PROJECT SPECIAL PROVISIONS GEOENVIRONMENTAL

CONTAMINATED SOIL (7/1/2020)

The Contractor's attention is directed to the fact that soil contaminated with petroleum hydrocarbon compounds exist within the project area. The known areas of contamination are indicated on corresponding plans sheets. Information relating to these contaminated areas, sample locations, and investigation reports will be available at the following web address by navigating to the correct letting year and month then selecting, "Plans and Proposals", "U-4424", "Individual Sheets/520 GeoEnvironmental":

http://dotw-xfer01.dot.state.nc.us/dsplan/

Petroleum contaminated soil may be encountered during any earthwork activities on the project. The Contractor shall only excavate those soils that the Engineer designates necessary to complete a particular task. The Engineer shall determine if soil is contaminated based on petroleum odors and unusual soil staining. Contaminated soil not required to be excavated is to remain in place and undisturbed. Undisturbed soil shall remain in place, whether contaminated or not. The Contractor shall transport all contaminated soil excavated from the project to a facility licensed to accept contaminated soil.

In the event that a stockpile is needed, the stockpile shall be created within the property boundaries of the source material and in accordance with the Diagram for Temporary Containment and Treatment of Petroleum-Contaminated Soil per North Carolina Department of Environmental Quality's Division of Waste Management UST Section GUIDELINES FOR EX SITU PETROLEUM CONTAMINATED SOIL REMEDIATION. If the volume of contaminated material exceeds available space on site, the Contractor shall obtain a permit from the NCDEQ UST Section's Regional Office for off-site temporary storage. The Contractor shall provide copies of disposal manifests completed per the disposal facilities requirements and weigh tickets to the Engineer.

If groundwater is encountered and dewatering is required in areas of known contamination then the contractor shall containerize the groundwater in vessels provided by the Department. The Department will be responsible for the sampling and disposal of the water. Handling contaminated ground water will be incidental to the project.

Measurement and Payment:

The quantity of contaminated soil hauled, and disposed of shall be the actual number of tons of material, which has been acceptably transported and weighed with certified scales as documented by disposal manifests and weigh tickets. The quantity of contaminated soil, measured as provided above, shall be paid for at the contract unit price per ton for "Hauling and Disposal of Petroleum Contaminated Soil".

The above price and payment shall be full compensation for all work covered by this section, including, but not limited to stockpiling, loading, transportation, weighing, laboratory testing, disposal, equipment, decontamination of equipment, labor, and personal protective equipment.

Payment shall be made under:

Pay ItemPay UnitHauling and Disposal of Petroleum Contaminated SoilTon

John Pilipchuk

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7/1/2020

SEAL

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Pilipchuk



114 Edinburgh South Drive Suite 200 Cary, North Carolina 27511 919.827.0864 www.daa.com

June 17, 2020

Mr. Craig Haden North Carolina Department of Transportation Geotechnical Engineering Unit 1589 Mail Service Center Raleigh, North Carolina 27699-1589

Reference: Preliminary Site Assessment for the Joshua A. Myer Property (Parcel 46)

1810 W Wilson Street

Tarboro, Edgecombe County, North Carolina

State Project: U-4424 WBS Element 39062.1.2

DAA Project No. 20080204-010103

Dear Mr. Haden:

Draper Aden Associates (DAA) completed the Preliminary Site Assessment at the above-referenced property. DAA performed the work in accordance with the Technical and Cost proposal dated April 22, 2020, and the North Carolina Department of Transportation's (NCDOT's) Notice to Proceed dated April 23, 2019. Activities associated with the assessment consisted of conducting a geophysical investigation to identify whether an underground storage tank (UST) exists within the proposed right-of-way/easement (ROW/easement) and collecting soil and groundwater samples for laboratory analysis. The purpose of this report is to document the field activities, present the laboratory analyses, and provide recommendations regarding the property.

Location and Description

The Joshua A. Myers Property (Parcel #46) is located at 1810 W. Wilson Street in Tarboro, Edgecombe County, North Carolina. The property is situated on the west side of W. Wilson Street at the intersection with Simpson Drive (**Figure 1**). The property was a former automotive fueling station, but as of the date of the fieldwork, it houses a privately-owned garage beside a residence. Two buildings and a metal carport canopy are located at the site (**Figure 2**).

The garage, a single-story block structure with metal carport canopy in front, was the former fueling station. According to the property owner, one UST was located under the concrete former dispenser island in front of the building and has not been removed.

The NCDOT requested a Preliminary Site Assessment for the proposed ROW/easement because the property is a former automotive fueling station. The scope of work as defined in the Request for Technical and Cost Proposal was to evaluate the proposed right-of-way/easement with respect to the potential presence of known and unknown USTs within the proposed ROW/easement, and to assess whether subsurface contamination existed within the proposed ROW/easement. An estimate of the quantity of impacted soil was to be provided if impacted soils were encountered.

DAA reviewed the on-line North Carolina Department of Environmental Quality (NCDEQ) Incident Management database and no incident has been assigned to the site. DAA also examined the UST registration database to obtain UST ownership information; however, these files were not accessible online during the preparation of this report and a facility identification could not be obtained or verified.

Geophysical Survey

Prior to DAA's mobilization to the site for drilling, we conducted a geophysical survey within and near the proposed ROW/easement (i.e., study area) to determine potential presence of unknown UST(s). Areas near the ROW/easement were not within the scope of work, but the equipment traverses necessitated crossing the ROW/easement lines. The geophysical survey consisted of an electromagnetic survey using a Geonics EM61 time-domain electromagnetic (EM) induction meter to locate buried metallic objects, and ground penetrating radar (GPR) using a Noggin 250 with 250 MHz antennae specifically to locate USTs.

The geophysical team laid out a survey grid along the proposed right-of-way with the X-axis oriented approximately parallel to W Wilson Street and the Y-axis oriented approximately perpendicular to W. Wilson Street. **Figure 1** of the geophysical survey report in **Attachment A** shows the EM survey area.

The EM survey lines were spaced five feet apart and the instruments collected magnetic data continuously along each survey line with a data logger. After collection, DAA reviewed the data in the field with graphical user interface computer software. Following the electromagnetic survey, a GPR survey was conducted to further evaluate any notable metallic anomalies. GPR transects are shown on **Figures 5 through 7** of **Attachment A**.

DAA detected several anomalies in those areas accessible to the study area. The survey attributed all but one of the anomalies to visible cultural features or underground utilities. The geophysical data indicated a magnetic anomaly southwest of the carport that is adjacent to the proposed ROW/easement, and GPR signatures suggest a UST. Based on the lack of additional information or visual evidence of a tank, DAA classified the anomaly as a possible UST. **Attachment A** presents DAA's detailed report of findings and interpretations.

Site Assessment Activities

On May 18, 2020, DAA mobilized to the site to conduct a Geoprobe® direct-push investigation to evaluate subsurface soil and groundwater conditions within the proposed ROW/easement to a depth of 8 to 10 feet below ground surface (ft bgs). DAA advanced five direct-push probes (SB-1 through SB-5) at select locations throughout the proposed ROW/easement (**Figure 2**). The soil boring logs are included as **Attachment B**. The borings were located to evaluate the subsurface conditions in the study area (**Attachment C**).

The lithology encountered by the direct-push samples was generally consistent throughout the site. The ground surface was covered by approximately 6 inches of gravel or topsoil. Below this surface cover was a brown to light brown medium- to fine-grained sand. The lithology encountered in borings B-1 to B-3 included clayey sands and clay between 2 and 7 ft bgs. Bedrock was not encountered in any of the borings. DAA noted groundwater in the borings at a depth of approximately 5 to 7 ft bgs. Each boring was backfilled with a mix of bentonite (swelling clay to seal the boring) and drill cuttings to the surface after completion.

According to the 1985 Geologic Map of North Carolina, the site is within the Coastal Plain Physiographic Province in North Carolina. The strata indicated for this area is the Yorktown Formation, comprising fossiliferous clay and sand.

Continuous sampling using a Geoprobe® resulted in good recovery of soil samples from the direct-push holes. DAA collected, documented, and contained soil samples in four-foot long acetate sleeves inside the direct-push Macro-Core® sampler. The soils observed at the site are consistent with Yorktown Formation strata (see **Attachment B**)

Each of the sleeves was divided into two-foot long sections for soil sample screening. Soil from each two-foot interval was placed in a resealable plastic bag and the bag was set aside to allow time for volatilization of potential organic compounds to the bag headspace. A photoionization detector (PID) probe was inserted into the bag and the reading was recorded (**Table 1**).

Following completion of the soil sampling, the boring with the highest recorded PID field screen results or indications of potential contamination (odors, staining, etc.) was converted to a temporary groundwater monitoring well using the direct push screen point. Groundwater was encountered at a depth of 5 to 7 ft bgs and a probe screen was inserted to collect the groundwater sample using low-flow techniques.

DAA submitted for laboratory analysis one soil sample from each of the five borings at the depth interval with the highest PID reading measured at the time of collection (**Table 1**). The soil samples were submitted to REDLab in Wilmington, North Carolina, for analysis of total petroleum hydrocarbons (TPH) diesel range organics (DRO) and gasoline range organics (GRO) using ultraviolet fluorescence (UVF) methodology.

The groundwater sample was analyzed for volatile organic compounds (VOCs) using EPA Method 8260 and for semivolatile organic compounds (SVOCs) using EPA Method 8270. Contest Laboratories in East Longmeadow, Massachusetts conducted the VOC and SVOC analyses.

Analytical Results

Table 1 and **Figure 3** summarize the soil laboratory results for the five soil samples for TPH DRO/GRO. **Table 2** and **Figure 4** summarize the groundwater laboratory results, and **Attachment D** presents the complete laboratory reports.

One soil sample, SB-2, contained a detectable GRO concentration of 53.9 milligrams per kilogram (mg/kg). No other sample contained detectable GRO concentrations. Four soil samples (SB-2 through SB-5) contained detectable DRO concentrations ranging from 1.3 to 29 mg/kg. The action levels are 50 mg/kg for GRO and 100 mg/kg for DRO¹. None of the soil samples analyzed for this site contained DRO concentrations above their respective action levels.

The groundwater analytical results (**Table 2**) indicate the detection of several petroleum compounds. DAA compared these concentrations to the Groundwater Quality Standards established in 15A NCAC 2L (2L Standards). The compounds detected above the 2L Standards were associated with gasoline constituents. These included ethylbenzene, isopropylbenzene, naphthalene, n-propylbenzene, toluene, 1,2,4-trimethylbenzene, 1,3,5-trimethylbenzene, and total xylenes. Of the SVOC compounds detected, only 1-methylnaphthalenen and naphthalene were present are concentrations greater than the 2L Standards.

Contaminated Soil Volume Estimate

The GRO concentration in SB-2 nominally exceeds the action level and from a soil sample collected in the apparent smear zone of the groundwater. Based on the field screening results, the PID readings increased at the depth just above the groundwater. As such, groundwater contamination may be the likely source of the GRO detected in the soil sample SB-2. Therefore, no estimate of the volume of soil necessitating possible remediation was required for this report.

Conclusions and Recommendations

DAA conducted a Preliminary Site Assessment to evaluate the NCDOT proposed ROW/easement on the Joshua A. Myers Property (Parcel #46) located at 1810 W. Wilson Street in Tarboro, Edgecombe County, North Carolina. A geophysical survey indicated the presence of a possible UST within the proposed ROW/easement; however, no visual signs of a UST were noted.

Five soil borings and one temporary well screen point were advanced within the proposed ROW/easement to evaluate the subsurface soil and groundwater conditions within the site. One

¹ NCDEQ, Guidelines for North Carolina Action Limits for Total Petroleum Hydrocarbons (TPH), July 26, 2016,

of the soil samples analyzed contained a GRO concentration above the action level; however, the compounds are likely from groundwater contamination and no contaminated soil calculations were conducted. Groundwater contained several compounds, many above the 2L Standards.

Because compounds were detected above the action level in the soil and groundwater samples, DAA recommends that a copy of this report be submitted to the Division of Waste Management, UST Section, in the Raleigh Regional Office.

DAA appreciates the opportunity to work with the NCDOT on this project. If you have any questions, please contact us at (919) 827-0864.

Sincerely,

Draper Aden Associates

DocuSigned by:

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6/25/2020

Michael W. Branson, P.G.

Project Manager

Attachments

William D. Newcomb, P.G. Senior Hydrogeologist

Willen D. News

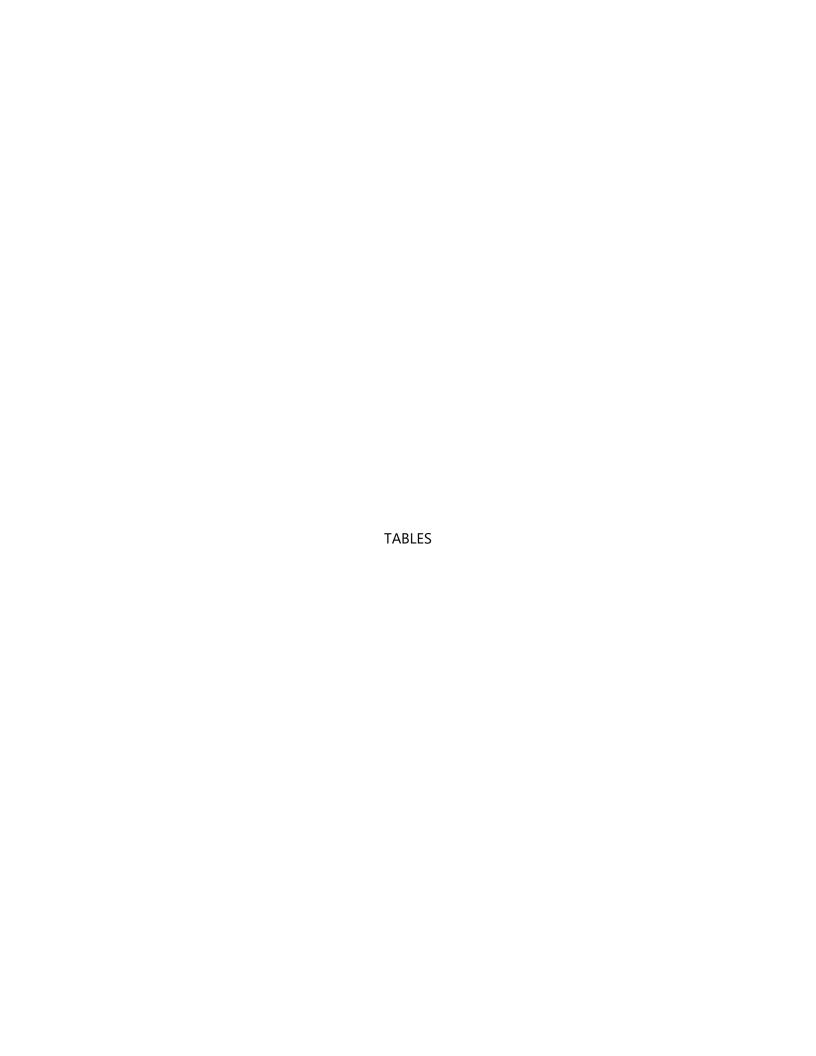


TABLE 1

SOIL FIELD SCREENING AND ANALYTICAL RESULTS JOSHUA A MYERS PROPRTY (PARCEL 46) TARBORO, EDGECOMBE COUNTY, NORTH CAROLINA STATE PROJECT: U-4424

WBS ELEMENT 39062.1.2 DAA PROJECT NO. 20080204-010103

SAMPLE ID	DEPTH (ft)	PID READING (ppm)	SAMPLE ID	ANALYTICAL RESULTS (mg/kg)	
		(ррііі)		UVF GRO	UVF DRO
	Action Level (mg/kg)			50	100
	0 - 2	0.8			
SB-1	2 - 4	0.8			
	4 - 6	2.2	SB-1	<0.68	<0.68
	6 - 8	1.9			
	0 - 2	1.9			
SB-2	2 - 4	2.5			
3D-2	4 - 6	3.3			
	6 - 8	1707	SB-2	53.9	29
	0 - 2	2.2			
SB-3	2 - 4	3			
30-3	4 - 6	3.2			
	6 - 8	24.5	SB-3	<0.65	3.8
	0 - 2	2.3			
SB-4	2 - 4	2.2			
SB-4	4 - 6	2.1			
	6 - 8	5.5	SB-4	<0.47	1.3
	0 - 2	3			
SB-5	2 - 4	3.5	SB-5	<0.75	2
	4 - 6	2.7			
	6 - 8	2.9			

- 1) ft feet
- 2) ppm parts per million
- 3) PID photoionization detector
- 4) mg/kg milligrams per kilogram
- 5) UVF DRO Diesel range organics by ultraviolet fluorescence (UVF)
- 6) UVF GRO Gasoline range organics by UVF
- 7) Action level for TPH based upon NCDEQ memo *Guidelines for North Carolina Action Limits for Total Petroleum Hydrocarbons* July 29, 2016. VOC action levels based on Maximum Soil Contaminant Concentrations.
- 8) Soil samples were collected on May 18, 2020.
- 9) **Bold** values are above the action level.

TABLE 2 GROUNDWATER ANALYTICAL RESULTS JOSHUA A MYERS PROPRTY (PARCEL 46) TARBORO, EDGECOMBE COUNTY, NORTH CAROLINA STATE PROJECT: U-4424 WBS ELEMENT 39062.1.2 DAA PROJECT NO. 20080204-010103

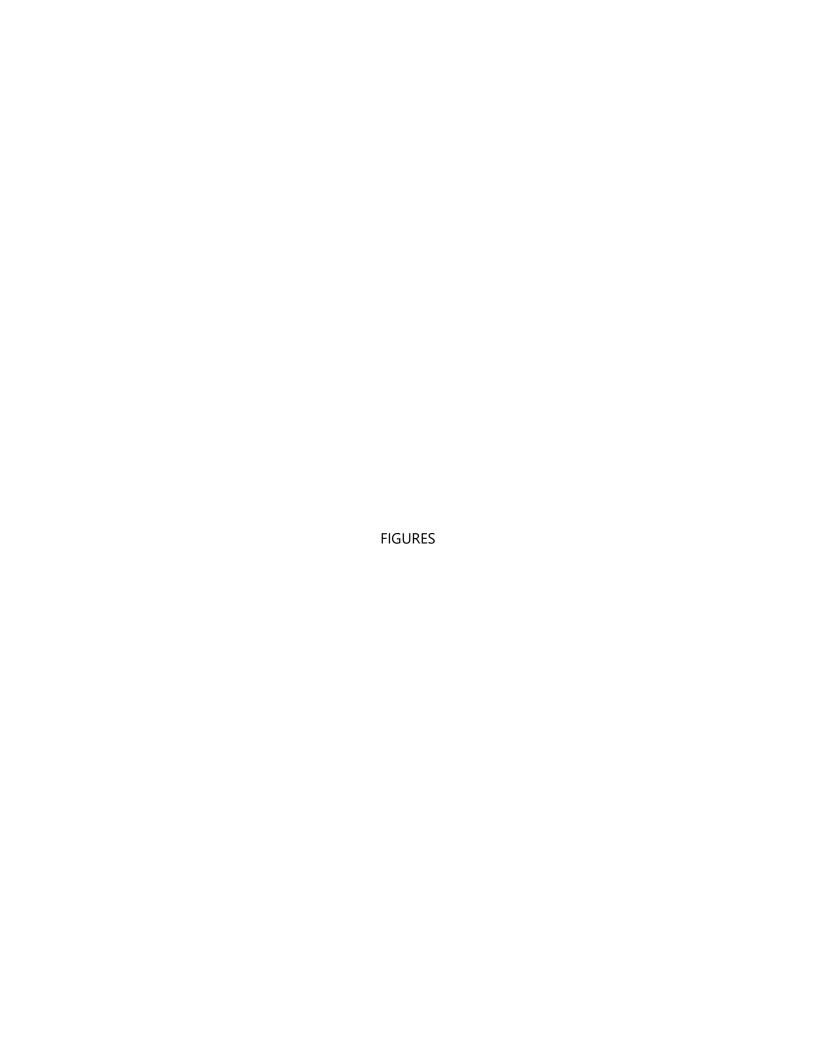
Sample ID	SB-2					
Analyte	15A NCAC 2L Standard (μg/L)	Concentration (μg/L)				
Volatile Organic Compounds						
Ethylbenzene	600	1600				
Isopropylbenzene	70	89				
Naphthalene	6	320				
n-Propylbenzene	70	130				
Toluene	600	3700				
1,2,4-Trimethylbenzene	400	920				
1,3,5-trimethylbenzene	400	220				
Xylenes	500	8000				
Semivolatile Organic Compounds						
Acenaphthene	80	0.075 J				
Acenaphthylene	200	0.037 J				
Fluorene	300	0.10 J				
1-Methylnaphthalene	1 (IMAC)	14				
2-Methylnaphthalene	30	30				
Naphthalene	6	190				
Phenanthrene	200	0.082				

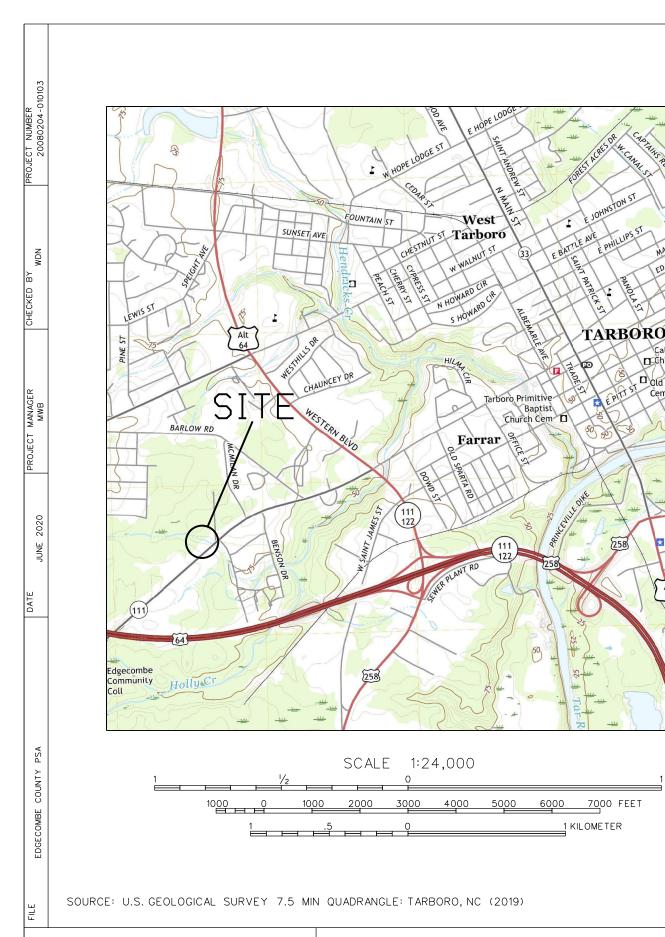
¹⁾ µg/L - micrograms per liter

²⁾ Groundwater sample was collected on May 18, 2020.

³⁾ J - Estimated value between the method detection limit and the reporting limit.

⁴⁾ IMAC - Interim Maximum Allowable Concentration







114 EDINBURGH SOUTH DRIVE CARY, NORTH CAROLINA 27511 TEL: (919) 873-1060 FAX: (919) 873-1074

VICINITY MAP

JOSHUA A MYERS PROPERTY 1810 W WILSON STREET TARBORO, NORTH CAROLINA

FIGURE

MARTIN LUTHER KING

EDMONDSON AVE

Calvary Episcopal

111 122

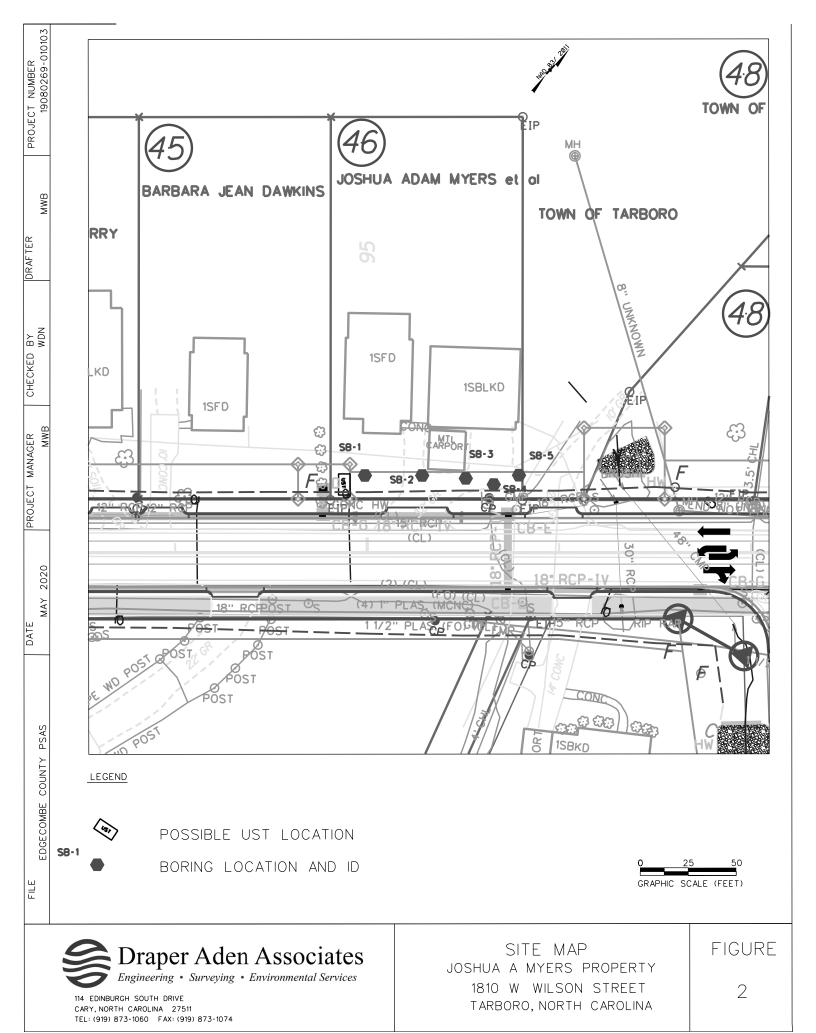
WALSTON ST

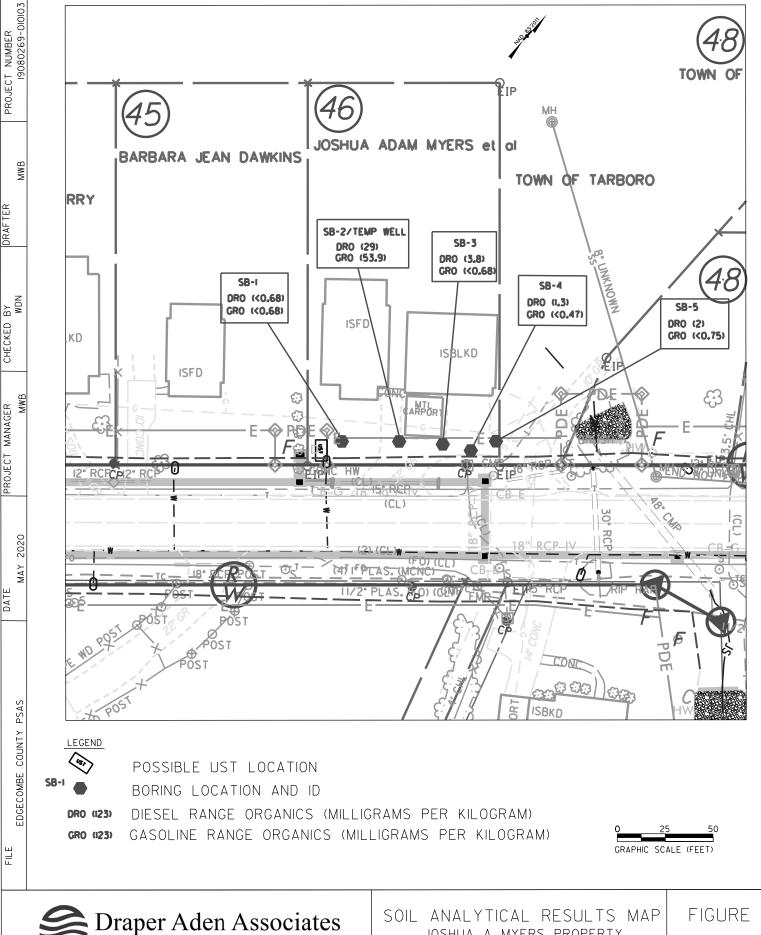
1 MILE

BEASLE

Old Town

Cem





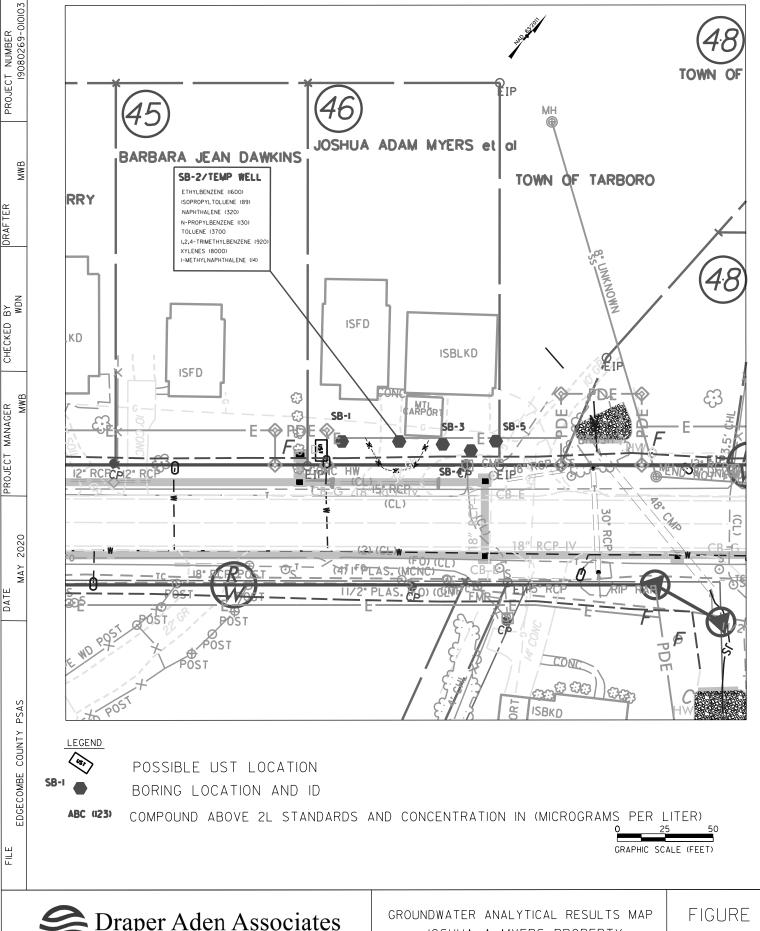
Draper Aden Associates

Engineering · Surveying · Environmental Services

CARY, NORTH CAROLINA 275II TEL: (919) 873-1060 FAX: (919) 873-1074 JOSHUA A MYERS PROPERTY

1810 W WILSON STREET

TARBORO, NORTH CAROLINA



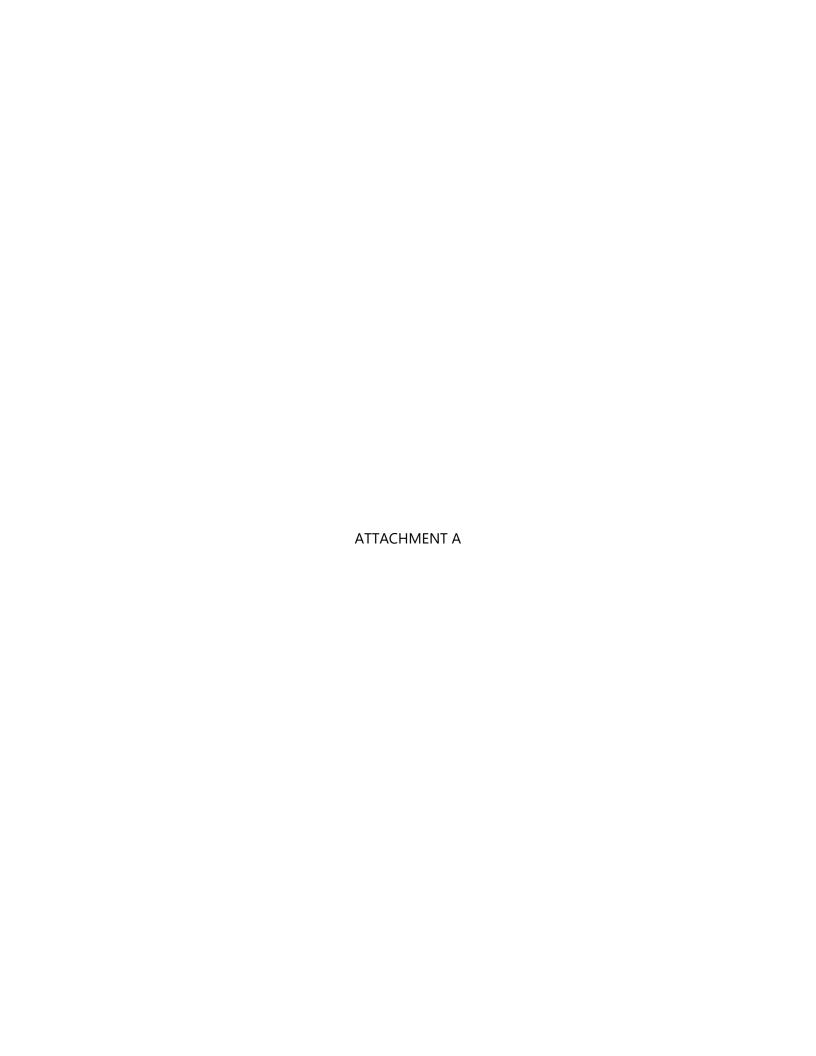


II4 EDINBURGH SOUTH DRIVE CARY, NORTH CAROLINA 275II TEL: (919) 873-1060 FAX: (919) 873-1074 GROUNDWATER ANALYTICAL RESULTS MAP

JOSHUA A MYERS PROPERTY

1810 W WILSON STREET

TARBORO, NORTH CAROLINA



Geophysical Study For Possible USTs

1810 West Wilson Street Tarboro, North Carolina



North Carolina Department of Transportation 1589 Mail Service Center Raleigh, NC 27699-1589 37918

May 28, 2020

DAA Project Number: 20080204-010203





2206 South Main Street Blacksburg, Virginia 24060 540.552.0444 www.daa.com

May 28, 2020

Mr. John Pilipchuck, P.E Geotechnical Engineering Unit N.C. Department of Transportation 1589 Mail Service Center Raleigh, NC 27699-1589 37918

RE: Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, North Carolina Draper Aden Associates Project No. 20080204-010203

Dear Mr. Pilipchuck:

Draper Aden Associates has completed the geophysical study at 1810 West Wilson Street near Tarboro, North Carolina. The objective of this study was to assist in determining if any underground storage tanks (USTs) may be present beneath the study area. To meet this objective, a combination of ground penetrating radar (GPR) and electromagnetic induction (EM) techniques were utilized. The following report documents our methodologies and findings.

We value our professional relationship with N.C. Department of Transportation and hope that you will contact us with any similar needs in the future. If you have any questions regarding this report, or if we can be of any further service to you please do not hesitate to contact us.

Sincerely,

Draper Aden Associates

Johanna Vaughan, P.G.

Geologist

Francis Douglas Pinckney, P.E.

Team Leader/Senior Project Engineer Geotechnical and Construction Services

> Blacksburg • Charlottesville • Manassas • Newport News • Richmond • Virginia Beach Fayetteville • Raleigh

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FIGURES

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1.0 EXECUTIVE SUMMARY

Draper Aden Associates (DAA) was retained by N.C. Department of Transportation to conduct a geophysical study at 1810 West Wilson Street near Tarboro, North Carolina. The objective of this study was to assist in determining if any underground storage tanks (USTs) may be present beneath the study area. To meet this objective, a combination of ground penetrating radar (GPR) and EM61 electromagnetic induction techniques were utilized.

The EM61 data were collected on May 4, 2020 in grid fashion with sub-parallel traverses spaced approximately four feet apart, tracked by a GPS unit capable of sub-foot accuracy. The contoured results from the EM61 data reveal one anomaly located approximately 45 feet southwest of the southwest corner of the carport which may represent a possible UST, identified as Anomaly 1.

GPR data were collected on May 4, 2020 utilizing the same grid as the EM61 data collection, with sub-parallel traverses spaced approximately two feet apart, tracked by a GPS unit capable of sub-foot accuracy. Analysis of the GPR data in cross-section revealed broad hyperbolic reflectors consistent with a possible UST in the same vicinity as EM Anomaly 1, located approximately 2.5 feet below the ground surface. Plan-view analysis of the GPR data reveals a weak anomaly in the depth slice from 2.5 to 3.0 feet depth that correlates well with EM Anomaly 1.

The combined analysis of the EM61 and GPR data reveals good correlation between the two methods, with a coincident anomaly observed in the data from each method (Anomaly 1). It is uncertain if Anomaly 1 represents a UST, but of the collected geophysical data, this feature is most likely to represent a potential UST.

This study was conducted by registered professional geologists with extensive experience in the collection, processing, and interpretation of geophysical data. It should be noted, however, that all geophysical methods are interpretive, and additional invasive exploration would be required to verify or refute the interpretations within this report.

2.0 INTRODUCTION

Draper Aden Associates (DAA) was retained by N.C. Department of Transportation to conduct a geophysical study at 1810 West Wilson Street near Tarboro, North Carolina (Figure 1). The objective of this study was to assist in determining if any underground storage tanks (USTs) may be present beneath the study area. To meet this objective, a combination of ground penetrating radar (GPR) and electromagnetic induction (EM) techniques were utilized. The following report documents our methodologies and findings.

The tasks involved in this study included:

- 1. Collection, processing, and interpretation of EM61 data;
- 2. Collection, processing, and interpretation of GPR data;
- 3. Preparation of this document to detail our methods and findings.

3.0 ELECTROMAGNETIC INDUCTION (EM) STUDY

3.1 EM Field Methods

The instrument used for this investigation was the EM61 manufactured by Geonics, LTD. The EM61 data were collected on May 4, 2020 in grid fashion with sub-parallel traverses spaced approximately four feet apart (Figure 2). The distribution of the EM61 data was tracked by a global positioning system (GPS) unit capable of sub-foot accuracy. The collected data were subsequently contoured laterally and analyzed for evidence of any possible USTs.

3.2 EM61 Results

The contoured results from the EM61 data are presented in Figure 3, overlain onto Google Earth aerial imagery. The data reveals a "hotspot" (isolated areas of elevated EM response) which may represent a possible UST. This feature is identified as Anomaly 1 and located approximately 45 feet southwest of the southwest corner of the carport. It is uncertain if the EM anomaly represent a UST, but of the collected data, this feature is most likely to possibly represent a UST. Since the EM61 instrument is particularly sensitive to buried metallic materials or objects, we consider it likely that the anomaly is metallic in composition.

4.0 GPR STUDY

4.1 GPR Field Methods

The instrument used for this investigation was the Noggin 250 manufactured by Sensors and Software, Inc. in Ontario, Canada, which utilizes a 250 MHz antenna mounted on a moveable cart. GPR data were collected on May 4, 2020 utilizing the same grid as the EM61 data collection, with sub-parallel traverses spaced approximately two feet apart, tracked by a GPS unit capable of sub-foot accuracy (Figure 4).

4.2 GPR Results

The GPR data were analyzed as vertical cross-sections and as depth slices, or plan-view maps of the GPR response from various depth intervals for evidence of possible USTs. Figure 5 depicts two sample GPR cross-sections from the collected data which contain broad hyperbolic reflectors consistent with a possible UST located approximately 2.5 feet below the ground surface. These features correlate to Anomaly 1 identified in the EM61 data.

Figures 6 and Figure 7 illustrate the plan-view GPR response in 6-inch-thick depth intervals spanning from 1.0 to 6.0 feet depth. An anomaly in the depth slice from 2.5 to 3.0 feet depth correlates well with EM Anomaly 1. Numerous other areas of elevated GPR response observed throughout the depth slices may represent miscellaneous buried objects, materials, or conditions, such as intermittent clay layers or zones of wet soil. However, due to a general lack of elevated EM response in other areas (beyond the vicinity of Anomaly 1), these other areas of anomalous GPR response are considered likely to not be metallic in composition.

5.0 CONCLUSIONS

The combined analysis of the EM61 and GPR data reveals good correlation between the two methods, with a coincident anomaly observed in the data from each method (Anomaly 1).

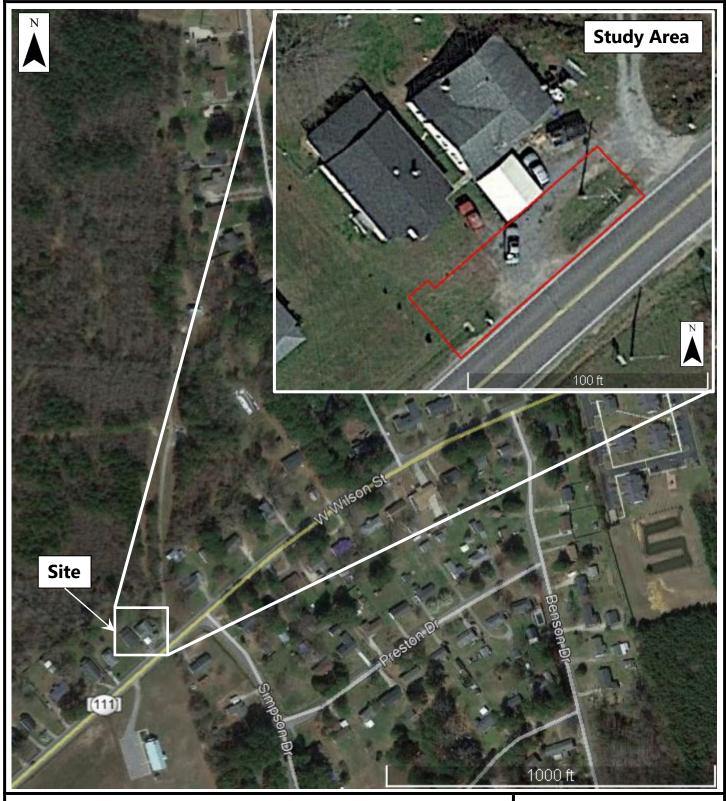
It is uncertain if Anomaly 1 represents a UST, but of the collected geophysical data, this feature is most likely to represent a potential UST. Since the EM61 instrument is particularly sensitive to buried metallic materials or objects, we consider it likely that Anomaly 1 is metallic in composition. Furthermore, the location of Anomaly1 is characterized in profile view of the GPR data as a broad hyperbolic reflector, which is consistent with a possible UST. The results of the study are summarized in Figure 8, depicting the location of Anomaly 1.

6.0 LIMITATIONS

This study was conducted by registered professional geologists with extensive experience in the collection, processing, and interpretation of geophysical data. It should be noted, however, that

all geophysical methods are interpretive, and additional invasive exploration would be required to verify or refute the interpretations within this report.

7.0 FIGURES



Site Location

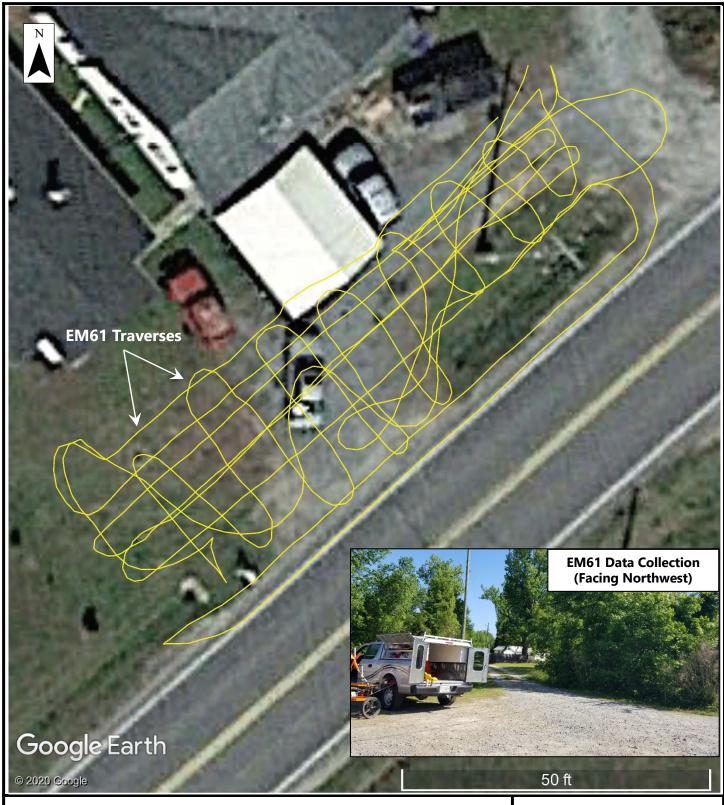
Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC

PROJECT: 20080204-010203

Draper Aden Associates

Engineering * Surveying * Environmental Services

2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291 Richmond, VA Charlottesville, VA Hampton Roads, VA Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE:05/15/2020 **FIGURE**



EM61 Traverse Map

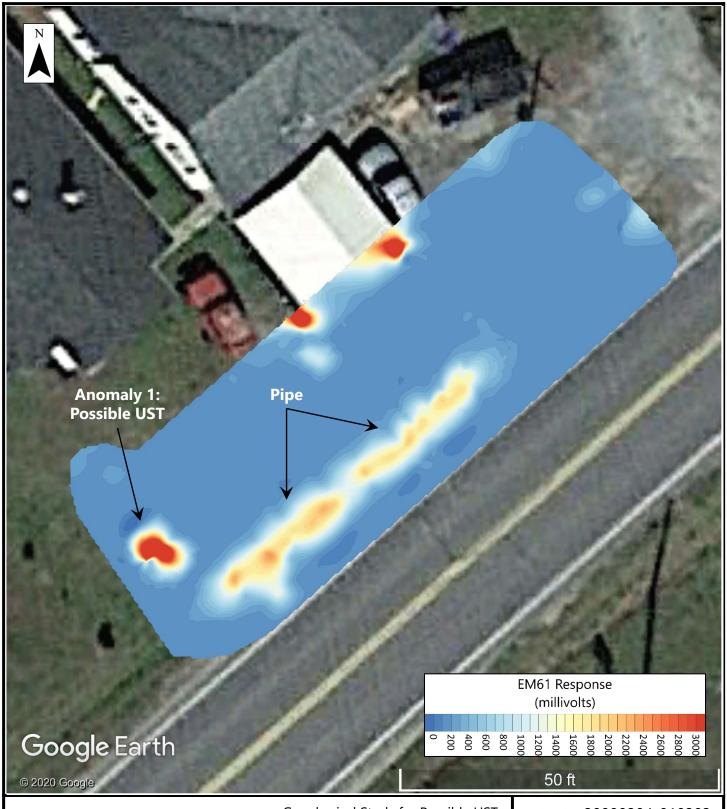
Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC

PROJECT: 20080204-010203

Draper Aden Associates

Engineering • Surveying • Environmental Services

2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291 Richmond, VA Charlottesville, VA Hampton Roads, VA Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE:05/15/2020 **FIGURE**



Contoured EM61 Results

Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC

PROJECT: 20080204-010203

Draper Aden Associates

Engineering * Surveying * Environmental Services

2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291 Richmond, VA Charlottesville, VA Hampton Roads, VA Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE: 05/15/2020 **FIGURE**



GPR Traverse Map

Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC

PROJECT: 20080204-010203

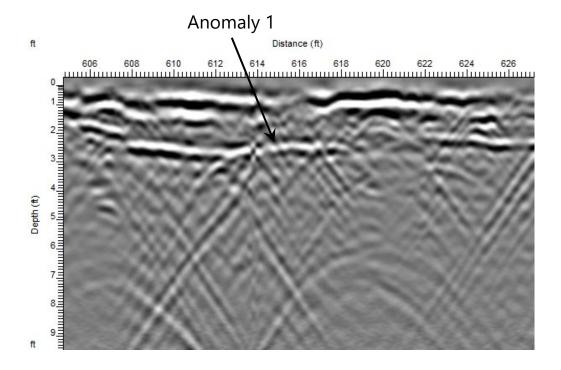


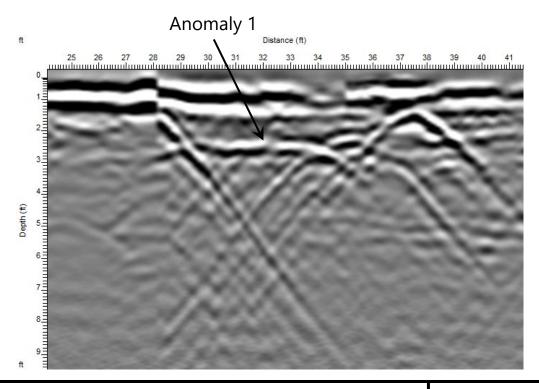
Draper Aden Associates

Engineering • Surveying • Environmental Services

2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291 Richmond, VA Charlottesville, VA Hampton Roads, VA Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE:05/15/2020 **FIGURE**

Southwest Portion of the Study Area





Sample GPR Cross-sections From the Collected Data

Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC PROJECT: 20080204-010203

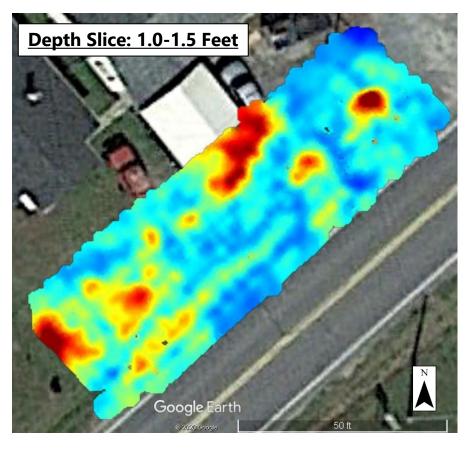


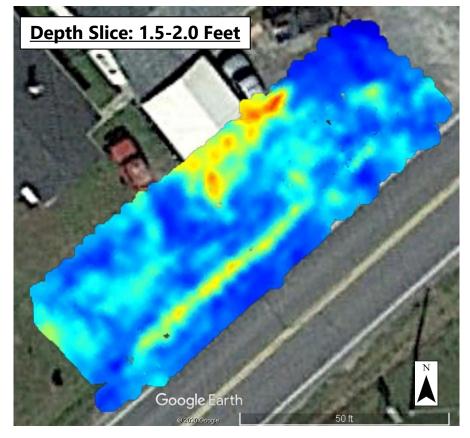
Draper Aden Associates

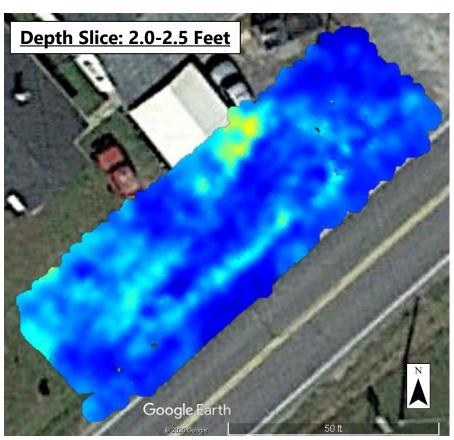
Engineering • Surveying • Environmental Services

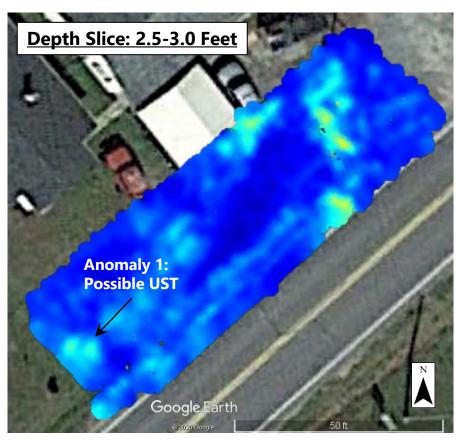
2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291 Richmond, VA Charlottesville, VA Hampton Roads, VA Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP

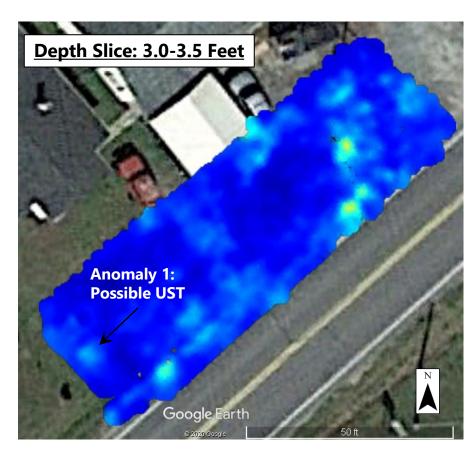
CHECKED: CMP DATE: 05/15/2020 **FIGURE**

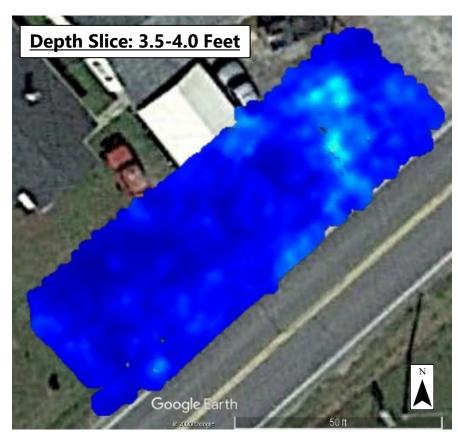












GPR Depth Slices from 1.0 Feet to 4.0 Feet Depth

Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC PROJECT: 20080204-010203

Draper Aden Associates

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2206 South Main Street Richmond VA Rale

2206 South Main Street Richmond, VA
Blacksburg, VA 24060 Charlottesville, VA
540-552-0444 Fax: 540-552-0291 Hampton Roads, VA

Pices

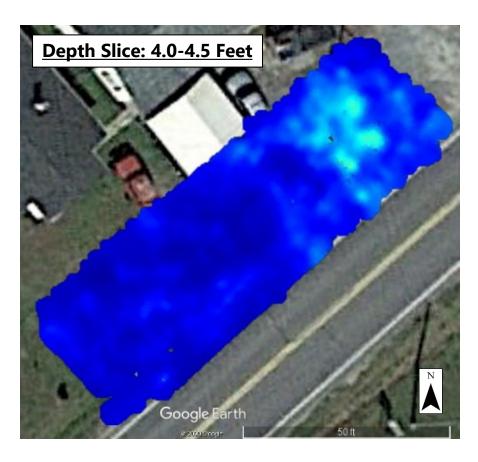
Raleigh, NC

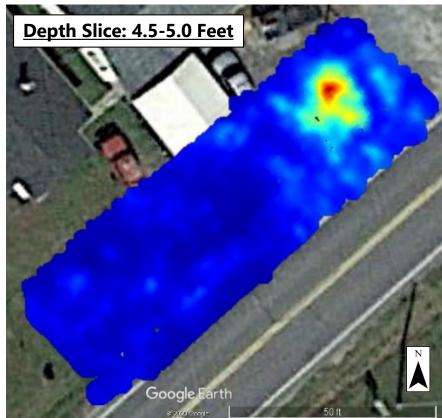
Fayetteville, NC

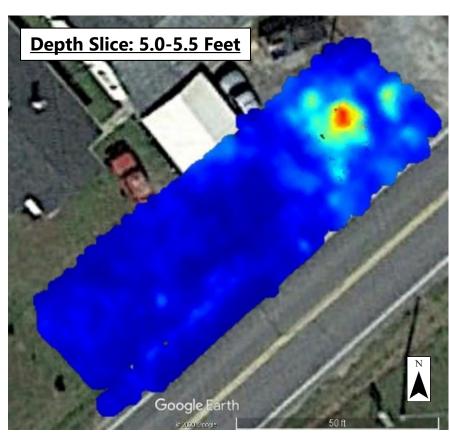
Northern Virginia

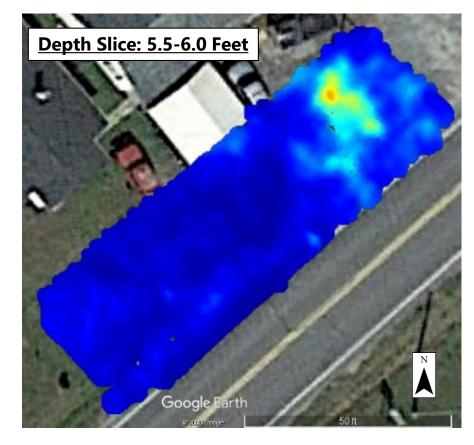
Virginia Beach, VA

DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE: 05/15/2020 FIGURE









Geophysical Study for Possible USTs 1810 West Wilson Street, Tarboro, NC PROJECT: 20080204-010203

GPR Depth Slices from 4.0 Feet to 6.0 Feet Depth



2206 South Main Street
Blacksburg, VA 24060
540-552-0444 Fax: 540-552-0291

Richmond, VA
Charlottesville, VA
Hampton Roads, VA

VA Raleigh, NC
ille, VA Fayetteville, NC
oads, VA Northern Virginia
Virginia Beach, VA

DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE: 05/15/2020 FIGURE



Geophysical Study for Possible USTs

1810 West Wilson Street, Tarboro, NC

Summary of GPR and EM61 Results



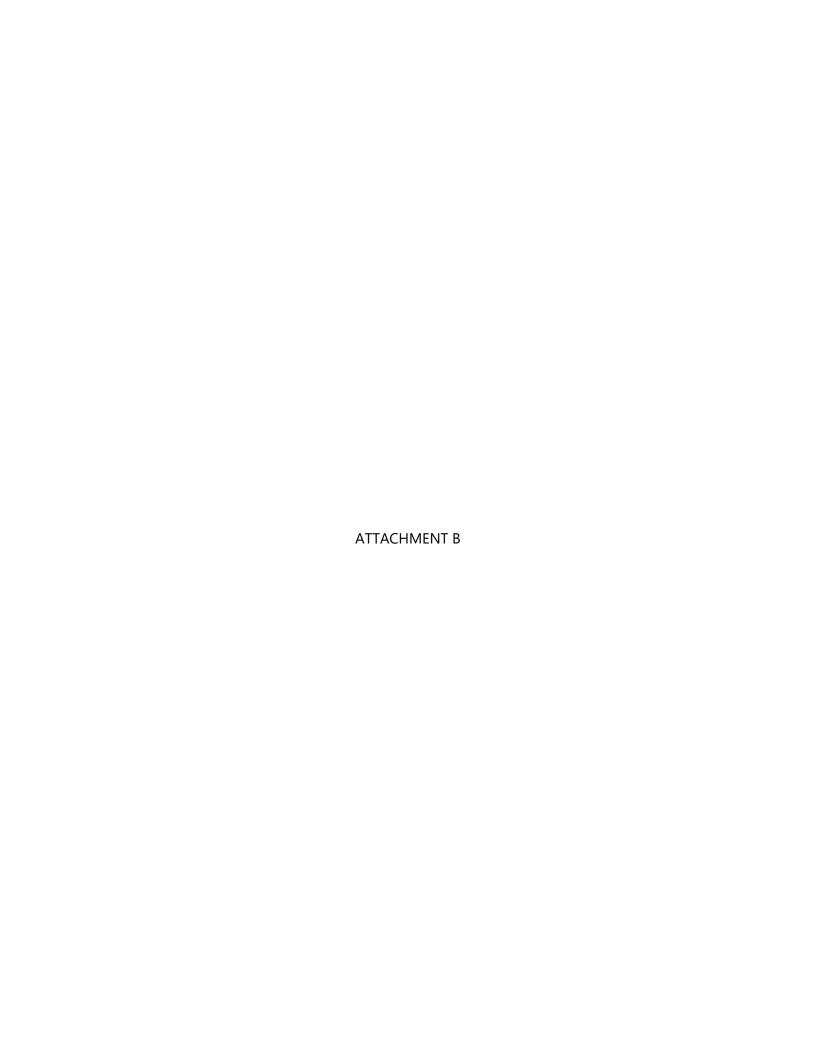
Draper Aden Associates

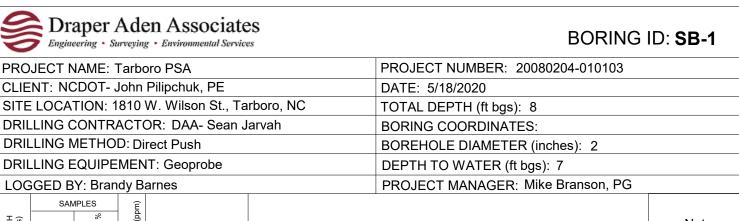
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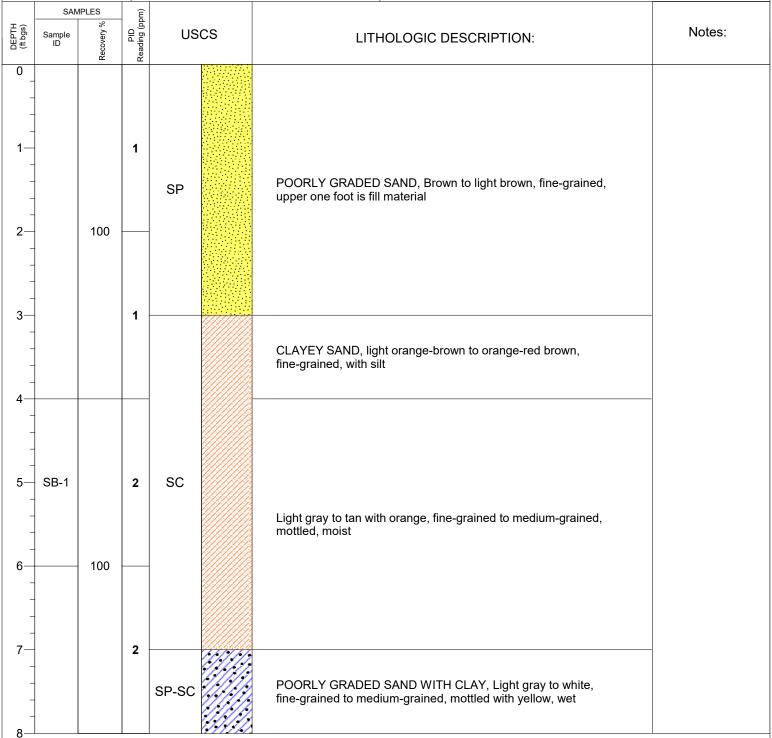
2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291 Richmond, VA Charlottesville, VA Hampton Roads, VA Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP

CHECKED: CMP DATE: 05/15/2020 **FIGURE**

PROJECT: 20080204-010203





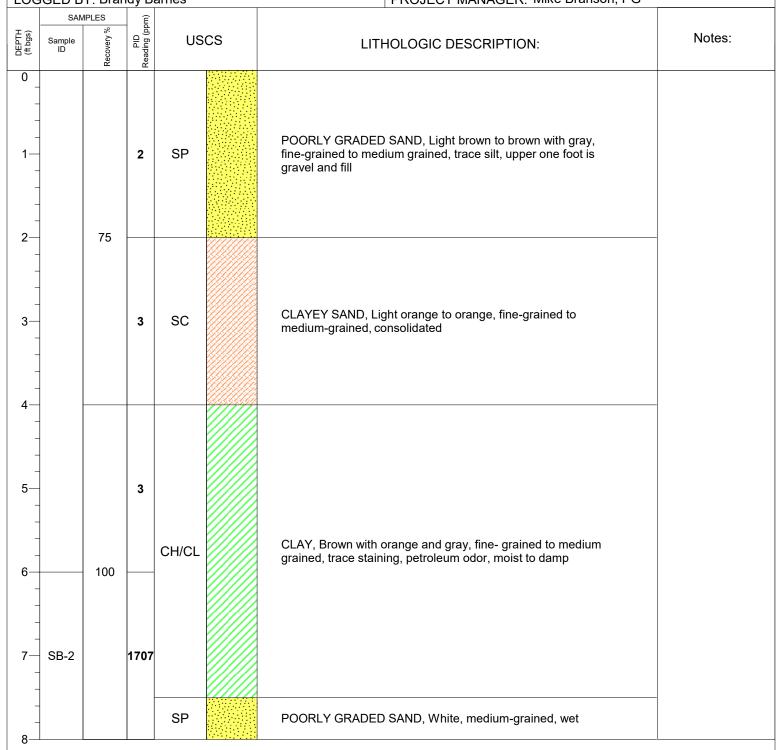


End of Borehole at 8 feet



BORING ID: SB-2

PROJECT NAME: Tarboro PSA	PROJECT NUMBER: 20080204-010103
CLIENT: NCDOT- John Pilipchuk, PE	DATE: 5/18/2020
SITE LOCATION: 1810 W. Wilson St., Tarboro, NC	TOTAL DEPTH (ft bgs): 8
DRILLING CONTRACTOR: DAA- Sean Jarvah	BORING COORDINATES:
DRILLING METHOD: Direct Push	BOREHOLE DIAMETER (inches): 2
DRILLING EQUIPEMENT: Geoprobe	DEPTH TO WATER (ft bgs): 7
LOGGED BY: Brandy Barnes	PROJECT MANAGER: Mike Branson, PG

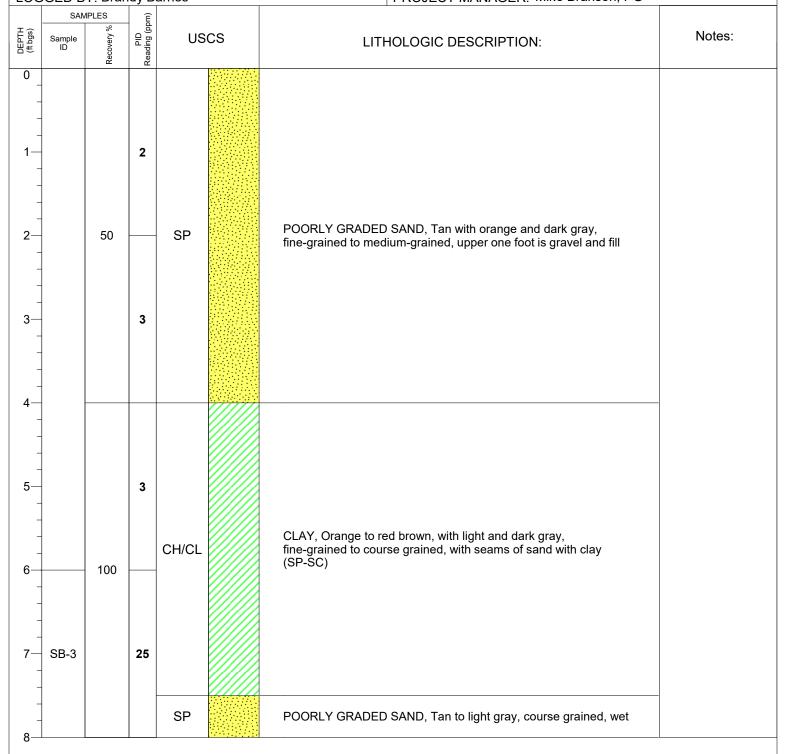


End of Borehole at 8 feet

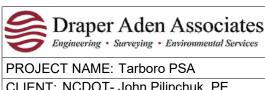


BORING ID: SB-3

PROJECT NAME: Tarboro PSA	PROJECT NUMBER: 20080204-010103
CLIENT: NCDOT- John Pilipchuk, PE	DATE: 5/18/2020
SITE LOCATION: 1810 W. Wilson St., Tarboro, NC	TOTAL DEPTH (ft bgs): 8
DRILLING CONTRACTOR: DAA- Sean Jarvah	BORING COORDINATES:
DRILLING METHOD: Direct Push	BOREHOLE DIAMETER (inches): 2
DRILLING EQUIPEMENT: Geoprobe	DEPTH TO WATER (ft bgs): 7.5
LOGGED BY: Brandy Barnes	PROJECT MANAGER: Mike Branson, PG

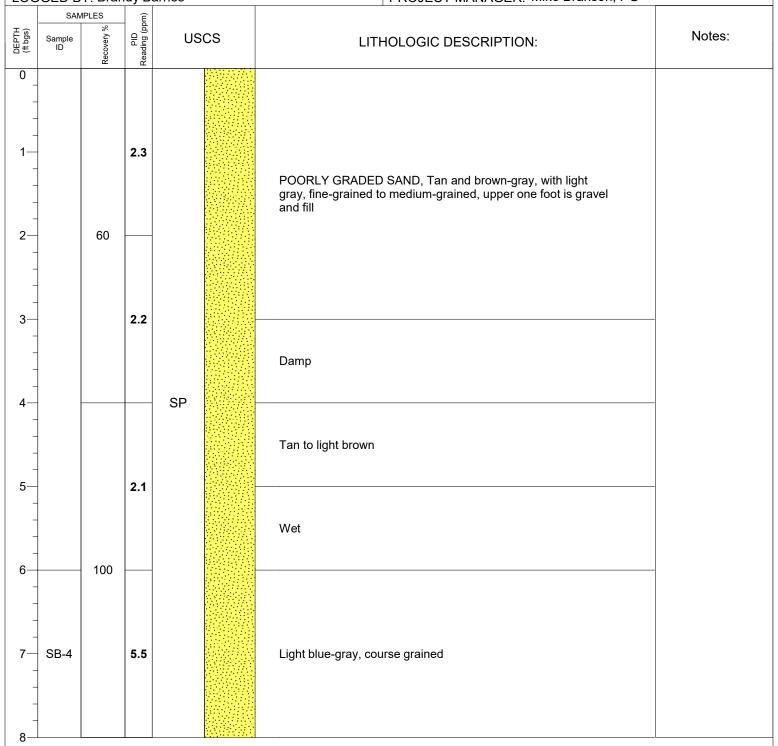


End of Borehole at 8 feet

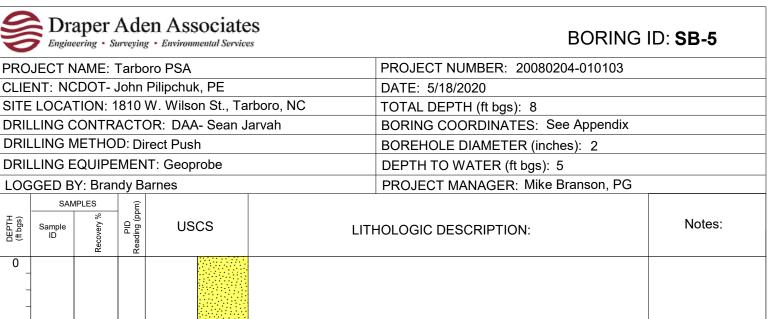


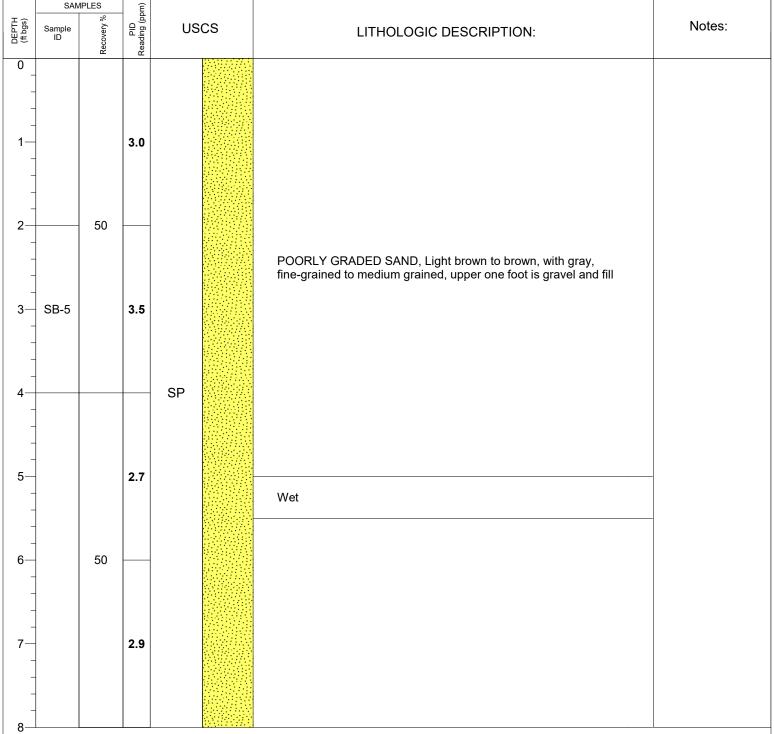
BORING ID: SB-4

PROJECT NAME: Tarboro PSA	PROJECT NUMBER: 20080204-010103
CLIENT: NCDOT- John Pilipchuk, PE	DATE: 5/18/2020
SITE LOCATION: 1810 W. Wilson St., Tarboro, NC	TOTAL DEPTH (ft bgs): 8
DRILLING CONTRACTOR: DAA- Sean Jarvah	BORING COORDINATES:
DRILLING METHOD: Direct Push	BOREHOLE DIAMETER (inches): 2
DRILLING EQUIPEMENT: Geoprobe	DEPTH TO WATER (ft bgs): 5
LOGGED BY: Brandy Barnes	PROJECT MANAGER: Mike Branson, PG

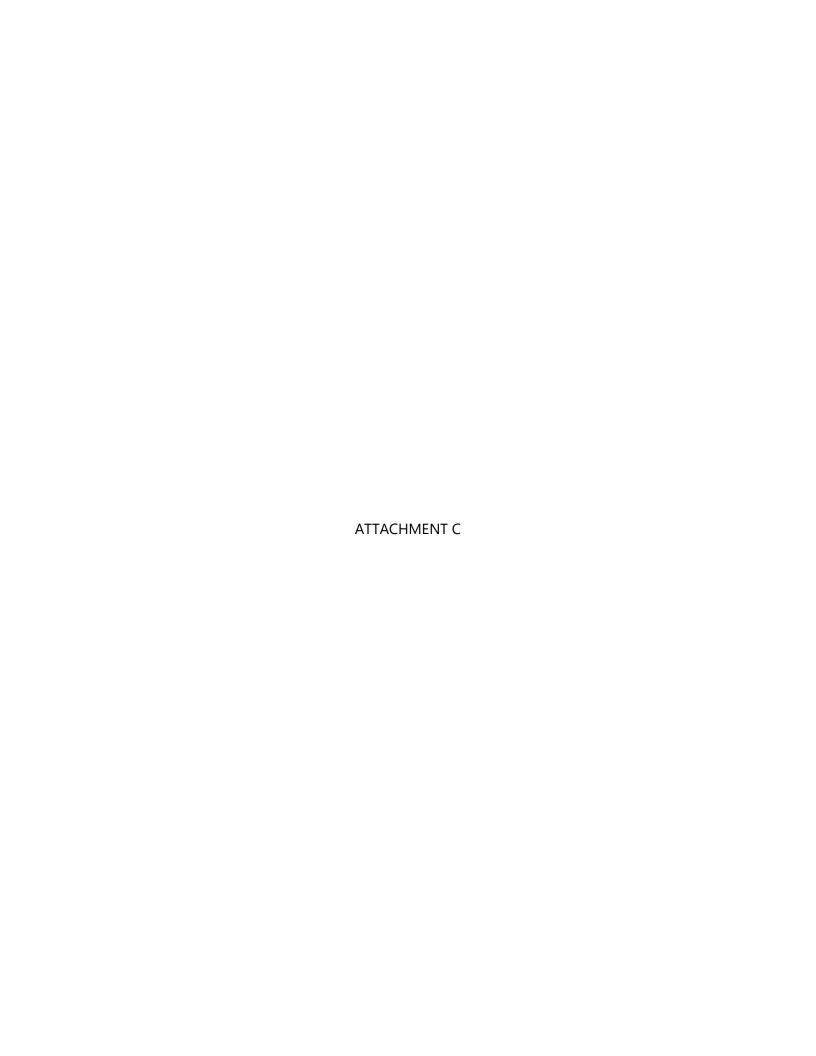


End of Borehole at 8 feet



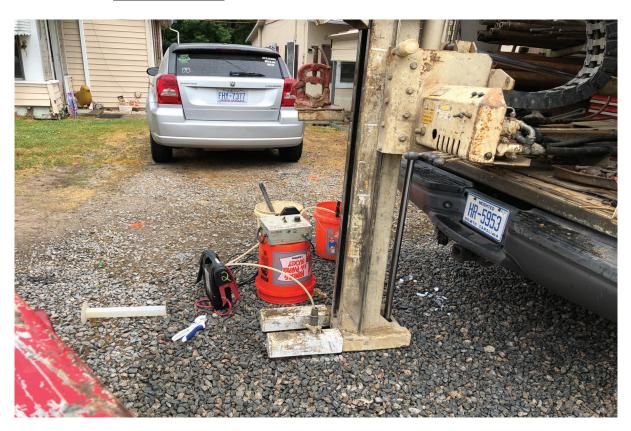


End of Borehole at 8 feet





PHOTOGRAPH 1. Location of SB-1.



PHOTOGRAPH 2. Location of SB-2.



Tarboro Preliminary Site Assessment
Parcel 46
1810 W. Wilson Street, Tarboro, NC
DAA PN: 20080204-010103



PHOTOGRAPH 3. Location of SB-3.



PHOTOGRAPH 4. Location of SB-4.



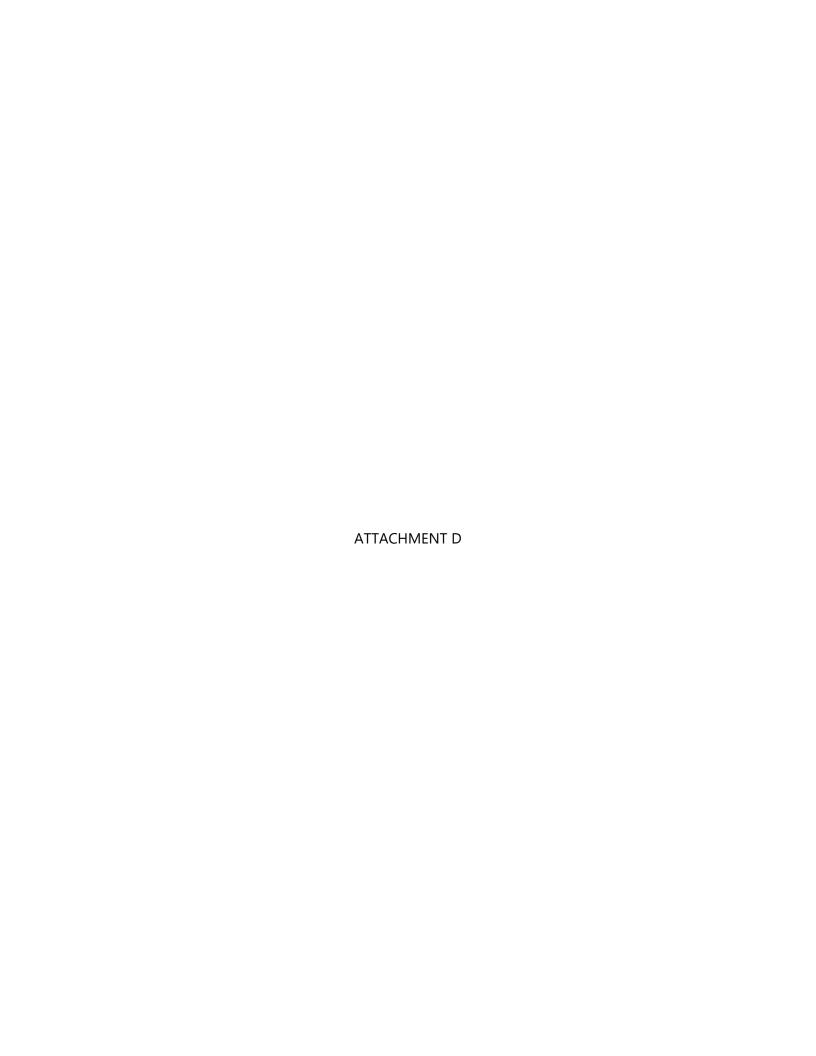


PHOTOGRAPH 5. Location of SB-5



PHOTOGRAPH 6. Soil identification and sampling area.











Hydrocarbon Analysis Results

Client: DAA Address: 114 Edinburgh S. Dr.

Cary, NC 27511

Samples taken Samples extracted Samples analysed

Monday, May 18, 2020 Monday, May 18, 2020 Wednesday, May 20, 2020

Contact: Michael Branson Operator Harry Wooten

Project: Tarboro PSA 20080204-010103

													F03640
Matrix	Sample ID	Dilution used	BTEX (C6 - C9)	GRO (C5 - C10)	DRO (C10 - C35)	TPH (C5 - C35)	Total Aromatics (C10-C35)	16 EPA PAHs	ВаР		Ratios		HC Fingerprint Match
										% light	% mid	% heavy	
S	SB-1	27.2	<0.68	<0.68	<0.68	<0.68	<0.14	<0.22	<0.027	0	0	0	PHC not detected,(BO)
S	SB-2	52.5	18.5	53.9	29	82.9	11.8	<0.42	<0.053	97.6	2.2	0.2	Deg Gas 82.2%,(FCM)
S	SB-3	25.9	<0.65	<0.65	3.8	3.8	0.83	<0.21	<0.026	0	90.2	9.8	Deg.Diesel 76.3%,(FCM)
S	SB-4	18.9	<0.47	<0.47	1.3	1.3	0.54	<0.15	<0.019	0	100	0	Deg.Light PHC 77%,(FCM)
S	SB-5	30.1	<0.75	<0.75	2	2	0.97	<0.24	<0.03	0	86.8	13.2	V.Deg.PHC 90.6%,(FCM)

Initial Calibrator QC check

OK

Final FCM QC Check OK

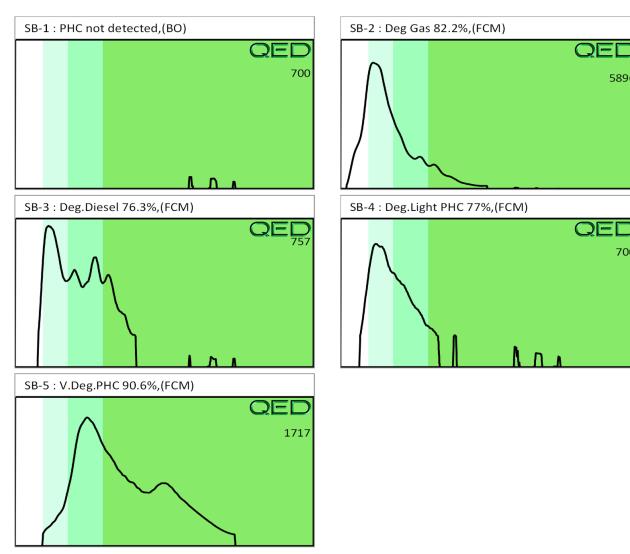
98.9 %

Results generated by a QED HC-1 analyser. Concentration values in mg/kg for soil samples and mg/L for water samples. Soil values are not corrected for moisture or stone content Fingerprints provide a tentative hydrocarbon identification. The abbreviations are:- FCM = Results calculated using Fundamental Calibration Mode: % = confidence for sample fingerprint match to library (SBS) or (LBS) = Site Specific or Library Background Subtraction applied to result: (PFM) = Poor Fingerprint Match: (T) = Turbid: (P) = Particulate present

5896

700

Tarboro PSA 20080204-010103 Project:



Ref. No HOZ						2000	Nemida	
	20 JMwy 2020 1 C100 Date/Time	d by	Accepted by	2/19/5		shed by	Belinquished by	 2
RED Lab USE ONLY	Date/Time	d by	Accepted by			shed by	Relinquished by	-
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V-1-1-1		4	88	×	×		Shoi	
43.7		1.1	8	X	×		1030	
Q. H. H	2	1	88	X	×		11010	
52.3 44.2 8.1		-45	88	R	-	4	5/18/26/09:45	
Total Wt. Tare Wt. Sample Wt.	Sample ID	" Ne.	GC Initials	Analysis Type UVF GC	TAT Requested 4 Hour 48 Hour	TAT Re 24 Hour	Sample Collection Date/Time	
trans DCE, TCE, and PCE. Specify target analytes in the space provided below.	TICAL REQUEST FORM	ND ANALY	CHAIN OF CUSTODY AND ANALYTICAL R	CHAIN	Sam s	40 g	Collected by:	
Analyses are for BTEX and Chlorinated	DIAGNOSTICS	NMENTAL D	RAPID ENVIRONMENTAL DIAGNOSTICS				Phone #:	
total BTEX, GRO, DRO, TPH, PAH total				01010	mbronsca@daa.com	S MO	Email:	
Each UVF sample will be analyzed for	U		7	10208002	3.E	Social Control	Project Ref.:	
MARBIONC Bidg, Suite 2003	TM TM		347	M.Bra	MC 270	365	Ontact:	
5598 Marvin K Moss Lane	70	Darce	1	s. pr.	nagas.	1488	Address:	
RED Lab IIC						DA P	lient Name:	

A PROPERTY OF



May 28, 2020

Mike Branson Draper Aden Associates 114 Edinburgh Drive South, Suite 200 Cary, NC 27519

Project Location: Tarboro, NC - Parcel 46

Client Job Number:

Project Number: 20080204-010103

Keny K. Mille

Laboratory Work Order Number: 20E0837

Enclosed are results of analyses for samples received by the laboratory on May 19, 2020. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Kerry K. McGee Project Manager

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Draper Aden Associates 114 Edinburgh Drive South, Suite 200 Cary, NC 27519 ATTN: Mike Branson

REPORT DATE: 5/28/2020

PURCHASE ORDER NUMBER:

PROJECT NUMBER: 20080204-010103

ANALYTICAL SUMMARY

WORK ORDER NUMBER: 20E0837

The results of analyses performed on the following samples submitted to the CON-TEST Analytical Laboratory are found in this report.

PROJECT LOCATION: Tarboro, NC - Parcel 46

FIELD SAMPLE # LAB ID: MATRIX SAMPLE DESCRIPTION TEST SUB LAB

SB-2 20E0837-01 Ground Water SW-846 8260D
SW-846 8270E

Page 3 of 30



CASE NARRATIVE SUMMARY

All reported results are within defined laboratory quality control objectives unless listed below or otherwise qualified in this report. For method 8260D:Elevated reporting limit for sample 20E0837-01 due to the high concentration of target compounds.



SW-846 8260D

Qualifications:

L-04

Laboratory fortified blank/laboratory control sample recovery and duplicate recovery are outside of control limits. Reported value for this compound is likely to be biased on the low side. Analyte & Samples(s) Qualified:

Trichlorofluoromethane (Freon 11)

20E0837-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1

RL-11

Elevated reporting limit due to high concentration of target compounds.

Analyte & Samples(s) Qualified:

20E0837-01[SB-2]

V-05

Continuing calibration verification (CCV) did not meet method specifications and was biased on the low side for this compound.

Analyte & Samples(s) Qualified:

1,1,1,2-Tetrachloroethane

20E0837-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

Bromoform

20E0837-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

Dichlorodifluoromethane (Freon 12

20E0837-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

tert-Butyl Alcohol (TBA)

20E0837-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

Trichlorofluoromethane (Freon 11)

20E0837-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

SW-846 8270E

Oualifications:

V-04

Initial calibration did not meet method specifications. Compound was calibrated using a response factor where %RSD is outside of method specified criteria. Reported result is estimated.

Analyte & Samples(s) Qualified:

Benzidine

20E0837-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

V-05

Continuing calibration verification (CCV) did not meet method specifications and was biased on the low side for this compound.

Analyte & Samples(s) Qualified:

20E0837-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

Benzidine

20E0837-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

V-06

Continuing calibration verification (CCV) did not meet method specifications and was biased on the high side for this compound.

Analyte & Samples(s) Qualified:

Benzoic Acid

20E0837-01[SB-2], S048793-CCV1

V-34

Initial calibration verification (ICV) did not meet method specifications and was biased on the low side for this compound. Reported result is estimated

Analyte & Samples(s) Qualified:

20E0837-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

20E0837-01[SB-2], S048793-CCV1

The results of analyses reported only relate to samples submitted to the Con-Test Analytical Laboratory for testing.

I certify that the analyses listed above, unless specifically listed as subcontracted, if any, were performed under my direction according to the approved methodologies listed in this document, and that based upon my inquiry of those individuals immediately responsible for obtaining the information, the material contained in this report is, to the best of my knowledge and belief, accurate and complete.

Tod E. Kopyscinski Laboratory Director



Project Location: Tarboro, NC - Parcel 46 Sample Description: Work Order: 20E0837

Date Received: 5/19/2020
Field Sample #: SB-2

Sampled: 5/18/2020 11:35

Sample ID: 20E0837-01
Sample Matrix: Ground Water

Sample Flags: RL-11			Volatilo	e Organic Co	mpounds by G	C/MS				
Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst
Acetone	ND	2000	150	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Acrylonitrile	ND	200	21	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Amyl Methyl Ether (TAME)	ND	20	5.6	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Benzene	ND	40	7.2	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromobenzene	ND	40	6.0	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromochloromethane	ND	40	13	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromodichloromethane	ND	20	6.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromoform	ND	40	18	$\mu g/L$	40	V-05	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromomethane	ND	80	55	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
2-Butanone (MEK)	ND	800	78	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Butyl Alcohol (TBA)	ND	800	170	$\mu g/L$	40	V-05	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
n-Butylbenzene	ND	40	8.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
sec-Butylbenzene	ND	40	6.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Butylbenzene	ND	40	6.8	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Butyl Ethyl Ether (TBEE)	ND	20	6.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Carbon Disulfide	ND	200	180	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Carbon Tetrachloride	ND	40	4.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chlorobenzene	ND	40	6.0	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chlorodibromomethane	ND	20	8.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chloroethane	ND	80	14	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chloroform	ND	80	6.8	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chloromethane	ND	80	18	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
2-Chlorotoluene	ND	40	4.8	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
4-Chlorotoluene	ND	40	5.6	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dibromo-3-chloropropane (DBCP)	ND	200	21	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dibromoethane (EDB)	ND	20	7.6	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Dibromomethane	ND	40	15	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dichlorobenzene	ND	40	6.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,3-Dichlorobenzene	ND	40	4.8	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,4-Dichlorobenzene	ND	40	5.2	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
trans-1,4-Dichloro-2-butene	ND	80	12	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Dichlorodifluoromethane (Freon 12)	ND	80	10	μg/L	40	V-05	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1-Dichloroethane	ND	40	6.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dichloroethane	ND	40	16	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1-Dichloroethylene	ND	40	13	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
cis-1,2-Dichloroethylene	ND	40	5.2	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
trans-1,2-Dichloroethylene	ND	40	12	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dichloropropane	ND	40	8.0	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,3-Dichloropropane	ND	20	4.4	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
2,2-Dichloropropane	ND	40	8.0	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1-Dichloropropene	ND	80	6.4	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
cis-1,3-Dichloropropene	ND	20	5.2	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
trans-1,3-Dichloropropene	ND	20	9.2	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Diethyl Ether	ND	80	14	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF

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Project Location: Tarboro, NC - Parcel 46 Sample Description: Work Order: 20E0837

Date Received: 5/19/2020
Field Sample #: SB-2

Sampled: 5/18/2020 11:35

Sample ID: 20E0837-01
Sample Matrix: Ground Water

Sample Flags: RL-11

Volatile Organic Compounds by GC/MS

Sample Plags, RL-11			701	tthe Organic Com	pounds by C	JC/M5				
								Date	Date/Time	
Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Prepared	Analyzed	Analys
Diisopropyl Ether (DIPE)	ND	20	6.8	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,4-Dioxane	ND	2000	900	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Ethylbenzene	1600	40	5.2	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Hexachlorobutadiene	ND	24	19	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
2-Hexanone (MBK)	ND	400	61	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Isopropylbenzene (Cumene)	89	80	6.8	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
p-Isopropyltoluene (p-Cymene)	ND	80	8.0	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Methyl tert-Butyl Ether (MTBE)	ND	40	10	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Methylene Chloride	ND	200	14	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
4-Methyl-2-pentanone (MIBK)	ND	400	67	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Naphthalene	320	200	12	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
n-Propylbenzene	130	40	5.2	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Styrene	ND	80	4.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1,1,2-Tetrachloroethane	ND	40	11	$\mu g/L$	40	V-05	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1,2,2-Tetrachloroethane	ND	20	8.8	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Tetrachloroethylene	ND	40	7.2	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Tetrahydrofuran	ND	400	20	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Toluene	3700	40	5.6	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2,3-Trichlorobenzene	ND	200	23	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2,4-Trichlorobenzene	ND	40	16	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,3,5-Trichlorobenzene	ND	40	12	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1,1-Trichloroethane	ND	40	8.0	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1,2-Trichloroethane	ND	40	6.4	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Trichloroethylene	ND	40	9.6	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Trichlorofluoromethane (Freon 11)	ND	80	13	μg/L	40	V-05, L-04	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2,3-Trichloropropane	ND	80	10	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	ND	40	13	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2,4-Trimethylbenzene	920	40	7.2	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,3,5-Trimethylbenzene	220	40	5.6	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Vinyl Chloride	ND	80	18	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
m+p Xylene	5600	80	12	μg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
o-Xylene	2400	40	6.8	$\mu g/L$	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Surrogates		% Reco	very	Recovery Limits	3	Flag/Qual				
1,2-Dichloroethane-d4		121		70-130					5/21/20 19:49	
Toluene-d8		95.9		70-130					5/21/20 19:49	
4-Bromofluorobenzene		102		70-130					5/21/20 19:49	



Project Location: Tarboro, NC - Parcel 46 Sample Description: Work Order: 20E0837

Date Received: 5/19/2020
Field Sample #: SB-2

Sampled: 5/18/2020 11:35

Sample ID: 20E0837-01
Sample Matrix: Ground Water

Semivolatile C	Organic Com	pounds by GC/MS	
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Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst
Acenaphthene (SIM)	0.075	0.31	0.034	μg/L	1	J	SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Acenaphthylene (SIM)	0.037	0.21	0.036	μg/L	1	J	SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Acetophenone	ND	10	0.40	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Aniline	ND	5.1	0.77	μg/L	1	V-05, V-34	SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Anthracene (SIM)	ND	0.21	0.033	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Benzidine	ND	21	17	μg/L	1	V-04, V-05	SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Benzo(a)anthracene (SIM)	ND	0.051	0.016	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Benzo(a)pyrene (SIM)	ND	0.10	0.012	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Benzo(b)fluoranthene (SIM)	ND	0.051	0.015	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Benzo(g,h,i)perylene (SIM)	ND	0.51	0.018	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Benzo(k)fluoranthene (SIM)	ND	0.21	0.012	$\mu g/L$	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Benzoic Acid	ND	10	5.5	μg/L	1	V-06	SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Bis(2-chloroethoxy)methane	ND	10	0.48	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Bis(2-chloroethyl)ether	ND	10	0.53	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Bis(2-chloroisopropyl)ether	ND	10	0.75	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Bis(2-Ethylhexyl)phthalate	ND	10	0.53	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Bromophenylphenylether	ND	10	0.30	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Butylbenzylphthalate	ND	10	0.30	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Carbazole	ND	10	0.29	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Chloroaniline	ND	10	0.35	$\mu g/L$	1	V-34	SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Chloro-3-methylphenol	ND	10	0.49	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Chloronaphthalene	ND	10	0.47	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Chlorophenol	ND	10	0.39	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Chlorophenylphenylether	ND	10	0.32	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Chrysene (SIM)	ND	0.21	0.015	$\mu g/L$	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Dibenz(a,h)anthracene (SIM)	ND	0.10	0.017	$\mu g/L$	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Dibenzofuran	ND	5.1	0.27	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Di-n-butylphthalate	ND	10	0.47	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2-Dichlorobenzene	ND	5.1	0.47	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,3-Dichlorobenzene	ND	5.1	0.47	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,4-Dichlorobenzene	ND	5.1	0.39	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
3,3-Dichlorobenzidine	ND	10	0.37	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dichlorophenol	ND	10	0.31	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Diethylphthalate	ND	10	0.23	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dimethylphenol	ND	10	0.82	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Dimethylphthalate	ND	10	0.31	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4,6-Dinitro-2-methylphenol	ND	10	2.0	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dinitrophenol	ND	10	1.7	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dinitrotoluene	ND	10	0.34	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,6-Dinitrotoluene	ND	10	0.35	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Di-n-octylphthalate	ND	10	0.54	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2-Diphenylhydrazine/Azobenzene	ND	10	0.38	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Fluoranthene (SIM)	ND	0.51	0.026	$\mu g/L$	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Fluorene (SIM)	0.10	1.0	0.035	$\mu g/L$	1	J	SW-846 8270E	5/21/20	5/28/20 1:32	IMR

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Date/Time

Date



39 Spruce Street * East Longmeadow, MA 01028 * FAX 413/525-6405 * TEL. 413/525-2332

Project Location: Tarboro, NC - Parcel 46 Sample Description: Work Order: 20E0837

Date Received: 5/19/2020
Field Sample #: SB-2

Sampled: 5/18/2020 11:35

Sample ID: 20E0837-01
Sample Matrix: Ground Water

Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Prepared	Analyzed	Analyst
Hexachlorobenzene	ND	10	0.44	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Hexachlorobutadiene	ND	10	0.61	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Hexachlorocyclopentadiene	ND	10	4.9	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Hexachloroethane	ND	10	0.54	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Indeno(1,2,3-cd)pyrene (SIM)	ND	0.10	0.018	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Isophorone	ND	10	0.31	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1-Methylnaphthalene	14	5.1	0.29	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Methylnaphthalene	30	5.1	0.27	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Methylphenol	ND	10	0.47	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
3/4-Methylphenol	ND	10	0.21	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Naphthalene	190	21	1.8	μg/L	4		SW-846 8270E	5/21/20	5/27/20 20:55	BGL
2-Nitroaniline	ND	10	0.41	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
3-Nitroaniline	ND	10	0.42	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Nitroaniline	ND	10	0.51	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Nitrobenzene	ND	10	0.42	μg/L μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Nitrophenol	ND	10	0.43	μg/L μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Nitrophenol	ND ND	10	0.43		1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
N-Nitrosodimethylamine				μg/L						
•	ND	10	1.9	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
N-Nitrosodiphenylamine/Diphenylamine	ND	10	0.30	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
N-Nitrosodi-n-propylamine	ND	10	0.53	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Pentachloronitrobenzene	ND	10	1.5	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Pentachlorophenol	ND	10	0.34	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Phenanthrene (SIM)	0.082	0.051	0.031	μg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Phenol	ND	10	0.20	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Pyrene (SIM)	ND	1.0	0.024	$\mu g/L$	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Pyridine	ND	5.1	2.9	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2,4,5-Tetrachlorobenzene	ND	10	0.34	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2,4-Trichlorobenzene	ND	5.1	0.57	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4,5-Trichlorophenol	ND	10	0.49	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4,6-Trichlorophenol	ND	10	0.34	$\mu g/L$	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Surrogates		% Reco	very	Recovery Limits		Flag/Qual				
2-Fluorophenol		35.1		15-110					5/27/20 13:13	
2-Fluorophenol		43.5		15-110					5/27/20 20:55	
Phenol-d6		27.6		15-110					5/27/20 13:13	
Phenol-d6		33.3		15-110					5/27/20 20:55	
Nitrobenzene-d5		63.9		30-130					5/28/20 1:32	
Nitrobenzene-d5		63.0		30-130					5/27/20 13:13	
Nitrobenzene-d5		71.2		30-130					5/27/20 20:55	
2-Fluorobiphenyl		64.1		30-130					5/27/20 13:13	
2-Fluorobiphenyl 2-Fluorobiphenyl		61.1 76.2		30-130 30-130					5/28/20 1:32 5/27/20 20:55	
2,4,6-Tribromophenol		71.0		15-110					5/27/20 20.33	
2,4,6-Tribromophenol		77.3		15-110					5/27/20 20:55	
p-Terphenyl-d14		64.5		30-130					5/28/20 1:32	
p-Terphenyl-d14		73.2		30-130					5/27/20 13:13	
p-Terphenyl-d14		82.3		30-130					5/27/20 13.13	
r - r				100					20.00	



Sample Extraction Data

Prep Method: SW-846 5030B Analytical Method: SW-846 8260D

Lab Number [Field ID]	Batch	Initial [mL]	Final [mL]	Date
20E0837-01 [SB-2]	B258487	0.125	5.00	05/21/20

Prep Method: SW-846 3510C Analytical Method: SW-846 8270E

Lab Number [Field ID]	Batch	Initial [mL]	Final [mL]	Date
20E0837-01 [SB-2]	B258532	975	1.00	05/21/20
20E0837-01RE1 [SB-2]	B258532	975	1.00	05/21/20

Prep Method: SW-846 3510C Analytical Method: SW-846 8270E

Lab Number [Field ID]	Batch	Initial [mL]	Final [mL]	Date
20E0837-01 [SB-2]	B258763	975	1.00	05/21/20

RPD

%REC



39 Spruce Street * East Longmeadow, MA 01028 * FAX 413/525-6405 * TEL. 413/525-2332

QUALITY CONTROL

Spike

Source

Volatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Limit	Units	Level	Result	%REC	%REC Limits	RPD	Limit	Notes
Batch B258487 - SW-846 5030B										
Blank (B258487-BLK1)				Prepared &	Analyzed: 05	/21/20				
Acetone	ND	50	$\mu \text{g/L}$							
Acrylonitrile	ND	5.0	μg/L							
tert-Amyl Methyl Ether (TAME)	ND	0.50	μg/L							
Benzene	ND	1.0	μg/L							
Bromobenzene	ND	1.0	μg/L							
Bromochloromethane	ND	1.0	μg/L							
Bromodichloromethane	ND	0.50	μg/L							
Bromoform	ND	1.0	μg/L							V-05
Bromomethane	ND	2.0	μg/L							
2-Butanone (MEK)	ND	20	μg/L							
tert-Butyl Alcohol (TBA)	ND	20	μg/L							V-05
n-Butylbenzene	ND	1.0	μg/L							
sec-Butylbenzene	ND	1.0	μg/L							
tert-Butylbenzene	ND	1.0	μg/L							
tert-Butyl Ethyl Ether (TBEE)	ND	0.50	μg/L							
Carbon Disulfide	ND	5.0	μg/L							
Carbon Tetrachloride	ND	1.0	μg/L							
Chlorobenzene	ND	1.0	μg/L							
Chlorodibromomethane	ND	0.50	μg/L							
Chloroethane	ND	2.0	μg/L							
Chloroform	ND	2.0	μg/L							
Chloromethane	ND ND	2.0	μg/L							
2-Chlorotoluene	ND ND	1.0	μg/L							
4-Chlorotoluene	ND ND	1.0	μg/L							
1,2-Dibromo-3-chloropropane (DBCP)	ND ND	5.0	μg/L							
1,2-Dibromoethane (EDB)		0.50	μg/L μg/L							
Dibromomethane (BBB)	ND	1.0	μg/L							
1,2-Dichlorobenzene	ND	1.0	μg/L μg/L							
1,3-Dichlorobenzene	ND	1.0	μg/L μg/L							
1,4-Dichlorobenzene	ND	1.0	μg/L μg/L							
trans-1,4-Dichloro-2-butene	ND	2.0	μg/L μg/L							
Dichlorodifluoromethane (Freon 12)	ND	2.0								V 05
1,1-Dichloroethane	ND	1.0	μg/L							V-05
1,2-Dichloroethane	ND		μg/L							
	ND	1.0	μg/L							
1,1-Dichloroethylene cis-1,2-Dichloroethylene	ND	1.0	μg/L							
	ND	1.0 1.0	μg/L							
trans-1,2-Dichloroethylene 1,2-Dichloropropane	ND	1.0	μg/L							
	ND		μg/L							
1,3-Dichloropropane	ND	0.50	μg/L							
2,2-Dichloropropane	ND	1.0	μg/L							
1,1-Dichloropropene	ND	2.0	μg/L							
cis-1,3-Dichloropropene	ND	0.50	μg/L							
trans-1,3-Dichloropropene	ND	0.50	μg/L							
Diethyl Ether	ND	2.0	μg/L							
Diisopropyl Ether (DIPE)	ND	0.50	μg/L							
1,4-Dioxane	ND	50	μg/L							
Ethylbenzene	ND	1.0	μg/L							
Hexachlorobutadiene	ND	0.60	μg/L							
2-Hexanone (MBK)	ND	10	μg/L							
Isopropylbenzene (Cumene)	ND	1.0	$\mu g \! / \! L$							
p-Isopropyltoluene (p-Cymene)	ND	1.0	$\mu g \! / \! L$							
Methyl tert-Butyl Ether (MTBE)	ND	1.0	μg/L							



QUALITY CONTROL

Spike

Source

%REC

RPD

Volatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
atch B258487 - SW-846 5030B										
lank (B258487-BLK1)				Prepared & A	Analyzed: 05	/21/20				
lethylene Chloride	ND	5.0	μg/L							
Methyl-2-pentanone (MIBK)	ND	10	$\mu g/L$							
aphthalene	ND	2.0	$\mu \text{g/L}$							
Propylbenzene	ND	1.0	μg/L							
yrene	ND	1.0	μg/L							
1,1,2-Tetrachloroethane	ND	1.0	μg/L							V-05
1,2,2-Tetrachloroethane	ND	0.50	μg/L							
etrachloroethylene	ND	1.0	μg/L							
trahydrofuran	ND	10	μg/L							
luene	ND	1.0	μg/L							
2,3-Trichlorobenzene	ND	5.0	μg/L							
2,4-Trichlorobenzene	ND	1.0	μg/L							
3,5-Trichlorobenzene	ND	1.0	μg/L							
1,1-Trichloroethane	ND	1.0	μg/L							
1,2-Trichloroethane	ND	1.0	μg/L							
richloroethylene	ND ND	1.0	μg/L							
richlorofluoromethane (Freon 11)	ND ND	2.0	μg/L μg/L							L-04, V-05
2,3-Trichloropropane	ND ND	2.0	μg/L μg/L							L-04, ¥-03
1,2-Trichloro-1,2,2-trifluoroethane (Freon	ND ND	1.0	μg/L μg/L							
3) 2,4-Trimethylbenzene	N.D.	1.0	ua/I							
3,5-Trimethylbenzene	ND	1.0	μg/L							
· ·	ND		μg/L							
nyl Chloride	ND	2.0	μg/L							
+p Xylene	ND	2.0	μg/L							
Xylene	ND	1.0	μg/L							
rrogate: 1,2-Dichloroethane-d4	27.5		$\mu g/L$	25.0		110	70-130			
rrogate: Toluene-d8	23.6		$\mu g/L$	25.0		94.5	70-130			
rrogate: 4-Bromofluorobenzene	25.2		μg/L	25.0		101	70-130			
CS (B258487-BS1)				Prepared & A	Analyzed: 05	/21/20				
cetone	89.2	50	μg/L	100		89.2	70-160			
crylonitrile	11.2	5.0	μg/L	10.0		112	70-130			
			/T	400		83.6	70-130			
	8.36	0.50	μg/L	10.0			70-150			
	8.36 10.5	0.50 1.0	μg/L μg/L	10.0		105	70-130			
enzene										
enzene	10.5	1.0	$\mu g/L$	10.0		105	70-130			
enzene romobenzene romochloromethane	10.5 10.3	1.0 1.0	μg/L μg/L	10.0 10.0		105 103	70-130 70-130			
enzene romobenzene romochloromethane romodichloromethane	10.5 10.3 10.8	1.0 1.0 1.0	μg/L μg/L μg/L	10.0 10.0 10.0		105 103 108	70-130 70-130 70-130			V-05
enzene romobenzene romochloromethane romodichloromethane romoform	10.5 10.3 10.8 8.78	1.0 1.0 1.0 0.50	μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0		105 103 108 87.8	70-130 70-130 70-130 70-130			V-05
enzene romobenzene romochloromethane romodichloromethane romoform romomethane	10.5 10.3 10.8 8.78 8.87	1.0 1.0 1.0 0.50 1.0	μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7	70-130 70-130 70-130 70-130 70-130			V-05
enzene romobenzene romochloromethane romoform romomethane romomethane Butanone (MEK)	10.5 10.3 10.8 8.78 8.87 5.25	1.0 1.0 1.0 0.50 1.0 2.0	μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5	70-130 70-130 70-130 70-130 70-130 40-160			V-05
enzene comobenzene comochloromethane comodichloromethane comoform comomethane Butanone (MEK) rt-Butyl Alcohol (TBA)	10.5 10.3 10.8 8.78 8.87 5.25	1.0 1.0 1.0 0.50 1.0 2.0	μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110	70-130 70-130 70-130 70-130 70-130 40-160 40-160			
enzene comobenzene comochloromethane comodichloromethane comoform comomethane Butanone (MEK) ct-Butyl Alcohol (TBA) Butylbenzene	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7	1.0 1.0 1.0 0.50 1.0 2.0 20	µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110 68.8	70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160			
enzene comobenzene comochloromethane comodichloromethane comoform comomethane Butanone (MEK) ct-Butyl Alcohol (TBA) Butylbenzene c-Butylbenzene	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7	1.0 1.0 1.0 0.50 1.0 2.0 20 20	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 100 100		105 103 108 87.8 88.7 52.5 110 68.8 117	70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130			
enzene comobenzene comochloromethane comodichloromethane comomethane comomethane Butanone (MEK) ct-Butyl Alcohol (TBA) Butylbenzene cc-Butylbenzene ct-Butylbenzene	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0	1.0 1.0 0.50 1.0 2.0 20 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 100 100 10		105 103 108 87.8 88.7 52.5 110 68.8 117	70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130			
enzene comobenzene comodichloromethane comodichloromethane comodichloromethane comomethane Butanone (MEK) ct-Butyl Alcohol (TBA) Butylbenzene ct-Butylbenzene ct-Butylbenzene ct-Butyl Ether (TBEE)	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32	1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 0.50	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 100 100 10		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115	70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130			
enzene comobenzene comodichloromethane comodichloromethane comoform comomethane Butanone (MEK) ct-Butyl Alcohol (TBA) Butylbenzene ct-Butylbenzene ct-Butyl Ethyl Ether (TBEE) urbon Disulfide	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32 8.24	1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 0.50	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 100 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115 93.2 82.4	70-130 70-130 70-130 70-130 70-130 40-160 40-160 70-130 70-130 70-130 70-130			
enzene comobenzene comodichloromethane comodichloromethane comoform comomethane Butanone (MEK) ct-Butyl Alcohol (TBA) Butylbenzene ct-Butylbenzene ct-Butylbenzene ct-Butyl Ethyl Ether (TBEE) curbon Disulfide curbon Tetrachloride	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32 8.24 8.42	1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 0.50 5.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 100 100 10		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115 93.2 82.4 84.2	70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130			
enzene comobenzene comochloromethane comochloromethane comoform comomethane Butanone (MEK) rt-Butyl Alcohol (TBA) Butylbenzene c-Butylbenzene rt-Butyl Ethyl Ether (TBEE) arbon Disulfide arbon Tetrachloride allorobenzene	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32 8.24 8.42 9.77	1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 0.50 5.0 1.0 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115 93.2 82.4 84.2 97.7	70-130 70-130 70-130 70-130 70-130 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130			
enzene comobenzene comobenzene comochloromethane comodichloromethane comoform comomethane Butanone (MEK) tt-Butyl Alcohol (TBA) Butylbenzene ct-Butylbenzene tt-Butyl Ether (TBEE) urbon Disulfide urbon Tetrachloride allorobenzene tolorodibromomethane	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32 8.24 8.42 9.77 8.21	1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 1.0 1.0 0.50 5.0 1.0 0.50	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115 93.2 82.4 84.2 97.7 82.1	70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130 70-130			
enzene romobenzene romochloromethane romodichloromethane romoform romomethane Butanone (MEK) rt-Butyl Alcohol (TBA) Butylbenzene c-Butylbenzene rt-Butyl Ethyl Ether (TBEE) arbon Disulfide arbon Tetrachloride hlorodibromomethane	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32 8.24 8.42 9.77 8.21 7.43	1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 0.50 5.0 1.0 0.50 2.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115 93.2 82.4 84.2 97.7 82.1 74.3	70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130 70-130 70-130 70-130			
rt-Amyl Methyl Ether (TAME) enzene romobenzene romochloromethane romodichloromethane romomethane Butanone (MEK) rt-Butyl Alcohol (TBA) Butylbenzene ec-Butylbenzene rt-Butyl Ethyl Ether (TBEE) arbon Disulfide arbon Tetrachloride hlorodibromomethane hlorodirm hloroform	10.5 10.3 10.8 8.78 8.87 5.25 110 68.8 11.7 12.0 11.5 9.32 8.24 8.42 9.77 8.21	1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 1.0 1.0 0.50 5.0 1.0 0.50	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		105 103 108 87.8 88.7 52.5 110 68.8 117 120 115 93.2 82.4 84.2 97.7 82.1	70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130 70-130			



QUALITY CONTROL

Spike

Source

%REC

RPD

Volatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes	
Batch B258487 - SW-846 5030B											
LCS (B258487-BS1)				Prepared & A	Analyzed: 05	/21/20					
4-Chlorotoluene	10.9	1.0	$\mu g\!/\!L$	10.0		109	70-130				
1,2-Dibromo-3-chloropropane (DBCP)	9.15	5.0	μg/L	10.0		91.5	70-130				
1,2-Dibromoethane (EDB)	9.22	0.50	$\mu g \! / \! L$	10.0		92.2	70-130				
Dibromomethane	8.99	1.0	$\mu g/L$	10.0		89.9	70-130				
1,2-Dichlorobenzene	10.2	1.0	$\mu g/L$	10.0		102	70-130				
1,3-Dichlorobenzene	10.8	1.0	μg/L	10.0		108	70-130				
1,4-Dichlorobenzene	10.1	1.0	$\mu g/L$	10.0		101	70-130				
trans-1,4-Dichloro-2-butene	10.5	2.0	$\mu g/L$	10.0		105	70-130				
Dichlorodifluoromethane (Freon 12)	6.26	2.0	μg/L	10.0		62.6	40-160			V-05	†
1,1-Dichloroethane	10.8	1.0	$\mu g/L$	10.0		108	70-130				
1,2-Dichloroethane	8.15	1.0	$\mu g \! / \! L$	10.0		81.5	70-130				
1,1-Dichloroethylene	8.71	1.0	$\mu g/L$	10.0		87.1	70-130				
cis-1,2-Dichloroethylene	11.3	1.0	$\mu \text{g/L}$	10.0		113	70-130				
trans-1,2-Dichloroethylene	9.51	1.0	$\mu \text{g/L}$	10.0		95.1	70-130				
1,2-Dichloropropane	10.1	1.0	$\mu g/L$	10.0		101	70-130				
1,3-Dichloropropane	9.46	0.50	μg/L	10.0		94.6	70-130				
2,2-Dichloropropane	9.00	1.0	$\mu g/L$	10.0		90.0	40-130				†
1,1-Dichloropropene	9.29	2.0	μg/L	10.0		92.9	70-130				
cis-1,3-Dichloropropene	8.99	0.50	μg/L	10.0		89.9	70-130				
trans-1,3-Dichloropropene	8.97	0.50	μg/L	10.0		89.7	70-130				
Diethyl Ether	9.14	2.0	μg/L	10.0		91.4	70-130				
Diisopropyl Ether (DIPE)	10.5	0.50	μg/L	10.0		105	70-130				
1,4-Dioxane	125	50	μg/L	100		125	40-130				†
Ethylbenzene	10.6	1.0	μg/L	10.0		106	70-130				'
Hexachlorobutadiene	9.47	0.60	μg/L	10.0		94.7	70-130				
2-Hexanone (MBK)	96.2	10	μg/L	100		96.2	70-150				†
Isopropylbenzene (Cumene)		1.0	μg/L μg/L	10.0		103	70-100				'
p-Isopropyltoluene (p-Cymene)	10.3	1.0	μg/L μg/L	10.0		103	70-130				
Methyl tert-Butyl Ether (MTBE)	10.8	1.0	μg/L μg/L	10.0		80.7	70-130				
Methylene Chloride	8.07	5.0	μg/L μg/L	10.0		100	70-130				
4-Methyl-2-pentanone (MIBK)	10.0	10	μg/L μg/L	10.0		99.2	70-130				+
Naphthalene	99.2	2.0	μg/L μg/L								† †
•	10.8			10.0		108	40-130				1
n-Propylbenzene	10.4	1.0	μg/L	10.0		104	70-130				
Styrene	10.4	1.0	μg/L	10.0		104	70-130			****	
1,1,2-Tetrachloroethane	8.45	1.0	μg/L	10.0		84.5	70-130			V-05	
1,1,2,2-Tetrachloroethane	9.22	0.50	μg/L	10.0		92.2	70-130				
Tetrachloroethylene	9.18	1.0	μg/L	10.0		91.8	70-130				
Tetrahydrofuran	11.2	10	μg/L	10.0		112	70-130				
Toluene	9.73	1.0	μg/L	10.0		97.3	70-130				
1,2,3-Trichlorobenzene	9.65	5.0	μg/L	10.0		96.5	70-130				
1,2,4-Trichlorobenzene	10.4	1.0	μg/L	10.0		104	70-130				
1,3,5-Trichlorobenzene	9.75	1.0	μg/L	10.0		97.5	70-130				
1,1,1-Trichloroethane	9.20	1.0	μg/L	10.0		92.0	70-130				
1,1,2-Trichloroethane	9.18	1.0	μg/L	10.0		91.8	70-130				
Trichloroethylene	8.72	1.0	μg/L	10.0		87.2	70-130				
Trichlorofluoromethane (Freon 11)	6.45	2.0	μg/L	10.0		64.5 *	70-130			V-05, L-04	
1,2,3-Trichloropropane	9.56	2.0	$\mu g/L$	10.0		95.6	70-130				
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	9.13	1.0	μg/L	10.0		91.3	70-130				
1,2,4-Trimethylbenzene	10.6	1.0	μg/L	10.0		106	70-130				
1,3,5-Trimethylbenzene	10.0	1.0	$\mu g/L$	10.0		100	70-130				
Vinyl Chloride	6.50	2.0	$\mu g/L$	10.0		65.0	40-160				†



QUALITY CONTROL

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes	
Batch B258487 - SW-846 5030B						. ,		2			_
LCS (B258487-BS1)				Prepared &	Analyzed: 05/	/21/20					_
n+p Xylene	21.4	2.0	μg/L	20.0	maryzea. 03/	107	70-130				_
o-Xylene	10.7	1.0	μg/L μg/L	10.0		107	70-130				
		1.0									
Surrogate: 1,2-Dichloroethane-d4	24.7		μg/L	25.0		98.8	70-130				
Surrogate: Toluene-d8	24.6		μg/L	25.0		98.5	70-130				
Surrogate: 4-Bromofluorobenzene	26.9		μg/L	25.0		107	70-130				
LCS Dup (B258487-BSD1)				Prepared &	Analyzed: 05	/21/20					
Acetone	91.1	50	$\mu g/L$	100		91.1	70-160	2.08	25		
crylonitrile	10.4	5.0	$\mu g/L$	10.0		104	70-130	7.52	25		
ert-Amyl Methyl Ether (TAME)	8.33	0.50	$\mu g/L$	10.0		83.3	70-130	0.359	25		
Benzene	10.4	1.0	μg/L	10.0		104	70-130	0.960	25		
Bromobenzene	10.2	1.0	μg/L	10.0		102	70-130	0.195	25		
Bromochloromethane	10.5	1.0	$\mu g/L$	10.0		105	70-130	2.92	25		
Bromodichloromethane	8.94	0.50	$\mu g/L$	10.0		89.4	70-130	1.81	25		
romoform	8.36	1.0	μg/L	10.0		83.6	70-130	5.92	25	V-05	
romomethane	5.82	2.0	μg/L	10.0		58.2	40-160	10.3	25		
-Butanone (MEK)	111	20	μg/L	100		111	40-160	0.987	25		
ert-Butyl Alcohol (TBA)	64.0	20	μg/L	100		64.0	40-160	7.29	25	V-05	
Butylbenzene	11.7	1.0	μg/L	10.0		117	70-130	0.171	25		
ec-Butylbenzene	12.1	1.0	μg/L	10.0		121	70-130	1.16	25		
rt-Butylbenzene	11.6	1.0	μg/L	10.0		116	70-130	0.782	25		
rt-Butyl Ethyl Ether (TBEE)	9.64	0.50	μg/L μg/L	10.0		96.4	70-130	3.38	25		
arbon Disulfide	7.55	5.0	μg/L μg/L	10.0		75.5	70-130	8.74	25		
arbon Tetrachloride	7.33 8.48	1.0	μg/L μg/L	10.0		84.8	70-130	0.710	25		
hlorobenzene	8.48 9.61	1.0	μg/L μg/L	10.0		96.1	70-130	1.65	25		
hlorodibromomethane	9.61 8.24	0.50	μg/L μg/L	10.0		82.4	70-130	0.365	25		
hloroethane	8.24 7.52	2.0	μg/L μg/L	10.0		75.2	70-130	1.20	25		
hloroform		2.0	μg/L μg/L	10.0		108	70-130	1.20	25		
hloromethane	10.8	2.0	μg/L μg/L								
-Chlorotoluene	6.08	1.0		10.0		60.8	40-160	1.79	25 25		
-Chlorotoluene	10.8	1.0	μg/L μg/I	10.0		108	70-130	0.739	25 25		
	10.9		μg/L	10.0		109	70-130	0.276	25		
,2-Dibromo-3-chloropropane (DBCP)	9.06	5.0	μg/L	10.0		90.6	70-130	0.988	25		
,2-Dibromoethane (EDB)	9.22	0.50	μg/L ug/I	10.0		92.2	70-130	0.00	25		
Dishlorahanana	8.93	1.0	μg/L	10.0		89.3	70-130	0.670	25		
,2-Dichlorobenzene	10.3	1.0	μg/L	10.0		103	70-130	1.46	25 25		
,3-Dichlorobenzene	10.7	1.0	μg/L	10.0		107	70-130	1.21	25		
,4-Dichlorobenzene	10.3	1.0	μg/L	10.0		103	70-130	1.67	25		
ans-1,4-Dichloro-2-butene	9.18	2.0	μg/L	10.0		91.8	70-130	13.1	25		
Dichlorodifluoromethane (Freon 12)	6.23	2.0	μg/L	10.0		62.3	40-160	0.480	25	V-05	
1-Dichloroethane	11.1	1.0	μg/L	10.0		111	70-130	2.65	25		
2-Dichloroethane	8.22	1.0	μg/L	10.0		82.2	70-130	0.855	25		
1-Dichloroethylene	8.87	1.0	μg/L	10.0		88.7	70-130	1.82	25		
s-1,2-Dichloroethylene	11.2	1.0	μg/L	10.0		112	70-130	1.42	25		
ans-1,2-Dichloroethylene	9.45	1.0	μg/L	10.0		94.5	70-130	0.633	25		
2-Dichloropropane	9.71	1.0	μg/L	10.0		97.1	70-130	4.04	25		
3-Dichloropropane	9.45	0.50	μg/L	10.0		94.5	70-130	0.106	25		
,2-Dichloropropane	9.06	1.0	$\mu g \! / \! L$	10.0		90.6	40-130	0.664	25		
1-Dichloropropene	9.00	2.0	$\mu g/L$	10.0		90.0	70-130	3.17	25		
is-1,3-Dichloropropene	9.34	0.50	$\mu g/L$	10.0		93.4	70-130	3.82	25		
rans-1,3-Dichloropropene	8.87	0.50	$\mu g/L$	10.0		88.7	70-130	1.12	25		
Diethyl Ether	9.63	2.0	$\mu g/L$	10.0		96.3	70-130	5.22	25		
Diisopropyl Ether (DIPE)	10.7	0.50	μg/L	10.0		107	70-130	1.60	25		



QUALITY CONTROL

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes	
Batch B258487 - SW-846 5030B											
LCS Dup (B258487-BSD1)				Prepared &	Analyzed: 05/	/21/20					
1,4-Dioxane	127	50	μg/L	100		127	40-130	1.84	50		 † ‡
Ethylbenzene	10.4	1.0	$\mu \text{g/L}$	10.0		104	70-130	1.05	25		
Hexachlorobutadiene	9.30	0.60	$\mu g\!/\!L$	10.0		93.0	70-130	1.81	25		
2-Hexanone (MBK)	97.7	10	$\mu \text{g/L}$	100		97.7	70-160	1.60	25		†
Isopropylbenzene (Cumene)	10.2	1.0	$\mu \text{g/L}$	10.0		102	70-130	1.37	25		
p-Isopropyltoluene (p-Cymene)	10.7	1.0	$\mu \text{g/L}$	10.0		107	70-130	0.928	25		
Methyl tert-Butyl Ether (MTBE)	8.13	1.0	$\mu g\!/\!L$	10.0		81.3	70-130	0.741	25		
Methylene Chloride	9.70	5.0	$\mu g\!/\!L$	10.0		97.0	70-130	3.15	25		
4-Methyl-2-pentanone (MIBK)	102	10	$\mu g\!/\!L$	100		102	70-160	2.60	25		†
Naphthalene	11.1	2.0	$\mu g/L$	10.0		111	40-130	2.92	25		†
n-Propylbenzene	10.4	1.0	$\mu \text{g/L}$	10.0		104	70-130	0.673	25		
Styrene	10.4	1.0	$\mu g/L$	10.0		104	70-130	0.481	25		
1,1,1,2-Tetrachloroethane	8.44	1.0	$\mu \text{g/L}$	10.0		84.4	70-130	0.118	25	V-05	
1,1,2,2-Tetrachloroethane	9.22	0.50	$\mu g/L$	10.0		92.2	70-130	0.00	25		
Tetrachloroethylene	8.98	1.0	$\mu \text{g/L}$	10.0		89.8	70-130	2.20	25		
Tetrahydrofuran	10.6	10	$\mu \text{g/L}$	10.0		106	70-130	5.32	25		
Toluene	9.56	1.0	$\mu \text{g/L}$	10.0		95.6	70-130	1.76	25		
1,2,3-Trichlorobenzene	9.77	5.0	$\mu \text{g/L}$	10.0		97.7	70-130	1.24	25		
1,2,4-Trichlorobenzene	10.6	1.0	$\mu \text{g/L}$	10.0		106	70-130	1.62	25		
1,3,5-Trichlorobenzene	9.98	1.0	$\mu g/L$	10.0		99.8	70-130	2.33	25		
1,1,1-Trichloroethane	8.89	1.0	$\mu g/L$	10.0		88.9	70-130	3.43	25		
1,1,2-Trichloroethane	9.44	1.0	$\mu \text{g/L}$	10.0		94.4	70-130	2.79	25		
Trichloroethylene	9.17	1.0	$\mu g/L$	10.0		91.7	70-130	5.03	25		
Trichlorofluoromethane (Freon 11)	6.68	2.0	$\mu g/L$	10.0		66.8	* 70-130	3.50	25	L-04, V-05	
1,2,3-Trichloropropane	9.81	2.0	$\mu g/L$	10.0		98.1	70-130	2.58	25		
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	8.91	1.0	μg/L	10.0		89.1	70-130	2.44	25		
1,2,4-Trimethylbenzene	10.7	1.0	$\mu g/L$	10.0		107	70-130	0.282	25		
1,3,5-Trimethylbenzene	9.97	1.0	$\mu g \! / \! L$	10.0		99.7	70-130	0.500	25		
Vinyl Chloride	6.51	2.0	$\mu g \! / \! L$	10.0		65.1	40-160	0.154	25		†
m+p Xylene	21.3	2.0	$\mu g/L$	20.0		106	70-130	0.749	25		
o-Xylene	10.6	1.0	$\mu g/L$	10.0		106	70-130	1.41	25		
Surrogate: 1,2-Dichloroethane-d4	25.2		μg/L	25.0		101	70-130				_
Surrogate: Toluene-d8	24.6		$\mu g/L$	25.0		98.4	70-130				
Surrogate: 4-Bromofluorobenzene	26.3		$\mu g/L$	25.0		105	70-130				



QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch B258532 - SW-846 3510C										
Blank (B258532-BLK1)				Prepared: 05	5/21/20 Anal	yzed: 05/26/2	20			
Acetophenone	ND	10	μg/L					_		
Aniline	ND	5.0	$\mu \text{g/L}$							V-05
Benzidine	ND	20	$\mu \text{g/L}$							V-04, V-05
Benzoic Acid	ND	10	$\mu \text{g/L}$							
Bis(2-chloroethoxy)methane	ND	10	μg/L							
Bis(2-chloroethyl)ether	ND	10	μg/L							
Bis(2-chloroisopropyl)ether	ND	10	μg/L							
Bis(2-Ethylhexyl)phthalate	ND	10	μg/L							
-Bromophenylphenylether	ND	10	μg/L							
Butylbenzylphthalate	ND	10	μg/L							
Carbazole	ND	10	μg/L							_
-Chloroaniline	ND	10	μg/L							V-34
-Chloro-3-methylphenol	ND	10	μg/L							
-Chlorophonol	ND	10	μg/L							
-Chlorophenol -Chlorophenylphenylether	ND	10	μg/L							
-Chlorophenylphenylether Dibenzofuran	ND	10	μg/L							
Dienzoturan Di-n-butylphthalate	ND	5.0 10	μg/L μg/L							
,2-Dichlorobenzene	ND	5.0								
,3-Dichlorobenzene	ND	5.0	μg/L μg/L							
,4-Dichlorobenzene	ND ND	5.0	μg/L μg/L							
,3-Dichlorobenzidine	ND ND	10	μg/L μg/L							
,4-Dichlorophenol	ND ND	10	μg/L μg/L							
Diethylphthalate	ND ND	10	μg/L							
,4-Dimethylphenol	ND ND	10	μg/L μg/L							
Dimethylphthalate	ND ND	10	μg/L μg/L							
,6-Dinitro-2-methylphenol	ND	10	μg/L							
,4-Dinitrophenol	ND	10	μg/L							
,4-Dinitrotoluene	ND	10	μg/L							
,,6-Dinitrotoluene	ND	10	μg/L							
Di-n-octylphthalate	ND	10	μg/L							
,2-Diphenylhydrazine/Azobenzene	ND	10	μg/L							
Hexachlorobenzene	ND	10	μg/L							
Hexachlorobutadiene	ND	10	$\mu g/L$							
Hexachlorocyclopentadiene	ND	10	$\mu g/L$							
Hexachloroethane	ND	10	$\mu g \! / \! L$							
sophorone	ND	10	$\mu g/L$							
-Methylnaphthalene	ND	5.0	$\mu \text{g/L}$							
-Methylphenol	ND	10	$\mu g \! / \! L$							
/4-Methylphenol	ND	10	$\mu g \! / \! L$							
-Nitroaniline	ND	10	μg/L							
-Nitroaniline	ND	10	μg/L							
-Nitroaniline	ND	10	μg/L							
Vitrobenzene	ND	10	μg/L							
-Nitrophenol	ND	10	μg/L							
-Nitrophenol	ND	10	μg/L							
N-Nitrosodimethylamine	ND	10	μg/L							
N-Nitrosodiphenylamine/Diphenylamine	ND	10	μg/L							
N-Nitrosodi-n-propylamine	ND	10	μg/L							
Pentachloronitrobenzene	ND	10	μg/L							
Pentachlorophenol	ND	10 10	μg/L μg/L							



QUALITY CONTROL

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
eatch B258532 - SW-846 3510C										
Blank (B258532-BLK1)				Prepared: 05	5/21/20 Anal	yzed: 05/26/2	20			
yridine	ND	5.0	μg/L							
,2,4,5-Tetrachlorobenzene	ND	10	$\mu g/L$							
,2,4-Trichlorobenzene	ND	5.0	$\mu g/L$							
,4,5-Trichlorophenol	ND	10	μg/L							
,4,6-Trichlorophenol	ND	10	$\mu g/L$							
urrogate: 2-Fluorophenol	101		μg/L	200		50.3	15-110			
urrogate: Phenol-d6	70.3		$\mu g/L$	200		35.2	15-110			
urrogate: Nitrobenzene-d5	80.6		$\mu g/L$	100		80.6	30-130			
urrogate: 2-Fluorobiphenyl	80.5		$\mu g/L$	100		80.5	30-130			
urrogate: 2,4,6-Tribromophenol	169		$\mu g/L$	200		84.5	15-110			
urrogate: p-Terphenyl-d14	93.6		$\mu g/L$	100		93.6	30-130			
CS (B258532-BS1)				Prepared: 05	5/21/20 Anal	yzed: 05/26/2	20			
cetophenone	39.4	10	μg/L	50.0		78.7	40-140			
niline	31.7	5.0	μg/L	50.0		63.4	40-140			V-05
enzidine	30.5	20	μg/L	50.0		60.9	40-140			V-04, V-05
enzoic Acid	17.6	10	μg/L	50.0		35.2	10-130			,
is(2-chloroethoxy)methane	40.9	10	μg/L	50.0		81.8	40-140			
is(2-chloroethyl)ether	38.8	10	μg/L	50.0		77.5	40-140			
is(2-chloroisopropyl)ether	43.4	10	μg/L	50.0		86.9	40-140			
is(2-Ethylhexyl)phthalate	44.4	10	μg/L	50.0		88.7	40-140			
Bromophenylphenylether	40.1	10	μg/L	50.0		80.2	40-140			
utylbenzylphthalate	40.6	10	μg/L	50.0		81.2	40-140			
arbazole	40.7	10	μg/L	50.0		81.4	40-140			
Chloroaniline	35.8	10	μg/L	50.0		71.7	40-140			V-34
Chloro-3-methylphenol	38.5	10	μg/L	50.0		77.1	30-130			, , ,
Chloronaphthalene	34.0	10	μg/L	50.0		68.0	40-140			
Chlorophenol	35.1	10	μg/L	50.0		70.2	30-130			
Chlorophenylphenylether	39.1	10	μg/L	50.0		78.2	40-140			
bibenzofuran	40.6	5.0	μg/L μg/L	50.0		81.1	40-140			
i-n-butylphthalate	43.6	10	μg/L μg/L	50.0		87.3	40-140			
2-Dichlorobenzene	28.9	5.0	μg/L μg/L	50.0		57.9	40-140			
3-Dichlorobenzene	28.9	5.0	μg/L μg/L	50.0		54.3	40-140			
4-Dichlorobenzene	27.5	5.0	μg/L μg/L	50.0		55.0	40-140			
3-Dichlorobenzidine		10	μg/L μg/L	50.0		88.9	40-140			
4-Dichlorophenol	44.5 38.3	10	μg/L μg/L	50.0		76.6	30-130			
iethylphthalate		10	μg/L μg/L	50.0		80.7	40-140			
4-Dimethylphenol	40.4	10	μg/L μg/L	50.0		65.4	30-130			
imethylphthalate	32.7	10	μg/L μg/L	50.0		80.1	40-140			
6-Dinitro-2-methylphenol	40.0	10	μg/L μg/L	50.0		79.1	30-130			
4-Dinitrophenol	39.5	10	μg/L μg/L	50.0		79.1	30-130			
4-Dinitrophenol 4-Dinitrotoluene	36.2	10	μg/L μg/L	50.0		72.3	40-140			
6-Dinitrotoluene	38.5	10	μg/L μg/L	50.0		81.2	40-140			
i-n-octylphthalate	40.6	10	μg/L μg/L	50.0			40-140			
2-Diphenylhydrazine/Azobenzene	43.3	10	μg/L μg/L			86.6 90.2	40-140 40-140			
exachlorobenzene	45.1	10		50.0						
exachlorobenzene exachlorobutadiene	39.5		μg/L	50.0		79.0	40-140			
	28.0	10	μg/L	50.0		56.1	40-140			
exachlorocyclopentadiene	24.6	10	μg/L	50.0		49.3	30-140			
exachloroethane	26.3	10	μg/L	50.0		52.6	40-140			
ophorone	39.2	10	μg/L	50.0		78.3	40-140			
-Methylnaphthalene	34.4	5.0	μg/L	50.0		68.8	40-140			



QUALITY CONTROL

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes	
Batch B258532 - SW-846 3510C											_
LCS (B258532-BS1)				Prepared: 05	/21/20 Anal	yzed: 05/26/2	20				
3/4-Methylphenol	34.2	10	μg/L	50.0		68.3	30-130				
2-Nitroaniline	50.6	10	$\mu g/L$	50.0		101	40-140				
3-Nitroaniline	39.8	10	μg/L	50.0		79.7	40-140				
4-Nitroaniline	40.4	10	μg/L	50.0		80.9	40-140				
Nitrobenzene	36.2	10	μg/L	50.0		72.3	40-140				
2-Nitrophenol	38.0	10	μg/L	50.0		76.0	30-130				
4-Nitrophenol	20.4	10	μg/L	50.0		40.9	10-130				†
N-Nitrosodimethylamine	24.6	10	μg/L	50.0		49.2	40-140				
N-Nitrosodiphenylamine/Diphenylamine	44.6	10	μg/L	50.0		89.3	40-140				
N-Nitrosodi-n-propylamine	42.4	10	μg/L	50.0		84.8	40-140				
Pentachloronitrobenzene	40.3	10	μg/L	50.0		80.7	40-140				
Pentachlorophenol	35.0	10	μg/L μg/L	50.0		70.0	30-130				
Phenol		10	μg/L μg/L	50.0		36.6	20-130				†
Pyridine	18.3	5.0	μg/L μg/L	50.0		40.1	10-140				†
1,2,4,5-Tetrachlorobenzene	20.1	10									1
1,2,4-Trichlorobenzene	36.3		μg/L μg/I	50.0		72.6	40-140				
	30.8	5.0	μg/L ug/I	50.0		61.7	40-140				
2,4,5-Trichlorophenol	37.6	10	μg/L	50.0		75.3	30-130				
2,4,6-Trichlorophenol	38.6	10	μg/L	50.0		77.2	30-130				_
Surrogate: 2-Fluorophenol	101		$\mu g/L$	200		50.6	15-110				
Surrogate: Phenol-d6	75.4		$\mu g/L$	200		37.7	15-110				
Surrogate: Nitrobenzene-d5	78.2		$\mu g/L$	100		78.2	30-130				
Surrogate: 2-Fluorobiphenyl	82.5		$\mu g/L$	100		82.5	30-130				
Surrogate: 2,4,6-Tribromophenol	163		μg/L	200		81.3	15-110				
Surrogate: p-Terphenyl-d14	86.1		$\mu g/L$	100		86.1	30-130				
LCS Dup (B258532-BSD1)				Prepared: 05	/21/20 Anal	yzed: 05/26/2	20				
Acetophenone	36.3	10	μg/L	50.0		72.7	40-140	8.03	20		
Aniline	28.4	5.0	$\mu g/L$	50.0		56.7	40-140	11.1	50	V-05	
Benzidine	27.8	20	$\mu g/L$	50.0		55.6	40-140	9.16	20	V-04, V-05	
Benzoic Acid	17.1	10	$\mu g/L$	50.0		34.2	10-130	2.77	50		†
Bis(2-chloroethoxy)methane	36.7	10	μg/L	50.0		73.5	40-140	10.8	20		
Bis(2-chloroethyl)ether	34.6	10	μg/L	50.0		69.2	40-140	11.3	20		
Bis(2-chloroisopropyl)ether	37.6	10	μg/L	50.0		75.2	40-140	14.4	20		
Bis(2-Ethylhexyl)phthalate	41.2	10	μg/L μg/L	50.0		82.5	40-140	7.29	20		
4-Bromophenylphenylether	38.2	10	μg/L μg/L	50.0		76.4	40-140	4.83	20		
Butylbenzylphthalate	38.7	10	μg/L μg/L	50.0		77.3	40-140	4.92	20		
Carbazole	39.2	10	μg/L μg/L	50.0		78.4	40-140	3.75	20		
4-Chloroaniline	39.2 34.0	10	μg/L μg/L	50.0		68.1	40-140	5.73	20	V-34	
4-Chloro-3-methylphenol		10								v-34	
2-Chloronaphthalene	37.0		μg/L	50.0		74.0	30-130	4.11	20		
		10		50.0		63.3	40-140	7.15	20		
-	31.7	10	μg/L ug/I	50.0		660	20 120				
2-Chlorophenol	33.4	10	$\mu g/L$	50.0		66.9	30-130	4.84	20		
2-Chlorophenol 4-Chlorophenylphenylether	33.4 38.4	10 10	μg/L μg/L	50.0		76.8	40-140	1.76	20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran	33.4 38.4 38.9	10 10 5.0	μg/L μg/L μg/L	50.0 50.0		76.8 77.7	40-140 40-140	1.76 4.28	20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate	33.4 38.4 38.9 40.0	10 10 5.0 10	μg/L μg/L μg/L μg/L	50.0 50.0 50.0		76.8 77.7 80.0	40-140 40-140 40-140	1.76 4.28 8.65	20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene	33.4 38.4 38.9 40.0 29.4	10 10 5.0 10 5.0	μg/L μg/L μg/L μg/L μg/L	50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8	40-140 40-140 40-140 40-140	1.76 4.28 8.65 1.61	20 20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene 1,3-Dichlorobenzene	33.4 38.4 38.9 40.0	10 10 5.0 10 5.0 5.0	μg/L μg/L μg/L μg/L μg/L μg/L	50.0 50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8 56.1	40-140 40-140 40-140 40-140 40-140	1.76 4.28 8.65 1.61 3.26	20 20 20 20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene	33.4 38.4 38.9 40.0 29.4	10 10 5.0 10 5.0 5.0 5.0	µg/L µg/L µg/L µg/L µg/L µg/L	50.0 50.0 50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8 56.1 56.9	40-140 40-140 40-140 40-140 40-140 40-140	1.76 4.28 8.65 1.61 3.26 3.32	20 20 20 20 20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 3,3-Dichlorobenzidine	33.4 38.4 38.9 40.0 29.4 28.1	10 10 5.0 10 5.0 5.0	μg/L μg/L μg/L μg/L μg/L μg/L	50.0 50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8 56.1	40-140 40-140 40-140 40-140 40-140	1.76 4.28 8.65 1.61 3.26	20 20 20 20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 3,3-Dichlorobenzidine	33.4 38.4 38.9 40.0 29.4 28.1 28.4	10 10 5.0 10 5.0 5.0 5.0	µg/L µg/L µg/L µg/L µg/L µg/L	50.0 50.0 50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8 56.1 56.9	40-140 40-140 40-140 40-140 40-140 40-140	1.76 4.28 8.65 1.61 3.26 3.32	20 20 20 20 20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene	33.4 38.4 38.9 40.0 29.4 28.1 28.4 41.1	10 10 5.0 10 5.0 5.0 5.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L	50.0 50.0 50.0 50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8 56.1 56.9 82.2	40-140 40-140 40-140 40-140 40-140 40-140	1.76 4.28 8.65 1.61 3.26 3.32 7.83	20 20 20 20 20 20 20 20		
2-Chlorophenol 4-Chlorophenylphenylether Dibenzofuran Di-n-butylphthalate 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 3,3-Dichlorobenzidine 2,4-Dichlorophenol	33.4 38.4 38.9 40.0 29.4 28.1 28.4 41.1 36.3	10 10 5.0 10 5.0 5.0 5.0 10	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	50.0 50.0 50.0 50.0 50.0 50.0 50.0 50.0		76.8 77.7 80.0 58.8 56.1 56.9 82.2 72.5	40-140 40-140 40-140 40-140 40-140 40-140 30-130	1.76 4.28 8.65 1.61 3.26 3.32 7.83 5.50	20 20 20 20 20 20 20 20 20		



39 Spruce Street * East Longmeadow, MA 01028 * FAX 413/525-6405 * TEL. 413/525-2332

QUALITY CONTROL

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes	
Batch B258532 - SW-846 3510C											
LCS Dup (B258532-BSD1)				Prepared: 05	/21/20 Anal	yzed: 05/26/2	20				
4,6-Dinitro-2-methylphenol	37.7	10	μg/L	50.0		75.4	30-130	4.77	50		_ ‡
2,4-Dinitrophenol	35.7	10	$\mu g \! / \! L$	50.0		71.4	30-130	1.31	50		‡
2,4-Dinitrotoluene	38.0	10	$\mu g \! / \! L$	50.0		76.1	40-140	1.20	20		
2,6-Dinitrotoluene	39.2	10	$\mu g \! / \! L$	50.0		78.4	40-140	3.56	20		
Di-n-octylphthalate	40.2	10	$\mu g/L$	50.0		80.3	40-140	7.55	20		
,2-Diphenylhydrazine/Azobenzene	37.8	10	$\mu g \! / \! L$	50.0		75.6	40-140	17.7	20		
Hexachlorobenzene	37.6	10	$\mu g\!/\!L$	50.0		75.1	40-140	5.03	20		
Hexachlorobutadiene	28.1	10	$\mu g/L$	50.0		56.3	40-140	0.356	20		
Hexachlorocyclopentadiene	22.6	10	$\mu g \! / \! L$	50.0		45.3	30-140	8.42	50		† ‡
Hexachloroethane	27.0	10	$\mu g \! / \! L$	50.0		54.0	40-140	2.70	50		‡
sophorone	34.9	10	μg/L	50.0		69.7	40-140	11.6	20		
I-Methylnaphthalene	32.9	5.0	$\mu g \! / \! L$	50.0		65.7	40-140	4.61	20		
2-Methylphenol	34.0	10	μg/L	50.0		67.9	30-130	2.76	20		
3/4-Methylphenol	31.3	10	$\mu g/L$	50.0		62.5	30-130	8.87	20		
2-Nitroaniline	44.4	10	$\mu g/L$	50.0		88.7	40-140	13.2	20		
3-Nitroaniline	39.1	10	$\mu g/L$	50.0		78.2	40-140	1.93	20		
4-Nitroaniline	41.1	10	μg/L	50.0		82.1	40-140	1.55	20		
Nitrobenzene	32.7	10	μg/L	50.0		65.4	40-140	10.0	20		
2-Nitrophenol	35.6	10	μg/L	50.0		71.1	30-130	6.63	20		
4-Nitrophenol	20.0	10	μg/L	50.0		39.9	10-130	2.38	50		† ‡
N-Nitrosodimethylamine	23.9	10	μg/L	50.0		47.8	40-140	3.01	20		
N-Nitrosodiphenylamine/Diphenylamine	40.9	10	μg/L	50.0		81.8	40-140	8.74	20		
N-Nitrosodi-n-propylamine	37.1	10	μg/L	50.0		74.2	40-140	13.3	20		
Pentachloronitrobenzene	40.0	10	μg/L	50.0		79.9	40-140	0.971	20		
Pentachlorophenol	34.6	10	μg/L	50.0		69.2	30-130	1.15	50		‡
Phenol	16.7	10	μg/L	50.0		33.4	20-130	9.26	20		†
Pyridine	19.5	5.0	μg/L	50.0		39.0	10-140	2.73	50		† ‡
1,2,4,5-Tetrachlorobenzene	33.6	10	μg/L	50.0		67.2	40-140	7.73	20		
1,2,4-Trichlorobenzene	29.9	5.0	μg/L	50.0		59.9	40-140	2.93	20		
2,4,5-Trichlorophenol	36.4	10	μg/L	50.0		72.9	30-130	3.29	20		
2,4,6-Trichlorophenol	36.8	10	μg/L	50.0		73.7	30-130	4.61	50		‡
Surrogate: 2-Fluorophenol	101		μg/L	200		50.3	15-110				_
Surrogate: Phenol-d6	69.5		μg/L μg/L	200		34.8	15-110				
Surrogate: Nitrobenzene-d5	71.6		μg/L μg/L	100		71.6	30-130				
Surrogate: 2-Fluorobiphenyl	76.8		μg/L	100		76.8	30-130				
Surrogate: 2,4,6-Tribromophenol	170		μg/L μg/L	200		84.8	15-110				
Surrogate: p-Terphenyl-d14	82.9		μg/L	100		82.9	30-130				
Batch B258763 - SW-846 3510C											_
Blank (B258763-BLK1)				Prepared: 05	/21/20 Anal	yzed: 05/27/2	20				
Acenaphthene (SIM)	ND	0.30	μg/L								
Acenaphthylene (SIM)	ND	0.20	μg/L								
Anthracene (SIM)	ND	0.20	μg/L								
Benzo(a)anthracene (SIM)	ND	0.050	μg/L								
Benzo(a)pyrene (SIM)	ND	0.10	μg/L								
Benzo(b)fluoranthene (SIM)	ND	0.050	μg/L								
Benzo(g,h,i)perylene (SIM)	ND	0.50	μg/L								
Benzo(k)fluoranthene (SIM)	ND	0.20	μg/L								
Chrysene (SIM)	ND	0.20	μg/L								
Dibenz(a,h)anthracene (SIM)	ND	0.10	μg/L								
Dibenz(a,h)anthracene (SIM) Fluoranthene (SIM)	ND ND	0.10 0.50	μg/L μg/L								



QUALITY CONTROL

2-Methylnaphthalene (SIM) ND Naphthalene (SIM) ND Phenanthrene (SIM) ND Pyrene (SIM) ND Surrogate: Nitrobenzene-d5 86.0 Surrogate: 2-Fluorobiphenyl 76.3 Surrogate: p-Terphenyl-d14 74.9 LCS (B258763-BSI) Acenaphthene (SIM) 45.4 Acenaphthylene (SIM) 49.1 Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 57.9 Fluoranthene (SIM) 57.9 Fluoranthene (SIM) 48.4 Indeno(1,2,3-ed)pyrene (SIM) 43.4 Naphthalene (SIM) 43.4 Naphthalene (SIM) 43.4 Naphthalene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: Nitrobenzene-d5 87.1 Surrogate: Nitrobenzene-d5 87.1 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthylene (SIM) 43.9 Acenaphthylene (SIM) 43.9 Acenaphthylene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.5 Benzo(a)pyrene (SIM) 47.5 Benzo(a)pyrene (SIM) 47.5 Benzo(a)pyrene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2	1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µg µg µg µg µg 1.0 µş 4.0 µş 4.0 µş 1.0 µş	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	86.0 3 76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	30-130 30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Indeno(1,2,3-cd)pyrene (SIM) 2-Methylnaphthalene (SIM) ND Naphthalene (SIM) ND Phenanthrene (SIM) ND Pyrene (SIM) ND Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 LCS (B258763-BS1) Acenaphthene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Benzo(a)anthracene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: Sim) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: Dibenz(a,h)anthracene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: Sim) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: Sim) Acenaphthylene (SIM) Asphthalene (SIM) Acenaphthylene (SIM) Acenaphthyl	1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µg µg µg µg µg 1.0 µş 4.0 µş 4.0 µş 1.0 µş	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	d: 05/21/20 Analyzed)))))))))))))))	86.0 3 76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
C-Methylnaphthalene (SIM) ND	1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µg µg µg µg µg 1.0 µş 4.0 µş 4.0 µş 1.0 µş	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Naphthalene (SIM) ND Phenanthrene (SIM) ND Pyrene (SIM) ND Surrogate: Nitrobenzene-d5 86.0 Surrogate: 2-Fluorobiphenyl 76.3 Surrogate: p-Terphenyl-d14 74.9 LCS (B258763-BS1) 45.4 Acenaphthene (SIM) 44.2 Anthracene (SIM) 49.1 Benzo(a)anthracene (SIM) 51.6 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(b)fluoranthene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 57.9 Fluoranthene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 40.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 43.4 Naphthalene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 42.5 Acenaphthene (SIM	1.0	μg/L μg/L μg/L μg/L μg/L μg/L 100 μg/L 100 Prepare μg/L 50.	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Phenanthrene (SIM) Pyrene (SIM) ND Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 **CS (B258763-BS1) **Cenaphthene (SIM) Acenaphthylene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) S1.6 Benzo(b)fluoranthene (SIM) S2.5 Chrysene (SIM) Dibenz(a,h)anthracene (SIM) Fluoranthene (SIM) S1.2 Fluorene (SIM) Indeno(1,2,3-cd)pyrene (SIM) As Alladeno(1,2,3-cd)pyrene (SIM) Anphthalene (SIM) As Alladeno(1,2,3-cd) Phenanthrene (SIM) As Alladeno(1,2,3-cd) Pyrene (SIM) As Alladeno(1,2,3-cd) Pyrene (SIM) As Alladeno(1,2,3-cd) As Alladeno(1,2,3	1.0 µş µg µg µg µg µ4.0 µş 4.0 µş 1.0 µş 4.0 µş 1.0 µş	µg/L µg/L 100 µg/L 100 µg/L 100 µg/L 100 µg/L 100 Prepare µg/L 50.0 µg/L 50.0 µg/L 50.0 µg/L 50.0 µg/L 50.0 µg/L 50.0 µg/L 50.0 µg/L 50.0 µg/L 50.0	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Pyrene (SIM) Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 74.9 LCS (B258763-BS1) Acenaphthene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Anthracene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(a,h,i)perylene (SIM) Benzo(b,fluoranthene (SIM) Surrogate: SiM) Benzo(a,h)anthracene (SIM) S1.6 Benzo(k)fluoranthene (SIM) S2.5 Chrysene (SIM) Dibenz(a,h)anthracene (SIM) Fluoranthene (SIM) Fluorene (SIM) Fluoranthene (SIM) Ada, A Naphthalene (SIM) Aga, A Naphthalene (SIM) Prene (SIM) Pyrene (SIM) Asa, A Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: 2-Fluorobiphenyl Surrogate: 2-Fluorobiphenyl Acenaphthylene (SIM) Acen	1.0 µş µg µg µg µg µg µg 4.0 µş 2.0 µ	рву/L р	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl 76.3 Surrogate: 2-Fluorobiphenyl 76.3 Surrogate: p-Terphenyl-d14 74.9 LCS (B258763-BS1) Acenaphthene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: CSIM) Benzo(g,h,i)perylene (SIM) Benzo(k)fluoranthene (SIM) Surrogate: CSIM) Surrogate: CSIM) Surrogate: CSIM) Surrogate: CSIM) Surrogate: CSIM) Surrogate: Nitrobenzene-d5 Surrogate: Nitrobenzene-d5 Surrogate: P-Terphenyl-d14 LCS Dup (B258763-BSD1) Acenaphthylene (SIM) Andhracene (SIM) Andhracene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Art. Benzo(a)anthracene (SIM) Art. Benzo(a)pyrene (SIM) Art. Benzo(a)pyrene (SIM) Art. Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM)	6.0 µқ 4.0 µқ 4.0 µқ 1.0 µқ 2.0 µқ 2.0 µқ 2.0 µқ 2.0 µқ 2.0 µқ 2.0 µқ	μg/L 100 μg/L 100 μg/L 100 Prepare μg/L 50.0 μg/L 50.0 μg/L 50.0 μg/L 50.0 μg/L	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Surrogate: 2-Fluorobiphenyl 76.3 Surrogate: p-Terphenyl-d14 74.9 LCS (B258763-BS1) 45.4 Acenaphthene (SIM) 44.2 Anthracene (SIM) 49.1 Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 43.4 Naphthalene (SIM) 46.6 Pyrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 42.5 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 50.2 Benzo(a)pyrene (SIM) 50.2<	6.0 µş 4.0 µş 4.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 2.0 µş 1.0 µş 2.0 µş 2.0 µş 2.0 µş 2.0 µş 2.0 µş	μg/L 100 Prepare μg/L 50.0	d: 05/21/20 Analyzed) (d: 05/21/20 Analyzed)	76.3 3 74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	30-130 30-130 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Surrogate: p-Terphenyl-d14 74.9 LCS (B258763-BS1) 45.4 Acenaphthene (SIM) 44.2 Anthracene (SIM) 49.1 Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 50.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	6.0 µş 4.0 µş 4.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 2.0 µş 1.0 µş 2.0 µş 2.0 µş 2.0 µş 2.0 µş 2.0 µş	μg/L 100 Prepare μg/L 50.0	d: 05/21/20 Analyzed))))))))))))))))))	74.9 3 d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
LCS (B258763-BS1) Acenaphthene (SIM) 45.4 Acenaphthylene (SIM) 44.2 Anthracene (SIM) 49.1 Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)apyrene (SIM) 50.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	6.0 µş 4.0 µş 4.0 µş 1.0 µş 1.0 µş 1.0 µş 1.0 µş 10 µş 10 µş 2.0 µş 10 µş 2.0 µş 20 µş 20 µş	μg/L 50.0	d: 05/21/20 Analyzed	d: 05/27/20 90.8 4 88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Acenaphthene (SIM) 45.4 Acenaphthylene (SIM) 44.2 Anthracene (SIM) 49.1 Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 43.4 Naphthalene (SIM) 43.4 Naphthalene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2	4.0 µş 4.0 µş 4.0 µş 1.0 µş 1.0 µş 1.0 µs 1.0 µs 10 µs 4.0 µs 2.0 µs 10 µs 20 µs 20 µs 20 µs	μg/L 50.		90.8 4 98.3 4 98.3 4 97.1 4 103 4 107 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Acenaphthene (SIM) 45.4 Acenaphthylene (SIM) 44.2 Anthracene (SIM) 49.1 Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 43.4 Naphthalene (SIM) 43.4 Naphthalene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: Nitrobenzene-d5 87.1 Surrogate: P-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 50.2	4.0 µş 4.0 µş 4.0 µş 1.0 µş 1.0 µş 1.0 µs 1.0 µs 10 µs 4.0 µs 2.0 µs 10 µs 20 µs 20 µs 20 µs	μg/L 50.		90.8 4 98.3 4 98.3 4 97.1 4 103 4 107 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Acenaphthylene (SIM) Anthracene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) Benzo(g,h,i)perylene (SIM) Benzo(g,hi)perylene (SIM) Benzo(k)fluoranthene (SIM) Chrysene (SIM) Dibenz(a,h)anthracene (SIM) Fluoranthene (SIM) Fluoranthene (SIM) Fluoranthene (SIM) Fluoranthene (SIM) Indeno(1,2,3-cd)pyrene (SIM) 2-Methylnaphthalene (SIM) Naphthalene (SIM) Phenanthrene (SIM) Pyrene (SIM) Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 CCS Dup (B258763-BSD1) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: Simolate (Simolate) Surrogate: Simolate (Simolate) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM)	4.0 µş 4.0 µş 4.0 µş 1.0 µş 1.0 µş 1.0 µs 1.0 µs 10 µs 4.0 µs 2.0 µs 10 µs 20 µs 20 µs 20 µs	μg/L 50.1 μg/L 50.2		88.4 4 98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Anthracene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) Benzo(g,h,i)perylene (SIM) Benzo(k)fluoranthene (SIM) Benzo(k)fluoranthene (SIM) Chrysene (SIM) Dibenz(a,h)anthracene (SIM) Fluoranthene (SIM) Fluoranthene (SIM) Fluoranthene (SIM) Fluoranthene (SIM) Fluorene (SIM) Indeno(1,2,3-cd)pyrene (SIM) 2-Methylnaphthalene (SIM) Aphthalene (SIM) Phenanthrene (SIM) Phenanthrene (SIM) Pyrene (SIM) Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 LCS Dup (B258763-BSD1) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM)	4.0	μg/L 50.		98.3 4 97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Benzo(a)anthracene (SIM) 48.5 Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.5 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	1.0 µş 2.0 µş 1.0 µş 1.0 µş 4.0 µş 4.0 µş 2.0 µş 10 µş 20 µş 20 µş 20 µş	μg/L 50.		97.1 4 103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Benzo(a)pyrene (SIM) 51.6 Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	2.0	μg/L 50.		103 4 107 4 109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Benzo(b)fluoranthene (SIM) 53.7 Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthylene (SIM) 47.5 Acenaphthylene (SIM) 47.5 Benzo(a)anthracene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	1.0	μg/L 50.		109 4 105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8	40-140 40-140 40-140 40-140 40-140 40-140 40-140			
Benzo(g,h,i)perylene (SIM) 54.6 Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthylene (SIM) 47.5 Acenaphthylene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	4.0 µş 4.0 µş 4.0 µş 2.0 µş 10 µş 20 µş 20 µş 20 µş	$\begin{array}{ccc} \mu g/L & 50.0 \\ \end{array}$		105 4 94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140 40-140 40-140			
Benzo(k)fluoranthene (SIM) 52.5 Chrysene (SIM) 47.3 Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.5 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	 4.0 με 2.0 με 10 με 20 με 2.0 με 2.0 με 2.0 με 	μg/L 50.		94.6 4 116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140 40-140			
Dibenz(a,h)anthracene (SIM) 57.9 Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 47.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	2.0 µg 10 µg 20 µg 2.0 µg 2.0 µg 2.0 µg	μg/L 50.0		116 4 102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140 40-140			
Fluoranthene (SIM) 51.2 Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	10 μ _ξ 20 μ _ξ 2.0 μ _ξ 20 μ _ξ	μg/L 50.4 μg/L 50.4 μg/L 50.4 μg/L 50.4 μg/L 50.4 μg/L 50.4)))	102 4 96.9 4 121 4 86.8 4	40-140 40-140 40-140			
Fluorene (SIM) 48.4 Indeno(1,2,3-cd)pyrene (SIM) 60.7 2-Methylnaphthalene (SIM) 43.4 Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	 20 με 2.0 με 20 με 	μg/L 50. μg/L 50. μg/L 50. μg/L 50.)	96.9 4 121 4 86.8 4	40-140 40-140			
Indeno(1,2,3-cd)pyrene (SIM) 2-Methylnaphthalene (SIM) Naphthalene (SIM) Phenanthrene (SIM) Pyrene (SIM) Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 LCS Dup (B258763-BSD1) Acenaphthene (SIM) Acenaphthylene (SIM)	2.0 μ _ξ 20 μ _ξ	μg/L 50. μg/L 50. μg/L 50.)	121 4 86.8 4	40-140			
2-Methylnaphthalene (SIM) Naphthalene (SIM) Phenanthrene (SIM) Pyrene (SIM) Surrogate: Nitrobenzene-d5 Surrogate: 2-Fluorobiphenyl Surrogate: p-Terphenyl-d14 CCS Dup (B258763-BSD1) Acenaphthene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Acenaphthylene (SIM) Benzo(a)anthracene (SIM) Benzo(a)pyrene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Benzo(b)fluoranthene (SIM) Surrogate: 2-Fluorobiphenyl 90.5 87.1 87.1 87.5 87.1 87.5 88.1 89.5 80.2 80.2 80.2 80.2 80.2	20 με	μg/L 50. μg/L 50.)	86.8 4				
Naphthalene (SIM) 39.6 Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7		μg/L 50.			40-140			
Phenanthrene (SIM) 46.6 Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	• 0							
Pyrene (SIM) 45.4 Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	20 με		J	79.2 4	40-140			
Surrogate: Nitrobenzene-d5 87.1 Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	1.0 με	μg/L 50.)	93.2 4	40-140			
Surrogate: 2-Fluorobiphenyl 90.5 Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) 43.9 Acenaphthene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	20 με	μg/L 50.)	90.9 4	40-140			
Surrogate: p-Terphenyl-d14 73.7 LCS Dup (B258763-BSD1) Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	μд	μg/L 100)	87.1 3	30-130			
Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	μд	μg/L 100)	90.5	30-130			
Acenaphthene (SIM) 43.9 Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	μg	μg/L 100)	73.7 3	30-130			
Acenaphthylene (SIM) 42.5 Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7		Prepare	d: 05/21/20 Analyzed	d: 05/27/20				
Anthracene (SIM) 47.5 Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7	6.0 με	μg/L 50.)	87.7 4	40-140	3.49	20	
Benzo(a)anthracene (SIM) 47.2 Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7		μg/L 50.)	85.1 4	40-140	3.83	20	
Benzo(a)pyrene (SIM) 50.2 Benzo(b)fluoranthene (SIM) 52.7		μg/L 50.)	95.0 4	40-140	3.39	20	
Benzo(b)fluoranthene (SIM) 52.7		μg/L 50.)	94.5 4	40-140	2.71	20	
· · · · · · · · · · · · · · · · · · ·		μg/L 50.				2.75	20	
		μg/L 50.				1.92	20	
Benzo(g,h,i)perylene (SIM) 53.0		μg/L 50.				2.90	20	
Benzo(k)fluoranthene (SIM) 51.4		μg/L 50.				2.12	20	
Chrysene (SIM) 46.0		μg/L 50.				2.79	20	
Dibenz(a,h)anthracene (SIM) 56.8		μg/L 50.				1.99	20	
Fluoranthene (SIM) 49.4		μg/L 50.				3.58	20	
Fluorene (SIM) 47.3		μg/L 50.				2.34	20	
Indeno(1,2,3-cd)pyrene (SIM) 59.0		μg/L 50.0				2.94	20	
2-Methylnaphthalene (SIM) 42.2		μg/L 50.0				2.66	20	
Naphthalene (SIM) 39.2		μg/L 50.0				1.07	20	
Phenanthrene (SIM) 44.9		μg/L 50.0				3.67	20	
Pyrene (SIM) 44.5		μg/L 50.				2.14	20	
Surrogate: Nitrobenzene-d5 83.6	20 με	μ g/L 100)		30-130			
Surrogate: 2-Fluorobiphenyl 88.7 Surrogate: p-Terphenyl-d14 72.7	20 μg	μg/L 100		88.7 3	30-130			



FLAG/QUALIFIER SUMMARY

*	QC result is outside of established limits.
†	Wide recovery limits established for difficult compound.
‡	Wide RPD limits established for difficult compound.
#	Data exceeded client recommended or regulatory level
ND	Not Detected
RL	Reporting Limit is at the level of quantitation (LOQ)
DL	Detection Limit is the lower limit of detection determined by the MDL study
MCL	Maximum Contaminant Level
	Percent recoveries and relative percent differences (RPDs) are determined by the software using values in the calculation which have not been rounded.
	No results have been blank subtracted unless specified in the case narrative section.
J	Detected but below the Reporting Limit (lowest calibration standard); therefore, result is an estimated concentration (CLP J-Flag).
L-04	Laboratory fortified blank/laboratory control sample recovery and duplicate recovery are outside of control limits. Reported value for this compound is likely to be biased on the low side.
RL-11	Elevated reporting limit due to high concentration of target compounds.
V-04	Initial calibration did not meet method specifications. Compound was calibrated using a response factor where %RSD is outside of method specified criteria. Reported result is estimated.
V-05	Continuing calibration verification (CCV) did not meet method specifications and was biased on the low side for this compound.
V-06	Continuing calibration verification (CCV) did not meet method specifications and was biased on the high side for this compound.
V-34	Initial calibration verification (ICV) did not meet method specifications and was biased on the low side for this compound. Reported result is estimated.



CERTIFICATIONS

Analyte	Certifications	
SW-846 8260D in Water		
Acetone	NC	
Acrylonitrile	NC	
tert-Amyl Methyl Ether (TAME)	NC	
Benzene	NC	
Bromobenzene	NC	
Bromochloromethane	NC	
Bromodichloromethane	NC	
Bromoform	NC	
Bromomethane	NC	
2-Butanone (MEK)	NC	
tert-Butyl Alcohol (TBA)	NC	
n-Butylbenzene	NC	
sec-Butylbenzene	NC	
tert-Butylbenzene	NC	
tert-Butyl Ethyl Ether (TBEE)	NC	
Carbon Disulfide	NC	
Carbon Tetrachloride	NC	
Chlorobenzene	NC	
Chlorodibromomethane	NC	
Chloroethane	NC	
Chloroform	NC	
Chloromethane	NC	
2-Chlorotoluene	NC	
4-Chlorotoluene	NC	
1,2-Dibromo-3-chloropropane (DBCP)	NC	
1,2-Dibromoethane (EDB)	NC	
Dibromomethane	NC	
1,2-Dichlorobenzene	NC	
1,3-Dichlorobenzene	NC	
1,4-Dichlorobenzene	NC	
trans-1,4-Dichloro-2-butene	NC	
Dichlorodifluoromethane (Freon 12)	NC	
1,1-Dichloroethane	NC	
1,2-Dichloroethane	NC	
1,1-Dichloroethylene cis-1,2-Dichloroethylene	NC NC	
trans-1,2-Dichloroethylene	NC NC	
1,2-Dichloropropane	NC NC	
1,3-Dichloropropane	NC NC	
2,2-Dichloropropane	NC	
1,1-Dichloropropene	NC	
cis-1,3-Dichloropropene	NC	
trans-1,3-Dichloropropene	NC	
Diethyl Ether	NC	
Diisopropyl Ether (DIPE)	NC	
1,4-Dioxane	NC	
Ethylbenzene	NC	
		1



CERTIFICATIONS

Certifications
NC
NY,NH,ME,NC,VA
NY,NH,ME,NC,VA
CT,NY,NH,ME,NC,VA
NY,NH,ME,NC,VA
CT,NY,NH,ME,NC,VA
NC
CT,NY,NH,ME,NC,VA
CT,NY,NH,ME,NC,VA
CT,NY,NH,NC,VA
CT,NY,NH,ME,NC,VA
CT,NY,NH,ME,NC,VA



CERTIFICATIONS

Analyte	Certifications
SW-846 8270E in Soil	
Dibenzofuran	CT,NY,NH,ME,NC,VA
Di-n-butylphthalate	CT,NY,NH,ME,NC,VA
1,2-Dichlorobenzene	NY,NH,ME,NC,VA
1,3-Dichlorobenzene	NY,NH,ME,NC,VA
1,4-Dichlorobenzene	NY,NH,ME,NC,VA
3,3-Dichlorobenzidine	CT,NY,NH,ME,NC,VA
2,4-Dichlorophenol	CT,NY,NH,ME,NC,VA
Diethylphthalate	CT,NY,NH,ME,NC,VA
2,4-Dimethylphenol	CT,NY,NH,ME,NC,VA
Dimethylphthalate	CT,NY,NH,ME,NC,VA
4,6-Dinitro-2-methylphenol	CT,NY,NH,ME,NC,VA
2,4-Dinitrophenol	CT,NY,NH,ME,NC,VA
2,4-Dinitrotoluene	CT,NY,NH,ME,NC,VA
2,6-Dinitrotoluene	CT,NY,NH,ME,NC,VA
Di-n-octylphthalate	CT,NY,NH,ME,NC,VA
1,2-Diphenylhydrazine/Azobenzene	NY,NH,ME,NC,VA
Hexachlorobenzene	CT,NY,NH,ME,NC,VA
Hexachlorobutadiene	CT,NY,NH,ME,NC,VA
Hexachlorocyclopentadiene	CT,NY,NH,ME,NC,VA
Hexachloroethane	CT,NY,NH,ME,NC,VA
Isophorone	CT,NY,NH,ME,NC,VA
1-Methylnaphthalene	NC
2-Methylnaphthalene	CT,NY,NH,ME,NC,VA
2-Methylphenol	CT,NY,NH,ME,NC,VA
3/4-Methylphenol	CT,NY,NH,ME,NC,VA
Naphthalene	CT,NY,NH,ME,NC,VA
2-Nitroaniline	CT,NY,NH,ME,NC,VA
3-Nitroaniline	CT,NY,NH,ME,NC,VA
4-Nitroaniline	CT,NY,NH,ME,NC,VA
Nitrobenzene	CT,NY,NH,ME,NC,VA
2-Nitrophenol	CT,NY,NH,ME,NC,VA
4-Nitrophenol	CT,NY,NH,ME,NC,VA
N-Nitrosodimethylamine	CT,NY,NH,ME,NC,VA
N-Nitrosodi-n-propylamine	CT,NY,NH,ME,NC,VA
Pentachloronitrobenzene	NY,NC
Pentachlorophenol	CT,NY,NH,ME,NC,VA
Phenol	CT,NY,NH,ME,NC,VA
Pyridine	CT,NY,NH,ME,NC,VA
1,2,4,5-Tetrachlorobenzene	NY,NC
1,2,4-Trichlorobenzene	CT,NY,NH,ME,NC,VA
2,4,5-Trichlorophenol	CT,NY,NH,ME,NC,VA
2,4,6-Trichlorophenol	CT,NY,NH,ME,NC,VA
2-Fluorophenol	NC
SW-846 8270E in Water	
Acetophenone	NY,NC
Aniline	CT,NY,NC,ME,VA



CERTIFICATIONS

Analyte	Certifications
SW-846 8270E in Water	
Benzidine	CT,NY,NC,ME,NH,VA
Benzoic Acid	NY,NC,ME,NH,VA
Bis(2-chloroethoxy)methane	CT,NY,NC,ME,NH,VA
Bis(2-chloroethyl)ether	CT,NY,NC,ME,NH,VA
Bis(2-chloroisopropyl)ether	CT,NY,NC,ME,NH,VA
Bis(2-Ethylhexyl)phthalate	CT,NY,NC,ME,NH,VA
4-Bromophenylphenylether	CT,NY,NC,ME,NH,VA
Butylbenzylphthalate	CT,NY,NC,ME,NH,VA
Carbazole	NC
4-Chloroaniline	CT,NY,NC,ME,NH,VA
4-Chloro-3-methylphenol	CT,NY,NC,ME,NH,VA
2-Chloronaphthalene	CT,NY,NC,ME,NH,VA
2-Chlorophenol	CT,NY,NC,ME,NH,VA
4-Chlorophenylphenylether	CT,NY,NC,ME,NH,VA
Dibenzofuran	CT,NY,NC,ME,NH,VA
Di-n-butylphthalate	CT,NY,NC,ME,NH,VA
1,2-Dichlorobenzene	CT,NY,NC,ME,NH,VA
1,3-Dichlorobenzene	CT,NY,NC,ME,NH,VA
1,4-Dichlorobenzene	CT,NY,NC,ME,NH,VA
3,3-Dichlorobenzidine	CT,NY,NC,ME,NH,VA
2,4-Dichlorophenol	CT,NY,NC,ME,NH,VA
Diethylphthalate	CT,NY,NC,ME,NH,VA
2,4-Dimethylphenol	CT,NY,NC,ME,NH,VA
Dimethylphthalate	CT,NY,NC,ME,NH,VA
4,6-Dinitro-2-methylphenol	CT,NY,NC,ME,NH,VA
2,4-Dinitrophenol	CT,NY,NC,ME,NH,VA
2,4-Dinitrotoluene	CT,NY,NC,ME,NH,VA
2,6-Dinitrotoluene	CT,NY,NC,ME,NH,VA
Di-n-octylphthalate	CT,NY,NC,ME,NH,VA
1,2-Diphenylhydrazine/Azobenzene	NY,NC
Hexachlorobenzene	CT,NY,NC,ME,NH,VA
Hexachlorobutadiene	CT,NY,NC,ME,NH,VA
Hexachlorocyclopentadiene	CT,NY,NC,ME,NH,VA
Hexachloroethane	CT,NY,NC,ME,NH,VA
Isophorone	CT,NY,NC,ME,NH,VA
1-Methylnaphthalene	NC
2-Methylnaphthalene	CT,NY,NC,ME,NH,VA
2-Methylphenol	CT,NY,NC,NH,VA
3/4-Methylphenol	CT,NY,NC,NH,VA
Naphthalene	CT,NY,NC,ME,NH,VA
2-Nitroaniline	CT,NY,NC,ME,NH,VA
3-Nitroaniline	CT,NY,NC,ME,NH,VA
4-Nitroaniline	CT,NY,NC,ME,NH,VA
Nitrobenzene	CT,NY,NC,ME,NH,VA
2-Nitrophenol	CT,NY,NC,ME,NH,VA
4-Nitrophenol	CT,NY,NC,ME,NH,VA
N-Nitrosodimethylamine	CT,NY,NC,ME,NH,VA



CERTIFICATIONS

Certified Analyses included in this Report

2-Fluorophenol

Analyte Certifications

SW-846 8270E in Water

SW-846 8270E in Water	
N-Nitrosodi-n-propylamine	CT,NY,NC,ME,NH,VA
Pentachloronitrobenzene	NC
Pentachlorophenol	CT,NY,NC,ME,NH,VA
Phenol	CT,NY,NC,ME,NH,VA
Pyridine	CT,NY,NC,ME,NH,VA
1,2,4,5-Tetrachlorobenzene	NY,NC
1,2,4-Trichlorobenzene	CT,NY,NC,ME,NH,VA
2,4,5-Trichlorophenol	CT,NY,NC,ME,NH,VA
2,4,6-Trichlorophenol	CT,NY,NC,ME,NH,VA

 $The \ CON-TEST \ Environmental \ Laboratory \ operates \ under \ the \ following \ certifications \ and \ accreditations:$

NC

Code	Description	Number	Expires
AIHA	AIHA-LAP, LLC - ISO17025:2017	100033	03/1/2022
MA	Massachusetts DEP	M-MA100	06/30/2020
CT	Connecticut Department of Publile Health	PH-0567	09/30/2021
NY	New York State Department of Health	10899 NELAP	04/1/2021
NH-S	New Hampshire Environmental Lab	2516 NELAP	02/5/2021
RI	Rhode Island Department of Health	LAO00112	12/30/2020
NC	North Carolina Div. of Water Quality	652	12/31/2020
NJ	New Jersey DEP	MA007 NELAP	06/30/2020
FL	Florida Department of Health	E871027 NELAP	06/30/2020
VT	Vermont Department of Health Lead Laboratory	LL015036	07/30/2021
ME	State of Maine	2011028	06/9/2021
VA	Commonwealth of Virginia	460217	12/14/2020
NH-P	New Hampshire Environmental Lab	2557 NELAP	09/6/2020
VT-DW	Vermont Department of Health Drinking Water	VT-255716	06/12/2020
NC-DW	North Carolina Department of Health	25703	07/31/2020
PA	Commonwealth of Pennsylvania DEP	68-05812	06/30/2020

F68034

Phone: 413-525-2332

Par Cel EED http://www.contestlabs.com

CHAIN OF CUSTODY RECORD (North Carolina)

Doc # 379 Rev 1_03242017

Table of Contents 2 Preservation Codes: X = Sodium Hydroxide WW = Waste Water DW = Drinking Water B = Sodium Bisulfate GW = Ground Water ³Container Codes:
A = Amber Glass
G = Glass
P = Plastic SL = Sludge SOL = Solid O = Other (please S = Summa Canister T = Tediar Bag O = Other (please Thiosulfate O = Other (please Non Soxhlet S = Sulfuric Acid Matrix Codes: PCB ONLY H = HCL M = Methanol N = Nitric Acid Soxhiet Preservation Code Field Filtered Field Filtered Lab to Filter Lab to Filter Container Code ا و ST = Sterile # of Containers = Sodium S = Soil V = Vial define) A = Air define) l = lced define) Page ___ UST/Trust Fund Please use the following codes to indicate possible sample concentration NELAC and AMA-LAP, LLC Accredited Chromatogram
AIHA-LAP, LLC REC 39 Spruce Street East Longmeadow, MA 01028 H - High; M - Medium; L - Low; C - Clean; U - Unknown Program Information ANALYSIS REQUESTED within the Conc Code column above: IHSB Orphaned Landfill Other SWS Landfill State Lead DSCA Other: Õ. 28 \overline{O} 9 9 8 8 3 Matrix Municipality Brownfield 10-Day Grab 3-Day 4-Day School EXCEL CLP Like Data Pkg Required: Ending Composite PDF Government 1135 Email To: Due Date ormat: Fax To #: Federal 7-Day Other l-Day 2-Day GWPC MSCC SWSL HSB City 7 5/18/201 Project Entity Beginning Date Time SOCIOTE Email: info@contestlabs.com Date/Time: 5/18/15 9 parel Date/Time: (6.30 Date/Time: (155) 7 2 Client Sample ID / Description 5010 Fax: 413-525-6405 Date/Time: Date/Time: Date/Time: <u>ئ</u>ر ك Marre o Ada SB. 20080204-8 M. Ascasor ACT BUTCH Tarboros TACKSTO Sampled By: Korolly (Agnature) 2.6 Con-Test Quote Name/Number: MI COD-LEST by: (signature) Rejinature) (signature) (signature) ived by: (signature) Work Order# Con-Test Project Location: Project Manager: Project Number: Invoice Recipient; nquished Company Comments Address: Phone: Page 28 of 30

IMPORTANT!

We are continuing to respond to the impact of COVID-19 around the world. See our latest updates. For COVID-19-related recipient closures, you can redirect packages, Ask FedEx, or contact the shipper.



149454201820







DELIVERED

Signed for by: R.PIETRIAS

GET STATUS UPDATES OBTAIN PROOF OF DELIVERY

FROM		то	
Raleigh, NC US	The second of the second description of the second	EAST LONGMEADOW, MA US 1 1800	-

Multiple-piece Shipment

3 Piece shipment

TRACKING NUMBER	SHIPPER CITY, STATE	SHIP DATE	STATUS	DELIVERY DATE	DESTINATION/RECIPIENT CITY, STATE
149454201809 (master)	RALEIGH, NC	5/19/2020	Company Comments of the Commen	5/20/2020	East Longmeadow, MA
149454201810	RALEIGH, NC	5/19/2020	Comment Commen	5/20/2020	East Longmeadow, MA
149454201820	RALEIGH, NC	5/19/2020	Energy Contract Contr	5/20/2020	East Longmeadow, MA

Shipment Facts

TRACKING NUMBER	SERVICE	MASTER TRACKING NUMBER
149454201820	FedEx Priority Overnight	149454201809
WEIGHT	DIMENSIONS	DELIVERED TO
10 lbs / 4.54 kgs	25x14x14 in.	Receptionist/Front Desk
TOTAL PIECES	TOTAL SHIPMENT WEIGHT	TERMS
3	10 lbs / 4.54 kgs	Third Party
PACKAGING	SPECIAL HANDLING SECTION	STANDARD TRANSIT
Your Packaging	Deliver Weekday, Non Standard Packaging	⑦ E/20/2020 b - 10-20

5/20/2020 by 10:30 am 820&locale=en US&cntry code=en Page 29 of 30 I Have Not Confirmed Sample Container Numbers With Lab Staff Before Relinquishing Over Samples____



Doc# 277 Rev 5 2017

IZECE!	ved By	DAN	***************************************	Date	۶	26	· · · · · · · · · · · · · · · · · · ·	Time	10(4	
How were t	he samples	In Cooler		No Cooler			On Ice			····
	ived?			NO COOIEI	7700				_ No Ice	
		Direct from Sam	•				Ambient		_ Melted Ice	*
Were sam	ples within		By Gun#	<u> </u>		A	ctual Tem	1p- 28		
	ure? 2-6°C	T	By Blank #			A	ctual Tem	nn -		_
Was	s Custody Se	eal Intact?	- M	We	re Sam		ampered		<u> </u>	-
	s COC Relin						With Sa			_
		eaking/loose caps	s on any sam	oles?	6		· · · · · · · · · · · · · · · · · · ·	mpioo .		_
Is COC in ir	nk/ Legible?	T	,	-	noles re	ceive	d within h	olding time?		
Did COC		Client		Analysis	T			er Name		_
pertinent In	formation?	Project		ID's			•	Dates/Times		_
Are Sample	labels filled	out and legible?		•••					· ————	
	ab to Filters?		\overline{F}		Who	was n	otified?			
Are there Ru	ushes?		MRKE				otified?			<u></u>
Are there Sh	ort Holds?		7				otified?			irw
ls there eno	ugh Volume'	?	·T				-			-
ls there Hea	dspace whe	re applicable?	- Commence Control	i	MS/MS	D?	F	ing or	^	
							mples rec		+	
Proper Medi	a contantor					no sai	umes rec	wren /		
	anks receive		£				inples req	Juirea?	* <u> </u>	-
Were trip bla		d?	F		On CO		f f	•	<u> </u>	-
Were trip bla Do all sampl	anks received les have the	d? proper pH?	NA	(£	uired? Base		
Were trip bla	anks received les have the	d? proper pH? Containers:	1	Acid	On CO		#	Base	Anala	
Were trip bla Do all sampl Vials	anks received les have the	d? proper pH? Containers: 1 Liter Amb.	A MANAGEMENT OF THE PROPERTY O	Acid1 Liter F	On CO		£	Base 16 oz		*
Were trip bla Do all sampl Vials Unp-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb.	1	Acid 1 Liter F 500 mL	On CO		£	Base 16 oz 8oz Am	ıb/Clear	***
Were trip bla Do all sampl Vials Unp- HCL-	anks received les have the	d? proper pH? Containers: 1 Liter Amb.	1	Acid 1 Liter F 500 mL I 250 mL I	Plastic Plastic Plastic		F	Base 16 oz 8oz Am 4oz Am	b/Clear b/Clear	*
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb.	1	1 Liter F 500 mL I 250 mL I Col./Bac	Plastic Plastic Plastic Plastic cteria		F	Base 16 oz 8oz Am 4oz Am 2oz Am	b/Clear b/Clear b/Clear	#
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint	1	Acid 1 Liter F 500 mL I 250 mL I Col./Bac	Plastic Plastic Plastic Cteria		#	Base 16 oz 8oz Am 4oz Am 2oz Am Enc	b/Clear b/Clear b/Clear	*
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass	1	1 Liter F 500 mL I 250 mL I Col./Bai Other P Plastic	Plastic Plastic Plastic Cteria Plastic Bag		#	Base 16 oz 8oz Am 4oz Am 2oz Am	b/Clear b/Clear b/Clear	***
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit	1	Acid 1 Liter F 500 mL I 250 mL I Col./Bac Other P Plastic Ziplo	Plastic Plastic Plastic cteria Plastic Bag		#	Base 16 oz 8oz Am 4oz Am 2oz Am Enc	b/Clear b/Clear b/Clear	***************************************
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate	2	1 Liter F 500 mL I 250 mL I Col./Bai Other P Plastic	Plastic Plastic Plastic cteria Plastic Bag		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc	b/Clear b/Clear b/Clear	
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers:	1	Acid 1 Liter F 500 mL I 250 mL I Col./Ba Other P Plastic Ziplo	Plastic Plastic Plastic cteria Plastic Bag ock		#	Base 16 oz 8oz Am 4oz Am 2oz Am Enc	b/Clear b/Clear b/Clear core	***
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers: 1 Liter Amb.	2	1 Liter F 500 mL I 250 mL I Col./Bac Other P Plastic Ziplo Unused M	Plastic Plastic Plastic cteria Plastic Bag ock		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc Frozen:	b/Clear b/Clear b/Clear core	
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric- Vials Unp-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers:	2	1 Liter F 500 mL I 250 mL I Col./Bac Other P Plastic Ziplo Unused M 1 Liter P 500 mL F	Plastic Plastic Cteria Plastic Bag ck ledia Plastic		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc Frozen:	b/Clear b/Clear b/Clear core Amb. b/Clear	***
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric- Vials Jnp- HCL-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb.	2	Acid 1 Liter F 500 mL I 250 mL I Col./Bac Other P Plastic Ziplo Unused M 1 Liter P 500 mL F 250 mL F	Plastic Plastic Cteria Plastic Bag ock edia Plastic Plastic		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc Frozen: 16 oz 8oz Am 4oz Am	Amb. b/Clear	
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric- Vials Jnp- HCL- Meoh-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Col./Bacteria	2	Acid 1 Liter F 500 mL I 250 mL I Col./Ba Other P Plastic Ziplo Unused M 1 Liter P 500 mL F 250 mL F	Plastic Plastic Cteria Plastic Bag Ck Plastic Plastic Plastic Plastic Plastic Plastic Plastic		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc Frozen: 16 oz 8oz Am 4oz Am 2oz Am 2oz Am	Amb. b/Clear b/Clear core	
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric- Vials Unp- HCL- Meoh- Bisulfate- Bisulfate- Bisulfate-	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb.	2	Acid 1 Liter F 500 mL I 250 mL I Col./Bac Other P Plastic Ziplo Unused M 1 Liter P 500 mL F 250 mL F Flashp Other C	Plastic Plastic Cteria Plastic Bag ock Plastic		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc Frozen: 16 oz 8oz Am 4oz Am 2oz Am Enc	Amb. b/Clear b/Clear core	
Were trip bla Do all sampl Vials Unp- HCL- Meoh- Bisulfate- DI- Thiosulfate- Sulfuric- Vials Unp- HCL- Meoh- Bisulfate- DI- DI- DI- DI- DI- DI- DI- DI- DI- DI	anks received les have the	d? proper pH? Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Flashpoint Other Glass SOC Kit Perchlorate Containers: 1 Liter Amb. 500 mL Amb. 250 mL Amb. Col./Bacteria Other Plastic	2	Acid 1 Liter F 500 mL I 250 mL I Col./Ba Other P Plastic Ziplo Unused M 1 Liter P 500 mL F 250 mL F	Plastic Plastic Control Plastic		*	Base 16 oz 8oz Am 4oz Am 2oz Am Enc Frozen: 16 oz 8oz Am 4oz Am 2oz Am 2oz Am	Amb. b/Clear b/Clear core	