7/1/11	4297	7/3/11	15:01	SW3540C8082A

Sample#: 21886-013

Sample ID: D-Floor Covering

Matrix: Solid

Sampled: 6/28/11 15:35

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	24 *	30-150	%	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A
decachlorobiphenyl SUR	25 *	30-150	%	1	JLZ	7/1/11	4297	7/8/11	11:07	SW3540C8082A

* The surrogate showed recovery outside the acceptance limits. Matrix interference suspected.

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-014

Sample ID: Transformer Pit

Matrix: Solid Percent Dry: 96.6% Results expressed on a dry weight basis.

Sampled: 6/28/11 12:55

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	43	30-150	%	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A
decachlorobiphenyl SUR	63	30-150	%	1	JLZ	7/1/11	4297	7/8/11	11:38	SW3540C8082A

Sample#: 21886-003

Sample ID: CA-TP-8

Matrix: Solid Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/28/11 10:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 220	220	ug/g	1	JLZ	7/1/11	4299	7/6/11	17:36	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	95	40-140	%	1	JLZ	7/1/11	4299	7/6/11	17:36	SW3550B8100m
o-terphenyl SUR	112	40-140	%	1	JLZ	7/1/11	4299	7/6/11	17:36	SW3550B8100m

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-004

Sample ID: CA-TP-10

Matrix: Solid Percent Dry: 88.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 11:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis			Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Lead	5.6	0.6	ug/g	1	BJS	7/6/11	4310	7/6/11	22:48	SW3051A6010C	

Sample#: 21886-005

Sample ID: CA-TP-11

Matrix: Solid Percent Dry: 97.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 12:31

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis			Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Lead	4.8	0.6	ug/g	1	BJS	7/6/11	4310	7/6/11	22:55	SW3051A6010C	

Sample#: 21886-006

Sample ID: CA-TP-13

Matrix: Solid Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis			Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	3.0	0.6	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Barium	23	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Chromium	36	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Copper	18	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Lead	7.2	0.6	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Mercury	< 0.08	0.08	ug/g	1	BJS	7/6/11	4308	7/7/11	9:59	SW7471B	
Nickel	7	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Silver	< 0.4	0.4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	
Zinc	21	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:03	SW3051A6010C	

Sample#: 21886-007

Sample ID: CA-TP-14

Matrix: Solid Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/28/11 15:50

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis			Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	1.4	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Barium	47	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Chromium	260	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Copper	14	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Lead	6.2	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Mercury	< 0.08	0.08	ug/g	1	BJS	7/6/11	4308	7/7/11	10:00	SW7471B	
Nickel	4	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Silver	< 0.5	0.5	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	
Zinc	17	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:10	SW3051A6010C	

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-008

Sample ID: CA-TP-15

Matrix: Solid Percent Dry: 75.2% Results expressed on a dry weight basis.

Sampled: 6/28/11 16:25

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	6.1	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Barium	22	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Chromium	9	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Copper	20	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Lead	28	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Mercury	< 0.10	0.10	ug/g	1	BJS	7/6/11	4308	7/7/11	10:02	SW7471B
Nickel	6	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Selenium	< 4	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Silver	< 0.5	0.5	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C
Zinc	23	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:18	SW3051A6010C

Sample#: 21886-009

Sample ID: CA-TP-16

Matrix: Solid Percent Dry: 98.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 17:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	1.2	0.6	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Barium	12	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Chromium	5	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Copper	10	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Lead	3.2	0.6	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Mercury	< 0.07	0.07	ug/g	1	BJS	7/6/11	4308	7/7/11	10:04	SW7471B
Nickel	4	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C
Zinc	9	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:25	SW3051A6010C

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-010

Sample ID: CA-TP-17

Matrix: Solid Percent Dry: 81.7% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	2.7	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Barium	11	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Chromium	9	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Copper	22	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Lead	4.8	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Mercury	< 0.07	0.07	ug/g	1	BJS	7/6/11	4308	7/7/11	10:06	SW7471B
Nickel	4	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Silver	< 0.5	0.5	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C
Zinc	13	3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:33	SW3051A6010C

Sample#: 21886-011

Sample ID: CA-TP-18

Matrix: Solid Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 6/28/11 14:48

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	2.7	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Barium	16	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Chromium	7	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Copper	13	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Lead	4.9	0.7	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Mercury	< 0.08	0.08	ug/g	1	BJS	7/6/11	4308	7/7/11	10:08	SW7471B
Nickel	6	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Selenium	< 4	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Silver	< 0.5	0.5	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C
Zinc	15	4	ug/g	1	BJS	7/6/11	4310	7/6/11	23:40	SW3051A6010C

Project ID: POLYCLAD 10001086

Job ID: 21886

Sample#: 21886-012

Sample ID: CA-TP-DUP-1

Matrix: Solid

Percent Dry: 93.7% Results expressed on a dry weight basis.

Sampled: 6/28/11

Parameter	Quant		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Arsenic	< 0.6	0.6	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Barium	92	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Chromium	300	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Copper	16	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Lead	7.0	0.6	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Mercury	< 0.09	0.09	ug/g	1	BJS	7/6/11	4308	7/7/11	10:13	SW7471B
Nickel	4	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C
Zinc	20	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:18	SW3051A6010C

Quality Control Report



124 Heritage Avenue Unit 10
Portsmouth, NH 03801

www.absoluteresourceassociates.com



Case Narrative

Lab # 21886

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 5 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

No exceptions noted.

Surrogate Recoveries

PCB: The surrogates for 21886-013 were below the acceptance criteria. Matrix interference suspected.

Laboratory Control Sample Results

VOC: The MLCS/D4295 did not meet the acceptance criteria for chloromethane. This compound showed high recovery. There is no impact to the data as this analyte was not detected in the associated samples. The MLCS/D4295 did not meet the acceptance criteria for dichlorodifluoromethane. The MLCSD4295 did not meet the acceptance criteria for bromomethane. The recovery was acceptable in the LCS. These compounds are known to be problematic in the method.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MB4295	dichlorodifluoromethane		<	0.1	ug/g				
		chloromethane		<	0.1	ug/g				
		vinyl chloride		<	0.1	ug/g				
		bromomethane		<	0.2	ug/g				
		chloroethane		<	0.1	ug/g				
		trichlorofluoromethane		<	0.1	ug/g				
		diethyl ether		<	0.5	ug/g				
		acetone		<	2.5	ug/g				
		1,1-dichloroethene		<	0.1	ug/g				
		methylene chloride		<	0.2	ug/g				
		carbon disulfide		<	0.1	ug/g				
		methyl t-butyl ether (MTBE)		<	0.1	ug/g				
		trans-1,2-dichloroethene		<	0.1	ug/g				
		isopropyl ether (DIPE)		<	0.1	ug/g				
		ethyl t-butyl ether (ETBE)		<	0.1	ug/g				
		1,1-dichloroethane		<	0.1	ug/g				
		t-butanol (TBA)		<	2.5	ug/g				
		2-butanone (MEK)		<	0.5	ug/g				
		2,2-dichloropropane		<	0.1	ug/g				
		cis-1,2-dichloroethene		<	0.1	ug/g				
		chloroform		<	0.1	ug/g				
		bromochloromethane		<	0.1	ug/g				
		tetrahydrofuran (THF)		<	0.5	ug/g				
		1,1,1-trichloroethane		<	0.1	ug/g				
		1,1-dichloropropene		<	0.1	ug/g				
		t-amyl-methyl ether (TAME)		<	0.1	ug/g				
		carbon tetrachloride		<	0.1	ug/g				
		1,2-dichloroethane		<	0.1	ug/g				
		benzene		<	0.1	ug/g				
		trichloroethene		<	0.1	ug/g				
		1,2-dichloropropane		<	0.1	ug/g				
		bromodichloromethane		<	0.1	ug/g				
		1,4-dioxane		<	2.5	ug/g				
		dibromomethane		<	0.1	ug/g				
		4-methyl-2-pentanone (MIBK)		<	0.5	ug/g				
		cis-1,3-dichloropropene		<	0.1	ug/g				
		toluene		<	0.1	ug/g				
		trans-1,3-dichloropropene		<	0.1	ug/g				
		2-hexanone		<	0.5	ug/g				
		1,1,2-trichloroethane		<	0.1	ug/g				
		1,3-dichloropropane		<	0.1	ug/g				
		tetrachloroethene		<	0.1	ug/g				
		dibromochloromethane		<	0.1	ug/g				
		1,2-dibromoethane (EDB)		<	0.1	ug/g				
		chlorobenzene		<	0.1	ug/g				
		1,1,1,2-tetrachloroethane		<	0.1	ug/g				
		ethylbenzene		<	0.1	ug/g				
		m&p-xylenes		<	0.1	ug/g				
		o-xylene		<	0.1	ug/g				
		styrene		<	0.1	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MB4295	bromoform		<	0.1	ug/g				
		isopropylbenzene		<	0.1	ug/g				
		1,1,2,2-tetrachloroethane		<	0.1	ug/g				
		1,2,3-trichloropropane		<	0.1	ug/g				
		n-propylbenzene		<	0.1	ug/g				
		bromobenzene		<	0.1	ug/g				
		1,3,5-trimethylbenzene		<	0.1	ug/g				
		2-chlorotoluene		<	0.1	ug/g				
		4-chlorotoluene		<	0.1	ug/g				
		tert-butylbenzene		<	0.1	ug/g				
		1,2,4-trimethylbenzene		<	0.1	ug/g				
		sec-butylbenzene		<	0.1	ug/g				
		1,3-dichlorobenzene		<	0.1	ug/g				
		4-isopropyltoluene		<	0.1	ug/g				
		1,4-dichlorobenzene		<	0.1	ug/g				
		1,2-dichlorobenzene		<	0.1	ug/g				
		n-butylbenzene		<	0.1	ug/g				
		1,2-dibromo-3-chloropropane		<	0.1	ug/g				
		1,2,4-trichlorobenzene		<	0.1	ug/g				
		1,3,5-trichlorobenzene		<	0.1	ug/g				
		hexachlorobutadiene		<	0.1	ug/g				
		naphthalene		<	0.2	ug/g				
		1,2,3-trichlorobenzene		<	0.1	ug/g				
		dibromofluoromethane SUR			84	%		78 114		
		toluene-D8 SUR			93	%		88 110		
		4-bromofluorobenzene SUR			92	%		86 115		
		a,a,a-trifluorotoluene SUR			93	%		70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCS4295	dichlorodifluoromethane		0.6	ug/g	1	59 *	70 130		
		chloromethane		2.2	ug/g	1	215 *	70 130		
		vinyl chloride		0.8	ug/g	1	83	70 130		
		bromomethane		0.7	ug/g	1	70	70 130		
		chloroethane		0.7	ug/g	1	70	70 130		
		trichlorofluoromethane		0.7	ug/g	1	73	70 130		
		diethyl ether		0.9	ug/g	1	94	70 130		
		acetone	<	2.5	ug/g	1	112			
		1,1-dichloroethene		0.8	ug/g	1	77	70 130		
		methylene chloride		0.9	ug/g	1	93	70 130		
		carbon disulfide		0.8	ug/g	1	78	70 130		
		methyl t-butyl ether (MTBE)		1.0	ug/g	1	103	70 130		
		trans-1,2-dichloroethene		0.9	ug/g	1	89	70 130		
		isopropyl ether (DIPE)		1.0	ug/g	1	95	70 130		
		ethyl t-butyl ether (ETBE)		1.0	ug/g	1	100	70 130		
		1,1-dichloroethane		1.0	ug/g	1	97	70 130		
		t-butanol (TBA)		5.3	ug/g	5	106	70 130		
		2-butanone (MEK)		1.0	ug/g	1	102	70 130		
		2,2-dichloropropane		1.1	ug/g	1	110	70 130		
		cis-1,2-dichloroethene		0.9	ug/g	1	94	70 130		
		chloroform		1.0	ug/g	1	98	70 130		
		bromochloromethane		1.0	ug/g	1	96	70 130		
		tetrahydrofuran (THF)		0.9	ug/g	1	91	70 130		
		1,1,1-trichloroethane		0.9	ug/g	1	92	70 130		
		1,1-dichloropropene		1.0	ug/g	1	99	70 130		
		t-amyl-methyl ether (TAME)		1.0	ug/g	1	105	70 130		
		carbon tetrachloride		1.0	ug/g	1	96	70 130		
		1,2-dichloroethane		1.1	ug/g	1	108	70 130		
		benzene		1.0	ug/g	1	98	70 130		
		trichloroethene		0.9	ug/g	1	92	70 130		
		1,2-dichloropropane		1.0	ug/g	1	104	70 130		
		bromodichloromethane		1.0	ug/g	1	103	70 130		
		1,4-dioxane	<	2.5	ug/g	2	103	70 130		
		dibromomethane		1.0	ug/g	1	102	70 130		
		4-methyl-2-pentanone (MIBK)		0.9	ug/g	1	87	70 130		
		cis-1,3-dichloropropene		1.0	ug/g	1	99	70 130		
		toluene		0.9	ug/g	1	93	70 130		
		trans-1,3-dichloropropene		1.0	ug/g	1	100	70 130		
		2-hexanone		0.8	ug/g	1	78	70 130		
		1,1,2-trichloroethane		1.0	ug/g	1	101	70 130		
		1,3-dichloropropane		1.2	ug/g	1	119	70 130		
		tetrachloroethene		1.0	ug/g	1	97	70 130		
		dibromochloromethane		1.1	ug/g	1	115	70 130		
		1,2-dibromoethane (EDB)		1.1	ug/g	1	112	70 130		
		chlorobenzene		1.1	ug/g	1	107	70 130		
		1,1,1,2-tetrachloroethane		1.1	ug/g	1	114	70 130		
		ethylbenzene		1.0	ug/g	1	99	70 130		
		m&p-xylenes		2.1	ug/g	2	107	70 130		
		o-xylene		1.1	ug/g	1	111	70 130		
		styrene		1.1	ug/g	1	113	70 130		
		bromoform		1.0	ug/g	1	98	70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCS4295	isopropylbenzene		0.9	ug/g	1	95	70 130		
		1,1,2,2-tetrachloroethane		1.1	ug/g	1	113	70 130		
		1,2,3-trichloropropane		1.1	ug/g	1	109	70 130		
		n-propylbenzene		1.0	ug/g	1	102	70 130		
		bromobenzene		1.1	ug/g	1	109	70 130		
		1,3,5-trimethylbenzene		1.0	ug/g	1	99	70 130		
		2-chlorotoluene		1.2	ug/g	1	121	70 130		
		4-chlorotoluene		1.0	ug/g	1	99	70 130		
		tert-butylbenzene		1.0	ug/g	1	96	70 130		
		1,2,4-trimethylbenzene		1.0	ug/g	1	102	70 130		
		sec-butylbenzene		0.9	ug/g	1	93	70 130		
		1,3-dichlorobenzene		1.0	ug/g	1	102	70 130		
		4-isopropyltoluene		0.9	ug/g	1	95	70 130		
		1,4-dichlorobenzene		1.0	ug/g	1	105	70 130		
		1,2-dichlorobenzene		1.1	ug/g	1	107	70 130		
		n-butylbenzene		0.9	ug/g	1	94	70 130		
		1,2-dibromo-3-chloropropane		1.2	ug/g	1	116	70 130		
		1,2,4-trichlorobenzene		0.9	ug/g	1	93	70 130		
		1,3,5-trichlorobenzene		0.9	ug/g	1	88	70 130		
		hexachlorobutadiene		0.9	ug/g	1	92	70 130		
		naphthalene		1.0	ug/g	1	100	70 130		
		1,2,3-trichlorobenzene		0.9	ug/g	1	92	70 130		
		dibromofluoromethane SUR		95	%			78 114		
		toluene-D8 SUR		94	%			88 110		
		4-bromofluorobenzene SUR		100	%			86 115		
		a,a,a-trifluorotoluene SUR		91	%			70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCSD4295	dichlorodifluoromethane		0.6	ug/g	1	60 *	70 130	2	30
		chloromethane		1.6	ug/g	1	160 *	70 130	29	30
		vinyl chloride		0.9	ug/g	1	86	70 130	4	30
		bromomethane	<	0.2	ug/g	1	21 *	70 130	107 *	30
		chloroethane		0.7	ug/g	1	71	70 130	1	30
		trichlorofluoromethane		0.8	ug/g	1	76	70 130	4	30
		diethyl ether		0.9	ug/g	1	92	70 130	2	30
		acetone	<	2.5	ug/g	1	104		8	30
		1,1-dichloroethene		0.8	ug/g	1	77	70 130	0	30
		methylene chloride		0.9	ug/g	1	93	70 130	0	30
		carbon disulfide		0.8	ug/g	1	80	70 130	2	30
		methyl t-butyl ether (MTBE)		1.0	ug/g	1	100	70 130	3	30
		trans-1,2-dichloroethene		0.9	ug/g	1	89	70 130	0	30
		isopropyl ether (DIPE)		0.9	ug/g	1	94	70 130	1	30
		ethyl t-butyl ether (ETBE)		1.0	ug/g	1	100	70 130	0	30
		1,1-dichloroethane		1.0	ug/g	1	95	70 130	2	30
		t-butanol (TBA)		4.7	ug/g	5	95	70 130	11	30
		2-butanone (MEK)		0.9	ug/g	1	90	70 130	12	30
		2,2-dichloropropane		1.0	ug/g	1	102	70 130	8	30
		cis-1,2-dichloroethene		0.9	ug/g	1	94	70 130	0	30
		chloroform		1.0	ug/g	1	95	70 130	3	30
		bromochloromethane		1.0	ug/g	1	95	70 130	1	30
		tetrahydrofuran (THF)		0.8	ug/g	1	75	70 130	19	30
		1,1,1-trichloroethane		0.9	ug/g	1	91	70 130	1	30
		1,1-dichloropropene		1.0	ug/g	1	96	70 130	3	30
		t-amyl-methyl ether (TAME)		1.0	ug/g	1	103	70 130	2	30
		carbon tetrachloride		0.9	ug/g	1	93	70 130	3	30
		1,2-dichloroethane		1.0	ug/g	1	104	70 130	4	30
		benzene		1.0	ug/g	1	96	70 130	2	30
		trichloroethene		0.9	ug/g	1	90	70 130	2	30
		1,2-dichloropropane		1.0	ug/g	1	103	70 130	1	30
		bromodichloromethane		1.0	ug/g	1	100	70 130	4	30
		1,4-dioxane	<	2.5	ug/g	2	88	70 130	16	30
		dibromomethane		1.0	ug/g	1	101	70 130	1	30
		4-methyl-2-pentanone (MIBK)		0.8	ug/g	1	85	70 130	2	30
		cis-1,3-dichloropropene		1.0	ug/g	1	96	70 130	3	30
		toluene		1.0	ug/g	1	97	70 130	4	30
		trans-1,3-dichloropropene		0.9	ug/g	1	95	70 130	5	30
		2-hexanone		0.7	ug/g	1	74	70 130	6	30
		1,1,2-trichloroethane		1.0	ug/g	1	99	70 130	2	30
		1,3-dichloropropane		1.1	ug/g	1	108	70 130	9	30
		tetrachloroethene		0.9	ug/g	1	94	70 130	3	30
		dibromochloromethane		1.0	ug/g	1	103	70 130	10	30
		1,2-dibromoethane (EDB)		1.1	ug/g	1	107	70 130	5	30
		chlorobenzene		1.0	ug/g	1	105	70 130	2	30
		1,1,1,2-tetrachloroethane		1.1	ug/g	1	106	70 130	8	30
		ethylbenzene		0.9	ug/g	1	91	70 130	9	30
		m&p-xylenes		2.0	ug/g	2	101	70 130	5	30
		o-xylene		1.0	ug/g	1	104	70 130	7	30
		styrene		1.1	ug/g	1	106	70 130	7	30
		bromoform		0.8	ug/g	1	85	70 130	15	30

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCSD4295	isopropylbenzene		0.9	ug/g	1	88	70 130	7	30
		1,1,2,2-tetrachloroethane		1.1	ug/g	1	108	70 130	5	30
		1,2,3-trichloropropane		1.1	ug/g	1	105	70 130	3	30
		n-propylbenzene		1.0	ug/g	1	101	70 130	1	30
		bromobenzene		1.1	ug/g	1	111	70 130	2	30
		1,3,5-trimethylbenzene		1.0	ug/g	1	99	70 130	1	30
		2-chlorotoluene		1.1	ug/g	1	109	70 130	11	30
		4-chlorotoluene		1.1	ug/g	1	110	70 130	11	30
		tert-butylbenzene		1.0	ug/g	1	96	70 130	0	30
		1,2,4-trimethylbenzene		1.0	ug/g	1	103	70 130	0	30
		sec-butylbenzene		0.9	ug/g	1	92	70 130	1	30
		1,3-dichlorobenzene		1.0	ug/g	1	103	70 130	0	30
		4-isopropyltoluene		0.9	ug/g	1	94	70 130	1	30
		1,4-dichlorobenzene		1.0	ug/g	1	103	70 130	1	30
		1,2-dichlorobenzene		1.1	ug/g	1	105	70 130	1	30
		n-butylbenzene		0.9	ug/g	1	93	70 130	1	30
		1,2-dibromo-3-chloropropane		1.0	ug/g	1	104	70 130	11	30
		1,2,4-trichlorobenzene		0.9	ug/g	1	90	70 130	3	30
		1,3,5-trichlorobenzene		0.9	ug/g	1	87	70 130	0	30
		hexachlorobutadiene		0.9	ug/g	1	93	70 130	1	30
		naphthalene		1.0	ug/g	1	96	70 130	3	30
		1,2,3-trichlorobenzene		0.9	ug/g	1	90	70 130	2	30
		dibromofluoromethane SUR		97	%			78 114		
		toluene-D8 SUR		99	%			88 110		
		4-bromofluorobenzene SUR		95	%			86 115		
		a,a,a-trifluorotoluene SUR		96	%			70 130		

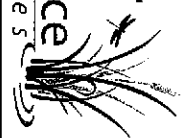
Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3540C8082A	BLK4297	PCB-1016		<	0.0	ug/g						
		PCB-1221		<	0.0	ug/g						
		PCB-1232		<	0.0	ug/g						
		PCB-1242		<	0.0	ug/g						
		PCB-1248		<	0.0	ug/g						
		PCB-1254		<	0.0	ug/g						
		PCB-1260		<	0.0	ug/g						
		tetrachloro-m-xylene SUR			65	%			30 150			
		decachlorobiphenyl SUR			84	%			30 150			
SW3540C8082A	LCS4297	PCB-1016			0.2	ug/g	0.2	79	40	140		
		PCB-1221		<	0.0	ug/g						
		PCB-1232		<	0.0	ug/g						
		PCB-1242		<	0.0	ug/g						
		PCB-1248		<	0.0	ug/g						
		PCB-1254		<	0.0	ug/g						
		PCB-1260			0.2	ug/g	0.2	95	40	140		
		tetrachloro-m-xylene SUR			59	%			30 150			
		decachlorobiphenyl SUR			86	%			30 150			
SW3540C8082A	LCSD4297	PCB-1016			0.1	ug/g	0.2	72	40	140	10	30
		PCB-1221		<	0.0	ug/g						
		PCB-1232		<	0.0	ug/g						
		PCB-1242		<	0.0	ug/g						
		PCB-1248		<	0.0	ug/g						
		PCB-1254		<	0.0	ug/g						
		PCB-1260			0.2	ug/g	0.2	92	40	140	3	30
		tetrachloro-m-xylene SUR			54	%			30 150			
		decachlorobiphenyl SUR			80	%			30 150			
SW3550B8100m	BLK4299	TPH C10-C36		<	200	ug/g						
		2-fluorobiphenyl SUR			92	%			40	140		
		o-terphenyl SUR			98	%			40	140		
SW3550B8100m	DUP4299	TPH C10-C36	21855-002		11000	ug/g				200	20	
		2-fluorobiphenyl SUR	21855-002		103	%			40	140		
		o-terphenyl SUR	21855-002		83	%			40	140		
SW3550B8100m	LCS4299	TPH C10-C36			2300	ug/g	2500	90	40	140		
		2-fluorobiphenyl SUR			105	%			40	140		
		o-terphenyl SUR			111	%			40	140		
SW3550B8100m	MS4299	TPH C10-C36	21855-002		12000	ug/g	3056.6	378	*	40	140	
		2-fluorobiphenyl SUR	21855-002		50	%			40	140		
		o-terphenyl SUR	21855-002		86	%			40	140		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3550B8270D	BLK4298	naphthalene		<	0.50	ug/g						
		2-methylnaphthalene		<	0.50	ug/g						
		acenaphthylene		<	0.50	ug/g						
		acenaphthene		<	0.50	ug/g						
		dibenzofuran		<	0.50	ug/g						
		fluorene		<	0.50	ug/g						
		phenanthrene		<	0.50	ug/g						
		anthracene		<	0.50	ug/g						
		fluoranthene		<	0.50	ug/g						
		pyrene		<	0.50	ug/g						
		benzo(a)anthracene		<	0.50	ug/g						
		chrysene		<	0.50	ug/g						
		benzo(b)fluoranthene		<	0.50	ug/g						
		benzo(k)fluoranthene		<	0.50	ug/g						
		benzo(a)pyrene		<	0.50	ug/g						
		indeno(1,2,3-cd)pyrene		<	0.50	ug/g						
		dibenzo(a,h)anthracene		<	0.50	ug/g						
		benzo(g,h,i)perylene		<	0.50	ug/g						
		2-fluorobiphenyl	SUR			88	%			43	116	
		o-terphenyl	SUR			86	%			33	141	
SW3550B8270D	DUP4298	naphthalene	21886-006	<	0.49	ug/g					30	
		2-methylnaphthalene	21886-006	<	0.49	ug/g					30	
		acenaphthylene	21886-006	<	0.49	ug/g					30	
		acenaphthene	21886-006	<	0.49	ug/g					30	
		dibenzofuran	21886-006	<	0.49	ug/g					30	
		fluorene	21886-006	<	0.49	ug/g					30	
		phenanthrene	21886-006	<	0.49	ug/g					30	
		anthracene	21886-006	<	0.49	ug/g					30	
		fluoranthene	21886-006	<	0.49	ug/g					30	
		pyrene	21886-006	<	0.49	ug/g					30	
		benzo(a)anthracene	21886-006	<	0.49	ug/g					30	
		chrysene	21886-006	<	0.49	ug/g					30	
		benzo(b)fluoranthene	21886-006	<	0.49	ug/g					30	
		benzo(k)fluoranthene	21886-006	<	0.49	ug/g					30	
		benzo(a)pyrene	21886-006	<	0.49	ug/g					30	
		indeno(1,2,3-cd)pyrene	21886-006	<	0.49	ug/g					30	
		dibenzo(a,h)anthracene	21886-006	<	0.49	ug/g					30	
		benzo(g,h,i)perylene	21886-006	<	0.49	ug/g					30	
		2-fluorobiphenyl	SUR	21886-006		86	%			43	116	
		o-terphenyl	SUR	21886-006		86	%			33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3550B8270D	LCS4298	naphthalene		3.4	ug/g	4	84	40 140			
		2-methylnaphthalene		3.6	ug/g	4	91	40 140			
		acenaphthylene		3.6	ug/g	4	91	40 140			
		acenaphthene		3.4	ug/g	4	85	40 140			
		dibenzofuran		<	0.50	ug/g					
		fluorene			3.5	ug/g	4	86	40 140		
		phenanthrene			3.2	ug/g	4	81	40 140		
		anthracene			3.6	ug/g	4	90	40 140		
		fluoranthene			3.5	ug/g	4	88	40 140		
		pyrene			4.3	ug/g	4	108	40 140		
		benzo(a)anthracene			3.9	ug/g	4	96	40 140		
		chrysene			3.8	ug/g	4	95	40 140		
		benzo(b)fluoranthene			4.4	ug/g	4	111	40 140		
		benzo(k)fluoranthene			3.3	ug/g	4	82	40 140		
		benzo(a)pyrene			3.5	ug/g	4	87	40 140		
		indeno(1,2,3-cd)pyrene			3.3	ug/g	4	83	40 140		
		dibenzo(a,h)anthracene			3.2	ug/g	4	79	40 140		
		benzo(g,h,i)perylene			3.3	ug/g	4	83	40 140		
		2-fluorobiphenyl SUR			87	%				43 116	
		o-terphenyl SUR			88	%				33 141	
SW3550B8270D	MS4298	naphthalene	21886-006	3.5	ug/g	4.08	84	40 140			
		2-methylnaphthalene	21886-006	3.5	ug/g	4.08	86	40 140			
		acenaphthylene	21886-006	3.9	ug/g	4.08	96	40 140			
		acenaphthene	21886-006	3.7	ug/g	4.08	91	40 140			
		dibenzofuran	21886-006	<	0.51	ug/g					
		fluorene	21886-006		3.6	ug/g	4.08	88	40 140		
		phenanthrene	21886-006		3.5	ug/g	4.08	86	40 140		
		anthracene	21886-006		3.5	ug/g	4.08	85	40 140		
		fluoranthene	21886-006		3.4	ug/g	4.08	85	40 140		
		pyrene	21886-006		4.3	ug/g	4.08	104	40 140		
		benzo(a)anthracene	21886-006		3.8	ug/g	4.08	93	40 140		
		chrysene	21886-006		3.9	ug/g	4.08	95	40 140		
		benzo(b)fluoranthene	21886-006		4.0	ug/g	4.08	97	40 140		
		benzo(k)fluoranthene	21886-006		3.5	ug/g	4.08	87	40 140		
		benzo(a)pyrene	21886-006		3.6	ug/g	4.08	87	40 140		
		indeno(1,2,3-cd)pyrene	21886-006		2.8	ug/g	4.08	68	40 140		
		dibenzo(a,h)anthracene	21886-006		2.6	ug/g	4.08	64	40 140		
		benzo(g,h,i)perylene	21886-006		2.7	ug/g	4.08	65	40 140		
		2-fluorobiphenyl SUR	21886-006		89	%				43 116	
		o-terphenyl SUR	21886-006		81	%				33 141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6010C	BLK4310	Silver		<	0.25	ug/g				
		Arsenic		<	0.50	ug/g				
		Barium		<	2.5	ug/g				
		Cadmium		<	0.20	ug/g				
		Chromium		<	2.5	ug/g				
		Copper		<	2.5	ug/g				
		Nickel		<	2.5	ug/g				
		Lead		<	0.50	ug/g				
		Selenium		<	2.5	ug/g				
		Zinc		<	2.5	ug/g				
SW3051A6010C	DUP4310	Lead	21906-011	4.9	ug/g				9	35
SW7471B	BLK4308	Mercury		<	0.02	ug/g				
SW7471B	CRM4308	Mercury		1.6	ug/g	1.1		0.49 1.76		
SW7471B	CRMD4308	Mercury		1.3	ug/g	1.1		0.49 1.76	20	20
SW7471B	MS4308	Mercury	21879-002	0.61	ug/g	0.406	110	75 125		
SW7471B	MSD4308	Mercury	21879-002	0.66	ug/g	0.406	121	75 125	7	35
SW7471B	REP4308	Mercury	21879-002							

Absolute Resource ASSOCIATES



124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001
absoluteresourceassociates.com

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

21886

Company Name: **CREDERE ASSOCIATES**
Company Address: **776 Main St. West**

Report To: **J. O'Donnell**

Phone #: **207-272-2225**

Invoice To: **T. Patten**

Project Name: **POLYCLAD**
Project #: **10081086**

Project Location: **MA ME VT**

Protocol: **RCPA SDWA NPDES OTHER**

Reporting Limits: **QAPP GW-1 S-1**

Quote # _____ Fund Pricing

Lab Sample ID	Field ID	# CONTAINERS	Matrix			Preservation Method							Sampling		
			WATER	SOLID	OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)	DATE	TIME	SAMPLER	
21886-01	CA-SS-1	1		X								06/28/11	1512	AW	
02	CA-SS-2	1										10/18	1628	AW	
03	CA-TP-8	2								X		11/05			
04	CA-TP-10	2								X		12/31			
05	CA-TP-11	2								X		1500			
06	CA-TP-13	2								X		1550			
07	CA-TP-14	2								X		1625			
08	CA-TP-15	2								X		1700			
09	CA-TP-16	2								X		1940			
10	CA-TP-17	2								X		1945			
11	CA-TP-18	2								X					

SPECIAL INSTRUCTIONS: **SPAWN FIELDS**

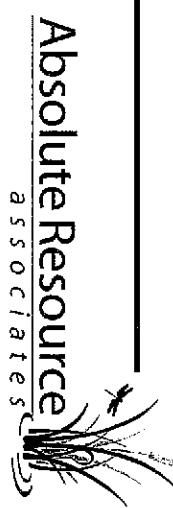
TAT REQUESTED
Priority (24 hr) *
Expedited (48 hr) *
Standard (10 Business Days)
*Date Needed _____

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.
 HARD COPY REQUIRED FAX (FAX#) _____
 OTHER (specify) _____

REPORTING INSTRUCTIONS
 PDF (e-mail address) **j.odonnell@credere.com**
 OTHER (specify) _____

RECEIVED ON ICE YES NO
TEMPERATURE **5** °C
Received by Laboratory: **J. O'Donnell**
Way Bill# _____
Date _____ Time _____

CUSTODY RECORD	Relinquished by: J. O'Donnell	Date: 6/28/11	Time: 110	Received by: J. O'Donnell	Date: 6/28/11	Time: 110
	Relinquished by: _____	Date: _____	Time: _____	Received by: _____	Date: _____	Time: _____



Absolute Resource Associates
a s s o c i a t e s

124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001
absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST**

21886

Company Name: **CAEDERE ASSOCIATES**

Company Address: **776 Main St, Westbrook, ME**

Project Name: **POLYCLAD**
Project #: **10001086**
Project Location: **(NH) MA ME VT**

Report To: **S. O'Donnell**

Protocol: **MCP**
Reporting: **QAPP**
Limits: **EPA DW**

Phone #: **207-272-2225**

Quote # _____
Fund Pricing

Invoice To: **T. Patten**

Other: _____
NH GREEN/ODD

Lab Sample ID	Field ID	# CONTAINERS	Matrix			Preservation Method							Sampling			SAMPLER	ANALYSIS REQUEST
			WATER	SOLID	OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)	DATE	TIME	TIME			
21886-12CA-TP-DW-1	Z	2	X	X				X				6/25/11	1535	590	X	<input checked="" type="checkbox"/> VOC 8260 <input type="checkbox"/> VOC 624 <input type="checkbox"/> VOC BTEX <input type="checkbox"/> MIBE, only <input type="checkbox"/> VOC 8021VT <input type="checkbox"/> VPH MADEP <input type="checkbox"/> MEGRO <input type="checkbox"/> GRO 8015 <input type="checkbox"/> VOC 524.2 <input type="checkbox"/> VOC 524.2 NH List <input type="checkbox"/> Gases-List: <input checked="" type="checkbox"/> TPH <input type="checkbox"/> DRO 8015 <input type="checkbox"/> MEDRO <input type="checkbox"/> EPH MADEP <input type="checkbox"/> TPH Fingerprint <input checked="" type="checkbox"/> 8270PAH <input type="checkbox"/> 8270ABN <input type="checkbox"/> 625 <input type="checkbox"/> EDB 504.1 <input checked="" type="checkbox"/> 8082 PCB <input type="checkbox"/> 8081 Pesticides <input type="checkbox"/> 608 Pest/PCB <input type="checkbox"/> O&G 1664 <input type="checkbox"/> Mineral O&G SM5520F <input type="checkbox"/> pH <input type="checkbox"/> BOD <input type="checkbox"/> Conductivity <input type="checkbox"/> Turbidity <input type="checkbox"/> TSS <input type="checkbox"/> TDS <input type="checkbox"/> TS <input type="checkbox"/> TVS <input type="checkbox"/> Alkalinity <input checked="" type="checkbox"/> RCRA Metals <input type="checkbox"/> Priority Pollutant Metals <input type="checkbox"/> TAL Metals <input checked="" type="checkbox"/> Total Metals-list: Pb <input checked="" type="checkbox"/> Dissolved Metals-list: Zn, Co, Ni <input type="checkbox"/> Ammonia <input type="checkbox"/> COD <input type="checkbox"/> TKN <input type="checkbox"/> TN <input type="checkbox"/> TON <input type="checkbox"/> T-Phosphorus <input type="checkbox"/> Phenols <input type="checkbox"/> Bacteria P/A <input type="checkbox"/> Bacteria MPN <input type="checkbox"/> Cyanide <input type="checkbox"/> Sulfide <input type="checkbox"/> Nitrate + Nitrite <input type="checkbox"/> Ortho P <input type="checkbox"/> Nitrate <input type="checkbox"/> Nitrite <input type="checkbox"/> Chloride <input type="checkbox"/> Sulfate <input type="checkbox"/> Bromide <input type="checkbox"/> Fluoride <input type="checkbox"/> Corrosivity <input type="checkbox"/> Reactive CN <input type="checkbox"/> Reactive S- <input type="checkbox"/> Ignitibility/FP <input type="checkbox"/> TCLP Metals <input type="checkbox"/> TCLP VOC <input type="checkbox"/> TCLP SVOC <input type="checkbox"/> TCLP Pesticide Subcontract: <input type="checkbox"/> TOC <input type="checkbox"/> Grain Size <input type="checkbox"/> TCLP Herbicides	
13 D-Tree	1	1	X	X								6/24/11	1255	590	X		
14 CA-3	1	1	X	X								6/24/11	1047	590	X		
16 Trip Blank	1	1															

TAT REQUESTED
Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)
*Date Needed _____

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

SPECIAL INSTRUCTIONS
BRAMFIELD

REPORTING INSTRUCTIONS
 HARD COPY REQUIRED FAX (FAX#) _____
 PDF (e-mail address) **y.odonnell@caedereilc.com** OTHER (specify) _____

RECEIVED ON ICE TEMPERATURE _____
YES NO
Date/Time: **6/29/11 11:01**

CUSTODY RECORD
OSD-01 Revision 12/23/10

Relinquished by: **[Signature]**
Date: **6/29/11**

Relinquished by: **[Signature]**
Date: _____

Received by Laboratory: **[Signature]**
Date: _____

Received by Laboratory: _____
Way Bill#: _____

Date: _____
Time: _____



ANALYTICAL REPORT

Lab Number:	L1110998
Client:	Absolute Resource Associates 124 Heritage Avenue Portsmouth, NH 03801
ATTN:	Jane Borkland
Phone:	(603) 436-2001
Project Name:	Not Specified
Project Number:	21886
Report Date:	07/28/11

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019
508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Alpha Sample ID	Client ID	Sample Location	Collection Date/Time
L1110998-01	21886-07 (CA-TP-14)	NH	06/28/11 15:50

Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

For additional information, please contact Client Services at 800-624-9220.

Sample Receipt

The samples were received at the laboratory above the required temperature range. The samples were transported to the laboratory in a cooler with ice and delivered directly from the sampling site.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:  Elizabeth Simmons

Title: Technical Director/Representative

Date: 07/28/11

INORGANICS & MISCELLANEOUS

Project Name: Not Specified

Lab Number: L1110998

Project Number: 21886

Report Date: 07/28/11

SAMPLE RESULTS

Lab ID: L1110998-01
 Client ID: 21886-07 (CA-TP-14)
 Sample Location: NH
 Matrix: Soil

Date Collected: 06/28/11 15:50
 Date Received: 07/21/11
 Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Analytical Method	Analyst
General Chemistry - Westborough Lab										
Solids, Total	92		%	0.10	NA	1	-	07/22/11 10:50	30,2540G	JC
Chromium, Hexavalent	4.5		mg/kg	0.87	--	1	07/24/11 14:00	07/24/11 18:00	1,7196A	JT



Project Name:
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Method Blank Analysis
Batch Quality Control

Parameter	Result Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Analytical Method	Analyst
General Chemistry - Westborough Lab for sample(s): 01 Batch: WG480802-1									
Chromium, Hexavalent	ND	mg/kg	0.80	--	1	07/24/11 14:00	07/24/11 18:00	1,7196A	JT

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Parameter	LCS %Recovery	Qual	LCS %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
General Chemistry - Westborough Lab Associated sample(s): 01 Batch: WG480802-2								
Chromium, Hexavalent	100		-		80-120	-		20



Matrix Spike Analysis
Batch Quality Control

Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	MSD Found	MSD %Recovery	MSD Qual	Recovery Limits	RPD Qual	RPD Limits
-----------	---------------	----------	----------	--------------	-----------	---------------	----------	-----------------	----------	------------

General Chemistry - Westborough Lab Associated sample(s): 01 QC Batch ID: WG480802-4 QC Sample: L1110998-01 Client ID: 21886-07 (CA-TP-14)

Chromium, Hexavalent	4.5	1150	1100	96	-	-	-	75-125	-	20
----------------------	-----	------	------	----	---	---	---	--------	---	----



Lab Duplicate Analysis

Batch Quality Control

Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Parameter	Native Sample	Duplicate Sample	Units	RPD	Qual	RPD Limits
General Chemistry - Westborough Lab Associated sample(s): 01 QC Batch ID: WG480548-1 QC Sample: L1110980-01 Client ID: DUP Sample						
Solids, Total	39	40	%	3		20
General Chemistry - Westborough Lab Associated sample(s): 01 QC Batch ID: WG480802-3 QC Sample: L1110998-01 Client ID: 21886-07 (CA-TP-14)						
Chromium, Hexavalent	4.5	4.3	mg/kg	5		20



Project Name: Not Specified

Lab Number: L1110998

Project Number: 21886

Report Date: 07/28/11

Sample Receipt and Container Information

Were project specific reporting limits specified? YES

Reagent H2O Preserved Vials Frozen on: NA

Cooler Information Custody Seal**Cooler**

A Absent

Container Information

Container ID	Container Type	Cooler	pH	Temp deg C	Pres	Seal	Analysis(*)
L1110998-01A	Amber 120ml unpreserved	A	N/A	8	Y	Absent	TS(7),HEXCR-7196(30)

*Values in parentheses indicate holding time in days



Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

GLOSSARY

Acronyms

EPA	- Environmental Protection Agency.
LCS	- Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
LCSD	- Laboratory Control Sample Duplicate: Refer to LCS.
LFB	- Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
MDL	- Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
MS	- Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.
MSD	- Matrix Spike Sample Duplicate: Refer to MS.
NA	- Not Applicable.
NC	- Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.
NI	- Not Ignitable.
RL	- Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
RPD	- Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.
SRM	- Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.

Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- A** - Spectra identified as "Aldol Condensation Product".
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than five times (5x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank.
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The RPD between the results for the two columns exceeds the method-specified criteria; however, the lower value has been reported due to obvious interference.
- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less

Report Format: Data Usability Report



Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

Data Qualifiers

than 5x the RL. (Metals only.)

R - Analytical results are from sample re-analysis.

RE - Analytical results are from sample re-extraction.

J - Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).

ND - Not detected at the reporting limit (RL) for the sample.

Project Name: Not Specified
Project Number: 21886

Lab Number: L1110998
Report Date: 07/28/11

REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - IIIA, 1997.
- 30 Standard Methods for the Examination of Water and Wastewater. APHA-AWWA-WPCF. 18th Edition. 1992.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised June 7, 2011 - Westboro Facility

The following list includes only those analytes/methods for which certification/approval is currently held.
For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0574. **NELAP Accredited Solid Waste/Soil.**

Drinking Water (Inorganic Parameters: Color, pH, Turbidity, Conductivity, Alkalinity, Chloride, Free Residual Chlorine, Fluoride, Calcium Hardness, Sulfate, Nitrate, Nitrite, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Dissolved Solids, Total Organic Carbon, Total Cyanide, Perchlorate. Organic Parameters: Volatile Organics 524.2, Total Trihalomethanes 524.2, 1,2-Dibromo-3-chloropropane (DBCP), Ethylene Dibromide (EDB), 1,4-Dioxane (Mod 8270). Microbiology Parameters: Total Coliform-MF mEndo (SM9222B), Total Coliform – Colilert (SM9223 P/A), E. Coli. – Colilert (SM9223 P/A), HPC – Pour Plate (SM9215B), Fecal Coliform – MF m-FC (SM9222D))

Wastewater/Non-Potable Water (Inorganic Parameters: Color, pH, Conductivity, Acidity, Alkalinity, Chloride, Total Residual Chlorine, Fluoride, Total Hardness, Silica, Sulfate, Sulfide, Ammonia, Kjeldahl Nitrogen, Nitrate, Nitrite, O-Phosphate, Total Phosphorus, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Titanium, Vanadium, Zinc, Total Residue (Solids), Total Dissolved Solids, Total Suspended Solids (non-filterable), BOD, CBOD, COD, TOC, Total Cyanide, Phenolics, Foaming Agents (MBAS), Bromide, Oil and Grease. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Acid Extractables (Phenols), Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, Polynuclear Aromatic Hydrocarbons, Haloethers, Chlorinated Hydrocarbons, Volatile Organics, TPH (HEM/SGT), Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH. Microbiology Parameters: Total Coliform – MF mEndo (SM9222B), Total Coliform – MTF (SM9221B), HPC – Pour Plate (SM9215B), Fecal Coliform – MF m-FC (SM9222D), Fecal Coliform – A-1 Broth (SM9221E).)

Solid Waste/Soil (Inorganic Parameters: pH, Sulfide, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Tin, Vanadium, Zinc, Total Cyanide, Ignitability, Phenolics, Corrosivity, TCLP Leach (1311), SPLP Leach (1312 metals only), Reactivity. Organic Parameters: PCBs, PCBs in Oil, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH, Dicamba, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Volatile Organics, Acid Extractables (Phenols), 3,3'-Dichlorobenzidine, Phthalates, Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Maine Department of Human Services Certificate/Lab ID: 2009024.

Drinking Water (Inorganic Parameters: SM9215B, 9222D, 9223B, EPA 180.1, 353.2, SM2130B, 2320B, 4500CI-D, 4500CN-C, 4500CN-E, 4500F-C, 4500H+B, 4500NO3-F, EPA 200.7, EPA 200.8, 245.1, EPA 300.0. Organic Parameters: 504.1, 524.2.)

Wastewater/Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664A, 350.1, 351.1, 353.2, 410.4, 420.1, SM2320B, 2510B, 2540C, 2540D, 426C, 4500CI-D, 4500CI-E, 4500CN-C, 4500CN-E, 4500F-B, 4500F-C, 4500H+B, 4500Norg-B, 4500Norg-C, 4500NH3-B, 4500NH3-G, 4500NH3-H, 4500NO3-F, 4500P-B, 4500P-E, 5210B, 5220D, 5310C, EPA 200.7, 200.8, 245.1. Organic Parameters: 608, 624, ME-DRO, ME-GRO, MA-EPH, MA-VPH.)

Solid Waste/Soil (Organic Parameters: ME-DRO, ME-GRO, MA-EPH, MA-VPH.)

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Drinking Water (Inorganic Parameters: (EPA 200.8 for: Sb,As,Ba,Be,Cd,Cr,Cu,Pb,Ni,Se,Tl) (EPA 200.7 for: Ba,Be,Ca,Cd,Cr,Cu,Na,Ni) 245.1, (300.0 for: Nitrate-N, Fluoride, Sulfate); (EPA 353.2 for: Nitrate-N, Nitrite-N); (SM4500NO3-F for: Nitrate-N and Nitrite-N); 4500F-C, 4500CN-CE, EPA 180.1, SM2130B, SM4500CI-D, 2320B, SM2540C, SM4500H-B. Organic Parameters: (EPA 524.2 for: Trihalomethanes, Volatile Organics); (504.1 for: 1,2-Dibromoethane, 1,2-Dibromo-3-Chloropropane), EPA 332. Microbiology Parameters: SM9215B; ENZ. SUB. SM9223; ColilertQT SM9223B; MF-SM9222D.)

Non-Potable Water (Inorganic Parameters: (EPA 200.8 for: Al,Sb,As,Be,Cd,Cr,Cu,Pb,Mn,Ni,Se,Ag,Tl,Zn); (EPA 200.7 for: Al,Sb,As,Be,Cd,Ca,Cr,Co,Cu,Fe,Pb,Mg,Mn,Mo,Ni,K,Se,Ag,Na,Sr,Ti,Tl, V,Zn); 245.1, SM4500H,B, EPA 120.1, SM2510B, 2540C, 2340B, 2320B, 4500CL-E, 4500F-BC, 426C, SM4500NH3-BH, (EPA 350.1 for: Ammonia-N), LACHAT 10-107-06-1-B for Ammonia-N, SM4500NO3-F, 353.2 for Nitrate-N, SM4500NH3-BC-NES, EPA 351.1, SM4500P-E, 4500P-B,E, 5220D, EPA 410.4, SM 5210B, 5310C, 4500CL-D, EPA 1664, SM14 510AC, EPA 420.1, SM4500-CN-CE, SM2540D.

Organic Parameters: (EPA 624 for Volatile Halocarbons, Volatile Aromatics),(608 for: Chlordane, Aldrin, Dieldrin, DDD, DDE, DDT, Heptachlor, Heptachlor Epoxide, PCBs-Water), (EPA 625 for SVOC Acid Extractables and SVOC Base/Neutral Extractables), 600/4-81-045-PCB-Oil. Microbiology Parameters: (ColilertQT SM9223B;Enterolert-QT: SM9222D-MF.)

New Hampshire Department of Environmental Services Certificate/Lab ID: 200307. *NELAP Accredited.*

Drinking Water (Inorganic Parameters: SM 9222B, 9223B, 9215B, EPA 200.7, 200.8, 245.2, 300.0, SM4500CN-E, 4500H+B, 4500NO3-F, 2320B, 2510B, 2540C, 4500F-C, 5310C, 2120B, EPA 332.0. Organic Parameters: 504.1, 524.2.)

Non-Potable Water (Inorganic Parameters: SM9222D, 9221B, 9222B, 9221E-EC, EPA 3005A, 200.7, 200.8, 245.1, 245.2, SW-846 6010B, 6020, 7196A, 7470A, SM3500-CR-D, EPA 120.1, 300.0, 350.1, 351.1, 353.2, 410.4, 420.1, 1664A, SW-846 9010, 9030, 9040B, 9050A, SM426C, SM2120B, 2310B, 2320B, 2540B, 2540D, 4500H+B, 4500CL-E, 4500CN-E, 4500NH3-H, 4500NO3-F, 4500NO2-B, 4500P-E, 4500-S2-D, 5210B, 5220D, 2510B, 2540C, 4500F-C, 5310C, 5540C, LACHAT 10-204-00-1-A, LACHAT 10-107-06-2-D. Organic Parameters: SW-846 3510C, 5030B, 8260B, 8270C, 8330, EPA 624, 625, 608, SW-846 8082, 8081A, 8151A.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6010B, 7196A, 7471A, 1010, 1030, 9010, 9012A, 9014, 9030B, 9040B, 9045C, 9050C, 9065,1311, 1312, 3005A, 3050B. Organic Parameters: SW-846 3540C, 3546, 3580A, 5030B, 5035, 8260B, 8270C, 8330, 8151A, 8015B, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA935. *NELAP Accredited.*

Drinking Water (Inorganic Parameters: SM9222B, 9221E, 9223B, 9215B, 4500CN-CE, 4500NO3-F, 4500F-C, EPA 300.0, 200.7, 200.8, 245.2, 2540C, SM2120B, 2320B, 2510B, 5310C, SM4500H-B. Organic Parameters: EPA 332, 504.1, 524.2.)

Non-Potable Water (Inorganic Parameters: SM5210B, EPA 410.4, SM5220D, 4500CI-E, EPA 300.0, SM2120B, SM4500F-BC, EPA 200.7, 351.1, LACHAT 10-107-06-2-D, EPA 353.2, SM4500NO3-F, 4500NO2-B, EPA 1664A, SM5310B, C or D, 4500-PE, EPA 420.1, SM510ABC, SM4500P-B5+E, 2540B, 2540C, 2540D, EPA 120.1, SM2510B, SM15 426C, 9222D, 9221B, 9221C, 9221E, 9222B, 9215B, 2310B, 2320B, 4500NH3-H, 4500-S D, EPA 350.1, 350.2, SW-846 1312, 6020, 7470A, 5540C, 4500H-B, EPA 200.8, SM3500Cr-D, 4500CN-CE, EPA 245.1, 245.2, SW-846 9040B, 3005A, EPA 6010B, 7196A, SW-846 9010B, 9030B. Organic Parameters: SW-846 8260B, 8270C, 8270C-SIM, 3510C, EPA 608, 624, 625, SW-846 3630C, 5030B, 8081A, 8082, 8151A, 8330, NJ OQA-QAM-025 Rev.7, NJ EPH.)

Solid & Chemical Materials (Inorganic Parameters: SW-846, 6010B, 7196A, 9010B, 9030B, 1010, 1030, 1311, 1312, 3005A, 3050B, 7471A, 9014, 9012A, 9040B, 9045C, 9050A, 9065. Organic Parameters: SW-846 8015B, 8081A, 8082, 8151A, 8330, 8260B, 8270C, 8270C-SIM, 3540C, 3545, 3546, 3550B, 3580A, 3630C, 5030B, 5035L, 5035H, NJ OQA-QAM-025 Rev.7, NJ EPH.)

New York Department of Health Certificate/Lab ID: 11148. *NELAP Accredited.*

Drinking Water (Inorganic Parameters: SM9223B, 9222B, 9215B, EPA 200.8, 200.7, 245.2, SM5310C, EPA 332.0, SM2320B, EPA 300.0, SM2120B, 4500CN-E, 4500F-C, 4500H-B, 4500NO3-F, 2540C, SM 2510B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: SM9221E, 9222D, 9221B, 9222B, 9215B, 5210B, 5310C, EPA 410.4, SM5220D, 2310B-4a, 2320B, EPA 200.7, 300.0, SM4500CL-E, 4500F-C, SM15 426C, EPA 350.1, SM4500NH3-BH, EPA 351.1, LACHAT 10-107-06-2, EPA 353.2, LACHAT 10-107-04-1-C, SM4500-NO3-F, 4500-NO2-B, 4500P-E, 2540C, 2540B, 2540D, EPA 200.8, EPA 6010B, 6020, EPA 7196A, SM3500Cr-D, EPA 245.1, 245.2, 7470A, SM2120B, LACHAT 10-204-00-1-A, EPA 9040B, SM4500-HB, EPA 1664A, EPA 420.1, SM14 510C, EPA 120.1, SM2510B, SM4500S-D, SM5540C, EPA 3005A, 9010B, 9030B.. Organic Parameters: EPA 624, 8260B, 8270C, 625, 608, 8081A, 8151A, 8330, 8082, EPA 3510C, 5030B.)

Solid & Hazardous Waste (Inorganic Parameters: 1010, 1030, EPA 6010B, 7196A, 7471A, 9012A, 9014, 9040B, 9045C, 9065, 9050, EPA 1311, 1312, 3005A, 3050B, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8015B, 8081A, 8151A, 8330, 8082, 3540C, 3545, 3546, 3580, 5030B, 5035.)

North Carolina Department of the Environment and Natural Resources Certificate/Lab ID : 666. Organic Parameters: MA-EPH, MA-VPH.

Pennsylvania Department of Environmental Protection Certificate/Lab ID : 68-03671. *NELAP Accredited.*

Drinking Water (Organic Parameters: EPA 524.2)

Non-Potable Water (Inorganic Parameters: EPA 1312. Organic Parameters: EPA 3510C, 5030B, 625, 624, 608, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Solid & Hazardous Waste (Inorganic Parameters: EPA 350.1, 1010, 1030, 1311, 1312, 3050B, 6010B, 7196A, 7471A, 9010B, 9012A, 9014, 9040B, 9045C, 9050, 9065, SM 4500NH3-H. Organic Parameters: 3540C, 3545, 3546, 3550B,

3580A, 3630C, 5035, 8015B, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Rhode Island Department of Health Certificate/Lab ID: LAO00065. *NELAP Accredited via NY-DOH.*

Refer to MA-DEP Certificate for Potable and Non-Potable Water.

Refer to NJ-DEP Certificate for Potable and Non-Potable Water.

Texas Commission on Environmental Quality Certificate/Lab ID: T104704476-09-1. *NELAP Accredited.*

Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664, 200.7, 200.8, 245.1, 245.2, 300.0, 350.1, 351.1, 353.2, 376.2, 410.4, 420.1, 6010, 6020, 7196, 7470, 9040, SM 2120B, 2310B, 2320B, 2510B, 2540B, 2540C, 2540D, 426C, 4500CL-E, 4500CN-E, 4500F-C, 4500H+B, 4500NH3-H, 4500NO2B, 4500P-E, 4500 S²⁻ D, 510C, 5210B, 5220D, 5310C, 5540C. Organic Parameters: EPA 608, 624, 625, 8081, 8082, 8151, 8260, 8270, 8330.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1311, 1312, 9012, 9014, 9040, 9045, 9050, 9065.)

Department of Defense Certificate/Lab ID: L2217.

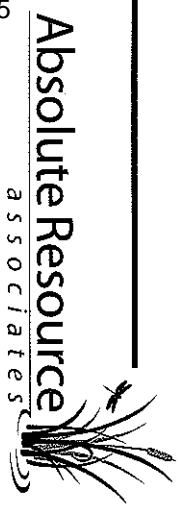
Drinking Water (Inorganic Parameters: SM 4500H-B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: EPA 200.7, 200.8, 6010B, 6020, 245.1, 245.2, 7470A, 9040B, 300.0, 332.0, 6860, 353.2, 410.4, 9060, 1664A, SM 4500CN-E, 4500H-B, 4500NO3-F, 5220D, 5310C, 2320B, 2540C, 3005A, 3015, 9010B, 9056. Organic Parameters: EPA 8260B, 8270C, 8330A, 625, 8082, 8081A, 3510C, 5030B, MassDEP EPH, MassDEP VPH.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 200.7, 6010B, 7471A, 9010, 9012A, 6860, 1311, 1312, 3050B, 7196A, 9010B, 3500-CR-D, 4500CN-CE, 2540G, Organic Parameters: EPA 8260B, 8270C, 8330A/B-prep, 8082, 8081A, 3540C, 3546, 3580A, 5035A, MassDEP EPH, MassDEP VPH.)

The following analytes are not included in our current NELAP/TNI Scope of Accreditation:

EPA 8260B: Freon-113, 1,2,4,5-Tetramethylbenzene, 4-Ethyltoluene. **EPA 8330A:** PETN, Picric Acid, Nitroglycerine, 2,6-DANT, 2,4-DANT. **EPA 8270C:** Methyl naphthalene, Dimethyl naphthalene, Total Methylnaphthalenes, Total Dimethylnaphthalenes, 1,4-Diphenylhydrazine (Azobenzene). **EPA 625:** 4-Chloroaniline, 4-Methylphenol. Total Phosphorus in a soil matrix, Chloride in a soil matrix, TKN in a soil matrix, NO₂ in a soil matrix, NO₃ in a soil matrix, SO₄ in a soil matrix.



Absolute Resource Associates
 124 Heritage Avenue #10
 Portsmouth, NH 03801
 603-436-2001
 absoluteresourceassociates.com

Alpha ALPHA Job # 2110998
 CHAIN-OF-CUSTODY RECORD
 AND ANALYSIS REQUEST
 PAGE 1 OF 1

Company Name: **See Above**
 Company Address: (Formerly Resource Labs)
 Project #: **218816**
 Project Location: **NH MA ME VT**
 Other:

Report to: **Sue Sylvester**
 Phone #: **603-077-1111**
 Protocol: RCRA SDWA NPDES
 MCP NHDES OTHER
 Reporting QAPP GW-1 S-1
 Limits: EPA DW Other
 Quote # **218816**
 PO # **218816**
 NH GREE/ODD
 Fund Pricing

Lab Sample ID	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		DATE	TIME	SAMPLER
			WATER	SOLID	OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)				
10998-1	218816-07	1		X									7/28/11	1550	
	(A-TP-14)														

TAT REQUESTED
 Priority (24 hr)*
 Expedited (48 hr)*
 Standard (10 Business Days)
 *Date Needed _____

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

REPORTING INSTRUCTIONS
 HARD COPY REQUIRED FAX (FAX#) _____
 PDF (e-mail address) **jennifer@absoluteresourceassociates.com**
Jane B OTHER (specify) _____

SPECIAL INSTRUCTIONS
 Please note holding time.

RECEIVED ON ICE YES NO
 TEMPERATURE _____ °C

CUSTODY RECORD
 OSD-01 Revision 12/23/10

Relinquished by: Jay	Date: 7/21/11	Time: 1:32	Received by: Scotts R	Date: 7/21/11	Time: 13:40
Relinquished by: _____	Date: _____	Time: _____	Received by: _____	Date: _____	Time: _____



Absolute Resource *associates*

124 Heritage Avenue #10 Portsmouth, NH 03801

Rick Vandenberg
CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: 10001086
Job ID: 21906
Date Received: 7/1/11

Project: POLYCLAD 10001086

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

Sue Sylvester
Principal, General Manager

Date of Approval: 7/14/2011
Total number of pages: 50

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-3 2-4	Solid	6/29/2011 12:59	21906-001	Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-2 6-8	Solid	6/29/2011 12:55	21906-002	Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-5 12-14	Solid	6/29/2011 17:10	21906-006	PCBs in soil by 8082 TPH in solids by 8100 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-1 26-28	Solid	6/29/2011 15:56	21906-007	TPH in solids by 8100 PAHs in solid by 8270 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-4 10-12	Solid	6/29/2011 17:40	21906-008	PCBs in soil by 8082 TPH in solids by 8100 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Selenium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-9 6'	Solid	6/30/2011 15:45	21906-009	TPH in solids by 8100 PAHs in solid by 8270 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-7A 10-12'	Solid	6/30/2011 17:30	21906-010	TPH in solids by 8100 PAHs in solid by 8270 Solid Digestion for ICP Analysis Lead in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-12 1'	Solid	6/30/2011 13:15	21906-011	Solid Digestion for ICP Analysis Lead in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solid by 8260 Petro & Haz Waste
CA-TP-DUP-2	Solid	6/30/2011 0:00	21906-012	TPH in solids by 8100 Percent Dry Matter for Sample Calc by SM2540B,G

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-002

Sample ID: CA-2 6-8

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 12:55

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
acetone	< 3	3	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
t-butanol (TBA)	< 3	3	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
2-butanone (MEK)	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
tetrahydrofuran (THF)	< 0.6	0.6	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,4-dioxane	< 3	3	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.5	0.5	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
2-hexanone	< 0.6	0.6	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-002

Sample ID: CA-2 6-8

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 12:55

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	79	78-114	%	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
toluene-D8 SUR	91	88-110	%	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
4-bromofluorobenzene SUR	95	86-115	%	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B
a,a,a-trifluorotoluene SUR	115	70-130	%	1	LMM	7/6/11	4307	7/8/11	15:29	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
acetone	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
t-butanol (TBA)	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
2-butanone (MEK)	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
tetrahydrofuran (THF)	< 0.7	0.7	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,4-dioxane	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.6	0.6	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
2-hexanone	< 0.7	0.7	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	82	78-114	%	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
toluene-D8 SUR	92	88-110	%	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
4-bromofluorobenzene SUR	95	86-115	%	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B
a,a,a-trifluorotoluene SUR	108	70-130	%	1	LMM	7/6/11	4307	7/8/11	16:00	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-007

Sample ID: CA-1 26-28

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/29/11 15:56

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
acetone	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
t-butanol (TBA)	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
2-butanone (MEK)	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
tetrahydrofuran (THF)	< 0.7	0.7	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,4-dioxane	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.6	0.6	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
2-hexanone	< 0.7	0.7	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-007

Sample ID: CA-1 26-28

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/29/11 15:56

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	80	78-114	%	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
toluene-D8 SUR	90	88-110	%	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
4-bromofluorobenzene SUR	100	86-115	%	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B
a,a,a-trifluorotoluene SUR	95	70-130	%	1	LMM	7/6/11	4307	7/8/11	18:06	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid

Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
chloromethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
vinyl chloride	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
bromomethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
chloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
trichlorofluoromethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
diethyl ether	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
acetone	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1-dichloroethene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
methylene chloride	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
carbon disulfide	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
trans-1,2-dichloroethene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
isopropyl ether (DIPE)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1-dichloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
t-butanol (TBA)	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
2-butanone (MEK)	< 0.5	0.5	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
2,2-dichloropropane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
cis-1,2-dichloroethene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
chloroform	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
bromochloromethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
tetrahydrofuran (THF)	< 0.9	0.9	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1,1-trichloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1-dichloropropene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
carbon tetrachloride	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2-dichloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
benzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
trichloroethene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2-dichloropropane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
bromodichloromethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,4-dioxane	< 4	4	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
dibromomethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.8	0.8	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
cis-1,3-dichloropropene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
toluene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
trans-1,3-dichloropropene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
2-hexanone	< 0.9	0.9	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1,2-trichloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,3-dichloropropane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
tetrachloroethene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
dibromochloromethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid

Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
chlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
ethylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
m&p-xylenes	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
o-xylene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
styrene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
bromoform	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
isopropylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2,3-trichloropropane	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
n-propylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
bromobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,3,5-trimethylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
2-chlorotoluene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
4-chlorotoluene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
tert-butylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2,4-trimethylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
sec-butylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,3-dichlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
4-isopropyltoluene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,4-dichlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2-dichlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
n-butylbenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2,4-trichlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,3,5-trichlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
hexachlorobutadiene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
1,2,3-trichlorobenzene	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	80	78-114	%	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
toluene-D8 SUR	93	88-110	%	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
4-bromofluorobenzene SUR	98	86-115	%	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B
a,a,a-trifluorotoluene SUR	108	70-130	%	1	LMM	7/6/11	4307	7/8/11	18:35	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-009

Sample ID: CA-TP-9 6'

Matrix: Solid

Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/30/11 15:45

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-009

Sample ID: CA-TP-9 6'

Matrix: Solid

Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/30/11 15:45

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	78	78-114	%	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
toluene-D8 SUR	93	88-110	%	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
4-bromofluorobenzene SUR	89	86-115	%	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B
a,a,a-trifluorotoluene SUR	98	70-130	%	1	LMM	7/6/11	4307	7/8/11	19:04	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-010

Sample ID: CA-TP-7A 10-12'

Matrix: Solid Percent Dry: 97.3% Results expressed on a dry weight basis.

Sampled: 6/30/11 17:30

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-010

Sample ID: CA-TP-7A 10-12'

Matrix: Solid Percent Dry: 97.3% Results expressed on a dry weight basis.

Sampled: 6/30/11 17:30

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	85	78-114	%	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
toluene-D8 SUR	91	88-110	%	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
4-bromofluorobenzene SUR	94	86-115	%	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B
a,a,a-trifluorotoluene SUR	98	70-130	%	1	LMM	7/6/11	4307	7/11/11	13:44	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-011

Sample ID: CA-TP-12 1'

Matrix: Solid Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/30/11 13:15

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
bromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
diethyl ether	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
methylene chloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
isopropyl ether (DIPE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
ethyl t-butyl ether (ETBE)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
t-butanol (TBA)	< 2	2	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
2-butanone (MEK)	< 0.2	0.2	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
tetrahydrofuran (THF)	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
t-amyl-methyl ether (TAME)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,4-dioxane	< 2	2	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.3	0.3	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
2-hexanone	< 0.4	0.4	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-011

Sample ID: CA-TP-12 1'

Matrix: Solid

Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/30/11 13:15

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,3,5-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
naphthalene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	84	78-114	%	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
toluene-D8 SUR	90	88-110	%	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
4-bromofluorobenzene SUR	96	86-115	%	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B
a,a,a-trifluorotoluene SUR	91	70-130	%	1	LMM	7/6/11	4307	7/11/11	14:16	SW5035A8260B

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
N-nitrosodimethylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
aniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
phenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-chlorophenol	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
bis(2-chloroethyl)ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
1,3-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
1,4-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
1,2-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzyl alcohol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
bis(2-chloroisopropyl) ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
hexachloroethane	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
N-nitroso-di-N-propylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
nitrobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
isophorone	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-nitrophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4-dimethylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
bis(2-chloroethoxy)methane	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4-dichlorophenol	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
1,2,4-trichlorobenzene	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
naphthalene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzoic acid	< 6	6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4-chloroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
hexachlorobutadiene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4-chloro-3-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-methylnaphthalene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
hexachlorocyclopentadiene	< 1	1	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4,6-trichlorophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4,5-trichlorophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-chloronaphthalene	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-nitroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
acenaphthylene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
dimethylphthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,6-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
acenaphthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
3-nitroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4-dinitrophenol	< 6	6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
dibenzofuran	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4-nitrophenol	< 2	2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
fluorene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
diethyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4-nitroaniline	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4,6-dinitro-2-methylphenol	< 2	2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
azobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
N-nitrosodiphenylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
4-bromophenyl phenyl ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
hexachlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
pentachlorophenol	< 1	1	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
phenanthrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
anthracene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
carbazole	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
di-n-butylphthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
fluoranthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzidine	< 4	4	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
pyrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
butyl benzyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzo(a)anthracene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
chrysene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
3,3'-dichlorobenzidine	< 4	4	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
bis(2-ethylhexyl)phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
di-n-octyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzo(b)fluoranthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzo(k)fluoranthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzo(a)pyrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
indeno(1,2,3-cd)pyrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
dibenzo(a,h)anthracene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
benzo(g,h,i)perylene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
Surrogate Recovery		Limits								
2-fluorophenol SUR	53	21-100	%	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
phenol-D5 SUR	60	10-102	%	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2,4,6-tribromophenol SUR	27	10-123	%	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
nitrobenzene-D5 SUR	47	35-114	%	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
2-fluorobiphenyl SUR	61	43-116	%	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D
p-terphenyl-D14 SUR	67	33-141	%	1	AJD	7/6/11	4314	7/9/11	16:04	SW3546/8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-007

Sample ID: CA-1 26-28

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/29/11 15:56

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
phenanthrene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
pyrene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
benzo(a)anthracene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
chrysene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
benzo(b)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
benzo(k)fluoranthene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
benzo(a)pyrene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
benzo(g,h,i)perylene	< 0.6	0.6	ug/g	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	78	43-116	%	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D
o-terphenyl SUR	82	33-141	%	1	AJD	7/7/11	4315	7/8/11	16:25	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid

Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
N-nitrosodimethylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
aniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
phenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-chlorophenol	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
bis(2-chloroethyl)ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
1,3-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
1,4-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
1,2-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzyl alcohol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
bis(2-chloroisopropyl) ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
hexachloroethane	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
N-nitroso-di-N-propylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
nitrobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
isophorone	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-nitrophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4-dimethylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
bis(2-chloroethoxy)methane	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4-dichlorophenol	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
1,2,4-trichlorobenzene	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
naphthalene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzoic acid	< 6	6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4-chloroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
hexachlorobutadiene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4-chloro-3-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-methylnaphthalene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
hexachlorocyclopentadiene	< 1	1	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4,6-trichlorophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4,5-trichlorophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-chloronaphthalene	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-nitroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
acenaphthylene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
dimethylphthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,6-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
acenaphthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
3-nitroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4-dinitrophenol	< 6	6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
dibenzofuran	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4-nitrophenol	< 2	2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
fluorene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
diethyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid

Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4-nitroaniline	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4,6-dinitro-2-methylphenol	< 2	2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
azobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
N-nitrosodiphenylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
4-bromophenyl phenyl ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
hexachlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
pentachlorophenol	< 1	1	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
phenanthrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
anthracene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
carbazole	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
di-n-butylphthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
fluoranthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzidine	< 4	4	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
pyrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
butyl benzyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzo(a)anthracene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
chrysene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
3,3'-dichlorobenzidine	< 4	4	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
bis(2-ethylhexyl)phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
di-n-octyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzo(b)fluoranthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzo(k)fluoranthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzo(a)pyrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
indeno(1,2,3-cd)pyrene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
dibenzo(a,h)anthracene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
benzo(g,h,i)perylene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
Surrogate Recovery		Limits								
2-fluorophenol SUR	48	21-100	%	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
phenol-D5 SUR	58	10-102	%	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2,4,6-tribromophenol SUR	29	10-123	%	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
nitrobenzene-D5 SUR	48	35-114	%	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
2-fluorobiphenyl SUR	64	43-116	%	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D
p-terphenyl-D14 SUR	71	33-141	%	1	AJD	7/6/11	4314	7/9/11	16:42	SW3546/8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-009

Sample ID: CA-TP-9 6'

Matrix: Solid

Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/30/11 15:45

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	81	43-116	%	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D
o-terphenyl SUR	86	33-141	%	1	AJD	7/7/11	4315	7/8/11	19:34	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-010

Sample ID: CA-TP-7A 10-12'

Matrix: Solid Percent Dry: 97.3% Results expressed on a dry weight basis.

Sampled: 6/30/11 17:30

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	77	43-116	%	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D
o-terphenyl SUR	77	33-141	%	1	AJD	7/7/11	4315	7/8/11	15:43	SW3550B8270D

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-003

Sample ID: CA-CC-1

Matrix: Solid Percent Dry: 98% Results expressed on a dry weight basis.

Sampled: 6/30/11 12:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
PCB-1221	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
PCB-1232	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
PCB-1242	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
PCB-1248	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
PCB-1254	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
PCB-1260	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	12:47	SW3540C8082A

DOR = Diluted out of range.

Note: Dilution due to matrix interference.

Sample#: 21906-004

Sample ID: CA-CC-DUP

Matrix: Solid Percent Dry: 78.9% Results expressed on a dry weight basis.

Sampled: 6/30/11 12:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
PCB-1221	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
PCB-1232	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
PCB-1242	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
PCB-1248	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
PCB-1254	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
PCB-1260	< 0.8	0.8	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	14:50	SW3540C8082A

DOR = Diluted out of range.

Note: Dilution due to matrix interference.

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-005

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 97.8% Results expressed on a dry weight basis.

Sampled: 6/30/11 12:15

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
PCB-1221	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
PCB-1232	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
PCB-1242	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
PCB-1248	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
PCB-1254	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
PCB-1260	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	13:18	SW3540C8082A

DOR = Diluted out of range.

Note: Dilution due to matrix interference.

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
PCB-1221	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
PCB-1232	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
PCB-1242	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
PCB-1248	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
PCB-1254	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
PCB-1260	< 0.7	0.7	ug/g	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	14:19	SW3540C8082A

DOR = Diluted out of range.

Note: Dilution due to matrix interference.

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
PCB-1221	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
PCB-1232	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
PCB-1242	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
PCB-1248	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
PCB-1254	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
PCB-1260	< 0.6	0.6	ug/g	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ	7/6/11	4305	7/11/11	13:48	SW3540C8082A

DOR = Diluted out of range.

Note: Dilution due to matrix interference.

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 250	250	ug/g	1	JLZ	7/7/11	4312	7/8/11	19:52	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	78	40-140	%	1	JLZ	7/7/11	4312	7/8/11	19:52	SW3550B8100m
o-terphenyl SUR	99	40-140	%	1	JLZ	7/7/11	4312	7/8/11	19:52	SW3550B8100m

Sample#: 21906-007

Sample ID: CA-1 26-28

Matrix: Solid Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 6/29/11 15:56

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 220	220	ug/g	1	JLZ	7/7/11	4312	7/8/11	20:09	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	74	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:09	SW3550B8100m
o-terphenyl SUR	99	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:09	SW3550B8100m

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 240	240	ug/g	1	JLZ	7/7/11	4312	7/8/11	20:25	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	74	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:25	SW3550B8100m
o-terphenyl SUR	93	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:25	SW3550B8100m

Sample#: 21906-009

Sample ID: CA-TP-9 6'

Matrix: Solid Percent Dry: 95.7% Results expressed on a dry weight basis.

Sampled: 6/30/11 15:45

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 200	200	ug/g	1	JLZ	7/7/11	4312	7/8/11	20:41	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	68	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:41	SW3550B8100m
o-terphenyl SUR	88	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:41	SW3550B8100m

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-010

Sample ID: CA-TP-7A 10-12'

Matrix: Solid Percent Dry: 97.3% Results expressed on a dry weight basis.

Sampled: 6/30/11 17:30

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 190	190	ug/g	1	JLZ	7/7/11	4312	7/8/11	20:58	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	68	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:58	SW3550B8100m
o-terphenyl SUR	87	40-140	%	1	JLZ	7/7/11	4312	7/8/11	20:58	SW3550B8100m

Sample#: 21906-012

Sample ID: CA-TP-DUP-2

Matrix: Solid Percent Dry: 94.7% Results expressed on a dry weight basis.

Sampled: 6/30/11 0:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
TPH C10-C36	< 200	200	ug/g	1	JLZ	7/7/11	4312	7/9/11	0:30	SW3550B8100m
Surrogate Recovery		Limits								
2-fluorobiphenyl SUR	80	40-140	%	1	JLZ	7/7/11	4312	7/9/11	0:30	SW3550B8100m
o-terphenyl SUR	103	40-140	%	1	JLZ	7/7/11	4312	7/9/11	0:30	SW3550B8100m

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-001

Sample ID: CA-3 2-4

Matrix: Solid Percent Dry: 95.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 12:59

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	4.4	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Barium	16	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Chromium	6	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Copper	16	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Lead	5.5	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Mercury	< 0.08	0.08	ug/g	1	BJS	7/6/11	4309	7/7/11	10:45	SW7471B
Nickel	7	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C
Zinc	21	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:41	SW3051A6010C

Sample#: 21906-002

Sample ID: CA-2 6-8

Matrix: Solid Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 12:55

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	1.5	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Barium	13	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Chromium	25	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Copper	12	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Lead	4.9	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Mercury	< 0.07	0.07	ug/g	1	BJS	7/6/11	4309	7/7/11	10:47	SW7471B
Nickel	5	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C
Zinc	23	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:49	SW3051A6010C

Project ID: POLYCLAD 10001086

Job ID: 21906

Sample#: 21906-006

Sample ID: CA-5 12-14

Matrix: Solid Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:10

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	2.2	0.7	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Barium	16	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Cadmium	< 0.3	0.3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Chromium	7	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Copper	22	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Lead	4.1	0.7	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Mercury	< 0.08	0.08	ug/g	1	BJS	7/6/11	4309	7/7/11	10:49	SW7471B
Nickel	6	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Silver	< 0.5	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C
Zinc	17	3	ug/g	1	BJS	7/6/11	4310	7/7/11	0:56	SW3051A6010C

Sample#: 21906-008

Sample ID: CA-4 10-12

Matrix: Solid Percent Dry: 81.4% Results expressed on a dry weight basis.

Sampled: 6/29/11 17:40

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	2.1	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Barium	15	3	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Chromium	6	3	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Copper	14	3	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Lead	4.0	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Mercury	< 0.07	0.07	ug/g	1	BJS	7/6/11	4309	7/7/11	10:51	SW7471B
Nickel	5	3	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Selenium	< 3	3	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C
Zinc	14	3	ug/g	1	BJS	7/6/11	4310	7/7/11	1:04	SW3051A6010C

Sample#: 21906-010

Sample ID: CA-TP-7A 10-12'

Matrix: Solid Percent Dry: 97.3% Results expressed on a dry weight basis.

Sampled: 6/30/11 17:30

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Lead	5.6	0.5	ug/g	1	BJS	7/6/11	4310	7/7/11	1:11	SW3051A6010C

Sample#: 21906-011

Sample ID: CA-TP-12 1'

Matrix: Solid Percent Dry: 91% Results expressed on a dry weight basis.

Sampled: 6/30/11 13:15

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Lead	4.5	0.6	ug/g	1	BJS	7/6/11	4310	7/7/11	1:19	SW3051A6010C

Quality Control Report



124 Heritage Avenue Unit 10
Portsmouth, NH 03801

www.absoluteresourceassociates.com



Case Narrative

Lab # 21906

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 1 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

No exceptions noted.

Surrogate Recoveries

PCB: The surrogates were diluted out of the calibration range in the following sample due to interferences caused by the matrix: 21906-003, -004, -005, -006, and -008.

Laboratory Control Sample Results

PCB: The relative percent difference between the LCS4305 and LCSD4305 was outside the acceptance criteria for PCB-1016 and PCB-1260. The percent recovery for these analytes in each QC parameter was within the acceptance criteria. No impact to the data suspected.

VOC: The MLCS/D4307 did not meet the acceptance criteria for dichlorodifluoromethane and vinyl chloride. These compounds are known to be problematic in the method.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MB4307	dichlorodifluoromethane		<	0.1	ug/g				
		chloromethane		<	0.1	ug/g				
		vinyl chloride		<	0.1	ug/g				
		bromomethane		<	0.2	ug/g				
		chloroethane		<	0.1	ug/g				
		trichlorofluoromethane		<	0.1	ug/g				
		diethyl ether		<	0.5	ug/g				
		acetone		<	2.5	ug/g				
		1,1-dichloroethene		<	0.1	ug/g				
		methylene chloride		<	0.2	ug/g				
		carbon disulfide		<	0.1	ug/g				
		methyl t-butyl ether (MTBE)		<	0.1	ug/g				
		trans-1,2-dichloroethene		<	0.1	ug/g				
		isopropyl ether (DIPE)		<	0.1	ug/g				
		ethyl t-butyl ether (ETBE)		<	0.1	ug/g				
		1,1-dichloroethane		<	0.1	ug/g				
		t-butanol (TBA)		<	2.5	ug/g				
		2-butanone (MEK)		<	0.5	ug/g				
		2,2-dichloropropane		<	0.1	ug/g				
		cis-1,2-dichloroethene		<	0.1	ug/g				
		chloroform		<	0.1	ug/g				
		bromochloromethane		<	0.1	ug/g				
		tetrahydrofuran (THF)		<	0.5	ug/g				
		1,1,1-trichloroethane		<	0.1	ug/g				
		1,1-dichloropropene		<	0.1	ug/g				
		t-amyl-methyl ether (TAME)		<	0.1	ug/g				
		carbon tetrachloride		<	0.1	ug/g				
		1,2-dichloroethane		<	0.1	ug/g				
		benzene		<	0.1	ug/g				
		trichloroethene		<	0.1	ug/g				
		1,2-dichloropropane		<	0.1	ug/g				
		bromodichloromethane		<	0.1	ug/g				
		1,4-dioxane		<	2.5	ug/g				
		dibromomethane		<	0.1	ug/g				
		4-methyl-2-pentanone (MIBK)		<	0.5	ug/g				
		cis-1,3-dichloropropene		<	0.1	ug/g				
		toluene		<	0.1	ug/g				
		trans-1,3-dichloropropene		<	0.1	ug/g				
		2-hexanone		<	0.5	ug/g				
		1,1,2-trichloroethane		<	0.1	ug/g				
		1,3-dichloropropane		<	0.1	ug/g				
		tetrachloroethene		<	0.1	ug/g				
		dibromochloromethane		<	0.1	ug/g				
		1,2-dibromoethane (EDB)		<	0.1	ug/g				
		chlorobenzene		<	0.1	ug/g				
		1,1,1,2-tetrachloroethane		<	0.1	ug/g				
		ethylbenzene		<	0.1	ug/g				
		m&p-xylenes		<	0.1	ug/g				
		o-xylene		<	0.1	ug/g				
		styrene		<	0.1	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MB4307	bromoform		<	0.1	ug/g				
		isopropylbenzene		<	0.1	ug/g				
		1,1,2,2-tetrachloroethane		<	0.1	ug/g				
		1,2,3-trichloropropane		<	0.1	ug/g				
		n-propylbenzene		<	0.1	ug/g				
		bromobenzene		<	0.1	ug/g				
		1,3,5-trimethylbenzene		<	0.1	ug/g				
		2-chlorotoluene		<	0.1	ug/g				
		4-chlorotoluene		<	0.1	ug/g				
		tert-butylbenzene		<	0.1	ug/g				
		1,2,4-trimethylbenzene		<	0.1	ug/g				
		sec-butylbenzene		<	0.1	ug/g				
		1,3-dichlorobenzene		<	0.1	ug/g				
		4-isopropyltoluene		<	0.1	ug/g				
		1,4-dichlorobenzene		<	0.1	ug/g				
		1,2-dichlorobenzene		<	0.1	ug/g				
		n-butylbenzene		<	0.1	ug/g				
		1,2-dibromo-3-chloropropane		<	0.1	ug/g				
		1,2,4-trichlorobenzene		<	0.1	ug/g				
		hexachlorobutadiene		<	0.1	ug/g				
		naphthalene		<	0.2	ug/g				
		1,2,3-trichlorobenzene		<	0.1	ug/g				
		dibromofluoromethane SUR			89	%		78 114		
		toluene-D8 SUR			93	%		88 110		
		4-bromofluorobenzene SUR			97	%		86 115		
		a,a,a-trifluorotoluene SUR			95	%		70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCS4307	dichlorodifluoromethane		0.6	ug/g	1	61 *	70 130		
		chloromethane		0.8	ug/g	1	76	70 130		
		vinyl chloride		0.5	ug/g	1	52 *	70 130		
		bromomethane		1.1	ug/g	1	106	70 130		
		chloroethane		0.9	ug/g	1	92	70 130		
		trichlorofluoromethane		0.8	ug/g	1	82	70 130		
		diethyl ether		1.0	ug/g	1	97	70 130		
		acetone	<	2.5	ug/g	1	89			
		1,1-dichloroethene		0.8	ug/g	1	76	70 130		
		methylene chloride		0.9	ug/g	1	94	70 130		
		carbon disulfide		0.8	ug/g	1	79	70 130		
		methyl t-butyl ether (MTBE)		1.0	ug/g	1	95	70 130		
		trans-1,2-dichloroethene		0.8	ug/g	1	84	70 130		
		isopropyl ether (DIPE)		1.0	ug/g	1	98	70 130		
		ethyl t-butyl ether (ETBE)		1.0	ug/g	1	96	70 130		
		1,1-dichloroethane		0.9	ug/g	1	89	70 130		
		t-butanol (TBA)		6.3	ug/g	5	127	70 130		
		2-butanone (MEK)		1.0	ug/g	1	98	70 130		
		2,2-dichloropropane		1.1	ug/g	1	107	70 130		
		cis-1,2-dichloroethene		0.9	ug/g	1	95	70 130		
		chloroform		1.0	ug/g	1	97	70 130		
		bromochloromethane		1.0	ug/g	1	95	70 130		
		tetrahydrofuran (THF)		1.0	ug/g	1	101	70 130		
		1,1,1-trichloroethane		0.9	ug/g	1	93	70 130		
		1,1-dichloropropene		1.0	ug/g	1	95	70 130		
		t-amyl-methyl ether (TAME)		0.9	ug/g	1	93	70 130		
		carbon tetrachloride		0.9	ug/g	1	91	70 130		
		1,2-dichloroethane		1.0	ug/g	1	97	70 130		
		benzene		1.0	ug/g	1	97	70 130		
		trichloroethene		0.9	ug/g	1	93	70 130		
		1,2-dichloropropane		1.0	ug/g	1	99	70 130		
		bromodichloromethane		1.0	ug/g	1	97	70 130		
		1,4-dioxane	<	2.5	ug/g	2	113	70 130		
		dibromomethane		1.0	ug/g	1	97	70 130		
		4-methyl-2-pentanone (MIBK)		1.0	ug/g	1	98	70 130		
		cis-1,3-dichloropropene		1.0	ug/g	1	102	70 130		
		toluene		1.0	ug/g	1	97	70 130		
		trans-1,3-dichloropropene		1.1	ug/g	1	109	70 130		
		2-hexanone		0.9	ug/g	1	91	70 130		
		1,1,2-trichloroethane		1.0	ug/g	1	100	70 130		
		1,3-dichloropropane		0.9	ug/g	1	89	70 130		
		tetrachloroethene		1.0	ug/g	1	97	70 130		
		dibromochloromethane		0.8	ug/g	1	81	70 130		
		1,2-dibromoethane (EDB)		0.9	ug/g	1	87	70 130		
		chlorobenzene		1.0	ug/g	1	100	70 130		
		1,1,1,2-tetrachloroethane		0.9	ug/g	1	95	70 130		
		ethylbenzene		0.9	ug/g	1	89	70 130		
		m&p-xylenes		2.0	ug/g	2	100	70 130		
		o-xylene		1.1	ug/g	1	106	70 130		
		styrene		0.9	ug/g	1	94	70 130		
		bromoform		1.1	ug/g	1	108	70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCS4307	isopropylbenzene		1.0	ug/g	1	101	70 130		
		1,1,2,2-tetrachloroethane		1.1	ug/g	1	112	70 130		
		1,2,3-trichloropropane		1.1	ug/g	1	108	70 130		
		n-propylbenzene		1.1	ug/g	1	107	70 130		
		bromobenzene		1.0	ug/g	1	101	70 130		
		1,3,5-trimethylbenzene		1.0	ug/g	1	105	70 130		
		2-chlorotoluene		1.1	ug/g	1	108	70 130		
		4-chlorotoluene		1.1	ug/g	1	107	70 130		
		tert-butylbenzene		1.1	ug/g	1	108	70 130		
		1,2,4-trimethylbenzene		1.0	ug/g	1	105	70 130		
		sec-butylbenzene		1.1	ug/g	1	110	70 130		
		1,3-dichlorobenzene		1.1	ug/g	1	112	70 130		
		4-isopropyltoluene		1.1	ug/g	1	111	70 130		
		1,4-dichlorobenzene		1.0	ug/g	1	101	70 130		
		1,2-dichlorobenzene		1.0	ug/g	1	102	70 130		
		n-butylbenzene		1.1	ug/g	1	110	70 130		
		1,2-dibromo-3-chloropropane		0.9	ug/g	1	89	70 130		
		1,2,4-trichlorobenzene		0.9	ug/g	1	88	70 130		
		hexachlorobutadiene		1.0	ug/g	1	96	70 130		
		naphthalene		0.9	ug/g	1	93	70 130		
		1,2,3-trichlorobenzene		1.0	ug/g	1	98	70 130		
		dibromofluoromethane SUR		93	%			78 114		
		toluene-D8 SUR		95	%			88 110		
		4-bromofluorobenzene SUR		108	%			86 115		
		a,a,a-trifluorotoluene SUR		97	%			70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCSD4307	dichlorodifluoromethane		0.6	ug/g	1	57 *	70 130	7	30
		chloromethane		0.8	ug/g	1	81	70 130	6	30
		vinyl chloride		0.5	ug/g	1	47 *	70 130	9	30
		bromomethane		1.2	ug/g	1	117	70 130	10	30
		chloroethane		0.9	ug/g	1	95	70 130	3	30
		trichlorofluoromethane		0.8	ug/g	1	83	70 130	1	30
		diethyl ether		1.0	ug/g	1	98	70 130	2	30
		acetone	<	2.5	ug/g	1	92		2	30
		1,1-dichloroethene		0.7	ug/g	1	74	70 130	3	30
		methylene chloride		0.9	ug/g	1	91	70 130	3	30
		carbon disulfide		0.8	ug/g	1	83	70 130	4	30
		methyl t-butyl ether (MTBE)		1.0	ug/g	1	100	70 130	5	30
		trans-1,2-dichloroethene		0.8	ug/g	1	81	70 130	3	30
		isopropyl ether (DIPE)		1.0	ug/g	1	101	70 130	3	30
		ethyl t-butyl ether (ETBE)		1.0	ug/g	1	99	70 130	3	30
		1,1-dichloroethane		0.9	ug/g	1	87	70 130	2	30
		t-butanol (TBA)		6.1	ug/g	5	122	70 130	4	30
		2-butanone (MEK)		1.0	ug/g	1	99	70 130	1	30
		2,2-dichloropropane		1.0	ug/g	1	105	70 130	2	30
		cis-1,2-dichloroethene		0.9	ug/g	1	94	70 130	1	30
		chloroform		1.0	ug/g	1	96	70 130	1	30
		bromochloromethane		0.9	ug/g	1	93	70 130	2	30
		tetrahydrofuran (THF)		1.0	ug/g	1	101	70 130	0	30
		1,1,1-trichloroethane		0.9	ug/g	1	90	70 130	3	30
		1,1-dichloropropene		1.0	ug/g	1	95	70 130	0	30
		t-amyl-methyl ether (TAME)		1.0	ug/g	1	99	70 130	6	30
		carbon tetrachloride		0.9	ug/g	1	87	70 130	4	30
		1,2-dichloroethane		1.0	ug/g	1	96	70 130	0	30
		benzene		1.0	ug/g	1	97	70 130	0	30
		trichloroethene		0.9	ug/g	1	92	70 130	1	30
		1,2-dichloropropane		1.0	ug/g	1	100	70 130	1	30
		bromodichloromethane		0.9	ug/g	1	94	70 130	4	30
		1,4-dioxane	<	2.5	ug/g	2	97	70 130	15	30
		dibromomethane		1.0	ug/g	1	100	70 130	2	30
		4-methyl-2-pentanone (MIBK)		1.0	ug/g	1	99	70 130	2	30
		cis-1,3-dichloropropene		1.0	ug/g	1	97	70 130	4	30
		toluene		1.0	ug/g	1	97	70 130	0	30
		trans-1,3-dichloropropene		1.1	ug/g	1	108	70 130	1	30
		2-hexanone		0.9	ug/g	1	94	70 130	3	30
		1,1,2-trichloroethane		1.0	ug/g	1	101	70 130	1	30
		1,3-dichloropropane		0.9	ug/g	1	92	70 130	4	30
		tetrachloroethene		1.0	ug/g	1	99	70 130	3	30
		dibromochloromethane		0.8	ug/g	1	79	70 130	3	30
		1,2-dibromoethane (EDB)		0.9	ug/g	1	89	70 130	2	30
		chlorobenzene		1.0	ug/g	1	101	70 130	1	30
		1,1,1,2-tetrachloroethane		0.9	ug/g	1	90	70 130	5	30
		ethylbenzene		0.9	ug/g	1	91	70 130	3	30
		m&p-xylenes		2.0	ug/g	2	102	70 130	2	30
		o-xylene		1.1	ug/g	1	108	70 130	1	30
		styrene		1.0	ug/g	1	95	70 130	1	30
		bromoform		1.0	ug/g	1	102	70 130	6	30

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCSD4307	isopropylbenzene		1.0	ug/g	1	100	70 130	1	30
		1,1,2,2-tetrachloroethane		1.1	ug/g	1	115	70 130	2	30
		1,2,3-trichloropropane		1.0	ug/g	1	105	70 130	3	30
		n-propylbenzene		1.1	ug/g	1	109	70 130	2	30
		bromobenzene		1.0	ug/g	1	103	70 130	1	30
		1,3,5-trimethylbenzene		1.0	ug/g	1	103	70 130	1	30
		2-chlorotoluene		1.1	ug/g	1	107	70 130	1	30
		4-chlorotoluene		1.1	ug/g	1	110	70 130	3	30
		tert-butylbenzene		1.1	ug/g	1	107	70 130	1	30
		1,2,4-trimethylbenzene		1.1	ug/g	1	108	70 130	3	30
		sec-butylbenzene		1.1	ug/g	1	111	70 130	1	30
		1,3-dichlorobenzene		1.2	ug/g	1	116	70 130	3	30
		4-isopropyltoluene		1.1	ug/g	1	112	70 130	1	30
		1,4-dichlorobenzene		1.0	ug/g	1	105	70 130	3	30
		1,2-dichlorobenzene		1.1	ug/g	1	106	70 130	4	30
		n-butylbenzene		1.1	ug/g	1	112	70 130	2	30
		1,2-dibromo-3-chloropropane		0.9	ug/g	1	85	70 130	4	30
		1,2,4-trichlorobenzene		0.9	ug/g	1	90	70 130	2	30
		hexachlorobutadiene		1.0	ug/g	1	103	70 130	8	30
		naphthalene		1.0	ug/g	1	99	70 130	6	30
		1,2,3-trichlorobenzene		1.0	ug/g	1	102	70 130	4	30
		dibromofluoromethane SUR		94	%			78 114		
		toluene-D8 SUR		96	%			88 110		
		4-bromofluorobenzene SUR		109	%			86 115		
		a,a,a-trifluorotoluene SUR		95	%			70 130		

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3540C8082A	BLK4305	PCB-1016		<	0.02	ug/g					
		PCB-1221		<	0.02	ug/g					
		PCB-1232		<	0.02	ug/g					
		PCB-1242		<	0.02	ug/g					
		PCB-1248		<	0.02	ug/g					
		PCB-1254		<	0.02	ug/g					
		PCB-1260		<	0.02	ug/g					
		tetrachloro-m-xylene SUR			51	%			30 150		
		decachlorobiphenyl SUR			62	%			30 150		
SW3540C8082A	LCS4305	PCB-1016			0.13	ug/g	0.2	64	40	140	
		PCB-1221		<	0.02	ug/g					
		PCB-1232		<	0.02	ug/g					
		PCB-1242		<	0.02	ug/g					
		PCB-1248		<	0.02	ug/g					
		PCB-1254		<	0.02	ug/g					
		PCB-1260			0.13	ug/g	0.2	64	40	140	
		tetrachloro-m-xylene SUR			48	%			30 150		
		decachlorobiphenyl SUR			57	%			30 150		
SW3540C8082A	LCSD4305	PCB-1016			0.18	ug/g	0.2	91	40	140	35 * 30
		PCB-1221		<	0.02	ug/g					
		PCB-1232		<	0.02	ug/g					
		PCB-1242		<	0.02	ug/g					
		PCB-1248		<	0.02	ug/g					
		PCB-1254		<	0.02	ug/g					
		PCB-1260			0.18	ug/g	0.2	90	40	140	34 * 30
		tetrachloro-m-xylene SUR			61	%			30 150		
		decachlorobiphenyl SUR			72	%			30 150		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	BLK4314	N-nitrosodimethylamine		<	0.2	ug/g				
		aniline		<	0.2	ug/g				
		phenol		<	0.2	ug/g				
		2-chlorophenol		<	0.5	ug/g				
		bis(2-chloroethyl)ether		<	0.2	ug/g				
		1,3-dichlorobenzene		<	0.2	ug/g				
		1,4-dichlorobenzene		<	0.2	ug/g				
		1,2-dichlorobenzene		<	0.2	ug/g				
		benzyl alcohol		<	0.2	ug/g				
		2-methylphenol		<	0.2	ug/g				
		bis(2-chloroisopropyl) ether		<	0.2	ug/g				
		hexachloroethane		<	0.2	ug/g				
		N-nitroso-di-N-propylamine		<	0.2	ug/g				
		4-methylphenol		<	0.2	ug/g				
		nitrobenzene		<	0.2	ug/g				
		isophorone		<	0.5	ug/g				
		2-nitrophenol		<	0.2	ug/g				
		2,4-dimethylphenol		<	0.2	ug/g				
		bis(2-chloroethoxy)methane		<	0.5	ug/g				
		2,4-dichlorophenol		<	0.5	ug/g				
		1,2,4-trichlorobenzene		<	0.5	ug/g				
		naphthalene		<	0.05	ug/g				
		benzoic acid		<	5.0	ug/g				
		4-chloroaniline		<	0.2	ug/g				
		hexachlorobutadiene		<	0.2	ug/g				
		4-chloro-3-methylphenol		<	0.2	ug/g				
		2-methylnaphthalene		<	0.05	ug/g				
		hexachlorocyclopentadiene		<	1.0	ug/g				
		2,4,6-trichlorophenol		<	0.2	ug/g				
		2,4,5-trichlorophenol		<	0.2	ug/g				
		2-chloronaphthalene		<	0.5	ug/g				
		2-nitroaniline		<	0.2	ug/g				
		acenaphthylene		<	0.05	ug/g				
		dimethylphthalate		<	0.5	ug/g				
		2,6-dinitrotoluene		<	0.2	ug/g				
		2,4-dinitrotoluene		<	0.2	ug/g				
		acenaphthene		<	0.05	ug/g				
		3-nitroaniline		<	0.2	ug/g				
		2,4-dinitrophenol		<	5.0	ug/g				
		dibenzofuran		<	0.05	ug/g				
		4-nitrophenol		<	1.0	ug/g				
		fluorene		<	0.05	ug/g				
		diethyl phthalate		<	0.5	ug/g				
		4-chlorophenyl phenyl ether		<	0.5	ug/g				
		4-nitroaniline		<	0.5	ug/g				
		4,6-dinitro-2-methylphenol		<	2.0	ug/g				
		azobenzene		<	0.2	ug/g				
		N-nitrosodiphenylamine		<	0.2	ug/g				
		4-bromophenyl phenyl ether		<	0.2	ug/g				
		hexachlorobenzene		<	0.2	ug/g				
		pentachlorophenol		<	1.0	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	BLK4314	phenanthrene		<	0.05	ug/g				
		anthracene		<	0.05	ug/g				
		carbazole		<	0.2	ug/g				
		di-n-butylphthalate		<	0.5	ug/g				
		fluoranthene		<	0.05	ug/g				
		benzidine		<	3.0	ug/g				
		pyrene		<	0.05	ug/g				
		butyl benzyl phthalate		<	0.5	ug/g				
		benzo(a)anthracene		<	0.05	ug/g				
		chrysene		<	0.05	ug/g				
		3,3'-dichlorobenzidine		<	3.0	ug/g				
		bis(2-ethylhexyl)phthalate		<	0.5	ug/g				
		di-n-octyl phthalate		<	0.2	ug/g				
		benzo(b)fluoranthene		<	0.05	ug/g				
		benzo(k)fluoranthene		<	0.05	ug/g				
		benzo(a)pyrene		<	0.02	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.05	ug/g				
		dibenzo(a,h)anthracene		<	0.05	ug/g				
		benzo(g,h,i)perylene		<	0.05	ug/g				
		2-fluorophenol SUR			43	%		21	100	
		phenol-D5 SUR			46	%		10	102	
		2,4,6-tribromophenol SUR			40	%		10	123	
		nitrobenzene-D5 SUR			38	%		35	114	
		2-fluorobiphenyl SUR			49	%		43	116	
		p-terphenyl-D14 SUR			66	%		33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCS4314	N-nitrosodimethylamine		<	0.2	ug/g				
		aniline		<	0.2	ug/g				
		phenol			5.6	ug/g	10	56	30	130
		2-chlorophenol			6.5	ug/g	10	65	30	130
		bis(2-chloroethyl)ether		<	0.2	ug/g				
		1,3-dichlorobenzene		<	0.2	ug/g				
		1,4-dichlorobenzene		<	0.2	ug/g				
		1,2-dichlorobenzene		<	0.2	ug/g				
		benzyl alcohol		<	0.2	ug/g				
		2-methylphenol		<	0.2	ug/g				
		bis(2-chloroisopropyl) ether		<	0.2	ug/g				
		hexachloroethane		<	0.2	ug/g				
		N-nitroso-di-N-propylamine		<	0.2	ug/g				
		4-methylphenol		<	0.2	ug/g				
		nitrobenzene		<	0.2	ug/g				
		isophorone		<	0.5	ug/g				
		2-nitrophenol		<	0.2	ug/g				
		2,4-dimethylphenol		<	0.2	ug/g				
		bis(2-chloroethoxy)methane		<	0.5	ug/g				
		2,4-dichlorophenol		<	0.5	ug/g				
		1,2,4-trichlorobenzene		<	0.5	ug/g				
		naphthalene			2.4	ug/g	4	59	40	140
		benzoic acid		<	5.0	ug/g				
		4-chloroaniline		<	0.2	ug/g				
		hexachlorobutadiene		<	0.2	ug/g				
		4-chloro-3-methylphenol			6.6	ug/g	10	66	30	130
		2-methylnaphthalene			2.45	ug/g	4	61	40	140
		hexachlorocyclopentadiene		<	1.0	ug/g				
		2,4,6-trichlorophenol		<	0.2	ug/g				
		2,4,5-trichlorophenol		<	0.2	ug/g				
		2-chloronaphthalene		<	0.5	ug/g				
		2-nitroaniline		<	0.2	ug/g				
		acenaphthylene			2.9	ug/g	4	72	40	140
		dimethylphthalate		<	0.5	ug/g				
		2,6-dinitrotoluene		<	0.2	ug/g				
		2,4-dinitrotoluene		<	0.2	ug/g				
		acenaphthene			2.8	ug/g				
		3-nitroaniline		<	0.2	ug/g				
		2,4-dinitrophenol		<	5.0	ug/g				
		dibenzofuran		<	0.05	ug/g				
		4-nitrophenol			6.4	ug/g	10	64	30	130
		fluorene			3.0	ug/g	4	76	40	140
		diethyl phthalate		<	0.5	ug/g				
		4-chlorophenyl phenyl ether		<	0.5	ug/g				
		4-nitroaniline		<	0.5	ug/g				
		4,6-dinitro-2-methylphenol		<	2.0	ug/g				
		azobenzene		<	0.2	ug/g				
		N-nitrosodiphenylamine		<	0.2	ug/g				
		4-bromophenyl phenyl ether		<	0.2	ug/g				
		hexachlorobenzene		<	0.2	ug/g				
		pentachlorophenol			4.3	ug/g	10	43	30	130

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCS4314	phenanthrene		2.9	ug/g	4	72	40 140		
		anthracene		3.0	ug/g	4	75	40 140		
		carbazole	<	0.2	ug/g					
		di-n-butylphthalate	<	0.5	ug/g					
		fluoranthene		3.3	ug/g	4	81	40 140		
		benzidine	<	3.0	ug/g					
		pyrene		3.5	ug/g	4	87	40 140		
		butyl benzyl phthalate	<	0.5	ug/g					
		benzo(a)anthracene		3.3	ug/g	4	82	40 140		
		chrysene		3.3	ug/g	4	82	40 140		
		3,3'-dichlorobenzidine	<	3.0	ug/g					
		bis(2-ethylhexyl)phthalate	<	0.5	ug/g					
		di-n-octyl phthalate	<	0.2	ug/g					
		benzo(b)fluoranthene		3.4	ug/g	4	85	40 140		
		benzo(k)fluoranthene		3.0	ug/g	4	76	40 140		
		benzo(a)pyrene		3.3	ug/g	4	82	40 140		
		indeno(1,2,3-cd)pyrene		3.2	ug/g	4	81	40 140		
		dibenzo(a,h)anthracene		3.2	ug/g	4	79	40 140		
		benzo(g,h,i)perylene		3.2	ug/g	4	79	40 140		
		2-fluorophenol SUR		53	%			21 100		
		phenol-D5 SUR		57	%			10 102		
		2,4,6-tribromophenol SUR		52	%			10 123		
		nitrobenzene-D5 SUR		50	%			35 114		
		2-fluorobiphenyl SUR		64	%			43 116		
		p-terphenyl-D14 SUR		71	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3546/8270D	LCS4314	N-nitrosodimethylamine		<	0.2	ug/g						
		aniline		<	0.2	ug/g						
		phenol			5.4	ug/g	10	54	30	130	4	30
		2-chlorophenol			5.9	ug/g	10	59	30	130	9	30
		bis(2-chloroethyl)ether		<	0.2	ug/g						
		1,3-dichlorobenzene		<	0.2	ug/g						
		1,4-dichlorobenzene		<	0.2	ug/g						
		1,2-dichlorobenzene		<	0.2	ug/g						
		benzyl alcohol		<	0.2	ug/g						
		2-methylphenol		<	0.2	ug/g						
		bis(2-chloroisopropyl) ether		<	0.2	ug/g						
		hexachloroethane		<	0.2	ug/g						
		N-nitroso-di-N-propylamine		<	0.2	ug/g						
		4-methylphenol		<	0.2	ug/g						
		nitrobenzene		<	0.2	ug/g						
		isophorone		<	0.5	ug/g						
		2-nitrophenol		<	0.2	ug/g						
		2,4-dimethylphenol		<	0.2	ug/g						
		bis(2-chloroethoxy)methane		<	0.5	ug/g						
		2,4-dichlorophenol		<	0.5	ug/g						
		1,2,4-trichlorobenzene		<	0.5	ug/g						
		naphthalene			2.2	ug/g	4	56	40	140	6	30
		benzoic acid		<	5.0	ug/g						
		4-chloroaniline		<	0.2	ug/g						
		hexachlorobutadiene		<	0.2	ug/g						
		4-chloro-3-methylphenol			6.4	ug/g	10	64	30	130	2	30
		2-methylnaphthalene			2.32	ug/g	4	58	40	140	5	30
		hexachlorocyclopentadiene		<	1.0	ug/g						
		2,4,6-trichlorophenol		<	0.2	ug/g						
		2,4,5-trichlorophenol		<	0.2	ug/g						
		2-chloronaphthalene		<	0.5	ug/g						
		2-nitroaniline		<	0.2	ug/g						
		acenaphthylene			2.9	ug/g	4	72	40	140	0	30
		dimethylphthalate		<	0.5	ug/g						
		2,6-dinitrotoluene		<	0.2	ug/g						
		2,4-dinitrotoluene		<	0.2	ug/g						
		acenaphthene			2.7	ug/g						
		3-nitroaniline		<	0.2	ug/g						
		2,4-dinitrophenol		<	5.0	ug/g						
		dibenzofuran		<	0.05	ug/g						
		4-nitrophenol			6.2	ug/g	10	62	30	130	3	30
		fluorene			3.0	ug/g	4	75	40	140	1	30
		diethyl phthalate		<	0.5	ug/g						
		4-chlorophenyl phenyl ether		<	0.5	ug/g						
		4-nitroaniline		<	0.5	ug/g						
		4,6-dinitro-2-methylphenol		<	2.0	ug/g						
		azobenzene		<	0.2	ug/g						
		N-nitrosodiphenylamine		<	0.2	ug/g						
		4-bromophenyl phenyl ether		<	0.2	ug/g						
		hexachlorobenzene		<	0.2	ug/g						
		pentachlorophenol			3.6	ug/g	10	36	30	130	17	30

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit			
SW3546/8270D	LCSD4314	phenanthrene		2.9	ug/g	4	72	40 140	0	30			
		anthracene		3.0	ug/g	4	76	40 140	0	30			
		carbazole		<	0.2	ug/g							
		di-n-butylphthalate		<	0.5	ug/g							
		fluoranthene			3.4	ug/g	4	84	40 140	3	30		
		benzidine		<	3.0	ug/g							
		pyrene			3.5	ug/g	4	88	40 140	2	30		
		butyl benzyl phthalate		<	0.5	ug/g							
		benzo(a)anthracene			3.4	ug/g	4	84	40 140	2	30		
		chrysene			3.4	ug/g	4	84	40 140	2	30		
		3,3'-dichlorobenzidine		<	3.0	ug/g							
		bis(2-ethylhexyl)phthalate		<	0.5	ug/g							
		di-n-octyl phthalate		<	0.2	ug/g							
		benzo(b)fluoranthene			3.5	ug/g	4	87	40 140	3	30		
		benzo(k)fluoranthene			3.2	ug/g	4	80	40 140	5	30		
		benzo(a)pyrene			3.3	ug/g	4	84	40 140	1	30		
		indeno(1,2,3-cd)pyrene			3.2	ug/g	4	81	40 140	0	30		
		dibenzo(a,h)anthracene			3.2	ug/g	4	79	40 140	0	30		
		benzo(g,h,i)perylene			3.2	ug/g	4	80	40 140	1	30		
		2-fluorophenol SUR			49	%				21 100			
		phenol-D5 SUR			51	%				10 102			
		2,4,6-tribromophenol SUR			49	%				10 123			
		nitrobenzene-D5 SUR			46	%				35 114			
		2-fluorobiphenyl SUR			63	%				43 116			
		p-terphenyl-D14 SUR			70	%				33 141			
		SW3550B8100m	BLK4312	TPH C10-C36		<	200	ug/g					
				2-fluorobiphenyl SUR			68	%			40 140		
				o-terphenyl SUR			88	%			40 140		
SW3550B8100m	DUP4312	2-fluorobiphenyl SUR	21913-001		115	%			40 140				
		o-terphenyl SUR	21913-001		104	%			40 140				
SW3550B8100m	LCS4312	TPH C10-C36			1500	ug/g	2500	59	40 140				
		2-fluorobiphenyl SUR			73	%			40 140				
		o-terphenyl SUR			89	%			40 140				
SW3550B8100m	MS4312	2-fluorobiphenyl SUR	21913-001		78	%			40 140				
		o-terphenyl SUR	21913-001		108	%			40 140				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3550B8270D	BLK4315	naphthalene		<	0.50	ug/g						
		2-methylnaphthalene		<	0.50	ug/g						
		acenaphthylene		<	0.50	ug/g						
		acenaphthene		<	0.50	ug/g						
		dibenzofuran		<	0.50	ug/g						
		fluorene		<	0.50	ug/g						
		phenanthrene		<	0.50	ug/g						
		anthracene		<	0.50	ug/g						
		fluoranthene		<	0.50	ug/g						
		pyrene		<	0.50	ug/g						
		benzo(a)anthracene		<	0.50	ug/g						
		chrysene		<	0.50	ug/g						
		benzo(b)fluoranthene		<	0.50	ug/g						
		benzo(k)fluoranthene		<	0.50	ug/g						
		benzo(a)pyrene		<	0.50	ug/g						
		indeno(1,2,3-cd)pyrene		<	0.50	ug/g						
		dibenzo(a,h)anthracene		<	0.50	ug/g						
		benzo(g,h,i)perylene		<	0.50	ug/g						
		2-fluorobiphenyl	SUR			78	%			43	116	
		o-terphenyl	SUR			79	%			33	141	
SW3550B8270D	LCS4315	naphthalene			3.1	ug/g	4	78	40	140		
		2-methylnaphthalene			3.6	ug/g	4	90	40	140		
		acenaphthylene			3.5	ug/g	4	88	40	140		
		acenaphthene			3.3	ug/g	4	82	40	140		
		dibenzofuran		<	0.50	ug/g						
		fluorene			3.4	ug/g	4	84	40	140		
		phenanthrene			3.2	ug/g	4	81	40	140		
		anthracene			3.4	ug/g	4	84	40	140		
		fluoranthene			3.3	ug/g	4	83	40	140		
		pyrene			3.8	ug/g	4	95	40	140		
		benzo(a)anthracene			3.6	ug/g	4	90	40	140		
		chrysene			3.8	ug/g	4	95	40	140		
		benzo(b)fluoranthene			3.9	ug/g	4	96	40	140		
		benzo(k)fluoranthene			3.2	ug/g	4	79	40	140		
		benzo(a)pyrene			3.3	ug/g	4	82	40	140		
		indeno(1,2,3-cd)pyrene			3.1	ug/g	4	78	40	140		
		dibenzo(a,h)anthracene			2.8	ug/g	4	71	40	140		
		benzo(g,h,i)perylene			3.3	ug/g	4	83	40	140		
		2-fluorobiphenyl	SUR			76	%			43	116	
		o-terphenyl	SUR			78	%			33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6010C	BLK4310	Silver		<	0.25	ug/g				
		Arsenic		<	0.50	ug/g				
		Barium		<	2.5	ug/g				
		Cadmium		<	0.20	ug/g				
		Chromium		<	2.5	ug/g				
		Copper		<	2.5	ug/g				
		Nickel		<	2.5	ug/g				
		Lead		<	0.50	ug/g				
		Selenium		<	2.5	ug/g				
		Zinc		<	2.5	ug/g				
SW3051A6010C	CRM4310	Silver		40	ug/g	38		25.1 51.9		
		Arsenic		390	ug/g	400		292 508		
		Barium		24	ug/g	25		0 51.3		
		Cadmium		16	ug/g	15		8.71 22		
		Chromium		13	ug/g	14		2.45 24.7		
		Copper		710	ug/g	730		592 866		
		Nickel		15	ug/g	17		6.2 27.5		
		Lead		5000	ug/g	5100		3753 6469		
		Selenium		6.2	ug/g	6.6		0 18.4		
		Zinc		3000	ug/g	3000		2447 3575		
SW3051A6010C	CRMD4310	Silver		41	ug/g	38		25.1 51.9	2	20
		Arsenic		400	ug/g	400		292 508	2	20
		Barium		26	ug/g	25		0 51.3	7	20
		Cadmium		16	ug/g	15		8.71 22	1	20
		Chromium		15	ug/g	14		2.45 24.7	15	20
		Copper		740	ug/g	730		592 866	4	20
		Nickel		16	ug/g	17		6.2 27.5	6	20
		Lead		5300	ug/g	5100		3753 6469	5	20
		Selenium		6.2	ug/g	6.6		0 18.4	1	20
		Zinc		3000	ug/g	3000		2447 3575	0	20
SW3051A6010C	DUP4310	Lead	21906-011	ERROR	ug/g				9	35
SW7471B	BLK4309	Mercury		<	0.02	ug/g				
SW7471B	CRM4309	Mercury		1.3	ug/g	1.1		0.49 1.76		
SW7471B	CRMD4309	Mercury		1.3	ug/g	1.1		0.49 1.76	1	20
SW7471B	DUP4309	Mercury	21894-009	1.4	ug/g				26	35
SW7471B	MS4309	Mercury	21894-009	1.1	ug/g	2.04	1	75 125		
SW7471B	MS4309	Mercury	21907-003	0.20	ug/g	0.172	113	75 125		

Company Name:

REDENE

Company Address:

776 Main St Westboro MA 01581

Report To:

RICK VAHDEMBERG

Phone #:

207-828-1272

Invoice To:

Same

Project Name:

POLYCLIN

Project #:

1000086

Project Location:

RD MA ME VT

Protocol:

RCRA SDWA NPDES MOP ~~NPDES~~ OTHER

Reporting Limits:

QAPP GW-1 S-1 EPA DW Other

Quote #

1000086

PO # 1000086

NH GREE/ODD Fund Pricing

ANALYSIS REQUEST

Lab Sample ID	Field ID	# CONTAINERS	Matrix			Preservation Method						Sampling		
			WATER	SOLID	OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)	DATE	TIME	SAMPLER*
02	CA-3 2-4	1	X								6/19/11	1259	2880	
03	CA-2 6-8	2	X								6/19/11	1255	2880	
04	CA-CC-1	1	X								6/13/11	1200	2880	
05	CA-CC-DUP	1	X								6/30/11	1200	2880	
06	CA-CC-2	1	X								6/30/11	1215	2880	
07	CA-S 12-H	2	X								6/29/11	1710	2880	
08	CA-1 26-28	2	X								6/29/11	1557	2880	
09	CA-4 10-12	2	X								6/29/11	1744	2880	
10	CA-TP-9 C'	2	X								6/30/11	1545	2880	
11	CA-TP-7A 10-12	2	X								6/30/11	1736	2880	
11	CA-TP-12 1'	2	X								6/30/11	1315	2880	

<input type="checkbox"/> VOC 8260	<input checked="" type="checkbox"/> VOC 8260 NHDES	<input type="checkbox"/> VOC 8260 MADEP
<input type="checkbox"/> VOC 624	<input type="checkbox"/> VOC BTEX	<input type="checkbox"/> MIBE, only
<input type="checkbox"/> VOC 8021VT	<input type="checkbox"/> VPH MADEP	<input type="checkbox"/> MEGRO
<input type="checkbox"/> GRO 8015	<input type="checkbox"/> VOC 524.2	<input type="checkbox"/> VOC 524.2 NH List
<input type="checkbox"/> TPH	<input type="checkbox"/> DRO 8015	<input type="checkbox"/> MEDRO
<input type="checkbox"/> EPH MADEP	<input type="checkbox"/> TPH Fingerprint	<input type="checkbox"/> 8270PAH
<input type="checkbox"/> 8270ABN	<input type="checkbox"/> 625	<input type="checkbox"/> EDB 504.1
<input checked="" type="checkbox"/> 8082 PCB	<input type="checkbox"/> 8081 Pesticides	<input type="checkbox"/> 608 Pest/PCB
<input type="checkbox"/> O&G 1664	<input type="checkbox"/> Mineral O&G SM5520F	<input type="checkbox"/> pH
<input type="checkbox"/> BOD	<input type="checkbox"/> Conductivity	<input type="checkbox"/> Turbidity
<input type="checkbox"/> TSS	<input type="checkbox"/> TDS	<input type="checkbox"/> TS
<input type="checkbox"/> TVS	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> RCRA Metals
<input type="checkbox"/> Priority Pollutant Metals	<input type="checkbox"/> TAL Metals	<input checked="" type="checkbox"/> Total Metals-list: Zn, Cu, Ni
<input type="checkbox"/> Dissolved Metals-list:	<input type="checkbox"/> Ammonia	<input type="checkbox"/> COD
<input type="checkbox"/> TKN	<input type="checkbox"/> TN	<input type="checkbox"/> TON
<input type="checkbox"/> T-Phosphorus	<input type="checkbox"/> Phenols	<input type="checkbox"/> Bacteria P/A
<input type="checkbox"/> Bacteria MPN	<input type="checkbox"/> Cyanide	<input type="checkbox"/> Sulfide
<input type="checkbox"/> Nitrate + Nitrite	<input type="checkbox"/> Ortho P	<input type="checkbox"/> Nitrate
<input type="checkbox"/> Nitrite	<input type="checkbox"/> Chloride	<input type="checkbox"/> Sulfate
<input type="checkbox"/> Bromide	<input type="checkbox"/> Fluoride	<input type="checkbox"/> Corrosivity
<input type="checkbox"/> Reactive CN	<input type="checkbox"/> Reactive S-	<input type="checkbox"/> Ignitibility/FP
<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC	<input type="checkbox"/> TCLP SVOC
<input type="checkbox"/> TCLP Pesticide	<input type="checkbox"/> Subcontract: TOC	<input type="checkbox"/> Grain Size
<input type="checkbox"/> TCLP Herbicides	<input type="checkbox"/> 8270 PAH	<input type="checkbox"/> Pb, total

TAT REQUESTED: Priority (24 hr)** Expedited (48 hr)** Standard (10 Business days)

Date Needed: 7/12/11

* See www.reslabs.com for sample acceptance policy and current accreditation lists.

SPECIAL INSTRUCTIONS: REPORTING INSTRUCTIONS: PDF (e-mail address) **RVAHDEMBERG@REDENE.COM**

HARD COPY REQUIRED FAX (FAX#) OTHER (specify)

RECEIVED ON ICE: YES NO TEMPERATURE: °C

Relinquished by:	Date	Time	Received by:	Date	Time
<i>[Signature]</i>	7/11/11	1021	<i>[Signature]</i>		
Relinquished by:	Date	Time	Received by:	Date	Time
Relinquished by:	Date	Time	Received by:	Date	Time



Absolute Resource *associates*

124 Heritage Avenue #10 Portsmouth, NH 03801

Jonathan O'Donnell
CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: None
Job ID: 21917
Date Received: 7/6/11

Project: POLYCLAD 10001086

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

Sue Sylvester
Principal, General Manager

Date of Approval: 7/14/2011
Total number of pages: 38

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-1	Water	7/5/2011 15:15	21917-001	VOCs in water by 8260 Petro & Haz Waste
CA-2	Water	7/5/2011 14:20	21917-002	Acid & Base/Neutral Extractables in water by 8270 Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Barium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Copper in water by 6010 Mercury in water by 7470 Nickel in water by 6010 Lead in water by 6010 Selenium in water by 6010 Zinc in water by 6010 VOCs in water by 8260 Petro & Haz Waste
CA-3	Water	7/5/2011 15:30	21917-003	Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Barium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Mercury in water by 7470 Lead in water by 6010 Selenium in water by 6010 VOCs in water by 8260 Petro & Haz Waste
MW-2	Water	7/5/2011 13:20	21917-004	Water Digestion for ICP Analysis Lead in water by 6010 VOCs in water by 8260 Petro & Haz Waste
SS-3	Solid	7/5/2011 12:00	21917-005	Acid & Base/Neutral Extractables in solid by 8270 Percent Dry Matter for Sample Calc by SM2540B,G

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-001

Sample ID: CA-1

Matrix: Water

Sampled: 7/5/11 15:15

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		
		Limit	Units	Factor	Analyst		Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
isopropyl ether (DIPE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
ethyl t-butyl ether (ETBE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
t-butanol (TBA)	< 30	30	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
t-amyl-methyl ether (TAME)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
bromodichloromethane	< 0.6	0.6	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,4-dioxane	< 50	50	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-001

Sample ID: CA-1

Matrix: Water

Sampled: 7/5/11 15:15

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		
		Limit	Units	Factor	Analyst		Batch	Date	Time
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
m&p-xylenes	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
o-xylene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
styrene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
bromoform	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
isopropylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
n-propylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
bromobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,3,5-trimethylbenzene	4	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
2-chlorotoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
4-chlorotoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
tert-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2,4-trimethylbenzene	9	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
sec-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
n-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,3,5-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
naphthalene	< 5	5	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
Surrogate Recovery		Limits							
dibromofluoromethane SUR	88	78-114	%	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
toluene-D8 SUR	92	88-110	%	1	LMM	1101353	7/7/11	15:21	SW5030B8260B
4-bromofluorobenzene SUR	93	86-115	%	1	LMM	1101353	7/7/11	15:21	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-002

Sample ID: CA-2

Matrix: Water

Sampled: 7/5/11 14:20

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		
		Limit	Units	Factor	Analyst		Batch	Date	Time
dichlorodifluoromethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
chloromethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
vinyl chloride	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
bromomethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
chloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
trichlorofluoromethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
diethyl ether	< 25	25	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
acetone	< 250	250	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1-dichloroethene	< 5	5	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
methylene chloride	< 25	25	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
carbon disulfide	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
methyl t-butyl ether (MTBE)	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
trans-1,2-dichloroethene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
isopropyl ether (DIPE)	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
ethyl t-butyl ether (ETBE)	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1-dichloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
t-butanol (TBA)	< 150	150	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
2-butanone (MEK)	< 50	50	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
2,2-dichloropropane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
cis-1,2-dichloroethene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
chloroform	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
bromochloromethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
tetrahydrofuran (THF)	< 50	50	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1,1-trichloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1-dichloropropene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
t-amyl-methyl ether (TAME)	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
carbon tetrachloride	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2-dichloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
benzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
trichloroethene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2-dichloropropane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
bromodichloromethane	< 3.0	3.0	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,4-dioxane	< 250	250	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
dibromomethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 50	50	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
cis-1,3-dichloropropene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
toluene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
trans-1,3-dichloropropene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
2-hexanone	< 50	50	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1,2-trichloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,3-dichloropropane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
tetrachloroethene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
dibromochloromethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-002

Sample ID: CA-2

Matrix: Water

Sampled: 7/5/11 14:20

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		Reference
		Limit	Units	Factor	Analyst		Batch	Date	
1,2-dibromoethane (EDB)	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
chlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1,1,2-tetrachloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
ethylbenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
m&p-xylenes	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
o-xylene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
styrene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
bromoform	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
isopropylbenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,1,2,2-tetrachloroethane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2,3-trichloropropane	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
n-propylbenzene	29	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
bromobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,3,5-trimethylbenzene	320	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
2-chlorotoluene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
4-chlorotoluene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
tert-butylbenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2,4-trimethylbenzene	710	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
sec-butylbenzene	43	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,3-dichlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
4-isopropyltoluene	100	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,4-dichlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2-dichlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
n-butylbenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2,4-trichlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,3,5-trichlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
hexachlorobutadiene	< 2.5	2.5	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
naphthalene	< 25	25	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
1,2,3-trichlorobenzene	< 10	10	ug/L	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
Surrogate Recovery		Limits							
dibromofluoromethane SUR	85	78-114	%	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
toluene-D8 SUR	92	88-110	%	5	LMM	1101353	7/7/11	23:16	SW5030B8260B
4-bromofluorobenzene SUR	99	86-115	%	5	LMM	1101353	7/7/11	23:16	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-003

Sample ID: CA-3

Matrix: Water

Sampled: 7/5/11 15:30

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		
		Limit	Units	Factor	Analyst		Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
isopropyl ether (DIPE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
ethyl t-butyl ether (ETBE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
t-butanol (TBA)	< 30	30	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
t-amyl-methyl ether (TAME)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
bromodichloromethane	< 0.6	0.6	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,4-dioxane	< 50	50	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-003

Sample ID: CA-3

Matrix: Water

Sampled: 7/5/11 15:30

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		
		Limit	Units	Factor	Analyst		Batch	Date	Time
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
m&p-xylenes	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
o-xylene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
styrene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
bromoform	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
isopropylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
n-propylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
bromobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
2-chlorotoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
4-chlorotoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
tert-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
sec-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
n-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,3,5-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
naphthalene	< 5	5	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
Surrogate Recovery		Limits							
dibromofluoromethane SUR	90	78-114	%	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
toluene-D8 SUR	93	88-110	%	1	LMM	1101353	7/7/11	15:52	SW5030B8260B
4-bromofluorobenzene SUR	91	86-115	%	1	LMM	1101353	7/7/11	15:52	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-004

Sample ID: MW-2

Matrix: Water

Sampled: 7/5/11 13:20

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis		
		Limit	Units	Factor	Analyst		Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
isopropyl ether (DIPE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
ethyl t-butyl ether (ETBE)	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
t-butanol (TBA)	< 30	30	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
t-amyl-methyl ether (TAME)	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
bromodichloromethane	< 0.6	0.6	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,4-dioxane	< 50	50	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-004

Sample ID: MW-2

Matrix: Water

Sampled: 7/5/11 13:20

Parameter	Result	Quant		Instr Dil'n		Prep Date	Analysis			Reference
		Limit	Units	Factor	Analyst		Batch	Date	Time	
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
chlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
ethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
m&p-xylenes	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
o-xylene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
styrene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
bromoform	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
isopropylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
n-propylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
bromobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
2-chlorotoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
4-chlorotoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
tert-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
sec-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
n-butylbenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,3,5-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
naphthalene	< 5	5	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
Surrogate Recovery		Limits								
dibromofluoromethane SUR	92	78-114	%	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
toluene-D8 SUR	96	88-110	%	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	
4-bromofluorobenzene SUR	93	86-115	%	1	LMM	1101353	7/7/11	16:24	SW5030B8260B	

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-002

Sample ID: CA-2

Matrix: Water

Sampled: 7/5/11 14:20

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
N-nitrosodimethylamine	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
aniline	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
phenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
2-chlorophenol	< 5	5	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
bis(2-chloroethyl)ether	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
1,3-dichlorobenzene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
1,4-dichlorobenzene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
1,2-dichlorobenzene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzyl alcohol	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2-methylphenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
bis(2-chloroisopropyl) ether	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
hexachloroethane	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
N-nitroso-di-N-propylamine	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
4-methylphenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
nitrobenzene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
isophorone	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2-nitrophenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
2,4-dimethylphenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
bis(2-chloroethoxy)methane	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2,4-dichlorophenol	< 5	5	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
1,2,4-trichlorobenzene	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
naphthalene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzoic acid	< 50	50	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
4-chloroaniline	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
hexachlorobutadiene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
4-chloro-3-methylphenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
2-methylnaphthalene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
hexachlorocyclopentadiene	< 100	100	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2,4,6-trichlorophenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
2,4,5-trichlorophenol	< 2	2	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
2-chloronaphthalene	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2-nitroaniline	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
acenaphthylene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
dimethylphthalate	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2,6-dinitrotoluene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2,4-dinitrotoluene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
acenaphthene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
3-nitroaniline	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2,4-dinitrophenol	< 50	50	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
dibenzofuran	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
4-nitrophenol	< 10	10	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
fluorene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
diethyl phthalate	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-002

Sample ID: CA-2

Matrix: Water

Sampled: 7/5/11 14:20

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
4-nitroaniline	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
4,6-dinitro-2-methylphenol	< 20	20	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
azobenzene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
N-nitrosodiphenylamine	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
4-bromophenyl phenyl ether	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
hexachlorobenzene	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
pentachlorophenol	< 10	10	ug/L	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
phenanthrene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
anthracene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
carbazole	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
di-n-butylphthalate	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
fluoranthene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzidine	< 300	300	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
pyrene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
butyl benzyl phthalate	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzo(a)anthracene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
chrysene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
3,3'-dichlorobenzidine	< 300	300	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
bis(2-ethylhexyl)phthalate	< 50	50	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
di-n-octyl phthalate	< 20	20	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzo(b)fluoranthene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzo(k)fluoranthene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzo(a)pyrene	< 2.0	2.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
indeno(1,2,3-cd)pyrene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
dibenzo(a,h)anthracene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
benzo(g,h,i)perylene	< 5.0	5.0	ug/L	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
Surrogate Recovery		Limits								
2-fluorophenol SUR	31	21-100	%	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
phenol-D5 SUR	20	10-102	%	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
2,4,6-tribromophenol SUR	70	10-123	%	1	AJD	7/8/11	4327	7/9/11	17:20	SW3510C8270D
nitrobenzene-D5 SUR	69	35-114	%	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
2-fluorobiphenyl SUR	72	43-116	%	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D
p-terphenyl-D14 SUR	72	33-141	%	10	AJD	7/8/11	4327	7/11/11	12:10	SW3510C8270D

Note: Dilution was necessary due to the high concentrations of non-target compounds.

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-005

Sample ID: SS-3

Matrix: Solid

Percent Dry: 85.2% Results expressed on a dry weight basis.

Sampled: 7/5/11 12:00

Parameter	Result	Limit	Units	Instr Dil'n		Prep		Analysis		
				Factor	Analyst	Date	Batch	Date	Time	Reference
N-nitrosodimethylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
aniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
phenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-chlorophenol	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
bis(2-chloroethyl)ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
1,3-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
1,4-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
1,2-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzyl alcohol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
bis(2-chloroisopropyl) ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
hexachloroethane	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
N-nitroso-di-N-propylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
nitrobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
isophorone	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-nitrophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4-dimethylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
bis(2-chloroethoxy)methane	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4-dichlorophenol	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
1,2,4-trichlorobenzene	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
naphthalene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzoic acid	< 6	6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4-chloroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
hexachlorobutadiene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4-chloro-3-methylphenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-methylnaphthalene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
hexachlorocyclopentadiene	< 1	1	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4,6-trichlorophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4,5-trichlorophenol	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-chloronaphthalene	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-nitroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
acenaphthylene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
dimethylphthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,6-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
acenaphthene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
3-nitroaniline	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4-dinitrophenol	< 6	6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
dibenzofuran	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4-nitrophenol	< 2	2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
fluorene	< 0.06	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
diethyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-005

Sample ID: SS-3

Matrix: Solid

Percent Dry: 85.2% Results expressed on a dry weight basis.

Sampled: 7/5/11 12:00

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4-nitroaniline	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4,6-dinitro-2-methylphenol	< 2	2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
azobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
N-nitrosodiphenylamine	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
4-bromophenyl phenyl ether	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
hexachlorobenzene	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
pentachlorophenol	< 1	1	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
phenanthrene	0.60	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
anthracene	0.12	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
carbazole	< 0.2	0.2	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
di-n-butylphthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
fluoranthene	1.5	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzidine	< 3	3	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
pyrene	1.3	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
butyl benzyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzo(a)anthracene	0.69	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
chrysene	0.80	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
3,3'-dichlorobenzidine	< 3	3	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
bis(2-ethylhexyl)phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
di-n-octyl phthalate	< 0.6	0.6	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzo(b)fluoranthene	0.82	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzo(k)fluoranthene	0.68	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzo(a)pyrene	0.71	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
indeno(1,2,3-cd)pyrene	0.34	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
dibenzo(a,h)anthracene	0.14	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
benzo(g,h,i)perylene	0.36	0.06	ug/g	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
Surrogate Recovery		Limits								
2-fluorophenol SUR	55	21-100	%	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
phenol-D5 SUR	57	10-102	%	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2,4,6-tribromophenol SUR	54	10-123	%	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
nitrobenzene-D5 SUR	44	35-114	%	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
2-fluorobiphenyl SUR	57	43-116	%	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D
p-terphenyl-D14 SUR	55	33-141	%	1	AJD	7/6/11	4314	7/9/11	18:35	SW3546/8270D

Project ID: POLYCLAD 10001086

Job ID: 21917

Sample#: 21917-002

Sample ID: CA-2

Matrix: Water

Sampled: 7/5/11 14:20

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	0.021	0.008	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Barium	0.13	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Chromium	0.13	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Copper	0.23	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Lead	0.045	0.008	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	BJS	7/12/11	4337	7/12/11	12:52	SW7470A
Nickel	< 0.05	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Selenium	< 0.05	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C
Zinc	0.12	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:18	SW3005A6010C

Sample#: 21917-003

Sample ID: CA-3

Matrix: Water

Sampled: 7/5/11 15:30

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Arsenic	< 0.008	0.008	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C
Barium	< 0.05	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	BJS	7/12/11	4337	7/12/11	12:49	SW7470A
Selenium	< 0.05	0.05	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	BJS	7/7/11	4318	7/7/11	17:25	SW3005A6010C

Sample#: 21917-004

Sample ID: MW-2

Matrix: Water

Sampled: 7/5/11 13:20

Parameter	Result	Quant		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Lead	0.008	0.008	mg/L	1	BJS	7/7/11	4318	7/7/11	17:32	SW3005A6010C

Quality Control Report



124 Heritage Avenue Unit 10
Portsmouth, NH 03801

www.absoluteresourceassociates.com



Case Narrative

Lab # 21917

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 3 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

VOC: The compound, hexachlorobutadiene, was detected in the BLK1101353 at 0.5ug/L. There is no impact to the data as this analyte was not detected in the associated field samples.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

VOC: The LCS1101353 did not meet the acceptance criteria for dichlorodifluoromethane and dibromochloromethane. These compounds are known to be problematic in the method. The LCSD1101353 did not meet the acceptance criteria for bromomethane. This compound showed high recovery. There is no impact to the data as this analyte was not detected in the associated samples.

ABN: The LCS/D4327 did not meet the acceptance criteria for phenol and 4-nitrophenol. These compounds are known to be problematic in the method.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

ABN: The Base/Neutral portion of the sample extract for 21917-002 required a re-analysis at a dilution due to interferences caused by high concentrations of non-target compounds present in the sample chromatogram.

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	BLK1101353	dichlorodifluoromethane		<	2	ug/L				
		chloromethane		<	2	ug/L				
		vinyl chloride		<	2	ug/L				
		bromomethane		<	2	ug/L				
		chloroethane		<	2	ug/L				
		trichlorofluoromethane		<	2	ug/L				
		diethyl ether		<	10	ug/L				
		acetone		<	50	ug/L				
		1,1-dichloroethene		<	1	ug/L				
		methylene chloride		<	5	ug/L				
		carbon disulfide		<	2	ug/L				
		methyl t-butyl ether (MTBE)		<	2	ug/L				
		trans-1,2-dichloroethene		<	2	ug/L				
		isopropyl ether (DIPE)		<	2	ug/L				
		ethyl t-butyl ether (ETBE)		<	2	ug/L				
		1,1-dichloroethane		<	2	ug/L				
		t-butanol (TBA)		<	30	ug/L				
		2-butanone (MEK)		<	10	ug/L				
		2,2-dichloropropane		<	2	ug/L				
		cis-1,2-dichloroethene		<	2	ug/L				
		chloroform		<	2	ug/L				
		bromochloromethane		<	2	ug/L				
		tetrahydrofuran (THF)		<	10	ug/L				
		1,1,1-trichloroethane		<	2	ug/L				
		1,1-dichloropropene		<	2	ug/L				
		t-amyl-methyl ether (TAME)		<	2	ug/L				
		carbon tetrachloride		<	2	ug/L				
		1,2-dichloroethane		<	2	ug/L				
		benzene		<	2	ug/L				
		trichloroethene		<	2	ug/L				
		1,2-dichloropropane		<	2	ug/L				
		bromodichloromethane		<	0.6	ug/L				
		1,4-dioxane		<	50	ug/L				
		dibromomethane		<	2	ug/L				
		4-methyl-2-pentanone (MIBK)		<	10	ug/L				
		cis-1,3-dichloropropene		<	2	ug/L				
		toluene		<	2	ug/L				
		trans-1,3-dichloropropene		<	2	ug/L				
		2-hexanone		<	10	ug/L				
		1,1,2-trichloroethane		<	2	ug/L				
		1,3-dichloropropane		<	2	ug/L				
		tetrachloroethene		<	2	ug/L				
		dibromochloromethane		<	2	ug/L				
		1,2-dibromoethane (EDB)		<	2	ug/L				
		chlorobenzene		<	2	ug/L				
		1,1,1,2-tetrachloroethane		<	2	ug/L				
		ethylbenzene		<	2	ug/L				
		m&p-xylenes		<	2	ug/L				
		o-xylene		<	2	ug/L				
		styrene		<	2	ug/L				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	BLK1101353	bromoform		<	2	ug/L				
		isopropylbenzene		<	2	ug/L				
		1,1,2,2-tetrachloroethane		<	2	ug/L				
		1,2,3-trichloropropane		<	2	ug/L				
		n-propylbenzene		<	2	ug/L				
		bromobenzene		<	2	ug/L				
		1,3,5-trimethylbenzene		<	2	ug/L				
		2-chlorotoluene		<	2	ug/L				
		4-chlorotoluene		<	2	ug/L				
		tert-butylbenzene		<	2	ug/L				
		1,2,4-trimethylbenzene		<	2	ug/L				
		sec-butylbenzene		<	2	ug/L				
		1,3-dichlorobenzene		<	2	ug/L				
		4-isopropyltoluene		<	2	ug/L				
		1,4-dichlorobenzene		<	2	ug/L				
		1,2-dichlorobenzene		<	2	ug/L				
		n-butylbenzene		<	2	ug/L				
		1,2-dibromo-3-chloropropane		<	2	ug/L				
		1,2,4-trichlorobenzene		<	2	ug/L				
		1,3,5-trichlorobenzene		<	2	ug/L				
		hexachlorobutadiene			0.5	ug/L				
		naphthalene		<	5	ug/L				
		1,2,3-trichlorobenzene		<	2	ug/L				
		dibromofluoromethane SUR			96	%		78 114		
		toluene-D8 SUR			96	%		88 110		
		4-bromofluorobenzene SUR			94	%		86 115		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCS1101353	dichlorodifluoromethane		14	ug/L	20	68 *	70 130		
		chloromethane		17	ug/L	20	85	70 130		
		vinyl chloride		18	ug/L	20	92	70 130		
		bromomethane		26	ug/L	20	130	70 130		
		chloroethane		19	ug/L	20	93	70 130		
		trichlorofluoromethane		17	ug/L	20	87	70 130		
		diethyl ether		19	ug/L	20	95	70 130		
		acetone	<	50	ug/L	20	98			
		1,1-dichloroethene		15	ug/L	20	77	70 130		
		methylene chloride		18	ug/L	20	91	70 130		
		carbon disulfide		17	ug/L	20	83	70 130		
		methyl t-butyl ether (MTBE)		19	ug/L	20	94	70 130		
		trans-1,2-dichloroethene		17	ug/L	20	85	70 130		
		isopropyl ether (DIPE)		19	ug/L	20	93	70 130		
		ethyl t-butyl ether (ETBE)		19	ug/L	20	93	70 130		
		1,1-dichloroethane		18	ug/L	20	88	70 130		
		t-butanol (TBA)		100	ug/L	100	103	70 130		
		2-butanone (MEK)		17	ug/L	20	85	70 130		
		2,2-dichloropropane		15	ug/L	20	75	70 130		
		cis-1,2-dichloroethene		19	ug/L	20	94	70 130		
		chloroform		19	ug/L	20	94	70 130		
		bromochloromethane		19	ug/L	20	96	70 130		
		tetrahydrofuran (THF)		19	ug/L	20	93	70 130		
		1,1,1-trichloroethane		17	ug/L	20	86	70 130		
		1,1-dichloropropene		19	ug/L	20	97	70 130		
		t-amyl-methyl ether (TAME)		18	ug/L	20	91	70 130		
		carbon tetrachloride		16	ug/L	20	82	70 130		
		1,2-dichloroethane		19	ug/L	20	95	70 130		
		benzene		19	ug/L	20	97	70 130		
		trichloroethene		19	ug/L	20	93	70 130		
		1,2-dichloropropane		19	ug/L	20	96	70 130		
		bromodichloromethane		17	ug/L	20	85	70 130		
		1,4-dioxane	<	50	ug/L	40	106	70 130		
		dibromomethane		19	ug/L	20	93	70 130		
		4-methyl-2-pentanone (MIBK)		17	ug/L	20	84	70 130		
		cis-1,3-dichloropropene		17	ug/L	20	85	70 130		
		toluene		19	ug/L	20	95	70 130		
		trans-1,3-dichloropropene		18	ug/L	20	89	70 130		
		2-hexanone		16	ug/L	20	80	70 130		
		1,1,2-trichloroethane		19	ug/L	20	95	70 130		
		1,3-dichloropropane		18	ug/L	20	90	70 130		
		tetrachloroethene		20	ug/L	20	100	70 130		
		dibromochloromethane		14	ug/L	20	69 *	70 130		
		1,2-dibromoethane (EDB)		17	ug/L	20	87	70 130		
		chlorobenzene		19	ug/L	20	97	70 130		
		1,1,1,2-tetrachloroethane		16	ug/L	20	79	70 130		
		ethylbenzene		18	ug/L	20	89	70 130		
		m&p-xylenes		40	ug/L	40	99	70 130		
		o-xylene		21	ug/L	20	105	70 130		
		styrene		18	ug/L	20	89	70 130		
		bromoform		16	ug/L	20	82	70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCS1101353	isopropylbenzene		19	ug/L	20	97	70 130		
		1,1,2,2-tetrachloroethane		21	ug/L	20	105	70 130		
		1,2,3-trichloropropane		20	ug/L	20	98	70 130		
		n-propylbenzene		21	ug/L	20	106	70 130		
		bromobenzene		20	ug/L	20	98	70 130		
		1,3,5-trimethylbenzene		20	ug/L	20	101	70 130		
		2-chlorotoluene		21	ug/L	20	106	70 130		
		4-chlorotoluene		21	ug/L	20	103	70 130		
		tert-butylbenzene		20	ug/L	20	102	70 130		
		1,2,4-trimethylbenzene		21	ug/L	20	103	70 130		
		sec-butylbenzene		21	ug/L	20	106	70 130		
		1,3-dichlorobenzene		21	ug/L	20	107	70 130		
		4-isopropyltoluene		21	ug/L	20	105	70 130		
		1,4-dichlorobenzene		20	ug/L	20	98	70 130		
		1,2-dichlorobenzene		21	ug/L	20	103	70 130		
		n-butylbenzene		20	ug/L	20	102	70 130		
		1,2-dibromo-3-chloropropane		14	ug/L	20	70	70 130		
		1,2,4-trichlorobenzene		18	ug/L	20	91	70 130		
		1,3,5-trichlorobenzene		19	ug/L	20	96	70 130		
		hexachlorobutadiene		20	ug/L	20	101	70 130		
		naphthalene		19	ug/L	20	97	70 130		
		1,2,3-trichlorobenzene		20	ug/L	20	100	70 130		
		dibromofluoromethane SUR		92	%			78 114		
		toluene-D8 SUR		96	%			88 110		
		4-bromofluorobenzene SUR		100	%			86 115		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCS1101353	dichlorodifluoromethane		14	ug/L	20	70	70 130	2	20
		chloromethane		17	ug/L	20	87	70 130	2	20
		vinyl chloride		19	ug/L	20	96	70 130	4	20
		bromomethane		26	ug/L	20	131 *	70 130	1	20
		chloroethane		19	ug/L	20	94	70 130	2	20
		trichlorofluoromethane		18	ug/L	20	91	70 130	4	20
		diethyl ether		20	ug/L	20	99	70 130	5	20
		acetone	<	50	ug/L	20	97		1	20
		1,1-dichloroethene		16	ug/L	20	82	70 130	6	20
		methylene chloride		18	ug/L	20	91	70 130	0	20
		carbon disulfide		17	ug/L	20	87	70 130	4	20
		methyl t-butyl ether (MTBE)		19	ug/L	20	97	70 130	3	20
		trans-1,2-dichloroethene		17	ug/L	20	86	70 130	1	20
		isopropyl ether (DIPE)		20	ug/L	20	99	70 130	6	20
		ethyl t-butyl ether (ETBE)		19	ug/L	20	95	70 130	2	20
		1,1-dichloroethane		18	ug/L	20	91	70 130	3	20
		t-butanol (TBA)		110	ug/L	100	110	70 130	7	20
		2-butanone (MEK)		18	ug/L	20	92	70 130	9	20
		2,2-dichloropropane		15	ug/L	20	76	70 130	1	20
		cis-1,2-dichloroethene		19	ug/L	20	95	70 130	1	20
		chloroform		20	ug/L	20	98	70 130	4	20
		bromochloromethane		20	ug/L	20	99	70 130	2	20
		tetrahydrofuran (THF)		20	ug/L	20	100	70 130	7	20
		1,1,1-trichloroethane		18	ug/L	20	92	70 130	7	20
		1,1-dichloropropene		20	ug/L	20	100	70 130	3	20
		t-amyl-methyl ether (TAME)		19	ug/L	20	94	70 130	3	20
		carbon tetrachloride		18	ug/L	20	91	70 130	10	20
		1,2-dichloroethane		20	ug/L	20	98	70 130	3	20
		benzene		20	ug/L	20	100	70 130	3	20
		trichloroethene		19	ug/L	20	93	70 130	0	20
		1,2-dichloropropane		20	ug/L	20	100	70 130	3	20
		bromodichloromethane		19	ug/L	20	93	70 130	8	20
		1,4-dioxane	<	50	ug/L	40	121	70 130	13	20
		dibromomethane		19	ug/L	20	97	70 130	4	20
		4-methyl-2-pentanone (MIBK)		18	ug/L	20	92	70 130	9	20
		cis-1,3-dichloropropene		18	ug/L	20	88	70 130	4	20
		toluene		20	ug/L	20	99	70 130	4	20
		trans-1,3-dichloropropene		19	ug/L	20	94	70 130	6	20
		2-hexanone		17	ug/L	20	84	70 130	5	20
		1,1,2-trichloroethane		20	ug/L	20	98	70 130	3	20
		1,3-dichloropropane		18	ug/L	20	92	70 130	3	20
		tetrachloroethene		20	ug/L	20	99	70 130	1	20
		dibromochloromethane		15	ug/L	20	76	70 130	10	20
		1,2-dibromoethane (EDB)		18	ug/L	20	88	70 130	2	20
		chlorobenzene		20	ug/L	20	100	70 130	2	20
		1,1,1,2-tetrachloroethane		17	ug/L	20	84	70 130	7	20
		ethylbenzene		18	ug/L	20	90	70 130	1	20
		m&p-xylenes		40	ug/L	40	101	70 130	1	20
		o-xylene		21	ug/L	20	105	70 130	0	20
		styrene		18	ug/L	20	92	70 130	3	20
		bromoform		18	ug/L	20	88	70 130	7	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCS1101353	isopropylbenzene		19	ug/L	20	97	70 130	1	20
		1,1,2,2-tetrachloroethane		22	ug/L	20	108	70 130	3	20
		1,2,3-trichloropropane		20	ug/L	20	99	70 130	1	20
		n-propylbenzene		21	ug/L	20	106	70 130	0	20
		bromobenzene		20	ug/L	20	100	70 130	2	20
		1,3,5-trimethylbenzene		20	ug/L	20	101	70 130	0	20
		2-chlorotoluene		21	ug/L	20	106	70 130	0	20
		4-chlorotoluene		21	ug/L	20	103	70 130	0	20
		tert-butylbenzene		21	ug/L	20	105	70 130	3	20
		1,2,4-trimethylbenzene		21	ug/L	20	103	70 130	0	20
		sec-butylbenzene		21	ug/L	20	105	70 130	1	20
		1,3-dichlorobenzene		22	ug/L	20	108	70 130	1	20
		4-isopropyltoluene		21	ug/L	20	105	70 130	0	20
		1,4-dichlorobenzene		20	ug/L	20	100	70 130	2	20
		1,2-dichlorobenzene		21	ug/L	20	104	70 130	1	20
		n-butylbenzene		20	ug/L	20	101	70 130	1	20
		1,2-dibromo-3-chloropropane		16	ug/L	20	82	70 130	15	20
		1,2,4-trichlorobenzene		19	ug/L	20	93	70 130	2	20
		1,3,5-trichlorobenzene		19	ug/L	20	95	70 130	1	20
		hexachlorobutadiene		20	ug/L	20	98	70 130	3	20
		naphthalene		20	ug/L	20	101	70 130	5	20
		1,2,3-trichlorobenzene		21	ug/L	20	103	70 130	3	20
		dibromofluoromethane SUR		94	%			78 114		
		toluene-D8 SUR		95	%			88 110		
		4-bromofluorobenzene SUR		104	%			86 115		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	BLK4327	N-nitrosodimethylamine		<	2	ug/L				
		aniline		<	2	ug/L				
		phenol		<	2	ug/L				
		2-chlorophenol		<	5	ug/L				
		bis(2-chloroethyl)ether		<	2	ug/L				
		1,3-dichlorobenzene		<	2	ug/L				
		1,4-dichlorobenzene		<	2	ug/L				
		1,2-dichlorobenzene		<	2	ug/L				
		benzyl alcohol		<	2	ug/L				
		2-methylphenol		<	2	ug/L				
		bis(2-chloroisopropyl) ether		<	2	ug/L				
		hexachloroethane		<	2	ug/L				
		N-nitroso-di-N-propylamine		<	2	ug/L				
		4-methylphenol		<	2	ug/L				
		nitrobenzene		<	2	ug/L				
		isophorone		<	5	ug/L				
		2-nitrophenol		<	2	ug/L				
		2,4-dimethylphenol		<	2	ug/L				
		bis(2-chloroethoxy)methane		<	5	ug/L				
		2,4-dichlorophenol		<	5	ug/L				
		1,2,4-trichlorobenzene		<	5	ug/L				
		naphthalene		<	0.5	ug/L				
		benzoic acid		<	50	ug/L				
		4-chloroaniline		<	2	ug/L				
		hexachlorobutadiene		<	2	ug/L				
		4-chloro-3-methylphenol		<	2	ug/L				
		2-methylnaphthalene		<	0.5	ug/L				
		hexachlorocyclopentadiene		<	10	ug/L				
		2,4,6-trichlorophenol		<	2	ug/L				
		2,4,5-trichlorophenol		<	2	ug/L				
		2-chloronaphthalene		<	5	ug/L				
		2-nitroaniline		<	2	ug/L				
		acenaphthylene		<	0.5	ug/L				
		dimethylphthalate		<	5	ug/L				
		2,6-dinitrotoluene		<	2	ug/L				
		2,4-dinitrotoluene		<	2	ug/L				
		acenaphthene		<	0.5	ug/L				
		3-nitroaniline		<	2	ug/L				
		2,4-dinitrophenol		<	50	ug/L				
		dibenzofuran		<	0.5	ug/L				
		4-nitrophenol		<	10	ug/L				
		fluorene		<	0.5	ug/L				
		diethyl phthalate		<	5	ug/L				
		4-chlorophenyl phenyl ether		<	5	ug/L				
		4-nitroaniline		<	5	ug/L				
		4,6-dinitro-2-methylphenol		<	20	ug/L				
		azobenzene		<	2	ug/L				
		N-nitrosodiphenylamine		<	2	ug/L				
		4-bromophenyl phenyl ether		<	2	ug/L				
		hexachlorobenzene		<	2	ug/L				
		pentachlorophenol		<	10	ug/L				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	BLK4327	phenanthrene		<	0.5	ug/L				
		anthracene		<	0.5	ug/L				
		carbazole		<	2	ug/L				
		di-n-butylphthalate		<	5	ug/L				
		fluoranthene		<	0.5	ug/L				
		benzidine		<	30	ug/L				
		pyrene		<	0.5	ug/L				
		butyl benzyl phthalate		<	5	ug/L				
		benzo(a)anthracene		<	0.5	ug/L				
		chrysene		<	0.5	ug/L				
		3,3'-dichlorobenzidine		<	30	ug/L				
		bis(2-ethylhexyl)phthalate		<	5	ug/L				
		di-n-octyl phthalate		<	2	ug/L				
		benzo(b)fluoranthene		<	0.5	ug/L				
		benzo(k)fluoranthene		<	0.5	ug/L				
		benzo(a)pyrene		<	0.2	ug/L				
		indeno(1,2,3-cd)pyrene		<	0.5	ug/L				
		dibenzo(a,h)anthracene		<	0.5	ug/L				
		benzo(g,h,i)perylene		<	0.5	ug/L				
		2-fluorophenol SUR			30	%		21	100	
		phenol-D5 SUR			19	%		10	102	
		2,4,6-tribromophenol SUR			54	%		10	123	
		nitrobenzene-D5 SUR			61	%		35	114	
		2-fluorobiphenyl SUR			73	%		43	116	
		p-terphenyl-D14 SUR			75	%		33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCS4327	N-nitrosodimethylamine		<	2	ug/L				
		aniline		<	2	ug/L				
		phenol			22	ug/L	100	22	*	30 130
		2-chlorophenol			64	ug/L	100	64		30 130
		bis(2-chloroethyl)ether		<	2	ug/L				
		1,3-dichlorobenzene		<	2	ug/L				
		1,4-dichlorobenzene		<	2	ug/L				
		1,2-dichlorobenzene		<	2	ug/L				
		benzyl alcohol		<	2	ug/L				
		2-methylphenol		<	2	ug/L				
		bis(2-chloroisopropyl) ether		<	2	ug/L				
		hexachloroethane		<	2	ug/L				
		N-nitroso-di-N-propylamine		<	2	ug/L				
		4-methylphenol		<	2	ug/L				
		nitrobenzene		<	2	ug/L				
		isophorone		<	5	ug/L				
		2-nitrophenol		<	2	ug/L				
		2,4-dimethylphenol		<	2	ug/L				
		bis(2-chloroethoxy)methane		<	5	ug/L				
		2,4-dichlorophenol		<	5	ug/L				
		1,2,4-trichlorobenzene		<	5	ug/L				
		naphthalene			26	ug/L	40	65		40 140
		benzoic acid		<	50	ug/L				
		4-chloroaniline		<	2	ug/L				
		hexachlorobutadiene		<	2	ug/L				
		4-chloro-3-methylphenol			64	ug/L	100	64		30 130
		2-methylnaphthalene			26	ug/L	40	65		40 140
		hexachlorocyclopentadiene		<	10	ug/L				
		2,4,6-trichlorophenol		<	2	ug/L				
		2,4,5-trichlorophenol		<	2	ug/L				
		2-chloronaphthalene		<	5	ug/L				
		2-nitroaniline		<	2	ug/L				
		acenaphthylene			28	ug/L	40	70		40 140
		dimethylphthalate		<	5	ug/L				
		2,6-dinitrotoluene		<	2	ug/L				
		2,4-dinitrotoluene		<	2	ug/L				
		acenaphthene			27	ug/L				
		3-nitroaniline		<	2	ug/L				
		2,4-dinitrophenol		<	50	ug/L				
		dibenzofuran		<	0.5	ug/L				
		4-nitrophenol			21	ug/L	100	21	*	30 130
		fluorene			29	ug/L	40	73		40 140
		diethyl phthalate		<	5	ug/L				
		4-chlorophenyl phenyl ether		<	5	ug/L				
		4-nitroaniline		<	5	ug/L				
		4,6-dinitro-2-methylphenol		<	20	ug/L				
		azobenzene		<	2	ug/L				
		N-nitrosodiphenylamine		<	2	ug/L				
		4-bromophenyl phenyl ether		<	2	ug/L				
		hexachlorobenzene		<	2	ug/L				
		pentachlorophenol			49	ug/L	100	49		30 130

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCS4327	phenanthrene		27	ug/L	40	66	40 140		
		anthracene		27	ug/L	40	68	40 140		
		carbazole	<	2	ug/L					
		di-n-butylphthalate	<	5	ug/L					
		fluoranthene		29	ug/L	40	73	40 140		
		benzidine	<	30	ug/L					
		pyrene		32	ug/L	40	80	40 140		
		butyl benzyl phthalate	<	5	ug/L					
		benzo(a)anthracene		29	ug/L	40	73	40 140		
		chrysene		30	ug/L	40	74	40 140		
		3,3'-dichlorobenzidine	<	30	ug/L					
		bis(2-ethylhexyl)phthalate	<	5	ug/L					
		di-n-octyl phthalate	<	2	ug/L					
		benzo(b)fluoranthene		29	ug/L	40	71	40 140		
		benzo(k)fluoranthene		30	ug/L	40	76	40 140		
		benzo(a)pyrene		30	ug/L	40	75	40 140		
		indeno(1,2,3-cd)pyrene		28	ug/L	40	71	40 140		
		dibenzo(a,h)anthracene		28	ug/L	40	70	40 140		
		benzo(g,h,i)perylene		28	ug/L	40	71	40 140		
		2-fluorophenol SUR		34	%				21 100	
		phenol-D5 SUR		23	%				10 102	
		2,4,6-tribromophenol SUR		64	%				10 123	
		nitrobenzene-D5 SUR		68	%				35 114	
		2-fluorobiphenyl SUR		78	%				43 116	
		p-terphenyl-D14 SUR		80	%				33 141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCSD4327	N-nitrosodimethylamine		<	2	ug/L				
		aniline		<	2	ug/L				
		phenol		21	ug/L	100	21 *	30 130	5	20
		2-chlorophenol		64	ug/L	100	64	30 130	0	20
		bis(2-chloroethyl)ether		<	2	ug/L				
		1,3-dichlorobenzene		<	2	ug/L				
		1,4-dichlorobenzene		<	2	ug/L				
		1,2-dichlorobenzene		<	2	ug/L				
		benzyl alcohol		<	2	ug/L				
		2-methylphenol		<	2	ug/L				
		bis(2-chloroisopropyl) ether		<	2	ug/L				
		hexachloroethane		<	2	ug/L				
		N-nitroso-di-N-propylamine		<	2	ug/L				
		4-methylphenol		<	2	ug/L				
		nitrobenzene		<	2	ug/L				
		isophorone		<	5	ug/L				
		2-nitrophenol		<	2	ug/L				
		2,4-dimethylphenol		<	2	ug/L				
		bis(2-chloroethoxy)methane		<	5	ug/L				
		2,4-dichlorophenol		<	5	ug/L				
		1,2,4-trichlorobenzene		<	5	ug/L				
		naphthalene		26	ug/L	40	64	40 140	1	20
		benzoic acid		<	50	ug/L				
		4-chloroaniline		<	2	ug/L				
		hexachlorobutadiene		<	2	ug/L				
		4-chloro-3-methylphenol		63	ug/L	100	63	30 130	0	20
		2-methylnaphthalene		25	ug/L	40	64	40 140	3	20
		hexachlorocyclopentadiene		<	10	ug/L				
		2,4,6-trichlorophenol		<	2	ug/L				
		2,4,5-trichlorophenol		<	2	ug/L				
		2-chloronaphthalene		<	5	ug/L				
		2-nitroaniline		<	2	ug/L				
		acenaphthylene		28	ug/L	40	70	40 140	0	20
		dimethylphthalate		<	5	ug/L				
		2,6-dinitrotoluene		<	2	ug/L				
		2,4-dinitrotoluene		<	2	ug/L				
		acenaphthene		27	ug/L					
		3-nitroaniline		<	2	ug/L				
		2,4-dinitrophenol		<	50	ug/L				
		dibenzofuran		<	0.5	ug/L				
		4-nitrophenol		19	ug/L	100	19 *	30 130	7	20
		fluorene		28	ug/L	40	70	40 140	3	20
		diethyl phthalate		<	5	ug/L				
		4-chlorophenyl phenyl ether		<	5	ug/L				
		4-nitroaniline		<	5	ug/L				
		4,6-dinitro-2-methylphenol		<	20	ug/L				
		azobenzene		<	2	ug/L				
		N-nitrosodiphenylamine		<	2	ug/L				
		4-bromophenyl phenyl ether		<	2	ug/L				
		hexachlorobenzene		<	2	ug/L				
		pentachlorophenol		51	ug/L	100	51	30 130	5	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3510C8270D	LCSD4327	phenanthrene		26	ug/L	40	66	40 140	1	20	
		anthracene		28	ug/L	40	69	40 140	2	20	
		carbazole	<	2	ug/L						
		di-n-butylphthalate	<	5	ug/L						
		fluoranthene		29	ug/L	40	71	40 140	2	20	
		benzidine	<	30	ug/L						
		pyrene		32	ug/L	40	81	40 140	1	20	
		butyl benzyl phthalate	<	5	ug/L						
		benzo(a)anthracene		29	ug/L	40	73	40 140	0	20	
		chrysene		29	ug/L	40	74	40 140	1	20	
		3,3'-dichlorobenzidine	<	30	ug/L						
		bis(2-ethylhexyl)phthalate	<	5	ug/L						
		di-n-octyl phthalate	<	2	ug/L						
		benzo(b)fluoranthene		29	ug/L	40	71	40 140	0	20	
		benzo(k)fluoranthene		29	ug/L	40	73	40 140	4	20	
		benzo(a)pyrene		29	ug/L	40	73	40 140	3	20	
		indeno(1,2,3-cd)pyrene		28	ug/L	40	70	40 140	1	20	
		dibenzo(a,h)anthracene		28	ug/L	40	70	40 140	1	20	
		benzo(g,h,i)perylene		28	ug/L	40	70	40 140	1	20	
		2-fluorophenol SUR		34	%				21 100		
		phenol-D5 SUR		22	%				10 102		
		2,4,6-tribromophenol SUR		66	%				10 123		
		nitrobenzene-D5 SUR		66	%				35 114		
		2-fluorobiphenyl SUR		71	%				43 116		
		p-terphenyl-D14 SUR		80	%				33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	BLK4314	N-nitrosodimethylamine		<	0.2	ug/g				
		aniline		<	0.2	ug/g				
		phenol		<	0.2	ug/g				
		2-chlorophenol		<	0.5	ug/g				
		bis(2-chloroethyl)ether		<	0.2	ug/g				
		1,3-dichlorobenzene		<	0.2	ug/g				
		1,4-dichlorobenzene		<	0.2	ug/g				
		1,2-dichlorobenzene		<	0.2	ug/g				
		benzyl alcohol		<	0.2	ug/g				
		2-methylphenol		<	0.2	ug/g				
		bis(2-chloroisopropyl) ether		<	0.2	ug/g				
		hexachloroethane		<	0.2	ug/g				
		N-nitroso-di-N-propylamine		<	0.2	ug/g				
		4-methylphenol		<	0.2	ug/g				
		nitrobenzene		<	0.2	ug/g				
		isophorone		<	0.5	ug/g				
		2-nitrophenol		<	0.2	ug/g				
		2,4-dimethylphenol		<	0.2	ug/g				
		bis(2-chloroethoxy)methane		<	0.5	ug/g				
		2,4-dichlorophenol		<	0.5	ug/g				
		1,2,4-trichlorobenzene		<	0.5	ug/g				
		naphthalene		<	0.05	ug/g				
		benzoic acid		<	5.0	ug/g				
		4-chloroaniline		<	0.2	ug/g				
		hexachlorobutadiene		<	0.2	ug/g				
		4-chloro-3-methylphenol		<	0.2	ug/g				
		2-methylnaphthalene		<	0.05	ug/g				
		hexachlorocyclopentadiene		<	1.0	ug/g				
		2,4,6-trichlorophenol		<	0.2	ug/g				
		2,4,5-trichlorophenol		<	0.2	ug/g				
		2-chloronaphthalene		<	0.5	ug/g				
		2-nitroaniline		<	0.2	ug/g				
		acenaphthylene		<	0.05	ug/g				
		dimethylphthalate		<	0.5	ug/g				
		2,6-dinitrotoluene		<	0.2	ug/g				
		2,4-dinitrotoluene		<	0.2	ug/g				
		acenaphthene		<	0.05	ug/g				
		3-nitroaniline		<	0.2	ug/g				
		2,4-dinitrophenol		<	5.0	ug/g				
		dibenzofuran		<	0.05	ug/g				
		4-nitrophenol		<	1.0	ug/g				
		fluorene		<	0.05	ug/g				
		diethyl phthalate		<	0.5	ug/g				
		4-chlorophenyl phenyl ether		<	0.5	ug/g				
		4-nitroaniline		<	0.5	ug/g				
		4,6-dinitro-2-methylphenol		<	2.0	ug/g				
		azobenzene		<	0.2	ug/g				
		N-nitrosodiphenylamine		<	0.2	ug/g				
		4-bromophenyl phenyl ether		<	0.2	ug/g				
		hexachlorobenzene		<	0.2	ug/g				
		pentachlorophenol		<	1.0	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	BLK4314	phenanthrene		<	0.05	ug/g				
		anthracene		<	0.05	ug/g				
		carbazole		<	0.2	ug/g				
		di-n-butylphthalate		<	0.5	ug/g				
		fluoranthene		<	0.05	ug/g				
		benzidine		<	3.0	ug/g				
		pyrene		<	0.05	ug/g				
		butyl benzyl phthalate		<	0.5	ug/g				
		benzo(a)anthracene		<	0.05	ug/g				
		chrysene		<	0.05	ug/g				
		3,3'-dichlorobenzidine		<	3.0	ug/g				
		bis(2-ethylhexyl)phthalate		<	0.5	ug/g				
		di-n-octyl phthalate		<	0.2	ug/g				
		benzo(b)fluoranthene		<	0.05	ug/g				
		benzo(k)fluoranthene		<	0.05	ug/g				
		benzo(a)pyrene		<	0.02	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.05	ug/g				
		dibenzo(a,h)anthracene		<	0.05	ug/g				
		benzo(g,h,i)perylene		<	0.05	ug/g				
		2-fluorophenol SUR			43	%		21	100	
		phenol-D5 SUR			46	%		10	102	
		2,4,6-tribromophenol SUR			40	%		10	123	
		nitrobenzene-D5 SUR			38	%		35	114	
		2-fluorobiphenyl SUR			49	%		43	116	
		p-terphenyl-D14 SUR			66	%		33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCS4314	N-nitrosodimethylamine		<	0.2	ug/g				
		aniline		<	0.2	ug/g				
		phenol			5.6	ug/g	10	56	30	130
		2-chlorophenol			6.5	ug/g	10	65	30	130
		bis(2-chloroethyl)ether		<	0.2	ug/g				
		1,3-dichlorobenzene		<	0.2	ug/g				
		1,4-dichlorobenzene		<	0.2	ug/g				
		1,2-dichlorobenzene		<	0.2	ug/g				
		benzyl alcohol		<	0.2	ug/g				
		2-methylphenol		<	0.2	ug/g				
		bis(2-chloroisopropyl) ether		<	0.2	ug/g				
		hexachloroethane		<	0.2	ug/g				
		N-nitroso-di-N-propylamine		<	0.2	ug/g				
		4-methylphenol		<	0.2	ug/g				
		nitrobenzene		<	0.2	ug/g				
		isophorone		<	0.5	ug/g				
		2-nitrophenol		<	0.2	ug/g				
		2,4-dimethylphenol		<	0.2	ug/g				
		bis(2-chloroethoxy)methane		<	0.5	ug/g				
		2,4-dichlorophenol		<	0.5	ug/g				
		1,2,4-trichlorobenzene		<	0.5	ug/g				
		naphthalene			2.4	ug/g	4	59	40	140
		benzoic acid		<	5.0	ug/g				
		4-chloroaniline		<	0.2	ug/g				
		hexachlorobutadiene		<	0.2	ug/g				
		4-chloro-3-methylphenol			6.6	ug/g	10	66	30	130
		2-methylnaphthalene			2.45	ug/g	4	61	40	140
		hexachlorocyclopentadiene		<	1.0	ug/g				
		2,4,6-trichlorophenol		<	0.2	ug/g				
		2,4,5-trichlorophenol		<	0.2	ug/g				
		2-chloronaphthalene		<	0.5	ug/g				
		2-nitroaniline		<	0.2	ug/g				
		acenaphthylene			2.9	ug/g	4	72	40	140
		dimethylphthalate		<	0.5	ug/g				
		2,6-dinitrotoluene		<	0.2	ug/g				
		2,4-dinitrotoluene		<	0.2	ug/g				
		acenaphthene			2.8	ug/g				
		3-nitroaniline		<	0.2	ug/g				
		2,4-dinitrophenol		<	5.0	ug/g				
		dibenzofuran		<	0.05	ug/g				
		4-nitrophenol			6.4	ug/g	10	64	30	130
		fluorene			3.0	ug/g	4	76	40	140
		diethyl phthalate		<	0.5	ug/g				
		4-chlorophenyl phenyl ether		<	0.5	ug/g				
		4-nitroaniline		<	0.5	ug/g				
		4,6-dinitro-2-methylphenol		<	2.0	ug/g				
		azobenzene		<	0.2	ug/g				
		N-nitrosodiphenylamine		<	0.2	ug/g				
		4-bromophenyl phenyl ether		<	0.2	ug/g				
		hexachlorobenzene		<	0.2	ug/g				
		pentachlorophenol			4.3	ug/g	10	43	30	130

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCS4314	phenanthrene		2.9	ug/g	4	72	40 140		
		anthracene		3.0	ug/g	4	75	40 140		
		carbazole	<	0.2	ug/g					
		di-n-butylphthalate	<	0.5	ug/g					
		fluoranthene		3.3	ug/g	4	81	40 140		
		benzidine	<	3.0	ug/g					
		pyrene		3.5	ug/g	4	87	40 140		
		butyl benzyl phthalate	<	0.5	ug/g					
		benzo(a)anthracene		3.3	ug/g	4	82	40 140		
		chrysene		3.3	ug/g	4	82	40 140		
		3,3'-dichlorobenzidine	<	3.0	ug/g					
		bis(2-ethylhexyl)phthalate	<	0.5	ug/g					
		di-n-octyl phthalate	<	0.2	ug/g					
		benzo(b)fluoranthene		3.4	ug/g	4	85	40 140		
		benzo(k)fluoranthene		3.0	ug/g	4	76	40 140		
		benzo(a)pyrene		3.3	ug/g	4	82	40 140		
		indeno(1,2,3-cd)pyrene		3.2	ug/g	4	81	40 140		
		dibenzo(a,h)anthracene		3.2	ug/g	4	79	40 140		
		benzo(g,h,i)perylene		3.2	ug/g	4	79	40 140		
		2-fluorophenol SUR		53	%			21 100		
		phenol-D5 SUR		57	%			10 102		
		2,4,6-tribromophenol SUR		52	%			10 123		
		nitrobenzene-D5 SUR		50	%			35 114		
		2-fluorobiphenyl SUR		64	%			43 116		
		p-terphenyl-D14 SUR		71	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3546/8270D	LCS4314	N-nitrosodimethylamine		<	0.2	ug/g						
		aniline		<	0.2	ug/g						
		phenol			5.4	ug/g	10	54	30	130	4	30
		2-chlorophenol			5.9	ug/g	10	59	30	130	9	30
		bis(2-chloroethyl)ether		<	0.2	ug/g						
		1,3-dichlorobenzene		<	0.2	ug/g						
		1,4-dichlorobenzene		<	0.2	ug/g						
		1,2-dichlorobenzene		<	0.2	ug/g						
		benzyl alcohol		<	0.2	ug/g						
		2-methylphenol		<	0.2	ug/g						
		bis(2-chloroisopropyl) ether		<	0.2	ug/g						
		hexachloroethane		<	0.2	ug/g						
		N-nitroso-di-N-propylamine		<	0.2	ug/g						
		4-methylphenol		<	0.2	ug/g						
		nitrobenzene		<	0.2	ug/g						
		isophorone		<	0.5	ug/g						
		2-nitrophenol		<	0.2	ug/g						
		2,4-dimethylphenol		<	0.2	ug/g						
		bis(2-chloroethoxy)methane		<	0.5	ug/g						
		2,4-dichlorophenol		<	0.5	ug/g						
		1,2,4-trichlorobenzene		<	0.5	ug/g						
		naphthalene			2.2	ug/g	4	56	40	140	6	30
		benzoic acid		<	5.0	ug/g						
		4-chloroaniline		<	0.2	ug/g						
		hexachlorobutadiene		<	0.2	ug/g						
		4-chloro-3-methylphenol			6.4	ug/g	10	64	30	130	2	30
		2-methylnaphthalene			2.32	ug/g	4	58	40	140	5	30
		hexachlorocyclopentadiene		<	1.0	ug/g						
		2,4,6-trichlorophenol		<	0.2	ug/g						
		2,4,5-trichlorophenol		<	0.2	ug/g						
		2-chloronaphthalene		<	0.5	ug/g						
		2-nitroaniline		<	0.2	ug/g						
		acenaphthylene			2.9	ug/g	4	72	40	140	0	30
		dimethylphthalate		<	0.5	ug/g						
		2,6-dinitrotoluene		<	0.2	ug/g						
		2,4-dinitrotoluene		<	0.2	ug/g						
		acenaphthene			2.7	ug/g						
		3-nitroaniline		<	0.2	ug/g						
		2,4-dinitrophenol		<	5.0	ug/g						
		dibenzofuran		<	0.05	ug/g						
		4-nitrophenol			6.2	ug/g	10	62	30	130	3	30
		fluorene			3.0	ug/g	4	75	40	140	1	30
		diethyl phthalate		<	0.5	ug/g						
		4-chlorophenyl phenyl ether		<	0.5	ug/g						
		4-nitroaniline		<	0.5	ug/g						
		4,6-dinitro-2-methylphenol		<	2.0	ug/g						
		azobenzene		<	0.2	ug/g						
		N-nitrosodiphenylamine		<	0.2	ug/g						
		4-bromophenyl phenyl ether		<	0.2	ug/g						
		hexachlorobenzene		<	0.2	ug/g						
		pentachlorophenol			3.6	ug/g	10	36	30	130	17	30

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCSD4314	phenanthrene		2.9	ug/g	4	72	40 140	0	30
		anthracene		3.0	ug/g	4	76	40 140	0	30
		carbazole	<	0.2	ug/g					
		di-n-butylphthalate	<	0.5	ug/g					
		fluoranthene		3.4	ug/g	4	84	40 140	3	30
		benzidine	<	3.0	ug/g					
		pyrene		3.5	ug/g	4	88	40 140	2	30
		butyl benzyl phthalate	<	0.5	ug/g					
		benzo(a)anthracene		3.4	ug/g	4	84	40 140	2	30
		chrysene		3.4	ug/g	4	84	40 140	2	30
		3,3'-dichlorobenzidine	<	3.0	ug/g					
		bis(2-ethylhexyl)phthalate	<	0.5	ug/g					
		di-n-octyl phthalate	<	0.2	ug/g					
		benzo(b)fluoranthene		3.5	ug/g	4	87	40 140	3	30
		benzo(k)fluoranthene		3.2	ug/g	4	80	40 140	5	30
		benzo(a)pyrene		3.3	ug/g	4	84	40 140	1	30
		indeno(1,2,3-cd)pyrene		3.2	ug/g	4	81	40 140	0	30
		dibenzo(a,h)anthracene		3.2	ug/g	4	79	40 140	0	30
		benzo(g,h,i)perylene		3.2	ug/g	4	80	40 140	1	30
		2-fluorophenol SUR		49	%			21 100		
		phenol-D5 SUR		51	%			10 102		
		2,4,6-tribromophenol SUR		49	%			10 123		
		nitrobenzene-D5 SUR		46	%			35 114		
		2-fluorobiphenyl SUR		63	%			43 116		
		p-terphenyl-D14 SUR		70	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3005A6010C	BLK4318	Silver		<	0.25	ug/g				
		Arsenic		<	0.50	ug/g				
		Barium		<	2.5	ug/g				
		Cadmium		<	0.20	ug/g				
		Chromium		<	2.5	ug/g				
		Copper		<	2.5	ug/g				
		Nickel		<	2.5	ug/g				
		Lead		<	0.50	ug/g				
		Selenium		<	2.5	ug/g				
		Zinc		<	2.5	ug/g				
SW3005A6010C	DUP4318	Arsenic	21885-001	<	0.008	mg/L				20
		Copper	21885-001	<	0.05	mg/L				20
		Lead	21885-001	<	0.01	mg/L				20
		Selenium	21885-001	<	0.05	mg/L				20
		Zinc	21885-001		0.18	mg/L			1	20
SW3005A6010C	LCS4318	Silver		0.25	mg/L	0.25	101	80 120		
		Arsenic		0.48	mg/L	0.5	96	80 120		
		Barium		0.49	mg/L	0.5	98	80 120		
		Cadmium		0.50	mg/L	0.5	100	80 120		
		Chromium		0.49	mg/L	0.5	98	80 120		
		Copper		0.49	mg/L	0.5	99	80 120		
		Nickel		0.49	mg/L	0.5	98	80 120		
		Lead		0.50	mg/L	0.5	100	80 120		
		Selenium		0.46	mg/L	0.5	92	80 120		
		Zinc		0.48	mg/L	0.5	96	80 120		
SW3005A6010C	LCSD4318	Silver		0.26	mg/L	0.25	104	80 120	3	20
		Arsenic		0.48	mg/L	0.5	97	80 120	1	20
		Barium		0.51	mg/L	0.5	101	80 120	3	20
		Cadmium		0.49	mg/L	0.5	98	80 120	1	20
		Chromium		0.51	mg/L	0.5	101	80 120	3	20
		Copper		0.50	mg/L	0.5	100	80 120	1	20
		Nickel		0.51	mg/L	0.5	101	80 120	3	20
		Lead		0.50	mg/L	0.5	100	80 120	0	20
		Selenium		0.47	mg/L	0.5	95	80 120	3	20
		Zinc		0.50	mg/L	0.5	99	80 120	3	20
SW3005A6010C	MS4318	Arsenic	21885-001	0.50	mg/L	0.5	99	70 130		
		Copper	21885-001	0.53	mg/L	0.5	99	70 130		
		Lead	21885-001	0.49	mg/L	0.5	97	70 130		
		Selenium	21885-001	0.47	mg/L	0.5	94	70 130		
		Zinc	21885-001	0.66	mg/L	0.5	97	70 130		
SW3005A6010C	MS4318	Lead	21919-006	0.49	mg/L	0.5	99	70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW7470A	BLK4337	Mercury		< 0.0002	mg/L					
SW7470A	LCS4337	Mercury		0.0023	mg/L	0.002	116	80 120		
SW7470A	LCSD4337	Mercury		0.0023	mg/L	0.002	115	80 120	0	20
SW7470A	MS4337	Mercury	21917-002	0.0023	mg/L	0.002	114	75 125		

Company Name: **REDGE ASSURERS, LLC**

Project Name: **POLYCLAD**

Company Address: **776 MAIN ST, WESTCAMB, ME**

Project #: **10001086**

Report To: **Jordan Obonell**

Project Location: **MA ME VT**

Phone #: **(207) 828-1272 EXT # 20**

Reporting Limits: **QAAP EPA DW GW-1 S-1**

Invoice To: **T. Patten**

Quote # NH GREEN/ODD Fund Pricing
 PO #

ANALYSIS REQUEST

- VOC 8260 VOC 8260 NHDES VOC 8260 MADEP
- VOC 624 VOC BTEX MIBE, only VOC 8021VT
- VPH MADEP MEGRO GRO 8015
- VOC 524.2 VOC 524.2 NH List
- TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
- 8270PAH 8270ABN 625 EDB 504.1 **SVOCs**
- 8082 PCB 8081 Pesticides 608 Past/PCB
- O&G 1664 Mineral O&G SM5520F
- pH BOD Conductivity Turbidity
- TSS TDS TS TVS Alkalinity
- RCRA Metals Priority Pollutants Metals TAL Metals
- Total Metals-list: **LEAD**
- Dissolved Metals-list: **TOTAL Zn, Cu, Ni (6010)**
- Ammonia COD TKN TN TON
- T-Phosphorus Phenols Bacteria P/A Bacteria MPN
- Cyanide Sulfide Nitrate + Nitrite Ortho P
- Nitrate Nitrite Chloride Sulfate Bromide Fluoride
- Corrosivity Reactive CN Reactive S- Ignitibility/FP
- TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
- Subcontract: TOC Grain Size TCLP Herbicides

Lab Sample ID	Field ID	# CONTAINERS	Matrix			Preservation Method							Sampling			
			WATER	SOLID	OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)	DATE	TIME	SAMPLER*		
CA-1	2	2	<input checked="" type="checkbox"/>			<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>						7-5	1515	50	X
CA-2	4	4	<input checked="" type="checkbox"/>			<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>						7-5	1420	50	X
CA-3	3	3	<input checked="" type="checkbox"/>			<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>						7-5	1530	50	X
CAW-2	3	3	<input checked="" type="checkbox"/>			<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>						7-5	1320	50	X
SS-3	1	1	<input checked="" type="checkbox"/>			<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>						7-5	1200	50	X

SPECIAL INSTRUCTIONS

BROWN FIELDS
 All metals run as total - nothing was field stored per Jonathan & Steve

TAT REQUESTED
 Priority (24 hr)
 Standard (48 hr)
 (10 Business Days)
 Date Needed **7/9

* See www.reslabs.com for sample acceptance policy and current accreditation lists.

REPORTING INSTRUCTIONS
 HARD COPY REQUIRED FAX (FAX#)
 PDF (e-mail address) **jdomeire@credalls.com**
 OTHER (specify)

RECEIVED ON CE YES NO
 TEMPERATURE **7** °C

CUSTODY RECORD

Relinquished by: **Jordan Obonell**
 Date: **7/6/11** Time: **1040**

Relinquished by: **Jordan Obonell**
 Date: **7/6/11** Time: **1040**

Received by Laboratory: **Jordan Obonell**
 Way Bill#:

Received by: **Jordan Obonell**
 Date: **7/6/11** Time: **1040**

Received by: **Jordan Obonell**
 Date: **7/6/11** Time: **1040**



Absolute Resource *associates*

124 Heritage Avenue #10 Portsmouth, NH 03801

CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: None
Job ID: 22020
Date Received: 7/18/11

Project: POLYCLAD 10001086

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

Cliff Chase
Partner, Technical Director

Date of Approval: 7/26/2011
Total number of pages: 3

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Project ID: POLYCLAD 10001086

Job ID: 22020

Sample#: 22020-001

Sample ID: CA-2

Matrix: Water

Sampled: 7/14/11 15:39

Parameter	Quant		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Arsenic	< 0.008	0.008	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C
Barium	< 0.05	0.05	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	BJS	7/19/11	4364	7/20/11	11:02	SW7470A
Selenium	< 0.05	0.05	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	BJS	7/20/11	4369	7/20/11	14:05	SW3005A6010C



Absolute Resource Associates
 124 Heritage Avenue #10
 Portsmouth, NH 03801
 603-436-2001
 absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD
 AND ANALYSIS REQUEST**

22020
 PAGE 1 OF 1

Company Name: **CREDERE ASSOCIATES LLC**
 Company Address: **776 Main St. Westbrook ME**
 Report To: **S. O'DONNELL**
 Phone #: **207-272-2225**
 Invoice To: **T. PATTEN**
 Project Name: **POLYCLAD**
 Project #: **10001086**
 Project Location: **ME VT**
 Protocol: **RCPRA MCP NHDES OTHER**
 Reporting: **QAPP GW-1 S-1**
 Limits: **EPA DW Other**
 Quote # _____
 NH GREEN/ODD Fund Pricing

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix	Preservation Method	DATE	TIME	SAMPLER	ANALYSIS REQUEST	RECEIVED ON ICE	TEMPERATURE
22020-01	CA-2	1	WATER <input checked="" type="checkbox"/> SOLID <input type="checkbox"/> OTHER <input type="checkbox"/>	HCl <input type="checkbox"/> HNO ₃ <input checked="" type="checkbox"/> H ₂ SO ₄ <input type="checkbox"/> NaOH <input type="checkbox"/> MeOH <input type="checkbox"/> OTHER (Specify) <input type="checkbox"/>	7/14/11	1539	SD	<input type="checkbox"/> VOC 8260 <input type="checkbox"/> VOC 8260 NHDES <input type="checkbox"/> VOC 8260 MADEP <input type="checkbox"/> VOC 624 <input type="checkbox"/> VOC BTEX <input type="checkbox"/> MIBE, only <input type="checkbox"/> VOC 8021VT <input type="checkbox"/> VPH MADEP <input type="checkbox"/> MEGRO <input type="checkbox"/> GRO 8015 <input type="checkbox"/> VOC 524.2 <input type="checkbox"/> VOC 524.2 NH List <input type="checkbox"/> Gases-List: <input type="checkbox"/> TPH <input type="checkbox"/> DRO 8015 <input type="checkbox"/> MEDRO <input type="checkbox"/> EPH MADEP <input type="checkbox"/> TPH Fingerprint <input type="checkbox"/> 8270PAH <input type="checkbox"/> 8270ABN <input type="checkbox"/> 625 <input type="checkbox"/> EDB 504.1 <input type="checkbox"/> 8082 PCB <input type="checkbox"/> 8081 Pesticides <input type="checkbox"/> 608 Pest/PCB <input type="checkbox"/> O&G 1664 <input type="checkbox"/> Mineral O&G SM5520F <input type="checkbox"/> pH <input type="checkbox"/> BOD <input type="checkbox"/> Conductivity <input type="checkbox"/> Turbidity <input type="checkbox"/> TSS <input type="checkbox"/> TDS <input type="checkbox"/> TS <input type="checkbox"/> TVS <input type="checkbox"/> Alkalinity <input checked="" type="checkbox"/> RCRA Metals <input type="checkbox"/> Priority Pollutant Metals <input type="checkbox"/> TAL Metals <input type="checkbox"/> Total Metals-list: <input type="checkbox"/> Dissolved Metals-list: <input type="checkbox"/> Ammonia <input type="checkbox"/> COD <input type="checkbox"/> TKN <input type="checkbox"/> TN <input type="checkbox"/> TON <input type="checkbox"/> T-Phosphorus <input type="checkbox"/> Phenols <input type="checkbox"/> Bacteria P/A <input type="checkbox"/> Bacteria MPN <input type="checkbox"/> Cyanide <input type="checkbox"/> Sulfide <input type="checkbox"/> Nitrate + Nitrite <input type="checkbox"/> Ortho P <input type="checkbox"/> Nitrate <input type="checkbox"/> Nitrite <input type="checkbox"/> Chloride <input type="checkbox"/> Sulfate <input type="checkbox"/> Bromide <input type="checkbox"/> Fluoride <input type="checkbox"/> Corrosivity <input type="checkbox"/> Reactive CN <input type="checkbox"/> Reactive S- <input type="checkbox"/> Ignitibility/FP <input type="checkbox"/> TCLP Metals <input type="checkbox"/> TCLP VOC <input type="checkbox"/> TCLP SVOC <input type="checkbox"/> TCLP Pesticide Subcontract: <input type="checkbox"/> TOC <input type="checkbox"/> Grain Size <input type="checkbox"/> TCLP Herbicides	YES <input type="checkbox"/> NO <input type="checkbox"/>	

TAT REQUESTED
 Priority (24 hr)*
 Expedited (48 hr)*
 Standard (10 Business Days)
 *Date Needed _____

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

REPORTING INSTRUCTIONS
 HARD COPY REQUIRED FAX (FAX#) _____
 PDF (e-mail address) **sdonnell@credere.com**
 OTHER (specify) _____

SPECIAL INSTRUCTIONS
BROWN FIELDS

Relinquished by Sampler: [Signature] Date: 7/14/11 Time: 2:00
 Relinquished by: [Signature] Date: 7/14/11 Time: 2:00
 Relinquished by: [Signature] Date: 7/14/11 Time: 2:00
 Received by Laboratory: [Signature] Date: 7/14/11 Time: 2:00
 May Bill#:

CUSTODY RECORD
 GSD-01 Revision 12/23/10

RECEIVED ON ICE YES NO
 TEMPERATURE _____ °C



EMSL Analytical, Inc.

7 Constitution Way, Suite 107, Woburn, MA 01801

Phone: (781) 933-8411 Fax: (781) 933-8412 Email: bostonlab@emsl.com

Attn: **Jonathan O'Donnell**
Crede Associates, LLC
776 Main Street
Westbrook, ME 04092

Customer ID: CRED25
Customer PO:
Received: 07/01/11 12:35 PM
EMSL Order: 131102954

Fax: (207) 887-1051 Phone: (204) 828-1272
Project: **10001087**

EMSL Proj:
Analysis Date: 7/9/2011


Test Report: Asbestos Analysis of Bulk Materials via EPA 600/R-93/116 Method using Polarized Light Microscopy

Sample	Description	Appearance	Non-Asbestos		Asbestos
			% Fibrous	% Non-Fibrous	% Type
D-1-A 131102954-0001	- Blue Floor Tile	Blue Non-Fibrous Homogeneous		95% Non-fibrous (other)	5% Chrysotile
D-1-B 131102954-0002	- Blue Floor Tile				Stop Positive (Not Analyzed)
D-1-C 131102954-0003	- Blue Floor Tile				Stop Positive (Not Analyzed)
D-2-A 131102954-0004	- White Speckled Floor Tile	White Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-2-B 131102954-0005	- White Speckled Floor Tile	White Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-2-C 131102954-0006	- White Speckled Floor Tile	White Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-3-A 131102954-0007	- Gray Floor Tile	Gray Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected

Initial report from 07/11/2011 12:12:20

Analyst(s)

Kevin Pine (10)



Renaldo Drakes, Laboratory Manager
or other approved signatory

Due to magnification limitations inherent in PLM, asbestos fibers in dimensions below the resolution capability of PLM may not be detected. Samples reported as <1% or none detected require additional testing by TEM to confirm asbestos quantities. The above test report relates only to the items tested and may not be reproduced in any form without the express approval of EMSL Analytical, Inc. EMSL's liability is limited to the cost of analysis. EMSL bears no responsibility for sample collection activities or analytical method limitations. In and use of test results are the responsibility of the client. Samples received in good condition unless otherwise noted.
Samples analyzed by EMSL Analytical, Inc. Woburn, MA NVLAP Lab Code 101147-0, CT PH-0315, MA AA000188, RI AAL-107T3 and VT AL357102



EMSL Analytical, Inc.

7 Constitution Way, Suite 107, Woburn, MA 01801

Phone: (781) 933-8411 Fax: (781) 933-8412 Email: bostonlab@emsl.com

Attn: **Jonathan O'Donnell**
Crede Associates, LLC
776 Main Street
Westbrook, ME 04092

Customer ID: CRED25
Customer PO:
Received: 07/01/11 12:35 PM
EMSL Order: 131102954

Fax: (207) 887-1051 Phone: (204) 828-1272
Project: **10001087**

EMSL Proj:
Analysis Date: 7/9/2011


Test Report: Asbestos Analysis of Bulk Materials via EPA 600/R-93/116 Method using Polarized Light Microscopy

Sample	Description	Appearance	Non-Asbestos		Asbestos
			% Fibrous	% Non-Fibrous	% Type
D-3-B <i>131102954-0008</i>	- Gray Floor Tile	Gray Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-3-C <i>131102954-0009</i>	- Gray Floor Tile	Gray Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-4-A <i>131102954-0010</i>	- Off White Floor Tile	White Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-4-B <i>131102954-0011</i>	- Off White Floor Tile	White Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected
D-4-C <i>131102954-0012</i>	- Off White Floor Tile	White Non-Fibrous Homogeneous		100% Non-fibrous (other)	None Detected

Initial report from 07/11/2011 12:12:20

Analyst(s)

Kevin Pine (10)



Renaldo Drakes, Laboratory Manager
or other approved signatory

Due to magnification limitations inherent in PLM, asbestos fibers in dimensions below the resolution capability of PLM may not be detected. Samples reported as <1% or none detected require additional testing by TEM to confirm asbestos quantities. The above test report relates only to the items tested and may not be reproduced in any form without the express approval of EMSL Analytical, Inc. EMSL's liability is limited to the cost of analysis. EMSL bears no responsibility for sample collection activities or analytical method limitations. In and use of test results are the responsibility of the client. Samples received in good condition unless otherwise noted.
Samples analyzed by EMSL Analytical, Inc. Woburn, MA NVLAP Lab Code 101147-0, CT PH-0315, MA AA000188, RI AAL-107T3 and VT AL357102



ANALYTICAL REPORT

Lab Number:	L1109811
Client:	Crede Associates, LLC 222 St. John Street Suite 314 Portland, ME 04102
ATTN:	Rick Vandenberg
Phone:	(207) 828-1272
Project Name:	POLYCLAD
Project Number:	10001087
Report Date:	07/08/11

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA030), NY (11627), CT (PH-0141), NH (2206), NJ (MA015), RI (LAO00299), ME (MA0030), PA (Registration #68-02089), LA NELAC (03090), FL NELAC (E87814), US Army Corps of Engineers.

320 Forbes Boulevard, Mansfield, MA 02048-1806
508-822-9300 (Fax) 508-822-3288 800-624-9220 - www.alphalab.com



Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Alpha Sample ID	Client ID	Sample Location	Collection Date/Time
L1109811-01	CA-SG-2	FRANKLIN, NH	06/28/11 14:30
L1109811-02	CA-SG-10	FRANKLIN, NH	06/28/11 14:40
L1109811-03	CA-SG-DUP	FRANKLIN, NH	06/28/11 14:30
L1109811-04	BLANK	FRANKLIN, NH	06/28/11 00:00

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

For additional information, please contact Client Services at 800-624-9220.

Volatile Organics in Air

The canister certification results are provided as an addendum.

L1109811-04 The RPD of the pre- and post-flow controller calibration check (46% RPD) was outside acceptable limits (< or = 20% RPD).

L1109811-01 through -03 have elevated detection limits due to the dilution required by the elevated concentrations of target compounds in the samples.

L1109811-01 through -03 The presence of Propylene could not be determined in these samples due to non-target compounds interfering with the identification and quantification of this compound.

Project Name: POLYCLAD
Project Number: 10001087


Lab Number: L1109811
Report Date: 07/08/11

Case Narrative (continued)

The WG477447-3 LCS recoveries for Freon-114 (61%) and trans-1,3-Dichloropropene (68%) are outside the 70%-130% acceptance limit. The LCS was within overall method allowances, therefore the analysis proceeded.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:



Kathleen O'Brien

Title: Technical Director/Representative

Date: 07/08/11

AIR

Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-01 D
 Client ID: CA-SG-2
 Sample Location: FRANKLIN, NH
 Matrix: Soil_Vapor
 Analytical Method: 48,TO-15
 Analytical Date: 07/07/11 23:01
 Analyst: RY

Date Collected: 06/28/11 14:30
 Date Received: 06/30/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Propylene	ND	5.00	--	ND	8.60	--		10
Dichlorodifluoromethane	47.6	2.00	--	235	9.89	--		10
Chloromethane	ND	2.00	--	ND	4.13	--		10
Freon-114	ND	2.00	--	ND	14.0	--		10
Vinyl chloride	ND	2.00	--	ND	5.11	--		10
1,3-Butadiene	ND	2.00	--	ND	4.42	--		10
Bromomethane	ND	2.00	--	ND	7.77	--		10
Chloroethane	ND	2.00	--	ND	5.28	--		10
Ethanol	3740	25.0	--	7050	47.1	--		10
Vinyl bromide	ND	2.00	--	ND	8.74	--		10
Acetone	274	10.0	--	651	23.8	--		10
Trichlorofluoromethane	4.20	2.00	--	23.6	11.2	--		10
Isopropanol	155	5.00	--	381	12.3	--		10
1,1-Dichloroethene	ND	2.00	--	ND	7.93	--		10
Methylene chloride	16.5	10.0	--	57.3	34.7	--		10
3-Chloropropene	ND	2.00	--	ND	6.26	--		10
Carbon disulfide	ND	2.00	--	ND	6.23	--		10
Freon-113	ND	2.00	--	ND	15.3	--		10
trans-1,2-Dichloroethene	ND	2.00	--	ND	7.93	--		10
1,1-Dichloroethane	ND	2.00	--	ND	8.09	--		10
Methyl tert butyl ether	ND	2.00	--	ND	7.21	--		10
Vinyl acetate	ND	2.00	--	ND	7.04	--		10
2-Butanone	10.2	2.00	--	30.1	5.90	--		10



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-01 D

Date Collected: 06/28/11 14:30

Client ID: CA-SG-2

Date Received: 06/30/11

Sample Location: FRANKLIN, NH

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
cis-1,2-Dichloroethene	ND	2.00	--	ND	7.93	--		10
Ethyl Acetate	ND	5.00	--	ND	18.0	--		10
Chloroform	ND	2.00	--	ND	9.77	--		10
Tetrahydrofuran	4.82	2.00	--	14.2	5.90	--		10
1,2-Dichloroethane	ND	2.00	--	ND	8.09	--		10
n-Hexane	8.78	2.00	--	30.9	7.05	--		10
1,1,1-Trichloroethane	ND	2.00	--	ND	10.9	--		10
Benzene	8.88	2.00	--	28.4	6.39	--		10
Carbon tetrachloride	ND	2.00	--	ND	12.6	--		10
Cyclohexane	2.22	2.00	--	7.64	6.88	--		10
1,2-Dichloropropane	ND	2.00	--	ND	9.24	--		10
Bromodichloromethane	ND	2.00	--	ND	13.4	--		10
1,4-Dioxane	ND	2.00	--	ND	7.21	--		10
Trichloroethene	ND	2.00	--	ND	10.7	--		10
2,2,4-Trimethylpentane	ND	2.00	--	ND	9.34	--		10
Heptane	3.85	2.00	--	15.8	8.20	--		10
cis-1,3-Dichloropropene	ND	2.00	--	ND	9.08	--		10
4-Methyl-2-pentanone	ND	2.00	--	ND	8.20	--		10
trans-1,3-Dichloropropene	ND	2.00	--	ND	9.08	--		10
1,1,2-Trichloroethane	ND	2.00	--	ND	10.9	--		10
Toluene	40.9	2.00	--	154	7.54	--		10
2-Hexanone	ND	2.00	--	ND	8.20	--		10
Dibromochloromethane	ND	2.00	--	ND	17.0	--		10
1,2-Dibromoethane	ND	2.00	--	ND	15.4	--		10
Tetrachloroethene	ND	2.00	--	ND	13.6	--		10
Chlorobenzene	ND	2.00	--	ND	9.21	--		10



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-01 D

Date Collected: 06/28/11 14:30

Client ID: CA-SG-2

Date Received: 06/30/11

Sample Location: FRANKLIN, NH

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Ethylbenzene	4.60	2.00	--	20.0	8.69	--		10
p/m-Xylene	15.0	4.00	--	65.2	17.4	--		10
Bromoform	ND	2.00	--	ND	20.7	--		10
Styrene	ND	2.00	--	ND	8.52	--		10
1,1,2,2-Tetrachloroethane	ND	2.00	--	ND	13.7	--		10
o-Xylene	4.13	2.00	--	17.9	8.69	--		10
4-Ethyltoluene	ND	2.00	--	ND	9.83	--		10
1,3,5-Trimethylbenzene	ND	2.00	--	ND	9.83	--		10
1,2,4-Trimethylbenzene	2.59	2.00	--	12.7	9.83	--		10
Benzyl chloride	ND	2.00	--	ND	10.4	--		10
1,3-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,4-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,2-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,2,4-Trichlorobenzene	ND	2.00	--	ND	14.8	--		10
Hexachlorobutadiene	ND	2.00	--	ND	21.3	--		10

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	93		60-140
Bromochloromethane	92		60-140
chlorobenzene-d5	95		60-140



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-02 D
 Client ID: CA-SG-10
 Sample Location: FRANKLIN, NH
 Matrix: Soil_Vapor
 Analytical Method: 48,TO-15
 Analytical Date: 07/07/11 23:37
 Analyst: RY

Date Collected: 06/28/11 14:40
 Date Received: 06/30/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Propylene	ND	5.00	--	ND	8.60	--		10
Dichlorodifluoromethane	3.24	2.00	--	16.0	9.89	--		10
Chloromethane	ND	2.00	--	ND	4.13	--		10
Freon-114	ND	2.00	--	ND	14.0	--		10
Vinyl chloride	ND	2.00	--	ND	5.11	--		10
1,3-Butadiene	ND	2.00	--	ND	4.42	--		10
Bromomethane	ND	2.00	--	ND	7.77	--		10
Chloroethane	ND	2.00	--	ND	5.28	--		10
Ethanol	3060	25.0	--	5760	47.1	--		10
Vinyl bromide	ND	2.00	--	ND	8.74	--		10
Acetone	89.0	10.0	--	211	23.8	--		10
Trichlorofluoromethane	18.9	2.00	--	106	11.2	--		10
Isopropanol	120	5.00	--	295	12.3	--		10
1,1-Dichloroethene	ND	2.00	--	ND	7.93	--		10
Methylene chloride	ND	10.0	--	ND	34.7	--		10
3-Chloropropene	ND	2.00	--	ND	6.26	--		10
Carbon disulfide	ND	2.00	--	ND	6.23	--		10
Freon-113	ND	2.00	--	ND	15.3	--		10
trans-1,2-Dichloroethene	ND	2.00	--	ND	7.93	--		10
1,1-Dichloroethane	ND	2.00	--	ND	8.09	--		10
Methyl tert butyl ether	ND	2.00	--	ND	7.21	--		10
Vinyl acetate	ND	2.00	--	ND	7.04	--		10
2-Butanone	11.1	2.00	--	32.7	5.90	--		10



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-02 D

Date Collected: 06/28/11 14:40

Client ID: CA-SG-10

Date Received: 06/30/11

Sample Location: FRANKLIN, NH

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
cis-1,2-Dichloroethene	ND	2.00	--	ND	7.93	--		10
Ethyl Acetate	ND	5.00	--	ND	18.0	--		10
Chloroform	ND	2.00	--	ND	9.77	--		10
Tetrahydrofuran	5.57	2.00	--	16.4	5.90	--		10
1,2-Dichloroethane	ND	2.00	--	ND	8.09	--		10
n-Hexane	7.17	2.00	--	25.3	7.05	--		10
1,1,1-Trichloroethane	ND	2.00	--	ND	10.9	--		10
Benzene	9.38	2.00	--	30.0	6.39	--		10
Carbon tetrachloride	ND	2.00	--	ND	12.6	--		10
Cyclohexane	2.24	2.00	--	7.71	6.88	--		10
1,2-Dichloropropane	ND	2.00	--	ND	9.24	--		10
Bromodichloromethane	ND	2.00	--	ND	13.4	--		10
1,4-Dioxane	ND	2.00	--	ND	7.21	--		10
Trichloroethene	ND	2.00	--	ND	10.7	--		10
2,2,4-Trimethylpentane	ND	2.00	--	ND	9.34	--		10
Heptane	4.67	2.00	--	19.1	8.20	--		10
cis-1,3-Dichloropropene	ND	2.00	--	ND	9.08	--		10
4-Methyl-2-pentanone	ND	2.00	--	ND	8.20	--		10
trans-1,3-Dichloropropene	ND	2.00	--	ND	9.08	--		10
1,1,2-Trichloroethane	ND	2.00	--	ND	10.9	--		10
Toluene	47.0	2.00	--	177	7.54	--		10
2-Hexanone	ND	2.00	--	ND	8.20	--		10
Dibromochloromethane	ND	2.00	--	ND	17.0	--		10
1,2-Dibromoethane	ND	2.00	--	ND	15.4	--		10
Tetrachloroethene	ND	2.00	--	ND	13.6	--		10
Chlorobenzene	ND	2.00	--	ND	9.21	--		10



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-02 D

Date Collected: 06/28/11 14:40

Client ID: CA-SG-10

Date Received: 06/30/11

Sample Location: FRANKLIN, NH

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Ethylbenzene	3.97	2.00	--	17.2	8.69	--		10
p/m-Xylene	12.7	4.00	--	55.2	17.4	--		10
Bromoform	ND	2.00	--	ND	20.7	--		10
Styrene	ND	2.00	--	ND	8.52	--		10
1,1,2,2-Tetrachloroethane	ND	2.00	--	ND	13.7	--		10
o-Xylene	3.26	2.00	--	14.2	8.69	--		10
4-Ethyltoluene	ND	2.00	--	ND	9.83	--		10
1,3,5-Trimethylbenzene	ND	2.00	--	ND	9.83	--		10
1,2,4-Trimethylbenzene	ND	2.00	--	ND	9.83	--		10
Benzyl chloride	ND	2.00	--	ND	10.4	--		10
1,3-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,4-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,2-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,2,4-Trichlorobenzene	ND	2.00	--	ND	14.8	--		10
Hexachlorobutadiene	ND	2.00	--	ND	21.3	--		10

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	89		60-140
Bromochloromethane	89		60-140
chlorobenzene-d5	91		60-140



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-03 D
 Client ID: CA-SG-DUP
 Sample Location: FRANKLIN, NH
 Matrix: Soil_Vapor
 Analytical Method: 48,TO-15
 Analytical Date: 07/08/11 00:14
 Analyst: RY

Date Collected: 06/28/11 14:30
 Date Received: 06/30/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Propylene	ND	5.00	--	ND	8.60	--		10
Dichlorodifluoromethane	54.1	2.00	--	268	9.89	--		10
Chloromethane	ND	2.00	--	ND	4.13	--		10
Freon-114	ND	2.00	--	ND	14.0	--		10
Vinyl chloride	ND	2.00	--	ND	5.11	--		10
1,3-Butadiene	ND	2.00	--	ND	4.42	--		10
Bromomethane	ND	2.00	--	ND	7.77	--		10
Chloroethane	ND	2.00	--	ND	5.28	--		10
Ethanol	4080	25.0	--	7690	47.1	--		10
Vinyl bromide	ND	2.00	--	ND	8.74	--		10
Acetone	282	10.0	--	670	23.8	--		10
Trichlorofluoromethane	4.82	2.00	--	27.1	11.2	--		10
Isopropanol	185	5.00	--	455	12.3	--		10
1,1-Dichloroethene	ND	2.00	--	ND	7.93	--		10
Methylene chloride	ND	10.0	--	ND	34.7	--		10
3-Chloropropene	ND	2.00	--	ND	6.26	--		10
Carbon disulfide	ND	2.00	--	ND	6.23	--		10
Freon-113	ND	2.00	--	ND	15.3	--		10
trans-1,2-Dichloroethene	ND	2.00	--	ND	7.93	--		10
1,1-Dichloroethane	ND	2.00	--	ND	8.09	--		10
Methyl tert butyl ether	ND	2.00	--	ND	7.21	--		10
Vinyl acetate	ND	2.00	--	ND	7.04	--		10
2-Butanone	10.3	2.00	--	30.4	5.90	--		10



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-03 D

Date Collected: 06/28/11 14:30

Client ID: CA-SG-DUP

Date Received: 06/30/11

Sample Location: FRANKLIN, NH

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
cis-1,2-Dichloroethene	ND	2.00	--	ND	7.93	--		10
Ethyl Acetate	ND	5.00	--	ND	18.0	--		10
Chloroform	ND	2.00	--	ND	9.77	--		10
Tetrahydrofuran	4.94	2.00	--	14.6	5.90	--		10
1,2-Dichloroethane	ND	2.00	--	ND	8.09	--		10
n-Hexane	9.28	2.00	--	32.7	7.05	--		10
1,1,1-Trichloroethane	ND	2.00	--	ND	10.9	--		10
Benzene	9.43	2.00	--	30.1	6.39	--		10
Carbon tetrachloride	ND	2.00	--	ND	12.6	--		10
Cyclohexane	2.59	2.00	--	8.92	6.88	--		10
1,2-Dichloropropane	ND	2.00	--	ND	9.24	--		10
Bromodichloromethane	ND	2.00	--	ND	13.4	--		10
1,4-Dioxane	ND	2.00	--	ND	7.21	--		10
Trichloroethene	ND	2.00	--	ND	10.7	--		10
2,2,4-Trimethylpentane	ND	2.00	--	ND	9.34	--		10
Heptane	4.57	2.00	--	18.7	8.20	--		10
cis-1,3-Dichloropropene	ND	2.00	--	ND	9.08	--		10
4-Methyl-2-pentanone	ND	2.00	--	ND	8.20	--		10
trans-1,3-Dichloropropene	ND	2.00	--	ND	9.08	--		10
1,1,2-Trichloroethane	ND	2.00	--	ND	10.9	--		10
Toluene	40.9	2.00	--	154	7.54	--		10
2-Hexanone	ND	2.00	--	ND	8.20	--		10
Dibromochloromethane	ND	2.00	--	ND	17.0	--		10
1,2-Dibromoethane	ND	2.00	--	ND	15.4	--		10
Tetrachloroethene	ND	2.00	--	ND	13.6	--		10
Chlorobenzene	ND	2.00	--	ND	9.21	--		10



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-03 D

Date Collected: 06/28/11 14:30

Client ID: CA-SG-DUP

Date Received: 06/30/11

Sample Location: FRANKLIN, NH

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Ethylbenzene	3.19	2.00	--	13.8	8.69	--		10
p/m-Xylene	10.4	4.00	--	45.2	17.4	--		10
Bromoform	ND	2.00	--	ND	20.7	--		10
Styrene	ND	2.00	--	ND	8.52	--		10
1,1,2,2-Tetrachloroethane	ND	2.00	--	ND	13.7	--		10
o-Xylene	2.68	2.00	--	11.6	8.69	--		10
4-Ethyltoluene	ND	2.00	--	ND	9.83	--		10
1,3,5-Trimethylbenzene	ND	2.00	--	ND	9.83	--		10
1,2,4-Trimethylbenzene	ND	2.00	--	ND	9.83	--		10
Benzyl chloride	ND	2.00	--	ND	10.4	--		10
1,3-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,4-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,2-Dichlorobenzene	ND	2.00	--	ND	12.0	--		10
1,2,4-Trichlorobenzene	ND	2.00	--	ND	14.8	--		10
Hexachlorobutadiene	ND	2.00	--	ND	21.3	--		10

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	89		60-140
Bromochloromethane	90		60-140
chlorobenzene-d5	91		60-140



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-04
 Client ID: BLANK
 Sample Location: FRANKLIN, NH
 Matrix: Soil_Vapor
 Analytical Method: 48,TO-15
 Analytical Date: 07/07/11 22:24
 Analyst: RY

Date Collected: 06/28/11 00:00
 Date Received: 06/30/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Propylene	ND	0.500	--	ND	0.860	--		1
Dichlorodifluoromethane	ND	0.200	--	ND	0.989	--		1
Chloromethane	ND	0.200	--	ND	0.413	--		1
Freon-114	ND	0.200	--	ND	1.40	--		1
Vinyl chloride	ND	0.200	--	ND	0.511	--		1
1,3-Butadiene	ND	0.200	--	ND	0.442	--		1
Bromomethane	ND	0.200	--	ND	0.777	--		1
Chloroethane	ND	0.200	--	ND	0.528	--		1
Ethanol	ND	2.50	--	ND	4.71	--		1
Vinyl bromide	ND	0.200	--	ND	0.874	--		1
Acetone	ND	1.00	--	ND	2.38	--		1
Trichlorofluoromethane	ND	0.200	--	ND	1.12	--		1
Isopropanol	ND	0.500	--	ND	1.23	--		1
1,1-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Methylene chloride	ND	1.00	--	ND	3.47	--		1
3-Chloropropene	ND	0.200	--	ND	0.626	--		1
Carbon disulfide	ND	0.200	--	ND	0.623	--		1
Freon-113	ND	0.200	--	ND	1.53	--		1
trans-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
1,1-Dichloroethane	ND	0.200	--	ND	0.809	--		1
Methyl tert butyl ether	ND	0.200	--	ND	0.721	--		1
Vinyl acetate	ND	0.200	--	ND	0.704	--		1
2-Butanone	ND	0.200	--	ND	0.590	--		1



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-04
 Client ID: BLANK
 Sample Location: FRANKLIN, NH

Date Collected: 06/28/11 00:00
 Date Received: 06/30/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
cis-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Ethyl Acetate	ND	0.500	--	ND	1.80	--		1
Chloroform	ND	0.200	--	ND	0.977	--		1
Tetrahydrofuran	ND	0.200	--	ND	0.590	--		1
1,2-Dichloroethane	ND	0.200	--	ND	0.809	--		1
n-Hexane	ND	0.200	--	ND	0.705	--		1
1,1,1-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Benzene	ND	0.200	--	ND	0.639	--		1
Carbon tetrachloride	ND	0.200	--	ND	1.26	--		1
Cyclohexane	ND	0.200	--	ND	0.688	--		1
1,2-Dichloropropane	ND	0.200	--	ND	0.924	--		1
Bromodichloromethane	ND	0.200	--	ND	1.34	--		1
1,4-Dioxane	ND	0.200	--	ND	0.721	--		1
Trichloroethene	ND	0.200	--	ND	1.07	--		1
2,2,4-Trimethylpentane	ND	0.200	--	ND	0.934	--		1
Heptane	ND	0.200	--	ND	0.820	--		1
cis-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
4-Methyl-2-pentanone	ND	0.200	--	ND	0.820	--		1
trans-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
1,1,2-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Toluene	ND	0.200	--	ND	0.754	--		1
2-Hexanone	ND	0.200	--	ND	0.820	--		1
Dibromochloromethane	ND	0.200	--	ND	1.70	--		1
1,2-Dibromoethane	ND	0.200	--	ND	1.54	--		1
Tetrachloroethene	ND	0.200	--	ND	1.36	--		1
Chlorobenzene	ND	0.200	--	ND	0.921	--		1



Project Name: POLYCLAD**Lab Number:** L1109811**Project Number:** 10001087**Report Date:** 07/08/11**SAMPLE RESULTS**

Lab ID: L1109811-04
 Client ID: BLANK
 Sample Location: FRANKLIN, NH

Date Collected: 06/28/11 00:00
 Date Received: 06/30/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Ethylbenzene	ND	0.200	--	ND	0.869	--		1
p/m-Xylene	ND	0.400	--	ND	1.74	--		1
Bromoform	ND	0.200	--	ND	2.07	--		1
Styrene	ND	0.200	--	ND	0.852	--		1
1,1,2,2-Tetrachloroethane	ND	0.200	--	ND	1.37	--		1
o-Xylene	ND	0.200	--	ND	0.869	--		1
4-Ethyltoluene	ND	0.200	--	ND	0.983	--		1
1,3,5-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
1,2,4-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
Benzyl chloride	ND	0.200	--	ND	1.04	--		1
1,3-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,4-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2,4-Trichlorobenzene	ND	0.200	--	ND	1.48	--		1
Hexachlorobutadiene	ND	0.200	--	ND	2.13	--		1

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	83		60-140
Bromochloromethane	85		60-140
chlorobenzene-d5	86		60-140



Project Name: POLYCLAD

Lab Number: L1109811

Project Number: 10001087

Report Date: 07/08/11

Method Blank Analysis Batch Quality Control

Analytical Method: 48,TO-15

Analytical Date: 07/07/11 19:52

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab for sample(s): 01-04 Batch: WG477447-4								
Propylene	ND	0.500	--	ND	0.860	--		1
Dichlorodifluoromethane	ND	0.200	--	ND	0.989	--		1
Chloromethane	ND	0.200	--	ND	0.413	--		1
Freon-114	ND	0.200	--	ND	1.40	--		1
Vinyl chloride	ND	0.200	--	ND	0.511	--		1
1,3-Butadiene	ND	0.200	--	ND	0.442	--		1
Bromomethane	ND	0.200	--	ND	0.777	--		1
Chloroethane	ND	0.200	--	ND	0.528	--		1
Ethanol	ND	2.50	--	ND	4.71	--		1
Vinyl bromide	ND	0.200	--	ND	0.874	--		1
Acetone	ND	1.00	--	ND	2.38	--		1
Trichlorofluoromethane	ND	0.200	--	ND	1.12	--		1
Isopropanol	ND	0.500	--	ND	1.23	--		1
1,1-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Methylene chloride	ND	1.00	--	ND	3.47	--		1
3-Chloropropene	ND	0.200	--	ND	0.626	--		1
Carbon disulfide	ND	0.200	--	ND	0.623	--		1
Freon-113	ND	0.200	--	ND	1.53	--		1
trans-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
1,1-Dichloroethane	ND	0.200	--	ND	0.809	--		1
Methyl tert butyl ether	ND	0.200	--	ND	0.721	--		1
Vinyl acetate	ND	0.200	--	ND	0.704	--		1
2-Butanone	ND	0.200	--	ND	0.590	--		1
cis-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Ethyl Acetate	ND	0.500	--	ND	1.80	--		1



Project Name: POLYCLAD

Lab Number: L1109811

Project Number: 10001087

Report Date: 07/08/11

Method Blank Analysis Batch Quality Control

Analytical Method: 48,TO-15

Analytical Date: 07/07/11 19:52

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab for sample(s): 01-04 Batch: WG477447-4								
Chloroform	ND	0.200	--	ND	0.977	--		1
Tetrahydrofuran	ND	0.200	--	ND	0.590	--		1
1,2-Dichloroethane	ND	0.200	--	ND	0.809	--		1
n-Hexane	ND	0.200	--	ND	0.705	--		1
1,1,1-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Benzene	ND	0.200	--	ND	0.639	--		1
Carbon tetrachloride	ND	0.200	--	ND	1.26	--		1
Cyclohexane	ND	0.200	--	ND	0.688	--		1
1,2-Dichloropropane	ND	0.200	--	ND	0.924	--		1
Bromodichloromethane	ND	0.200	--	ND	1.34	--		1
1,4-Dioxane	ND	0.200	--	ND	0.721	--		1
Trichloroethene	ND	0.200	--	ND	1.07	--		1
2,2,4-Trimethylpentane	ND	0.200	--	ND	0.934	--		1
Heptane	ND	0.200	--	ND	0.820	--		1
cis-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
4-Methyl-2-pentanone	ND	0.200	--	ND	0.820	--		1
trans-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
1,1,2-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Toluene	ND	0.200	--	ND	0.754	--		1
2-Hexanone	ND	0.200	--	ND	0.820	--		1
Dibromochloromethane	ND	0.200	--	ND	1.70	--		1
1,2-Dibromoethane	ND	0.200	--	ND	1.54	--		1
Tetrachloroethene	ND	0.200	--	ND	1.36	--		1
Chlorobenzene	ND	0.200	--	ND	0.921	--		1
Ethylbenzene	ND	0.200	--	ND	0.869	--		1



Project Name: POLYCLAD

Lab Number: L1109811

Project Number: 10001087

Report Date: 07/08/11

Method Blank Analysis Batch Quality Control

Analytical Method: 48,TO-15

Analytical Date: 07/07/11 19:52

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab for sample(s): 01-04 Batch: WG477447-4								
p/m-Xylene	ND	0.400	--	ND	1.74	--		1
Bromoform	ND	0.200	--	ND	2.07	--		1
Styrene	ND	0.200	--	ND	0.852	--		1
1,1,2,2-Tetrachloroethane	ND	0.200	--	ND	1.37	--		1
o-Xylene	ND	0.200	--	ND	0.869	--		1
4-Ethyltoluene	ND	0.200	--	ND	0.983	--		1
1,3,5-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
1,2,4-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
Benzyl chloride	ND	0.200	--	ND	1.04	--		1
1,3-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,4-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2,4-Trichlorobenzene	ND	0.200	--	ND	1.48	--		1
Hexachlorobutadiene	ND	0.200	--	ND	2.13	--		1

Lab Control Sample Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	LCS		LCSD		%Recovery		RPD	Qual	RPD Limits
	%Recovery	Qual	%Recovery	Qual	%Recovery	Limits			
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 Batch: WG477447-3									
Propylene	111	-	-	-	70-130	-	-	-	70-130
Dichlorodifluoromethane	102	-	-	-	70-130	-	-	-	70-130
Chloromethane	98	-	-	-	70-130	-	-	-	70-130
1,2-Dichloro-1,1,2,2-tetrafluoroethane	61	Q	-	-	70-130	-	-	-	70-130
Vinyl chloride	101	-	-	-	70-130	-	-	-	70-130
1,3-Butadiene	101	-	-	-	70-130	-	-	-	70-130
Bromomethane	102	-	-	-	70-130	-	-	-	70-130
Chloroethane	100	-	-	-	70-130	-	-	-	70-130
Ethyl Alcohol	91	-	-	-	70-130	-	-	-	70-130
Vinyl bromide	102	-	-	-	70-130	-	-	-	70-130
Acetone	85	-	-	-	70-130	-	-	-	70-130
Trichlorofluoromethane	106	-	-	-	70-130	-	-	-	70-130
iso-Propyl Alcohol	97	-	-	-	70-130	-	-	-	70-130
1,1-Dichloroethene	98	-	-	-	70-130	-	-	-	70-130
Methylene chloride	90	-	-	-	70-130	-	-	-	70-130
3-Chloropropene	88	-	-	-	70-130	-	-	-	70-130
Carbon disulfide	94	-	-	-	70-130	-	-	-	70-130
1,1,2-Trichloro-1,2,2-Trifluoroethane	105	-	-	-	70-130	-	-	-	70-130
trans-1,2-Dichloroethene	96	-	-	-	70-130	-	-	-	70-130
1,1-Dichloroethane	93	-	-	-	70-130	-	-	-	70-130
Methyl tert butyl ether	85	-	-	-	70-130	-	-	-	70-130

Lab Control Sample Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	LCS		LCSD		%Recovery		RPD	Qual	RPD Limits
	%Recovery	Qual	%Recovery	Qual	%Recovery	Limits			
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 Batch: WG477447-3									
Vinyl acetate	85	-	-	-	70-130	-	-	-	70-130
2-Butanone	90	-	-	-	70-130	-	-	-	70-130
cis-1,2-Dichloroethene	97	-	-	-	70-130	-	-	-	70-130
Ethyl Acetate	88	-	-	-	70-130	-	-	-	70-130
Chloroform	100	-	-	-	70-130	-	-	-	70-130
Tetrahydrofuran	80	-	-	-	70-130	-	-	-	70-130
1,2-Dichloroethane	95	-	-	-	70-130	-	-	-	70-130
n-Hexane	81	-	-	-	70-130	-	-	-	70-130
1,1,1-Trichloroethane	94	-	-	-	70-130	-	-	-	70-130
Benzene	90	-	-	-	70-130	-	-	-	70-130
Carbon tetrachloride	92	-	-	-	70-130	-	-	-	70-130
Cyclohexane	84	-	-	-	70-130	-	-	-	70-130
1,2-Dichloropropane	80	-	-	-	70-130	-	-	-	70-130
Bromodichloromethane	81	-	-	-	70-130	-	-	-	70-130
1,4-Dioxane	86	-	-	-	70-130	-	-	-	70-130
Trichloroethene	90	-	-	-	70-130	-	-	-	70-130
2,2,4-Trimethylpentane	82	-	-	-	70-130	-	-	-	70-130
Heptane	82	-	-	-	70-130	-	-	-	70-130
cis-1,3-Dichloropropene	83	-	-	-	70-130	-	-	-	70-130
4-Methyl-2-pentanone	85	-	-	-	70-130	-	-	-	70-130
trans-1,3-Dichloropropene	68	-	-	-	70-130	-	-	Q	70-130



Lab Control Sample Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	LCS		LCSD		%Recovery		RPD	Qual	RPD Limits
	%Recovery	Qual	%Recovery	Qual	%Recovery	Limits			
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 Batch: WG477447-3									
1,1,2-Trichloroethane	83	-	-	-	70-130	-	-	-	70-130
Toluene	88	-	-	-	70-130	-	-	-	70-130
2-Hexanone	100	-	-	-	70-130	-	-	-	70-130
Dibromochloromethane	93	-	-	-	70-130	-	-	-	70-130
1,2-Dibromoethane	96	-	-	-	70-130	-	-	-	70-130
Tetrachloroethene	100	-	-	-	70-130	-	-	-	70-130
Chlorobenzene	93	-	-	-	70-130	-	-	-	70-130
Ethylbenzene	87	-	-	-	70-130	-	-	-	70-130
p/m-Xylene	87	-	-	-	70-130	-	-	-	70-130
Bromoform	90	-	-	-	70-130	-	-	-	70-130
Styrene	94	-	-	-	70-130	-	-	-	70-130
1,1,2,2-Tetrachloroethane	94	-	-	-	70-130	-	-	-	70-130
o-Xylene	89	-	-	-	70-130	-	-	-	70-130
4-Ethyltoluene	92	-	-	-	70-130	-	-	-	70-130
1,3,5-Trimethylbenzene	94	-	-	-	70-130	-	-	-	70-130
1,2,4-Trimethylbenzene	102	-	-	-	70-130	-	-	-	70-130
Benzyl chloride	85	-	-	-	70-130	-	-	-	70-130
1,3-Dichlorobenzene	98	-	-	-	70-130	-	-	-	70-130
1,4-Dichlorobenzene	99	-	-	-	70-130	-	-	-	70-130
1,2-Dichlorobenzene	102	-	-	-	70-130	-	-	-	70-130
1,2,4-Trichlorobenzene	124	-	-	-	70-130	-	-	-	70-130



Lab Control Sample Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	LCS %Recovery	Qual	LCS %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 Batch: WG477447-3								
Hexachlorobutadiene	103	-	-	-	70-130	-	-	-



Lab Duplicate Analysis Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	Native Sample	Duplicate Sample	Units	RPD	Qual	RPD Limits
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 QC Batch ID: WG477447-5 QC Sample: L1109857-03 Client ID: DUP Sample						
Propylene	44.3	44.5	ppbV	0		25
Dichlorodifluoromethane	ND	ND	ppbV	NC		25
Chloromethane	1.09	1.09	ppbV	0		25
1,2-Dichloro-1,1,2,2-tetrafluoroethane	1.42	1.16	ppbV	20		25
Vinyl chloride	1.52	1.48	ppbV	3		25
1,3-Butadiene	ND	ND	ppbV	NC		25
Bromomethane	ND	ND	ppbV	NC		25
Chloroethane	ND	ND	ppbV	NC		25
Ethyl Alcohol	14.7	14.3	ppbV	3		25
Vinyl bromide	ND	ND	ppbV	NC		25
Acetone	119	121	ppbV	2		25
Trichlorofluoromethane	ND	ND	ppbV	NC		25
iso-Propyl Alcohol	ND	ND	ppbV	NC		25
1,1-Dichloroethene	ND	ND	ppbV	NC		25
Methylene chloride	ND	ND	ppbV	NC		25
3-Chloropropene	ND	ND	ppbV	NC		25
Carbon disulfide	1.28	1.38	ppbV	8		25
1,1,2-Trichloro-1,2,2-Trifluoroethane	ND	ND	ppbV	NC		25
trans-1,2-Dichloroethene	ND	ND	ppbV	NC		25



Lab Duplicate Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	Native Sample	Duplicate Sample	Units	RPD	RPD Limits
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 QC Batch ID: WG477447-5 QC Sample: L1109857-03 Client ID: DUP Sample					
1,1-Dichloroethane	ND	ND	ppbV	NC	25
Methyl tert butyl ether	ND	ND	ppbV	NC	25
Vinyl acetate	ND	ND	ppbV	NC	25
2-Butanone	11.2	11.2	ppbV	0	25
cis-1,2-Dichloroethene	1.66	1.66	ppbV	0	25
Ethyl Acetate	ND	ND	ppbV	NC	25
Chloroform	ND	ND	ppbV	NC	25
Tetrahydrofuran	ND	ND	ppbV	NC	25
1,2-Dichloroethane	ND	ND	ppbV	NC	25
n-Hexane	29.8	29.9	ppbV	0	25
1,1,1-Trichloroethane	ND	ND	ppbV	NC	25
Benzene	4.30	4.48	ppbV	4	25
Carbon tetrachloride	ND	ND	ppbV	NC	25
Cyclohexane	13.9	14.4	ppbV	4	25
1,2-Dichloropropane	ND	ND	ppbV	NC	25
Bromodichloromethane	ND	ND	ppbV	NC	25
1,4-Dioxane	ND	ND	ppbV	NC	25
Trichloroethene	ND	ND	ppbV	NC	25
2,2,4-Trimethylpentane	42.9	43.5	ppbV	1	25



Lab Duplicate Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	Native Sample	Duplicate Sample	Units	RPD	RPD Limits
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 QC Batch ID: WG477447-5 QC Sample: L1109857-03 Client ID: DUP Sample					
Heptane	8.78	8.87	ppbV	1	25
cis-1,3-Dichloropropene	ND	ND	ppbV	NC	25
4-Methyl-2-pentanone	ND	ND	ppbV	NC	25
trans-1,3-Dichloropropene	ND	ND	ppbV	NC	25
1,1,2-Trichloroethane	ND	ND	ppbV	NC	25
Toluene	2.54	2.49	ppbV	2	25
2-Hexanone	ND	ND	ppbV	NC	25
Dibromochloromethane	ND	ND	ppbV	NC	25
1,2-Dibromoethane	ND	ND	ppbV	NC	25
Tetrachloroethene	ND	ND	ppbV	NC	25
Chlorobenzene	ND	ND	ppbV	NC	25
Ethylbenzene	ND	ND	ppbV	NC	25
p/m-Xylene	ND	ND	ppbV	NC	25
Bromoform	ND	ND	ppbV	NC	25
Styrene	ND	ND	ppbV	NC	25
1,1,2,2-Tetrachloroethane	ND	ND	ppbV	NC	25
o-Xylene	ND	ND	ppbV	NC	25
4-Ethyltoluene	ND	ND	ppbV	NC	25
1,3,5-Trimethylbenzene	ND	ND	ppbV	NC	25



Lab Duplicate Analysis

Batch Quality Control

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Parameter	Native Sample	Duplicate Sample	Units	RPD	RPD Limits
Volatile Organics in Air (Low Level) - Mansfield Lab Associated sample(s): 01-04 QC Batch ID: WG477447-5 QC Sample: L1109857-03 Client ID: DUP					
Sample					
1,2,4-Trimethylbenzene	ND	ND	ppbV	NC	25
Benzyl chloride	ND	ND	ppbV	NC	25
1,3-Dichlorobenzene	ND	ND	ppbV	NC	25
1,4-Dichlorobenzene	ND	ND	ppbV	NC	25
1,2-Dichlorobenzene	ND	ND	ppbV	NC	25
1,2,4-Trichlorobenzene	ND	ND	ppbV	NC	25
Hexachlorobutadiene	ND	ND	ppbV	NC	25



Project Name: POLYCLAD

Project Number: 10001087

Serial_No:07081115:31
Lab Number: L1109811

Report Date: 07/08/11

Canister and Flow Controller Information

Samplenum	Client ID	Media ID	Media Type	Cleaning Batch ID	Initial Pressure (in. Hg)	Pressure on Receipt (in. Hg)	Flow Out mL/min	Flow In mL/min	% RSD
L1109811-01	CA-SG-2	0458	#30 SV		-	-	34	34	0
L1109811-01	CA-SG-2	469	2.7L Can	L1107923	-29.4	-4.2	-	-	-
L1109811-02	CA-SG-10	0412	#90 SV		-	-	36	39	8
L1109811-02	CA-SG-10	524	2.7L Can	L1107923	-29.4	-1.6	-	-	-
L1109811-03	CA-SG-DUP	0141	#16 AMB		-	-	33	33	0
L1109811-03	CA-SG-DUP	530	2.7L Can	L1107923	-29.4	-4.0	-	-	-
L1109811-04	BLANK	0293	#90 SV		-	-	32	20	46
L1109811-04	BLANK	554	2.7L Can	L1107923	-29.2	-29.2	-	-	-



Air Volatiles Can Certification

Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01
 Client ID: CAN 131 SHELF 2
 Sample Location:
 Matrix: Air
 Analytical Method: 48,TO-15
 Analytical Date: 06/15/11 16:38
 Analyst: RY

Date Collected: 06/03/11 00:00
 Date Received: 06/03/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Chlorodifluoromethane	ND	0.200	--	ND	0.707	--		1
Propylene	ND	0.500	--	ND	0.860	--		1
Propane	ND	0.200	--	ND	0.361	--		1
Dichlorodifluoromethane	ND	0.200	--	ND	0.989	--		1
Chloromethane	ND	0.200	--	ND	0.413	--		1
Freon-114	ND	0.200	--	ND	1.40	--		1
Methanol	ND	5.00	--	ND	6.55	--		1
Vinyl chloride	ND	0.200	--	ND	0.511	--		1
1,3-Butadiene	ND	0.200	--	ND	0.442	--		1
Butane	ND	0.200	--	ND	0.475	--		1
Bromomethane	ND	0.200	--	ND	0.777	--		1
Chloroethane	ND	0.200	--	ND	0.528	--		1
Ethanol	ND	2.50	--	ND	4.71	--		1
Dichlorofluoromethane	ND	0.200	--	ND	0.842	--		1
Vinyl bromide	ND	0.200	--	ND	0.874	--		1
Acrolein	ND	0.500	--	ND	1.15	--		1
Acetone	ND	1.00	--	ND	2.38	--		1
Acetonitrile	ND	0.200	--	ND	0.336	--		1
Trichlorofluoromethane	ND	0.200	--	ND	1.12	--		1
Isopropanol	ND	0.500	--	ND	1.23	--		1
Acrylonitrile	ND	0.200	--	ND	0.434	--		1
Pentane	ND	0.200	--	ND	0.590	--		1
Ethyl ether	ND	0.200	--	ND	0.606	--		1
1,1-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Tertiary butyl Alcohol	ND	0.500	--	ND	1.52	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Methylene chloride	ND	1.00	--	ND	3.47	--		1
3-Chloropropene	ND	0.200	--	ND	0.626	--		1
Carbon disulfide	ND	0.200	--	ND	0.623	--		1
Freon-113	ND	0.200	--	ND	1.53	--		1
trans-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
1,1-Dichloroethane	ND	0.200	--	ND	0.809	--		1
Methyl tert butyl ether	ND	0.200	--	ND	0.721	--		1
Vinyl acetate	ND	0.200	--	ND	0.704	--		1
2-Butanone	ND	0.200	--	ND	0.590	--		1
cis-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Ethyl Acetate	ND	0.500	--	ND	1.80	--		1
Chloroform	ND	0.200	--	ND	0.977	--		1
Tetrahydrofuran	ND	0.200	--	ND	0.590	--		1
2,2-Dichloropropane	ND	0.200	--	ND	0.924	--		1
1,2-Dichloroethane	ND	0.200	--	ND	0.809	--		1
n-Hexane	ND	0.200	--	ND	0.705	--		1
Diisopropyl ether	ND	0.200	--	ND	0.836	--		1
tert-Butyl Ethyl Ether	ND	0.200	--	ND	0.836	--		1
1,1,1-Trichloroethane	ND	0.200	--	ND	1.09	--		1
1,1-Dichloropropene	ND	0.200	--	ND	0.908	--		1
Benzene	ND	0.200	--	ND	0.639	--		1
Carbon tetrachloride	ND	0.200	--	ND	1.26	--		1
Cyclohexane	ND	0.200	--	ND	0.688	--		1
tert-Amyl Methyl Ether	ND	0.200	--	ND	0.836	--		1
Dibromomethane	ND	0.200	--	ND	1.42	--		1
1,2-Dichloropropane	ND	0.200	--	ND	0.924	--		1
Bromodichloromethane	ND	0.200	--	ND	1.34	--		1
1,4-Dioxane	ND	0.200	--	ND	0.721	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Trichloroethene	ND	0.200	--	ND	1.07	--		1
2,2,4-Trimethylpentane	ND	0.200	--	ND	0.934	--		1
Heptane	ND	0.200	--	ND	0.820	--		1
2,4,4-trimethyl-1-pentene	ND	0.500	--	ND	2.29	--		1
cis-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
4-Methyl-2-pentanone	ND	0.200	--	ND	0.820	--		1
2,4,4-trimethyl-2-pentene	ND	0.500	--	ND	2.29	--		1
trans-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
1,1,2-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Toluene	ND	0.200	--	ND	0.754	--		1
1,3-Dichloropropane	ND	0.200	--	ND	0.924	--		1
2-Hexanone	ND	0.200	--	ND	0.820	--		1
Dibromochloromethane	ND	0.200	--	ND	1.70	--		1
1,2-Dibromoethane	ND	0.200	--	ND	1.54	--		1
Butyl acetate	ND	0.500	--	ND	2.38	--		1
Octane	ND	0.200	--	ND	0.934	--		1
Tetrachloroethene	ND	0.200	--	ND	1.36	--		1
1,1,1,2-Tetrachloroethane	ND	0.200	--	ND	1.37	--		1
Chlorobenzene	ND	0.200	--	ND	0.921	--		1
Ethylbenzene	ND	0.200	--	ND	0.869	--		1
p/m-Xylene	ND	0.400	--	ND	1.74	--		1
Bromoform	ND	0.200	--	ND	2.07	--		1
Styrene	ND	0.200	--	ND	0.852	--		1
1,1,2,2-Tetrachloroethane	ND	0.200	--	ND	1.37	--		1
o-Xylene	ND	0.200	--	ND	0.869	--		1
1,2,3-Trichloropropane	ND	0.200	--	ND	1.20	--		1
Nonane	ND	0.200	--	ND	1.05	--		1
Isopropylbenzene	ND	0.200	--	ND	0.983	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								
Bromobenzene	ND	0.200	--	ND	0.793	--		1
2-Chlorotoluene	ND	0.200	--	ND	1.04	--		1
n-Propylbenzene	ND	0.200	--	ND	0.983	--		1
4-Chlorotoluene	ND	0.200	--	ND	1.04	--		1
4-Ethyltoluene	ND	0.200	--	ND	0.983	--		1
1,3,5-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
tert-Butylbenzene	ND	0.200	--	ND	1.10	--		1
1,2,4-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
Decane	ND	0.200	--	ND	1.16	--		1
Benzyl chloride	ND	0.200	--	ND	1.04	--		1
1,3-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,4-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
sec-Butylbenzene	ND	0.200	--	ND	1.10	--		1
p-Isopropyltoluene	ND	0.200	--	ND	1.10	--		1
1,2-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
n-Butylbenzene	ND	0.200	--	ND	1.10	--		1
1,2-Dibromo-3-chloropropane	ND	0.200	--	ND	1.93	--		1
Undecane	ND	0.200	--	ND	1.28	--		1
Dodecane	ND	0.200	--	ND	1.39	--		1
1,2,4-Trichlorobenzene	ND	0.200	--	ND	1.48	--		1
Naphthalene	ND	0.200	--	ND	1.05	--		1
1,2,3-Trichlorobenzene	ND	0.200	--	ND	1.48	--		1
Hexachlorobutadiene	ND	0.200	--	ND	2.13	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air (Low Level) - Mansfield Lab								

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	100		60-140
Bromochloromethane	98		60-140
chlorobenzene-d5	93		60-140



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01
 Client ID: CAN 131 SHELF 2
 Sample Location:
 Matrix: Air
 Analytical Method: 48,TO-15-SIM
 Analytical Date: 06/15/11 16:38
 Analyst: RY

Date Collected: 06/03/11 00:00
 Date Received: 06/03/11
 Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air by SIM - Mansfield Lab								
Dichlorodifluoromethane	ND	0.050	--	ND	0.247	--		1
Chloromethane	ND	0.500	--	ND	1.03	--		1
Freon-114	ND	0.050	--	ND	0.349	--		1
Vinyl chloride	ND	0.020	--	ND	0.051	--		1
1,3-Butadiene	ND	0.020	--	ND	0.044	--		1
Bromomethane	ND	0.020	--	ND	0.078	--		1
Chloroethane	ND	0.020	--	ND	0.053	--		1
Acetone	ND	2.00	--	ND	4.75	--		1
Trichlorofluoromethane	ND	0.050	--	ND	0.281	--		1
Acrylonitrile	ND	0.500	--	ND	1.08	--		1
1,1-Dichloroethene	ND	0.020	--	ND	0.079	--		1
Methylene chloride	ND	1.00	--	ND	3.47	--		1
Freon-113	ND	0.050	--	ND	0.383	--		1
Halothane	ND	0.050	--	ND	0.404	--		1
trans-1,2-Dichloroethene	ND	0.020	--	ND	0.079	--		1
1,1-Dichloroethane	ND	0.020	--	ND	0.081	--		1
Methyl tert butyl ether	ND	0.020	--	ND	0.072	--		1
2-Butanone	ND	0.500	--	ND	1.47	--		1
cis-1,2-Dichloroethene	ND	0.020	--	ND	0.079	--		1
Chloroform	ND	0.020	--	ND	0.098	--		1
1,2-Dichloroethane	ND	0.020	--	ND	0.081	--		1
1,1,1-Trichloroethane	ND	0.020	--	ND	0.109	--		1
Benzene	ND	0.100	--	ND	0.319	--		1
Carbon tetrachloride	ND	0.020	--	ND	0.126	--		1
1,2-Dichloropropane	ND	0.020	--	ND	0.092	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air by SIM - Mansfield Lab								
Bromodichloromethane	ND	0.020	--	ND	0.134	--		1
1,4-Dioxane	ND	0.100	--	ND	0.360	--		1
Trichloroethene	ND	0.020	--	ND	0.107	--		1
cis-1,3-Dichloropropene	ND	0.020	--	ND	0.091	--		1
4-Methyl-2-pentanone	ND	0.500	--	ND	2.05	--		1
trans-1,3-Dichloropropene	ND	0.020	--	ND	0.091	--		1
1,1,2-Trichloroethane	ND	0.020	--	ND	0.109	--		1
Toluene	ND	0.050	--	ND	0.188	--		1
Dibromochloromethane	ND	0.020	--	ND	0.170	--		1
1,2-Dibromoethane	ND	0.020	--	ND	0.154	--		1
Tetrachloroethene	ND	0.020	--	ND	0.136	--		1
1,1,1,2-Tetrachloroethane	ND	0.020	--	ND	0.137	--		1
Chlorobenzene	ND	0.020	--	ND	0.092	--		1
Ethylbenzene	ND	0.020	--	ND	0.087	--		1
p/m-Xylene	ND	0.040	--	ND	0.174	--		1
Bromoform	ND	0.020	--	ND	0.207	--		1
Styrene	ND	0.020	--	ND	0.085	--		1
1,1,2,2-Tetrachloroethane	ND	0.020	--	ND	0.137	--		1
o-Xylene	ND	0.020	--	ND	0.087	--		1
Isopropylbenzene	ND	0.500	--	ND	2.46	--		1
1,3,5-Trimethylbenzene	ND	0.020	--	ND	0.098	--		1
1,2,4-Trimethylbenzene	ND	0.020	--	ND	0.098	--		1
1,3-Dichlorobenzene	ND	0.020	--	ND	0.120	--		1
1,4-Dichlorobenzene	ND	0.020	--	ND	0.120	--		1
sec-Butylbenzene	ND	0.500	--	ND	2.74	--		1
p-Isopropyltoluene	ND	0.500	--	ND	2.74	--		1
1,2-Dichlorobenzene	ND	0.020	--	ND	0.120	--		1
n-Butylbenzene	ND	0.500	--	ND	2.74	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air by SIM - Mansfield Lab								
1,2,4-Trichlorobenzene	ND	0.050	--	ND	0.371	--		1
Naphthalene	ND	0.050	--	ND	0.262	--		1
1,2,3-Trichlorobenzene	ND	0.050	--	ND	0.371	--		1
Hexachlorobutadiene	ND	0.050	--	ND	0.533	--		1



Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**Air Canister Certification Results**

Lab ID: L1107923-01

Date Collected: 06/03/11 00:00

Client ID: CAN 131 SHELF 2

Date Received: 06/03/11

Sample Location:

Field Prep: Not Specified

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air by SIM - Mansfield Lab								

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-difluorobenzene	101		60-140
bromochloromethane	99		60-140
chlorobenzene-d5	96		60-140



AIR Petro Can Certification

Project Name: BATCH CANISTER CERTIFICATION**Lab Number:** L1107923**Project Number:** CANISTER QC BAT**Report Date:** 07/08/11**AIR CAN CERTIFICATION RESULTS**

Lab ID: L1107923-01
Client ID: CAN 131 SHELF 2
Sample Location: Not Specified
Matrix: Air
Analytical Method: 96,APH
Analytical Date: 06/10/11 19:04
Analyst: RY

Date Collected: 06/03/11 00:00
Date Received: 06/03/11
Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Petroleum Hydrocarbons in Air - Mansfield Lab						
1,3-Butadiene	ND		ug/m3	2.0	--	1
Methyl tert butyl ether	ND		ug/m3	2.0	--	1
Benzene	ND		ug/m3	2.0	--	1
Toluene	ND		ug/m3	2.0	--	1
C5-C8 Aliphatics, Adjusted	ND		ug/m3	12	--	1
Ethylbenzene	ND		ug/m3	2.0	--	1
p/m-Xylene	ND		ug/m3	4.0	--	1
o-Xylene	ND		ug/m3	2.0	--	1
Naphthalene	ND		ug/m3	2.0	--	1
C9-C12 Aliphatics, Adjusted	ND		ug/m3	14	--	1
C9-C10 Aromatics Total	ND		ug/m3	10	--	1

Project Name: POLYCLAD

Lab Number: L1109811

Project Number: 10001087

Report Date: 07/08/11

Sample Receipt and Container Information

Were project specific reporting limits specified? YES

Reagent H2O Preserved Vials Frozen on: NA

Cooler Information Custody Seal**Cooler**

NA Present/Intact

Container Information

Container ID	Container Type	Cooler	pH	Temp deg C	Pres	Seal	Analysis(*)
L1109811-01A	Canister - 2.7 Liter	NA	NA	NA	NA	Present/Intact	TO15-LL(30)
L1109811-02A	Canister - 2.7 Liter	NA	NA	NA	NA	Present/Intact	TO15-LL(30)
L1109811-03A	Canister - 2.7 Liter	NA	NA	NA	NA	Present/Intact	TO15-LL(30)
L1109811-04A	Canister - 2.7 Liter	NA	NA	NA	NA	Present/Intact	TO15-LL(30)

*Values in parentheses indicate holding time in days

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

GLOSSARY

Acronyms

EPA	- Environmental Protection Agency.
LCS	- Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
LCSD	- Laboratory Control Sample Duplicate: Refer to LCS.
LFB	- Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
MDL	- Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
MS	- Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.
MSD	- Matrix Spike Sample Duplicate: Refer to MS.
NA	- Not Applicable.
NC	- Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.
NI	- Not Ignitable.
RL	- Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
RPD	- Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.
SRM	- Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.

Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- A** - Spectra identified as "Aldol Condensation Product".
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than five times (5x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank.
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The RPD between the results for the two columns exceeds the method-specified criteria; however, the lower value has been reported due to obvious interference.
- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less

Report Format: Data Usability Report



Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

Data Qualifiers

than 5x the RL. (Metals only.)

R - Analytical results are from sample re-analysis.

RE - Analytical results are from sample re-extraction.

J - Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).

ND - Not detected at the reporting limit (RL) for the sample.

Project Name: POLYCLAD
Project Number: 10001087

Lab Number: L1109811
Report Date: 07/08/11

REFERENCES

- 48 Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air. Second Edition. EPA/625/R-96/010b, January 1999.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised March 23, 2011 – Mansfield Facility

The following list includes only those analytes/methods for which certification/approval is currently held. For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0141.

Wastewater/Non-Potable Water (Inorganic Parameters: pH, Turbidity, Conductivity, Alkalinity, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Vanadium, Zinc, Total Residue (Solids), Total Suspended Solids (non-filterable), Total Cyanide. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Acid Extractables, Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, PAHs, Haloethers, Chlorinated Hydrocarbons, Volatile Organics.)

Solid Waste/Soil (Inorganic Parameters: pH, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Organic Carbon, Total Cyanide, Corrosivity, TCLP 1311. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Volatile Organics, Acid Extractables, Benzidines, Phthalates, Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Florida Department of Health Certificate/Lab ID: E87814. *NELAP Accredited.*

Non-Potable Water (Inorganic Parameters: SM2320B, SM2540D, SM2540G.)

Solid & Chemical Materials (Inorganic Parameters: 6020, 7470, 7471, 9045. Organic Parameters: EPA 8260, 8270, 8082, 8081.)

Air & Emissions (EPA TO-15.)

Louisiana Department of Environmental Quality Certificate/Lab ID: 03090. *NELAP Accredited.*

Non-Potable Water (Inorganic Parameters: EPA 180.1, 245.7, 1631E, 3020, 6020A, 7470A, 9040, 9050A, SM2320B, 2540D, 2540G, 4500H-B, Organic Parameters: EPA 3510C, 3580A, 3630C, 3640A, 3660B, 3665A, 5030B, 8015D, 3570, 8081B, 8082A, 8260B, 8270C.)

Solid & Chemical Materials (Inorganic Parameters: EPA 1311, 3050, 3051A, 3060A, 6020A, 7196A, 7470A, 7471B, 7474, 9040B, 9045C, 9060. Organic Parameters: EPA 3540C, 3570B, 3580A, 3630C, 3640A, 3660, 3665A, 5035, 8015D, 8081B, 8082A, 8260B, 8270C.)

Biological Tissue (Inorganic Parameters: EPA 6020A. Organic Parameters: EPA 3570, 3510C, 3610B, 3630C, 3640A, 8270C.)

Air & Emissions (EPA TO-15.)

New Hampshire Department of Environmental Services Certificate/Lab ID: 2206. *NELAP Accredited.*

Non-Potable Water (Inorganic Parameters: EPA, 245.1, 245.7, 1631E, 180.1, 6020A, 7470A, 9040B, 9050A, SM2540D, 2540G, 4500H+B, 2320B. Organic Parameters: EPA 8081, 8082, 8260B, 8270C.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 1311, 1312, 3050B, 3051A, 3060A, 6020A, 7470A, 7471A, 9040B, 9045C, 7196A. Organic Parameters: SW-846 3540C, 3580, 3630C, 3640A, 3660B, 3665A, 5035, 8260B, 8270C, 8015D, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA015. *NELAP Accredited.*

Non-Potable Water (Inorganic Parameters: SW-846 1312, 3010, 3020A, 3015, SM2320B, EPA 200.8, SM2540D, 2540G, EPA 120.1, SM2510B, EPA 180.1, 245.1, 1631E, SW-846 7470A, 9040B, 6020, 9010B, 9014 Organic Parameters: SW-846 3510C, 3580A, 5030B, 5035L, 5035H, 3630C, 3640C, 3660B, 3665A, 8015B, 8081A, 8082, 8260B, 8270C)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6020, 9010B, 9014, 1311, 1312, 3050B, 3051, 3060A, 7196A, 7470A, 7471A, 9040B, 9045C, 9060. Organic Parameters: SW-846 3540C, 3570, 3580A, 5030B, 5035L, 5035H, 3630C, 3640A, 3660B, 3665A, 8081A, 8082, 8260B, 8270C, 8015B.)

Atmospheric Organic Parameters (EPA TO-15)

Biological Tissue (Inorganic Parameters: SW-846 6020 Organic Parameters: SW-846 8270C, 3510C, 3570, 3630C, 3640A)

New York Department of Health Certificate/Lab ID: 11627. **NELAP Accredited.**

Non-Potable Water (Inorganic Parameters: SM2320B, SM2540D, EPA 200.8, 6020, 1631E, 245.1, 9014, 9040B, 120.1, SM2510B, 4500CN-E, 4500H-B, EPA 376.2, 180.1, 9010B. Organic Parameters: EPA 8260B, 8270C, 8081A, 8082, 3510C, 5030B.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 6020, 7196A, 3060A, 7471A, 7474, 9014, 9040B, 9045C, 9010B. Organic Parameters: EPA 8260B, 8270C, 8081A, DRO 8015B, 8082, 1311, 1312, 3050B, 3580, 3570, 3051, 5035, 5030B.)

Air & Emissions (EPA TO-15.)

Rhode Island Department of Health Certificate/Lab ID: LAO00299. **NELAP Accredited via LA-DEQ.**

Refer to LA-DEQ Certificate for Non-Potable Water.

Texas Commission of Environmental Quality Certificate/Lab ID: T104704419-08-TX. **NELAP Accredited.**

Solid & Chemical Materials (Inorganic Parameters: EPA 6020, 7470, 7471, 1311, 7196, 9014, 9040, 9045, 9060. Organic Parameters: EPA 8015, 8270, 8260, 8081, 8082.)

Air (Organic Parameters: EPA TO-15)

Washington State Department of Ecology Certificate/Lab ID: C954. *Non-Potable Water* (Inorganic Parameters: SM2540D, 2510B, EPA 120.1, 180.1, 1631E, 245.7.)

Solid & Chemical Materials (Inorganic Parameters: EPA 9040, 9060, 6020, 7470, 7471, 7474. Organic Parameters: EPA 8081, 8082, 8015 Mod, 8270, 8260.)

U.S. Army Corps of Engineers

Department of Defense Certificate/Lab ID: L2217.01.

Non-Potable Water (Inorganic Parameters: EPA 6020A, SM4500H-B. Organic Parameters: 3020A, 3510C, 5030B, 8260B, 8270C, 8270C-ALK-PAH, 8082, 8081A, 8015D-SHC.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1311, 1312, 3050B, 6020A, 7471A, 9045C, 9060, SM 2540G, ASTM D422-63. Organic Parameters: EPA 3580A, 3570, 3540C, 5035A, 8260B, 8270C, 8270-ALK-PAH, 8082, 8081A, 8015D-SHC, 8015-DRO.

Air & Emissions (EPA TO-15.)

Analytes Not Accredited by NELAP

Certification is not available by NELAP for the following analytes: **8270C**: Biphenyl. **TO-15**: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene, 3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 2-Methylnaphthalene, 1-Methylnaphthalene.



ALPHA CHAIN OF CUSTODY

AIR ANALYSIS

PAGE 1 OF 1

320 Forbes Blvd, Mansfield, MA 02048
TEL: 508-822-9300 FAX: 508-822-3288

Client Information

Client: CREEDER

Address: 776 MAIN ST

WESTBORO MA 01581

Phone: 867-888-1272

Fax: 867-887-1051

Email: RIVANDEWELT@CREEDERLLC.COM

Other Project Specific Requirements/Comments:

Project Information

Project Name: POLYCLAD

Project Location: FARMINGTON, NH

Project #: 10001087

Project Manager: RICK VANDEWELT

ALPHA Quote #:

Turn-Around Time

Standard RUSH (only confirmed if pre-approved)

Date Due: Time:

Report Information - Data Deliverables

Date Rec'd in Lab:

FAX ADEX

Criteria Checker: (Default based on Regulatory Criteria Indicated)

Other Formats: EMAIL (standard pdf report)

Additional Deliverables:

Report to: (if different than Project Manager)

Billing Information

Same as Client info PO #:

ALPHA Job #: 1109811

Regulatory Requirements/Report Limits

State/Fed Program Criteria

ANALYSIS

All Columns Below Must Be Filled Out

ALPHA Lab ID (Lab Use Only)	Sample ID	COLLECTION				Sample Matrix*	Sampler's Initials	Can Size	I.D. Can	I.D. - Flow Controller	TO-14A by TO-15	TO-15	TO-15 SIM	APH	FIXED GASES	TO-13A	TO-4 / TO-10	Sample Comments (i.e. PID)
		Date	Start Time	End Time	Initial Vacuum													
1109811-01	CA-SG-2	6/28/11	1256	1430	-29.4	-4.22	SV	2.7	4169	0458	X						PID = 8 ppm	
	CA-SG-10		1306	1440	-29.4	-1.59	SV		524	0412	X						PID = 4 ppm	
	CA-SG-DUP		1256	1430	-29.4	-3.94	SV		530	0141	X						PID = 8 ppm	
	BLANK				-29.2	-29.2												

*SAMPLE MATRIX CODES

AA = Ambient Air (Indoor/Outdoor)
SV = Soil Vapor/Landfill Gas/SVE
Other = Please Specify

Container Type

Relinquished By:

Received By:

Date/Time:

[Signature]
UPS

6/29/11 9:00
6/30/11 13:35

[Signature]
UPS

6/30/11 13:35

Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguities are resolved. All samples submitted are subject to Alpha's Terms and Conditions. See reverse side.