



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 DRO<sup>1</sup>. 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-methylnaphthalene and naphthalene were present at 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

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<sup>1</sup> 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:

Mike Branson

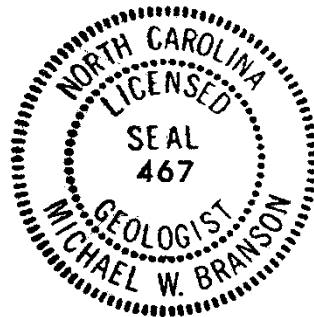
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6/25/2020

Michael W. Branson, P.G.  
Project Manager

Attachments



William D. Newcomb, P.G.  
Senior Hydrogeologist

## TABLES

**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)	
				UVF GRO	UVF DRO
Action Level (mg/kg)				50	100
SB-1	0 - 2	0.8			
	2 - 4	0.8			
	4 - 6	2.2	SB-1	<0.68	<0.68
	6 - 8	1.9			
SB-2	0 - 2	1.9			
	2 - 4	2.5			
	4 - 6	3.3			
	6 - 8	1707	SB-2	<b>53.9</b>	29
SB-3	0 - 2	2.2			
	2 - 4	3			
	4 - 6	3.2			
	6 - 8	24.5	SB-3	<0.65	3.8
SB-4	0 - 2	2.3			
	2 - 4	2.2			
	4 - 6	2.1			
	6 - 8	5.5	SB-4	<0.47	1.3
SB-5	0 - 2	3			
	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	<b>SB-2</b>	
Analyte	15A NCAC 2L Standard (µg/L)	Concentration (µg/L)
<b>Volatile Organic Compounds</b>		
Ethylbenzene	600	<b>1600</b>
Isopropylbenzene	70	<b>89</b>
Naphthalene	6	<b>320</b>
n-Propylbenzene	70	<b>130</b>
Toluene	600	<b>3700</b>
1,2,4-Trimethylbenzene	400	<b>920</b>
1,3,5-trimethylbenzene	400	220
Xylenes	500	<b>8000</b>
<b>Semivolatile Organic Compounds</b>		
Acenaphthene	80	0.075 J
Acenaphthylene	200	0.037 J
Fluorene	300	0.10 J
1-Methylnaphthalene	1 (IMAC)	<b>14</b>
2-Methylnaphthalene	30	30
Naphthalene	6	<b>190</b>
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



## FIGURES

PROJECT NUMBER  
20080204-010103

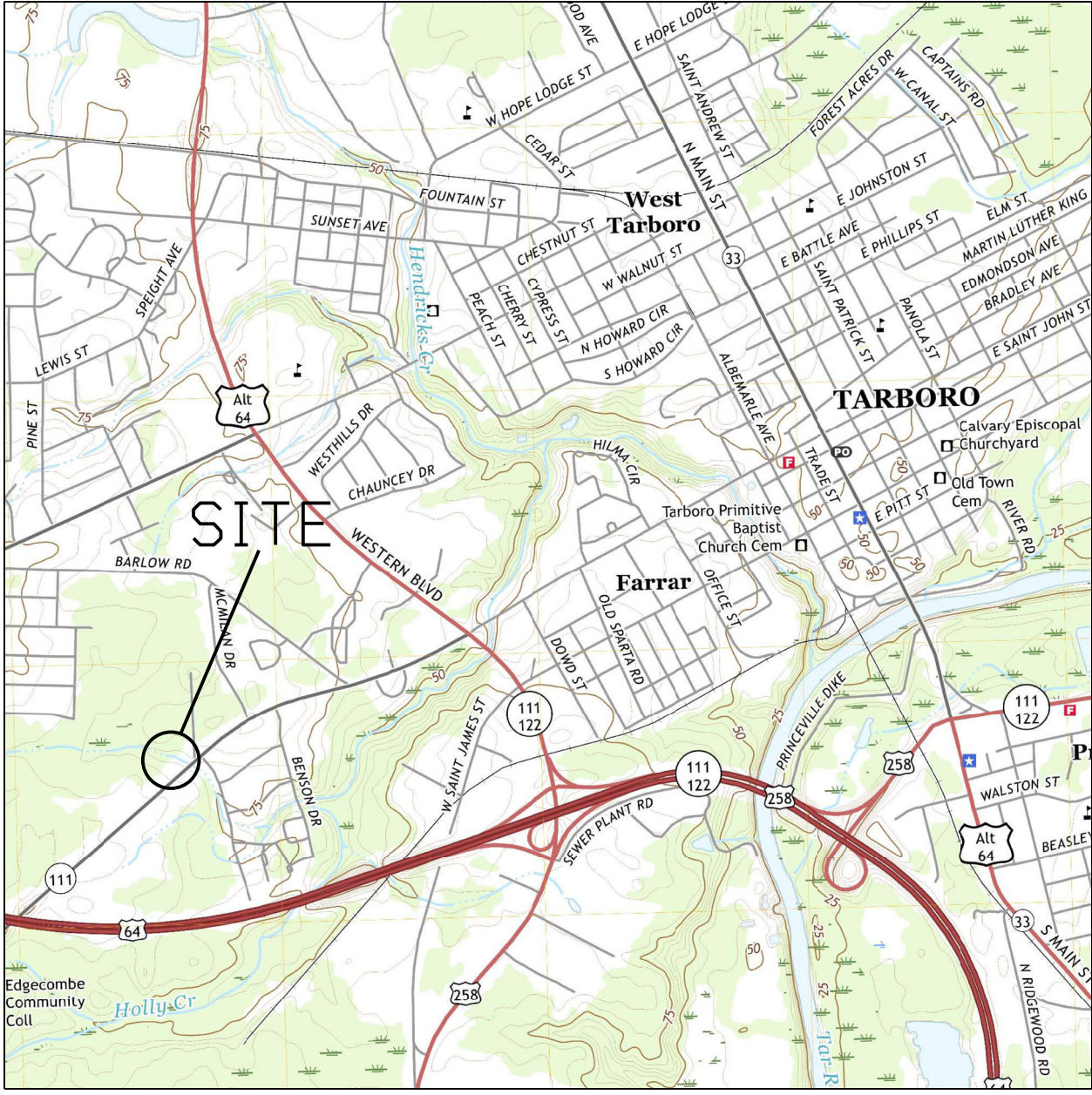
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PROJECT MANAGER  
MWB

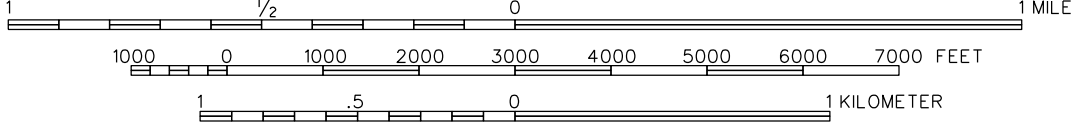
DATE  
JUNE 2020

EDGECOMBE COUNTY PSA

FILE



SCALE 1:24,000



SOURCE: U.S. GEOLOGICAL SURVEY 7.5 MIN QUADRANGLE: TARBORO, NC (2019)

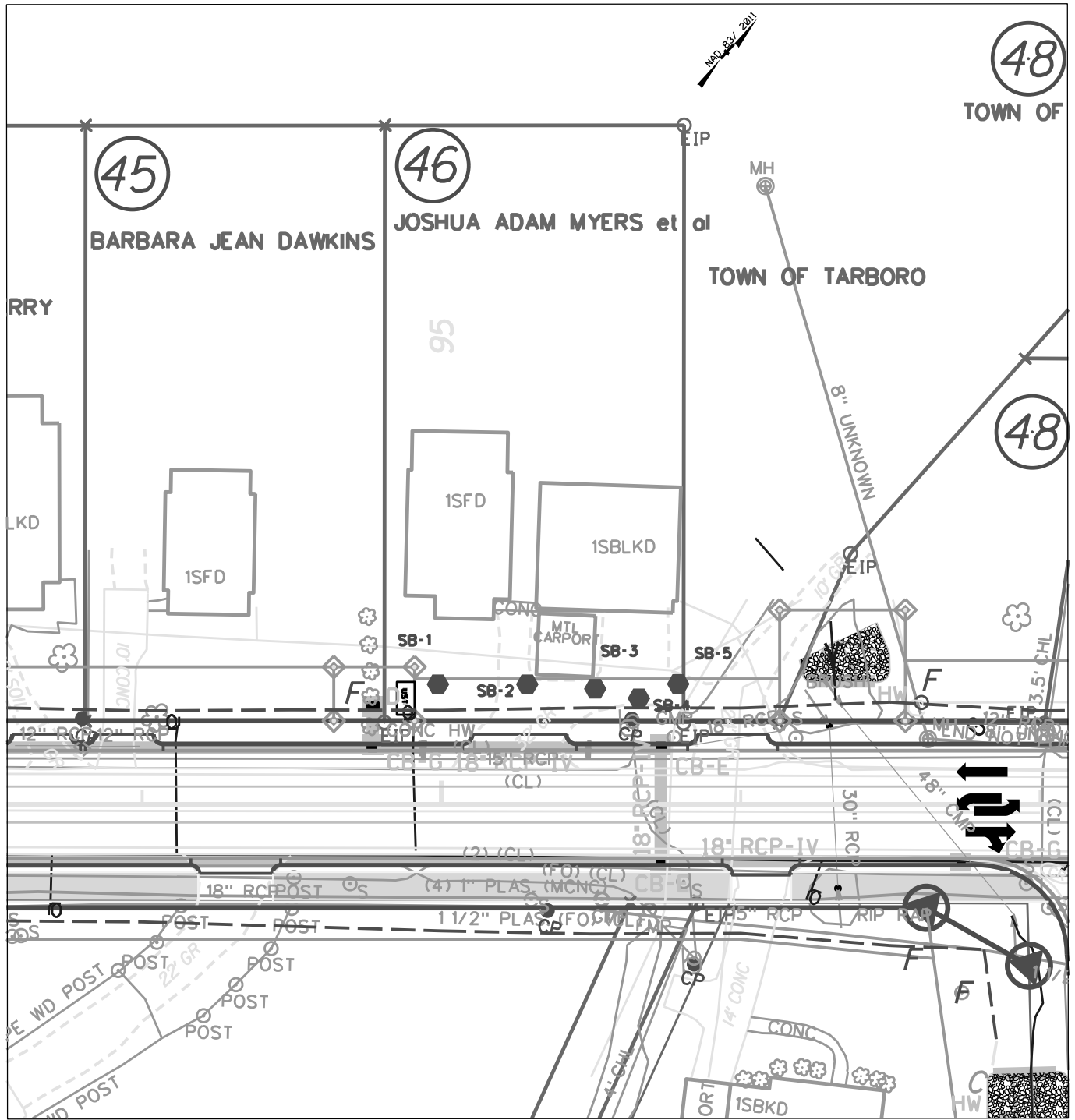


114 EDINBURGH SOUTH DRIVE  
CARY, NORTH CAROLINA 27511  
TEL: (919) 873-1060 FAX: (919) 873-1074



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

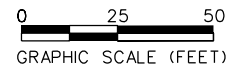
FIGURE  
1

PROJECT NUMBER 19080269-010103  
 DRAFTER MWB  
 CHECKED BY WDN  
 PROJECT MANAGER MWB  
 DATE MAY 2020  
 FILE EDGECOMBE COUNTY PSAS



LEGEND

-  POSSIBLE UST LOCATION
-  BORING LOCATION AND ID



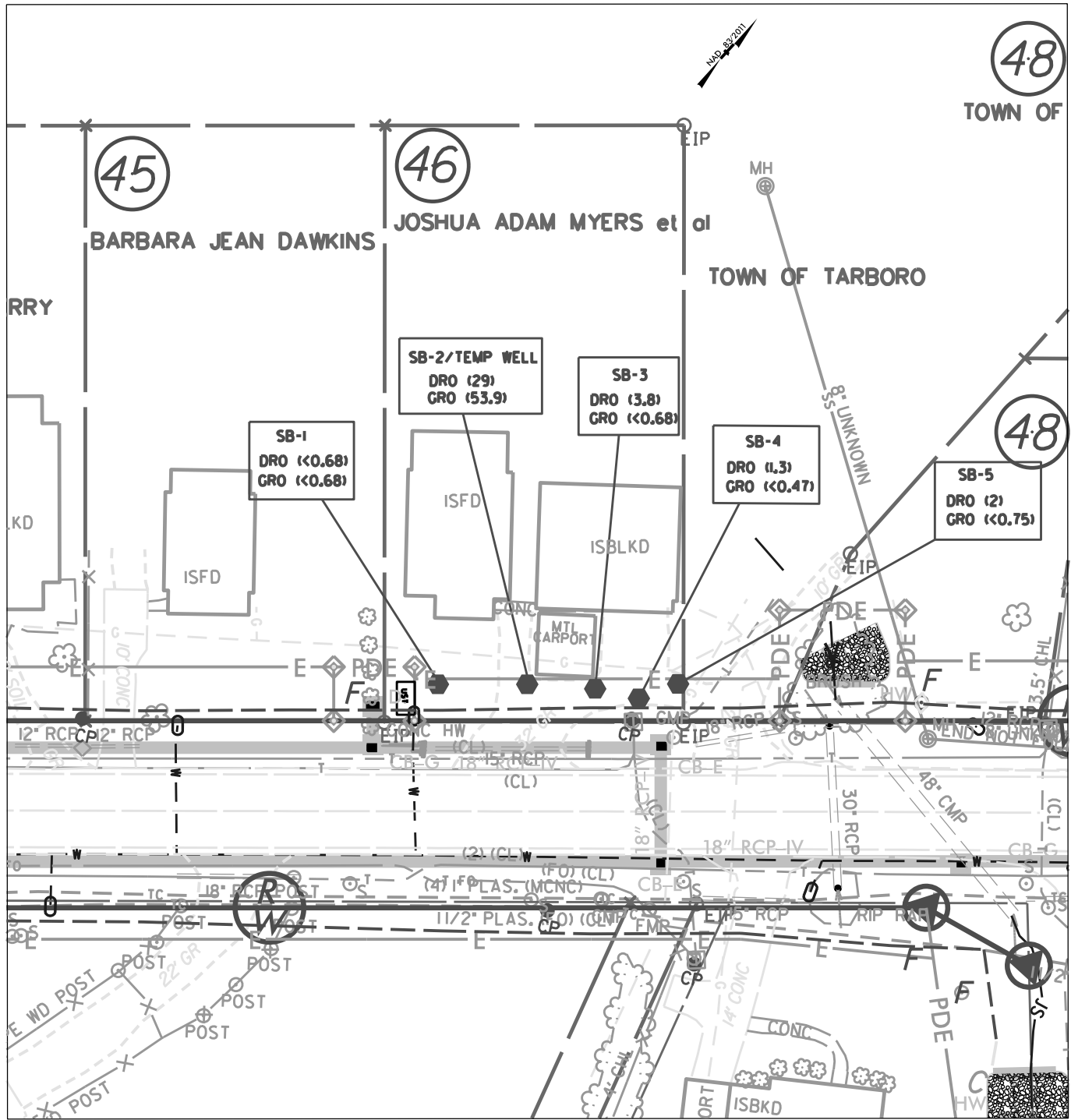
114 EDINBURGH SOUTH DRIVE  
 CARY, NORTH CAROLINA 27511  
 TEL: (919) 873-1060 FAX: (919) 873-1074

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

FIGURE

2

PROJECT NUMBER 19080269-010103  
 DRAFTER MWB  
 CHECKED BY WDN  
 PROJECT MANAGER MWB  
 DATE MAY 2020  
 FILE EDGECOMBE COUNTY PSAS



**LEGEND**

- POSSIBLE UST LOCATION
- SB-1** BORING LOCATION AND ID
- DRO (#23)** DIESEL RANGE ORGANICS (MILLIGRAMS PER KILOGRAM)
- GRO (#23)** GASOLINE RANGE ORGANICS (MILLIGRAMS PER KILOGRAM)

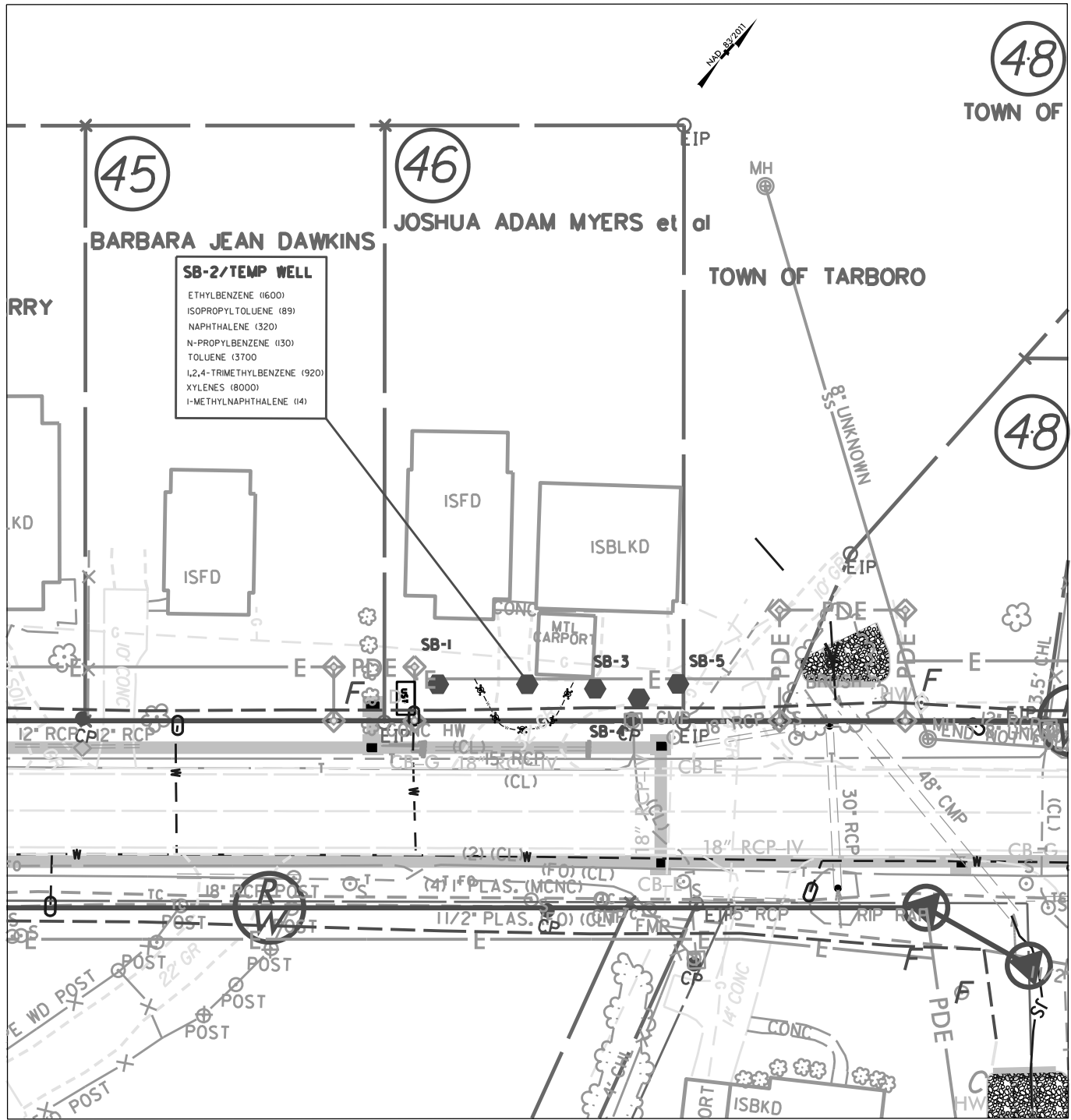


114 EDINBURGH SOUTH DRIVE  
 CARY, NORTH CAROLINA 27511  
 TEL: (919) 873-1060 FAX: (919) 873-1074

SOIL ANALYTICAL RESULTS MAP  
 JOSHUA A MYERS PROPERTY  
 1810 W WILSON STREET  
 TARBORO, NORTH CAROLINA

FIGURE  
 3

PROJECT NUMBER 19080269-010103  
 DRAFTER MWB  
 CHECKED BY WDN  
 PROJECT MANAGER MWB  
 DATE MAY 2020  
 FILE EDGECOMBE COUNTY PSAS



**LEGEND**

- POSSIBLE UST LOCATION
  - SB-1** BORING LOCATION AND ID
  - ABC (123)** COMPOUND ABOVE 2L STANDARDS AND CONCENTRATION IN (MICROGRAMS PER LITER)
- 0 25 50  
GRAPHIC SCALE (FEET)



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

GROUNDWATER ANALYTICAL RESULTS MAP  
 JOSHUA A MYERS PROPERTY  
 1810 W WILSON STREET  
 TARBORO, NORTH CAROLINA

FIGURE  
 4

ATTACHMENT A

# **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



**Draper Aden Associates**  
*Engineering • Surveying • Environmental Services*



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

A handwritten signature in blue ink, appearing to read "Johanna Vaughan".

Johanna Vaughan, P.G.  
Geologist



A handwritten signature in blue ink, appearing to read "F. Doug Pinckney".

Francis Douglas Pinckney, P.E.  
Team Leader/Senior Project Engineer  
Geotechnical and Construction Services



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- Figure 8. Summary of GPR and EM61 Results

## 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 Anomaly 1 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

*Geophysical Study for Possible USTs at 1810 West Wilson Street, Tarboro, North Carolina*

*DAA Project No. 20080204-010203*

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

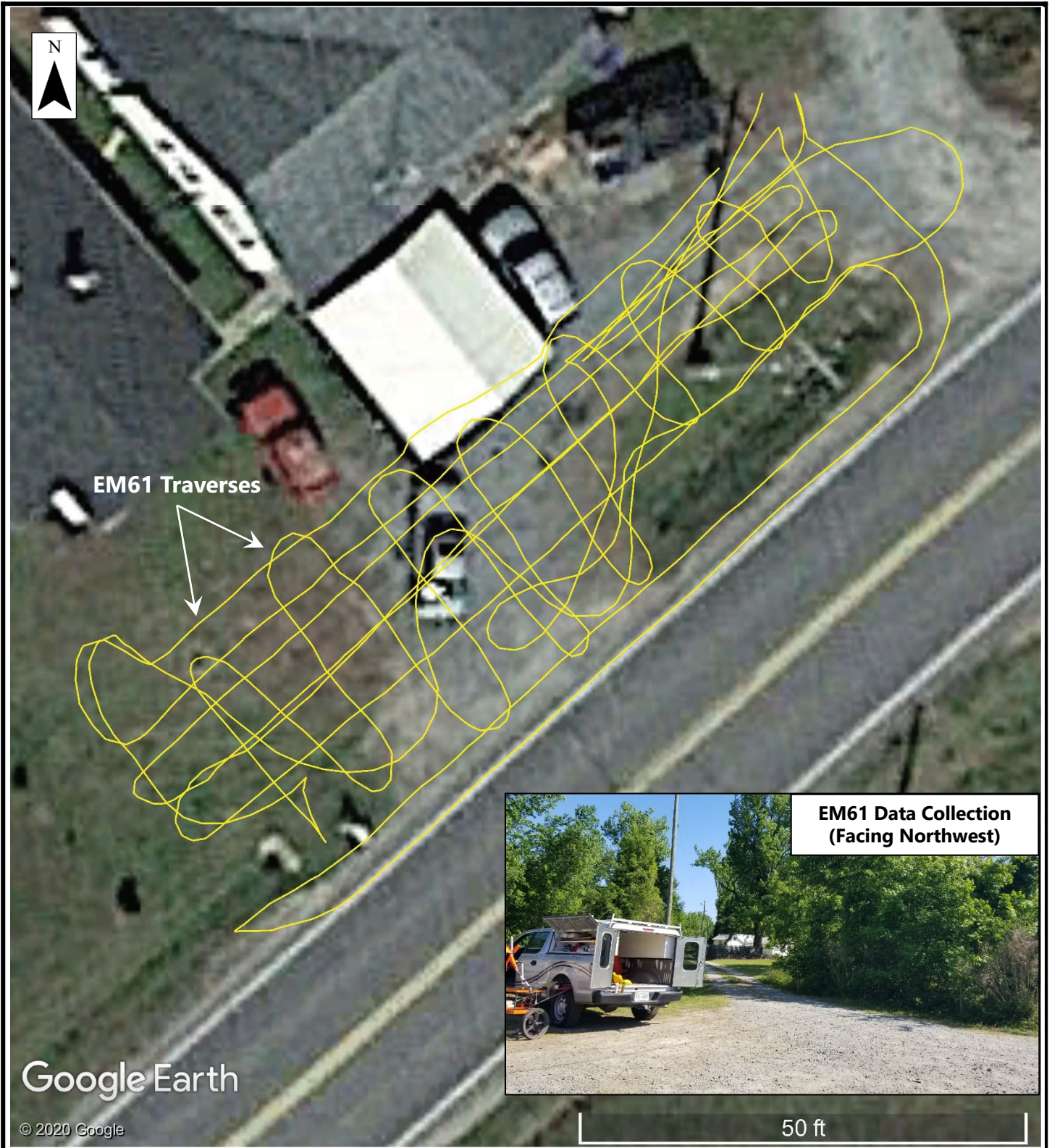
Richmond, VA  
 Charlottesville, VA  
 Hampton Roads, VA

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

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

FIGURE  
 1





**EM61 Traverse Map**

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

PROJECT: 20080204-010203



**Draper Aden Associates**

*Engineering ♦ Surveying ♦ Environmental Services*

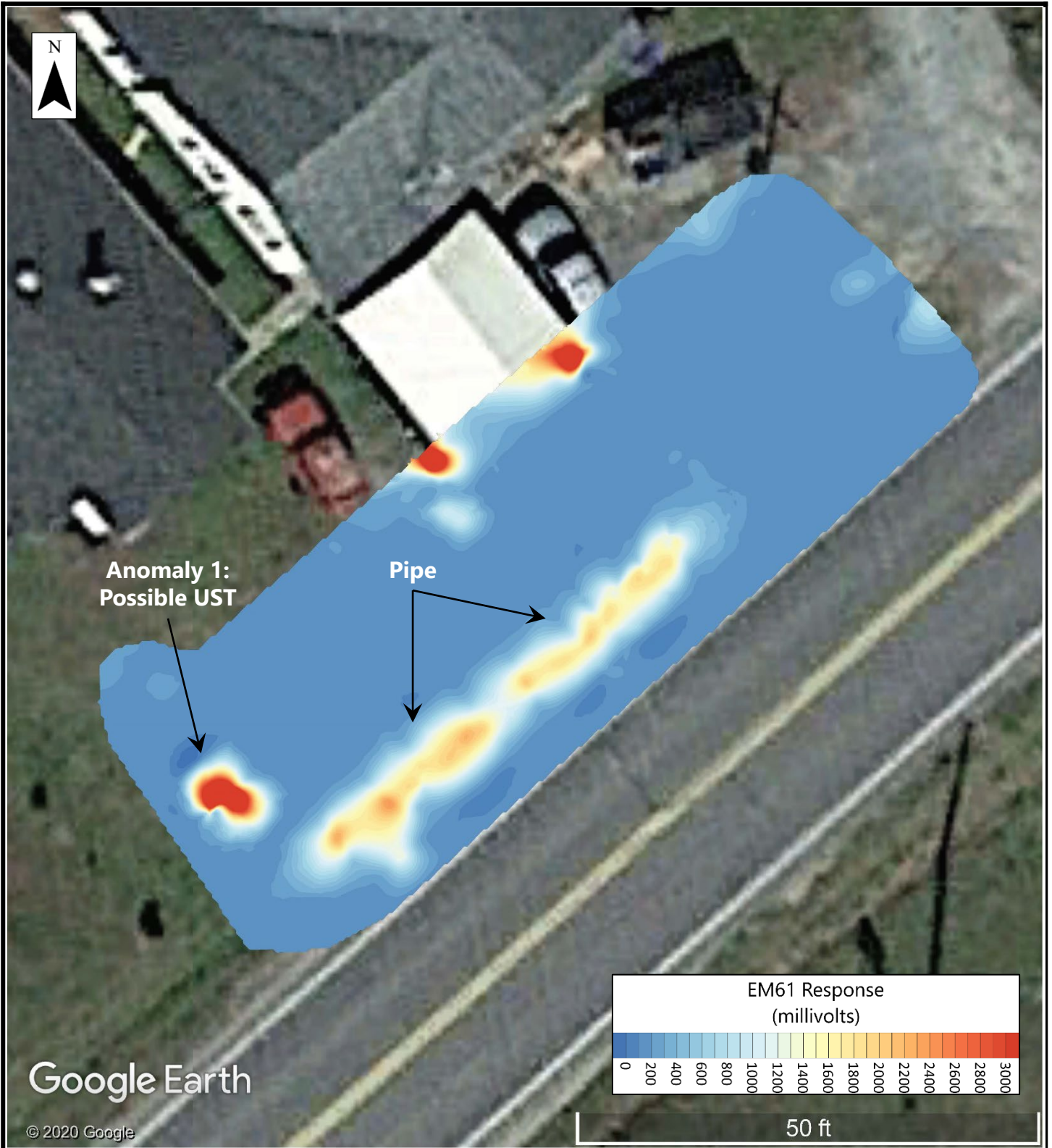
2206 South Main Street  
Blacksburg, VA 24060  
540-552-0444 Fax: 540-552-0291

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  
2



**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  
3



**GPR Traverse Map**

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

PROJECT: 20080204-010203



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Blacksburg, VA 24060  
540-552-0444 Fax: 540-552-0291

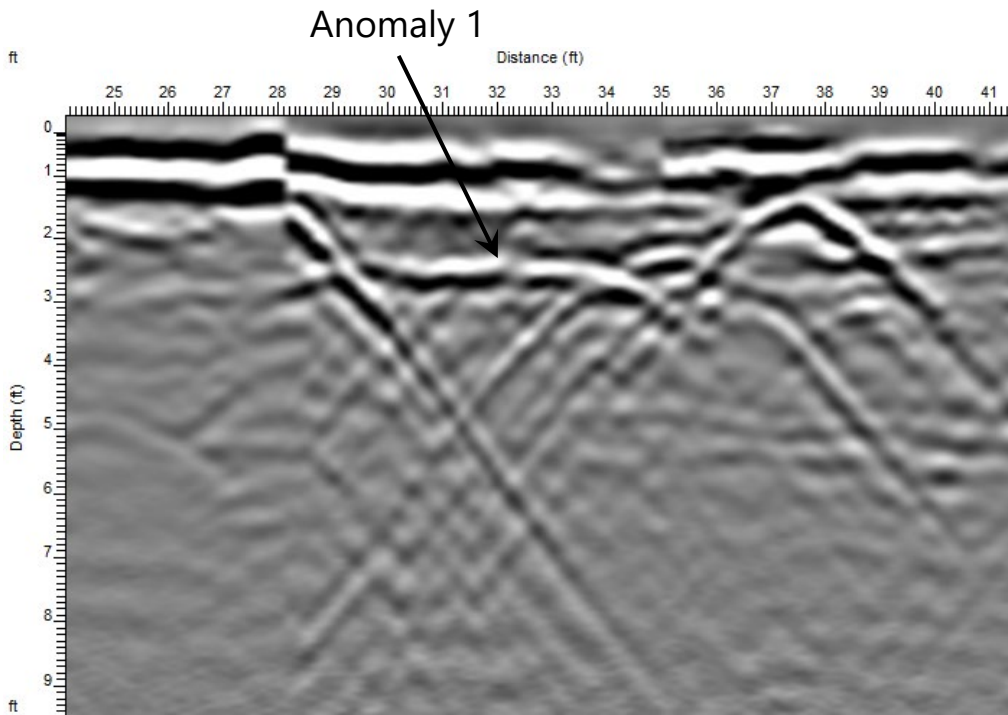
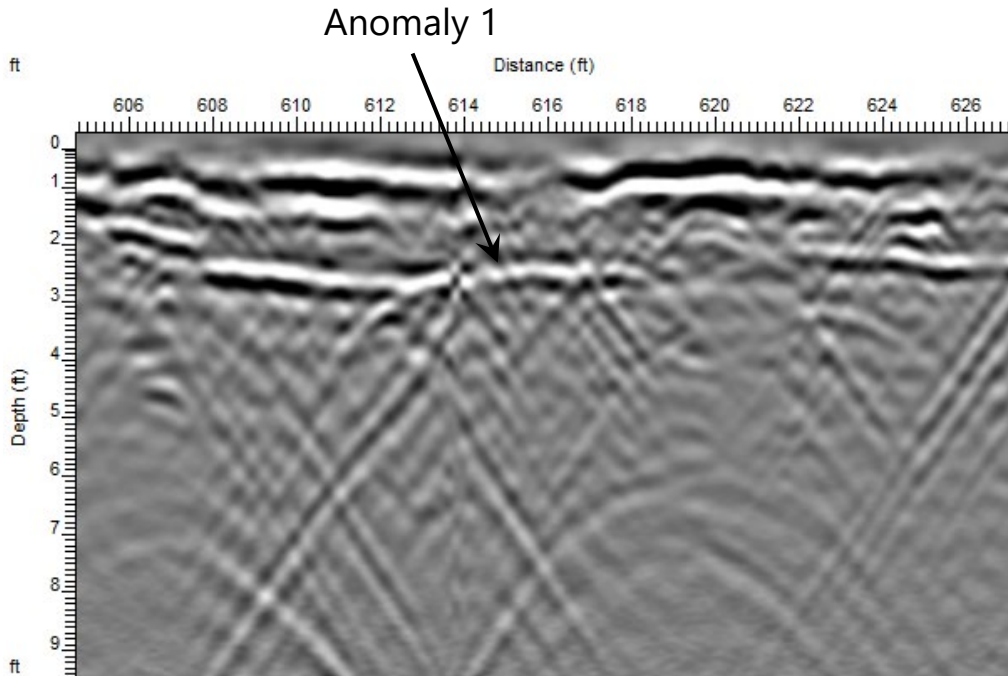
Richmond, VA  
Charlottesville, VA  
Hampton Roads, VA

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

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

FIGURE  
4

**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



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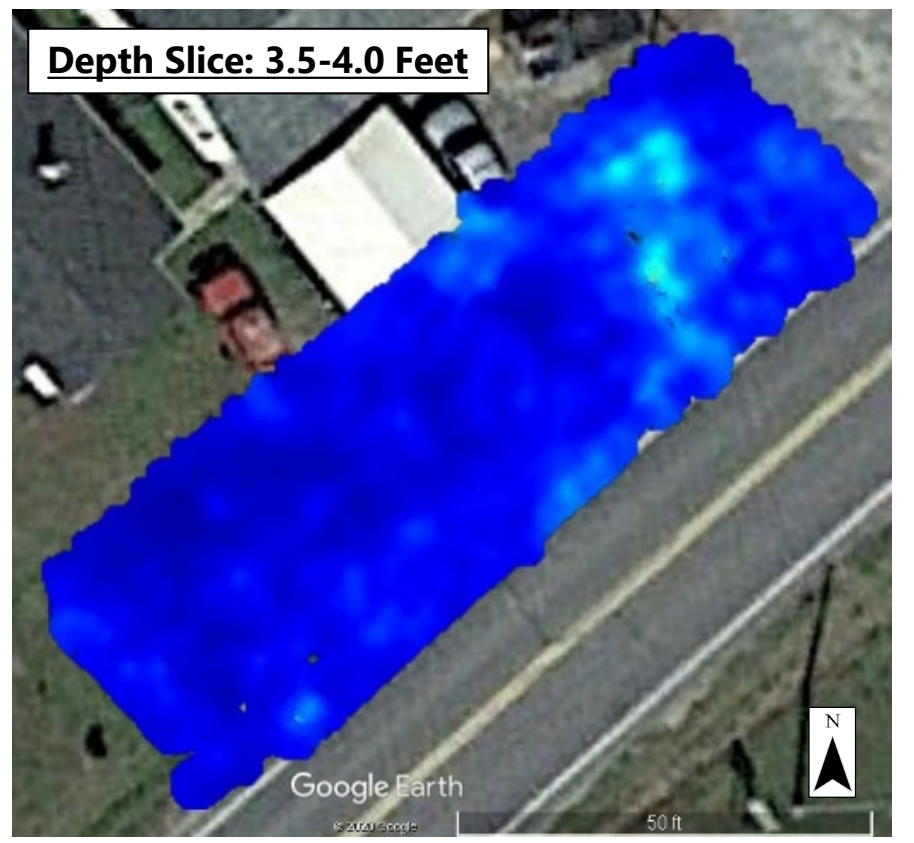
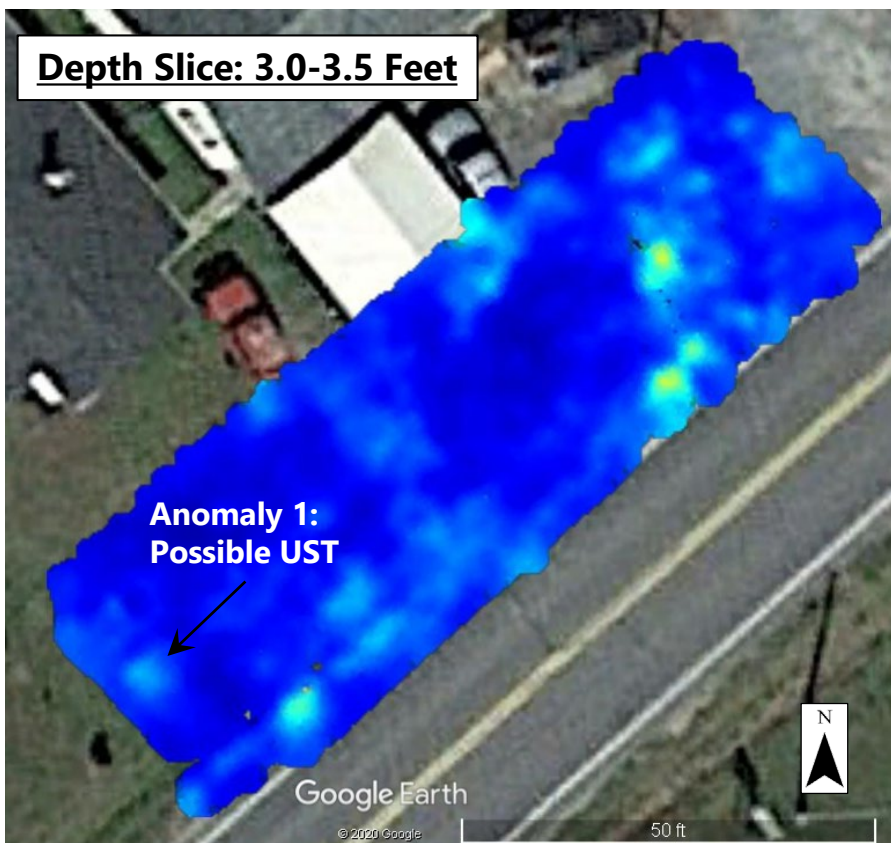
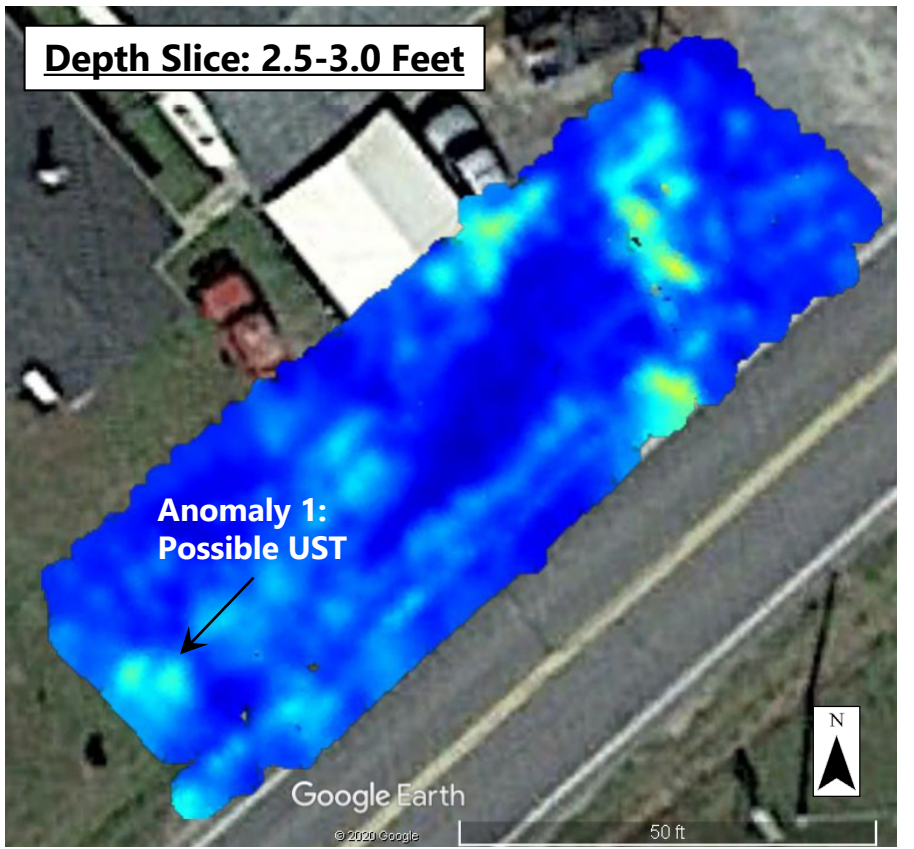
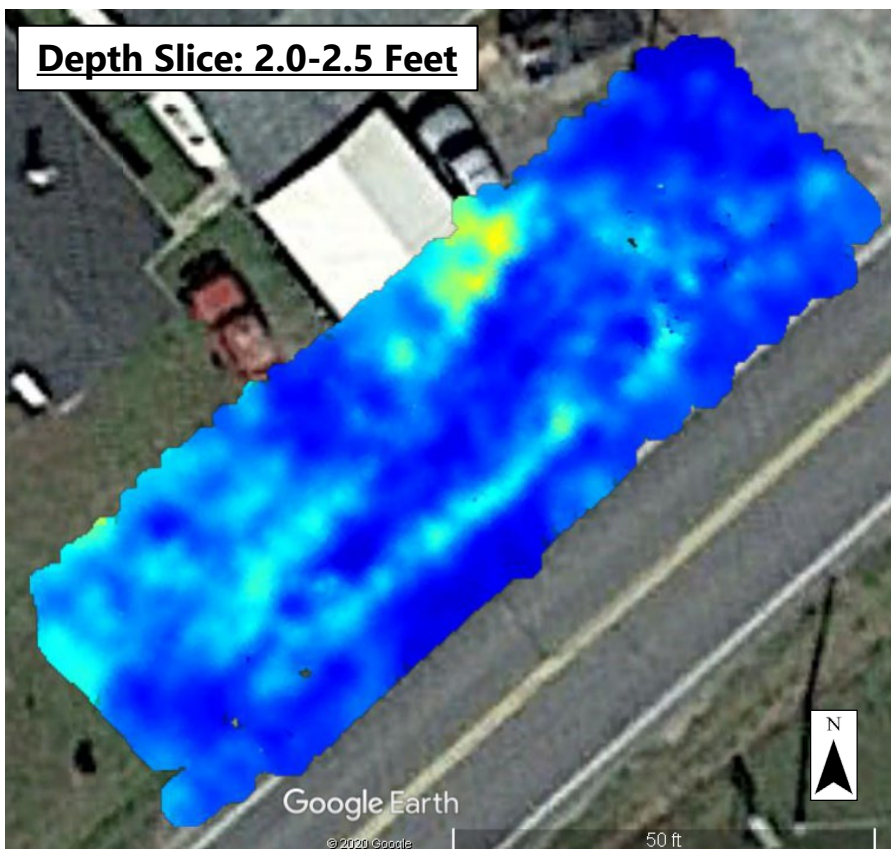
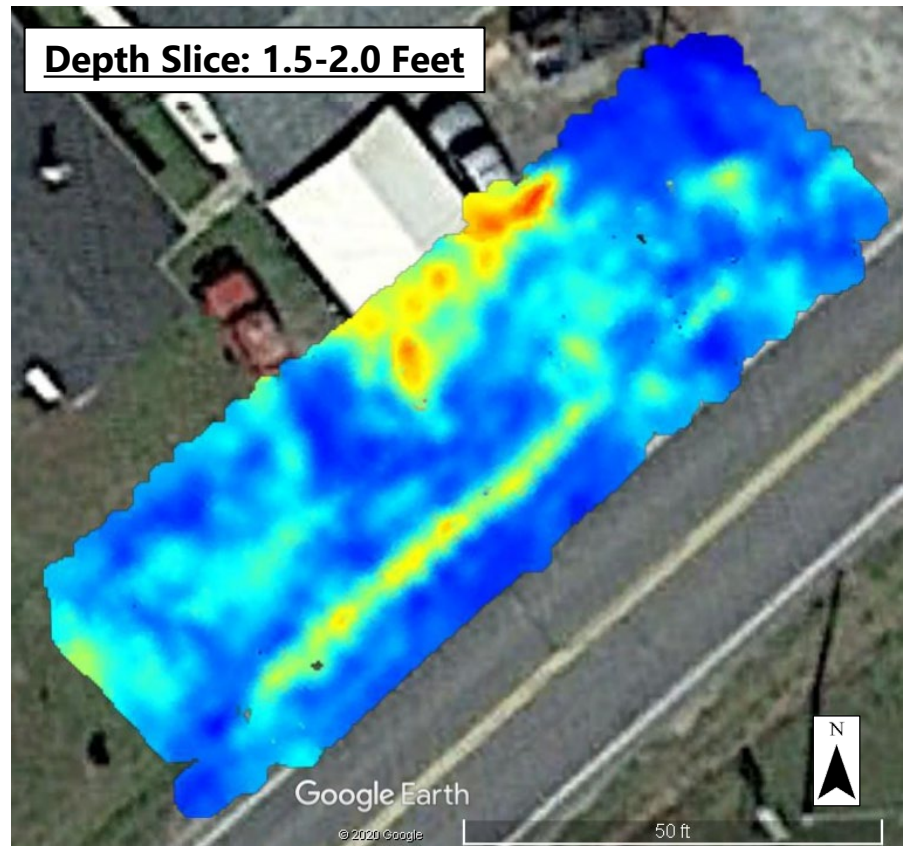
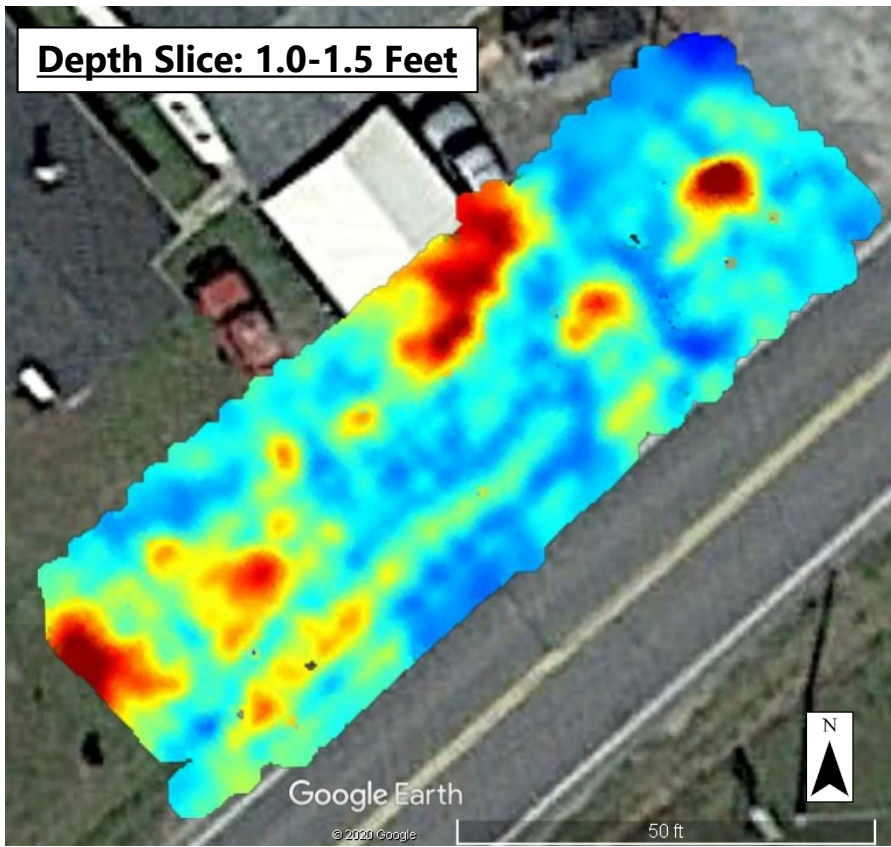
2206 South Main Street  
Blacksburg, VA 24060  
540-552-0444 Fax: 540-552-0291

Richmond, VA  
Charlottesville, VA  
Hampton Roads, VA

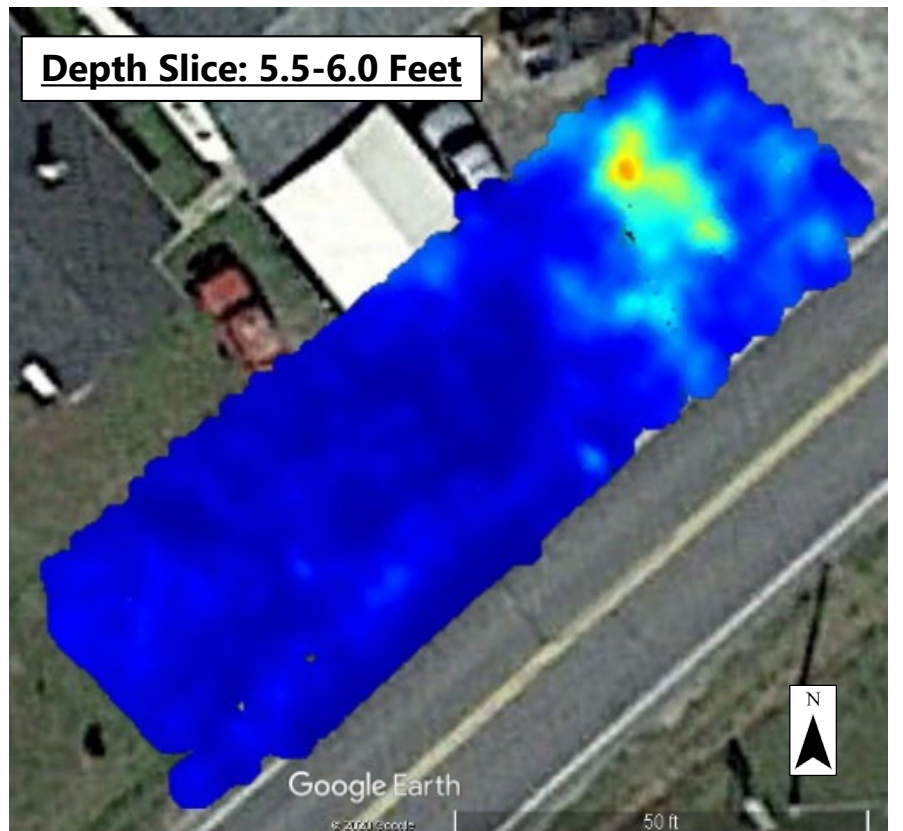
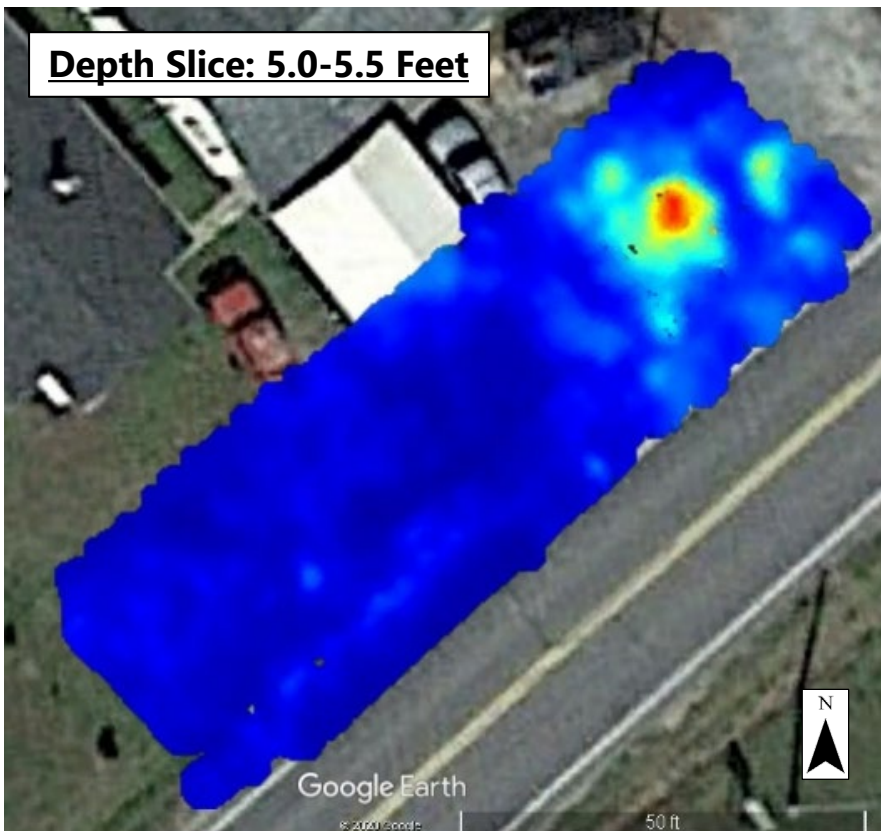
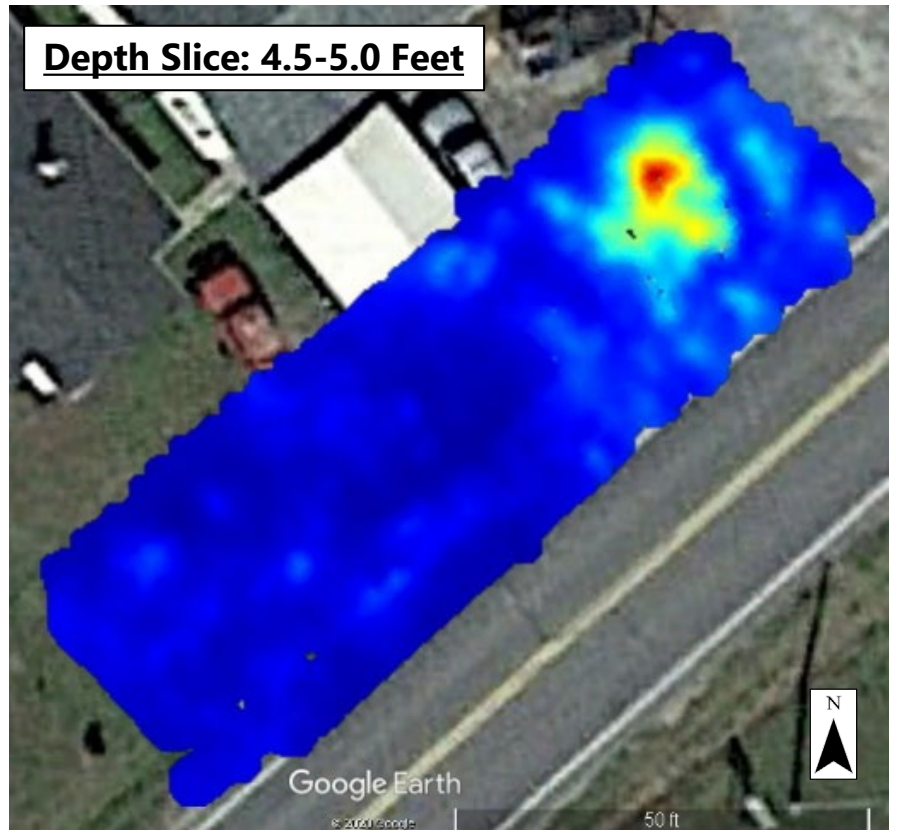
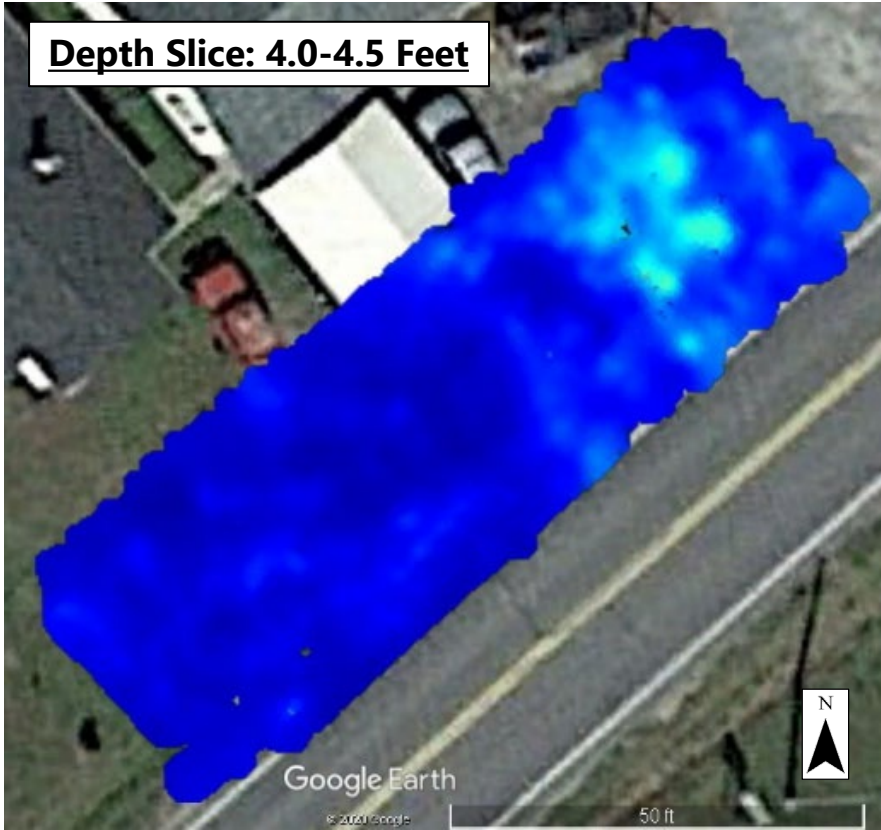
Raleigh, NC  
Fayetteville, NC  
Northern Virginia  
Virginia Beach, VA

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

**FIGURE  
5**




**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



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**

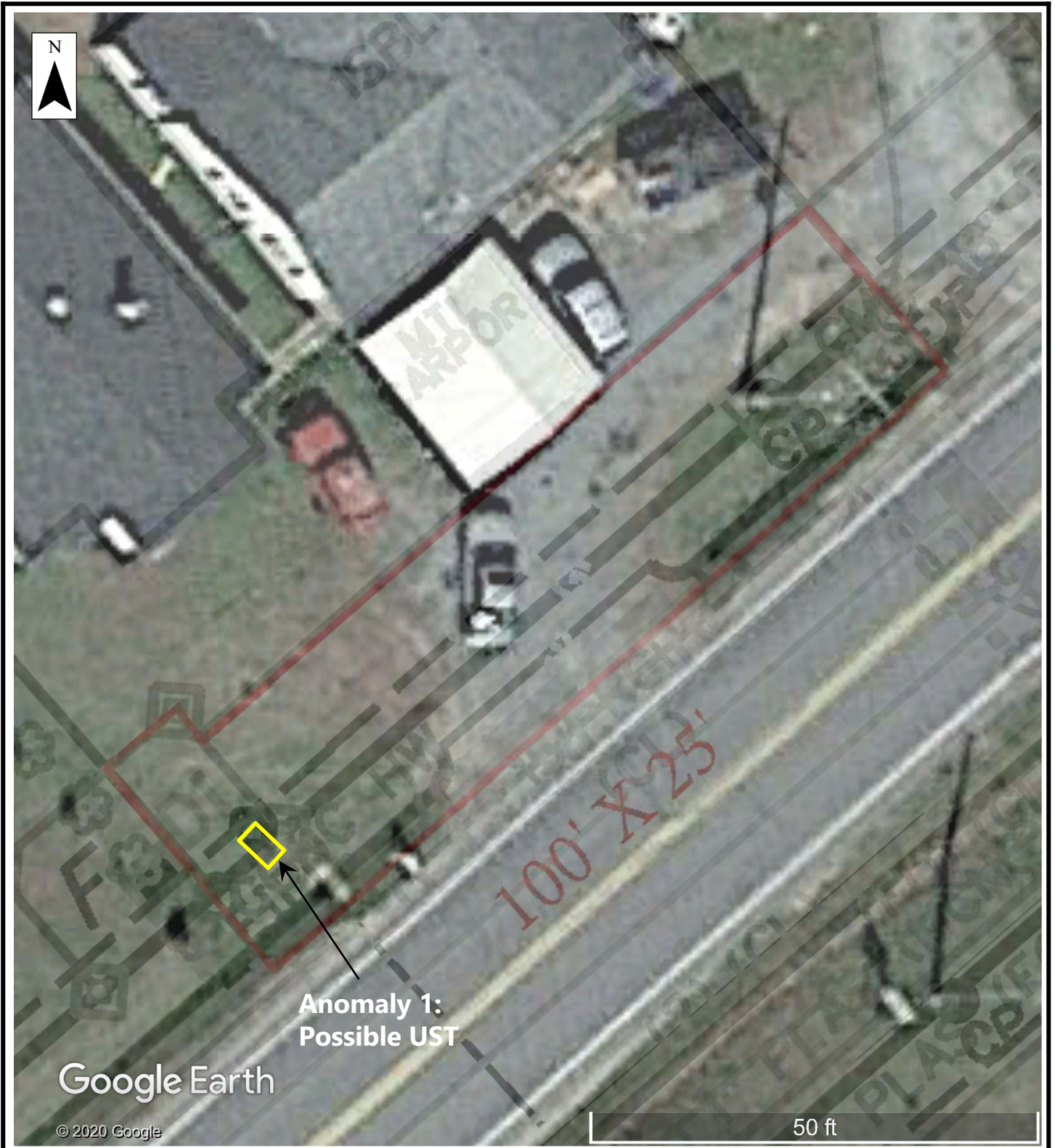
 **Draper Aden Associates**  
*Engineering • Surveying • Environmental Services*  
 2206 South Main Street  
 Blacksburg, VA 24060  
 540-552-0444 Fax: 540-552-0291

Richmond, VA  
 Charlottesville, VA  
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Raleigh, NC  
 Fayetteville, NC  
 Northern Virginia  
 Virginia Beach, VA

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

FIGURE  
7



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

PROJECT: 20080204-010203

**Summary of GPR and EM61 Results**



**Draper Aden Associates**

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

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

FIGURE  
 8

ATTACHMENT B



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

DEPTH (ft bgs)	SAMPLES		PID Reading (ppm)	USCS	LITHOLOGIC DESCRIPTION:	Notes:
	Sample ID	Recovery %				
0						
1			1	SP	POORLY GRADED SAND, Brown to light brown, fine-grained, upper one foot is fill material	
2		100				
3			1		CLAYEY SAND, light orange-brown to orange-red brown, fine-grained, with silt	
4						
5	SB-1		2	SC	Light gray to tan with orange, fine-grained to medium-grained, mottled, moist	
6		100				
7			2	SP-SC	POORLY GRADED SAND WITH CLAY, Light gray to white, fine-grained to medium-grained, mottled with yellow, wet	
8						

**End of Borehole at 8 feet**

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

DEPTH (ft bgs)	SAMPLES		PID Reading (ppm)	USCS	LITHOLOGIC DESCRIPTION:	Notes:
	Sample ID	Recovery %				
0						
1			2	SP	POORLY GRADED SAND, Light brown to brown with gray, fine-grained to medium grained, trace silt, upper one foot is gravel and fill	
2		75				
3			3	SC	CLAYEY SAND, Light orange to orange, fine-grained to medium-grained, consolidated	
4						
5			3			
6		100		CH/CL	CLAY, Brown with orange and gray, fine- grained to medium grained, trace staining, petroleum odor, moist to damp	
7	SB-2		1707			
8				SP	POORLY GRADED SAND, White, medium-grained, wet	

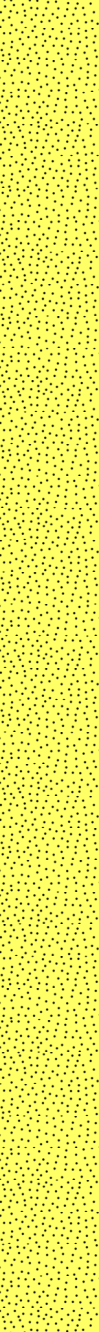
**End of Borehole at 8 feet**

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

DEPTH (ft bgs)	SAMPLES		PID Reading (ppm)	USCS	LITHOLOGIC DESCRIPTION:	Notes:
	Sample ID	Recovery %				
0						
1			2			
2		50		SP	POORLY GRADED SAND, Tan with orange and dark gray, fine-grained to medium-grained, upper one foot is gravel and fill	
3			3			
4						
5			3			
6		100		CH/CL	CLAY, Orange to red brown, with light and dark gray, fine-grained to course grained, with seams of sand with clay (SP-SC)	
7	SB-3		25			
8				SP	POORLY GRADED SAND, Tan to light gray, course grained, wet	

**End of Borehole at 8 feet**

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

DEPTH (ft bgs)	SAMPLES		PID Reading (ppm)	USCS	LITHOLOGIC DESCRIPTION:	Notes:
	Sample ID	Recovery %				
0						
1			2.3	SP	 POORLY GRADED SAND, Tan and brown-gray, with light gray, fine-grained to medium-grained, upper one foot is gravel and fill	
2		60				
3			2.2			Damp
4						Tan to light brown
5			2.1			Wet
6		100				
7	SB-4		5.5		Light blue-gray, course grained	
8	End of Borehole at 8 feet					

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: See Appendix
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

DEPTH (ft bgs)	SAMPLES		PID Reading (ppm)	USCS	LITHOLOGIC DESCRIPTION:	Notes:	
	Sample ID	Recovery %					
0							
1			3.0	SP	POORLY GRADED SAND, Light brown to brown, with gray, fine-grained to medium grained, upper one foot is gravel and fill		
2		50					
3	SB-5		3.5				
4							
5			2.7				Wet
6		50					
7			2.9				
8							

End of Borehole at 8 feet

ATTACHMENT C



**PHOTOGRAPH 1.** Location of SB-1.



**PHOTOGRAPH 2.** Location of SB-2.



**Draper Aden Associates**  
Engineering ^ Surveying ^ Environmental Services

**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.



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





**PHOTOGRAPH 5.** Location of SB-5



**PHOTOGRAPH 6.** Soil identification and sampling area.



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

ATTACHMENT D



### 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

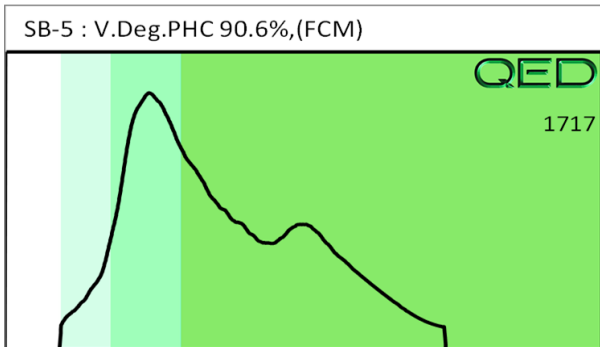
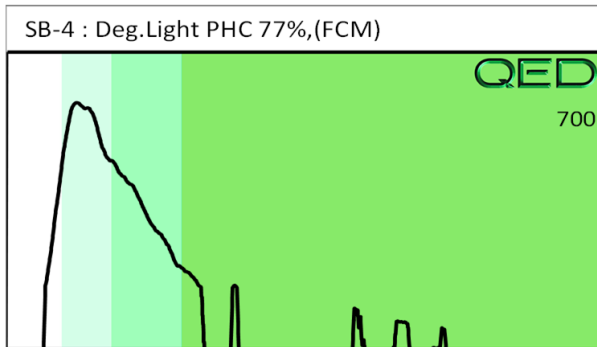
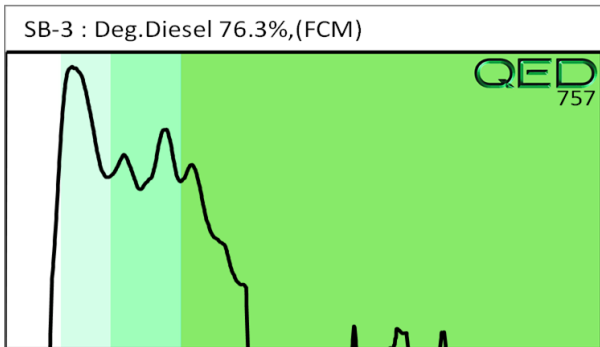
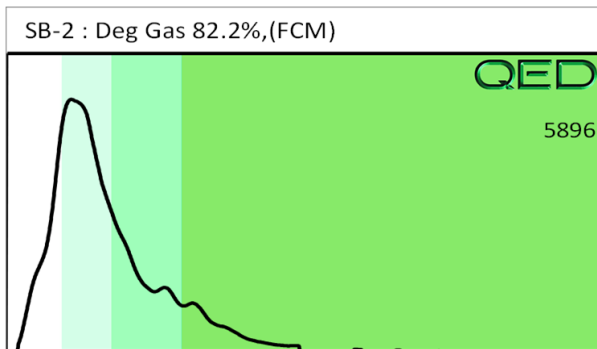
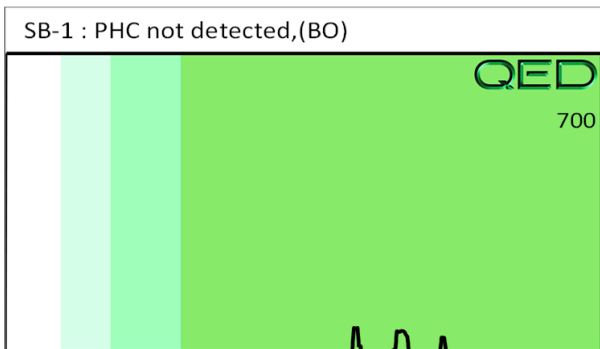
**Operator**

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	BaP	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



Client Name: **PAH**  
 Address: **114 Edinburgs S. br.**  
**Parcel 46**

Contact: **COV, NC, 27811**  
**Brandy Barnes**  
 Project Ref.: **Tobacco PSA 20080204-010103**

Email: **Embronsced@a.com**  
 Phone #: **519 261 0945**

Collected by: **Brandy Barnes**  
**CHAIN OF CUSTODY AND ANALYTICAL REQUEST FORM**



RED Lab, LLC  
 5598 Marvin K Moss Lane  
 MARBIONC Bldg, Suite 2003  
 Wilmington, NC 28409

Each UVF sample will be analyzed for total BTEX, GRO, DRO, TPH, PAH total aromatics and Bap. Standard GC Analyses are for BTEX and Chlorinated Solvents: VC, 1,1 DCE, 1,2 cis DCE, 1,2 trans DCE, TCE, and PCE. Specify target analytes in the space provided below.

Sample Collection Date/Time	TAT Requested	Analysis Type		Initials	Sample ID	Total Wt.	Tare Wt.	Sample Wt.
		48 Hour	UVF					
5/19/26 09:45	X	X		BB	SB-1	52.3	44.2	8.1
1610	X			BB	SB-2	52.0	44.0	8.0
1036	X	X		BB	SB-3	52.2	43.7	8.5
1045	X	X		BB	SB-4	52.6	44.5	8.1
1055	X	X		BB	SB-5	51.3	44.0	7.3

COMMENTS/REQUESTS:

TARGET GC/UVF ANALYTES:

Relinquished by: **1200** Accepted by: **5/19/26** Date/Time: **20 May 2020 12:00**

Relinquished by: **5** Accepted by: **402** Ref. No: **402**

RED Lab USE ONLY

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  
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,

A handwritten signature in black ink that reads "Kerry K. McGee". The signature is written in a cursive, flowing style.

Kerry K. McGee  
Project Manager

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---

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

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	



**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.

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

---

**Qualifications:****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:****Aniline**

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:****4-Chloroaniline**

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

**Aniline**

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.

A handwritten signature in black ink that reads "Tod Kopycinski". The signature is written in a cursive style with a large, sweeping initial "T".

Tod E. Kopycinski  
Laboratory Director

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

Sample Flags: RL-11

**Volatile Organic Compounds by GC/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	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromobenzene	ND	40	6.0	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromochloromethane	ND	40	13	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromodichloromethane	ND	20	6.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromoform	ND	40	18	µg/L	40	V-05	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Bromomethane	ND	80	55	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
2-Butanone (MEK)	ND	800	78	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Butyl Alcohol (TBA)	ND	800	170	µg/L	40	V-05	SW-846 8260D	5/21/20	5/21/20 19:49	MFF
n-Butylbenzene	ND	40	8.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
sec-Butylbenzene	ND	40	6.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Butylbenzene	ND	40	6.8	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
tert-Butyl Ethyl Ether (TBEE)	ND	20	6.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Carbon Disulfide	ND	200	180	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Carbon Tetrachloride	ND	40	4.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chlorobenzene	ND	40	6.0	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chlorodibromomethane	ND	20	8.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chloroethane	ND	80	14	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chloroform	ND	80	6.8	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Chloromethane	ND	80	18	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
2-Chlorotoluene	ND	40	4.8	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
4-Chlorotoluene	ND	40	5.6	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dibromo-3-chloropropane (DBCP)	ND	200	21	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dibromoethane (EDB)	ND	20	7.6	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Dibromomethane	ND	40	15	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2-Dichlorobenzene	ND	40	6.4	µg/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	µg/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

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

Sample Flags: RL-11

**Volatile Organic Compounds by GC/MS**

Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst
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	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Isopropylbenzene (Cumene)	89	80	6.8	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
p-Isopropyltoluene (p-Cymene)	ND	80	8.0	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Methyl tert-Butyl Ether (MTBE)	ND	40	10	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Methylene Chloride	ND	200	14	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
4-Methyl-2-pentanone (MIBK)	ND	400	67	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Naphthalene	320	200	12	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
n-Propylbenzene	130	40	5.2	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Styrene	ND	80	4.4	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,1,1,2-Tetrachloroethane	ND	40	11	µg/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	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Tetrachloroethylene	ND	40	7.2	µg/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	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
Trichloroethylene	ND	40	9.6	µg/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	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,2,4-Trimethylbenzene	920	40	7.2	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF
1,3,5-Trimethylbenzene	220	40	5.6	µg/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	µg/L	40		SW-846 8260D	5/21/20	5/21/20 19:49	MFF

Surrogates	% Recovery	Recovery Limits	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

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

Semivolatile Organic Compounds by GC/MS

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	µg/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	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Butylbenzylphthalate	ND	10	0.30	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Carbazole	ND	10	0.29	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Chloroaniline	ND	10	0.35	µg/L	1	V-34	SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Chloro-3-methylphenol	ND	10	0.49	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Chloronaphthalene	ND	10	0.47	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Chlorophenol	ND	10	0.39	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Chlorophenylphenylether	ND	10	0.32	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Chrysene (SIM)	ND	0.21	0.015	µg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Dibenz(a,h)anthracene (SIM)	ND	0.10	0.017	µg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Dibenzofuran	ND	5.1	0.27	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Di-n-butylphthalate	ND	10	0.47	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2-Dichlorobenzene	ND	5.1	0.47	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,3-Dichlorobenzene	ND	5.1	0.47	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,4-Dichlorobenzene	ND	5.1	0.39	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
3,3-Dichlorobenzidine	ND	10	0.37	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dichlorophenol	ND	10	0.31	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Diethylphthalate	ND	10	0.23	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dimethylphenol	ND	10	0.82	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Dimethylphthalate	ND	10	0.31	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4,6-Dinitro-2-methylphenol	ND	10	2.0	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dinitrophenol	ND	10	1.7	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4-Dinitrotoluene	ND	10	0.34	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,6-Dinitrotoluene	ND	10	0.35	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Di-n-octylphthalate	ND	10	0.54	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2-Diphenylhydrazine/Azobenzene	ND	10	0.38	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Fluoranthene (SIM)	ND	0.51	0.026	µg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Fluorene (SIM)	0.10	1.0	0.035	µg/L	1	J	SW-846 8270E	5/21/20	5/28/20 1:32	IMR

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

Semivolatile Organic Compounds by GC/MS

Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst
Hexachlorobenzene	ND	10	0.44	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Hexachlorobutadiene	ND	10	0.61	µg/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	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2-Nitrophenol	ND	10	0.43	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
4-Nitrophenol	ND	10	0.64	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
N-Nitrosodimethylamine	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	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
Pyrene (SIM)	ND	1.0	0.024	µg/L	1		SW-846 8270E	5/21/20	5/28/20 1:32	IMR
Pyridine	ND	5.1	2.9	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2,4,5-Tetrachlorobenzene	ND	10	0.34	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
1,2,4-Trichlorobenzene	ND	5.1	0.57	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4,5-Trichlorophenol	ND	10	0.49	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL
2,4,6-Trichlorophenol	ND	10	0.34	µg/L	1		SW-846 8270E	5/21/20	5/27/20 13:13	BGL

Surrogates	% Recovery	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	61.1	30-130	5/28/20 1:32
2-Fluorobiphenyl	76.2	30-130	5/27/20 20:55
2,4,6-Tribromophenol	71.0	15-110	5/27/20 13:13
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 20:55

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

**Sample Extraction Data****Prep Method: SW-846 5030B    Analytical Method: SW-846 8260D**

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

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

<b>Lab Number [Field ID]</b>	<b>Batch</b>	<b>Initial [mL]</b>	<b>Final [mL]</b>	<b>Date</b>
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**

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



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**QUALITY CONTROL**

**Volatile Organic Compounds by GC/MS - Quality Control**

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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**Batch B258487 - SW-846 5030B**

**Blank (B258487-BLK1)**

Prepared & Analyzed: 05/21/20

Acetone	ND	50	µg/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	2.0	µg/L							
2-Chlorotoluene	ND	1.0	µg/L							
4-Chlorotoluene	ND	1.0	µg/L							
1,2-Dibromo-3-chloropropane (DBCP)	ND	5.0	µg/L							
1,2-Dibromoethane (EDB)	ND	0.50	µg/L							
Dibromomethane	ND	1.0	µg/L							
1,2-Dichlorobenzene	ND	1.0	µg/L							
1,3-Dichlorobenzene	ND	1.0	µg/L							
1,4-Dichlorobenzene	ND	1.0	µg/L							
trans-1,4-Dichloro-2-butene	ND	2.0	µg/L							
Dichlorodifluoromethane (Freon 12)	ND	2.0	µg/L							V-05
1,1-Dichloroethane	ND	1.0	µg/L							
1,2-Dichloroethane	ND	1.0	µg/L							
1,1-Dichloroethylene	ND	1.0	µg/L							
cis-1,2-Dichloroethylene	ND	1.0	µg/L							
trans-1,2-Dichloroethylene	ND	1.0	µg/L							
1,2-Dichloropropane	ND	1.0	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	µg/L							
p-Isopropyltoluene (p-Cymene)	ND	1.0	µg/L							
Methyl tert-Butyl Ether (MTBE)	ND	1.0	µg/L							

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch B258487 - SW-846 5030B

Blank (B258487-BLK1)

Prepared & Analyzed: 05/21/20

Methylene Chloride	ND	5.0	µg/L							
4-Methyl-2-pentanone (MIBK)	ND	10	µg/L							
Naphthalene	ND	2.0	µg/L							
n-Propylbenzene	ND	1.0	µg/L							
Styrene	ND	1.0	µg/L							
1,1,1,2-Tetrachloroethane	ND	1.0	µg/L							V-05
1,1,2,2-Tetrachloroethane	ND	0.50	µg/L							
Tetrachloroethylene	ND	1.0	µg/L							
Tetrahydrofuran	ND	10	µg/L							
Toluene	ND	1.0	µg/L							
1,2,3-Trichlorobenzene	ND	5.0	µg/L							
1,2,4-Trichlorobenzene	ND	1.0	µg/L							
1,3,5-Trichlorobenzene	ND	1.0	µg/L							
1,1,1-Trichloroethane	ND	1.0	µg/L							
1,1,2-Trichloroethane	ND	1.0	µg/L							
Trichloroethylene	ND	1.0	µg/L							
Trichlorofluoromethane (Freon 11)	ND	2.0	µg/L							L-04, V-05
1,2,3-Trichloropropane	ND	2.0	µg/L							
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	ND	1.0	µg/L							
1,2,4-Trimethylbenzene	ND	1.0	µg/L							
1,3,5-Trimethylbenzene	ND	1.0	µg/L							
Vinyl Chloride	ND	2.0	µg/L							
m+p Xylene	ND	2.0	µg/L							
o-Xylene	ND	1.0	µg/L							
Surrogate: 1,2-Dichloroethane-d4	27.5		µg/L	25.0		110	70-130			
Surrogate: Toluene-d8	23.6		µg/L	25.0		94.5	70-130			
Surrogate: 4-Bromofluorobenzene	25.2		µg/L	25.0		101	70-130			

LCS (B258487-BS1)

Prepared & Analyzed: 05/21/20

Acetone	89.2	50	µg/L	100		89.2	70-160			†
Acrylonitrile	11.2	5.0	µg/L	10.0		112	70-130			
tert-Amyl Methyl Ether (TAME)	8.36	0.50	µg/L	10.0		83.6	70-130			
Benzene	10.5	1.0	µg/L	10.0		105	70-130			
Bromobenzene	10.3	1.0	µg/L	10.0		103	70-130			
Bromochloromethane	10.8	1.0	µg/L	10.0		108	70-130			
Bromodichloromethane	8.78	0.50	µg/L	10.0		87.8	70-130			
Bromoform	8.87	1.0	µg/L	10.0		88.7	70-130			V-05
Bromomethane	5.25	2.0	µg/L	10.0		52.5	40-160			†
2-Butanone (MEK)	110	20	µg/L	100		110	40-160			†
tert-Butyl Alcohol (TBA)	68.8	20	µg/L	100		68.8	40-160			V-05 †
n-Butylbenzene	11.7	1.0	µg/L	10.0		117	70-130			
sec-Butylbenzene	12.0	1.0	µg/L	10.0		120	70-130			
tert-Butylbenzene	11.5	1.0	µg/L	10.0		115	70-130			
tert-Butyl Ethyl Ether (TBEE)	9.32	0.50	µg/L	10.0		93.2	70-130			
Carbon Disulfide	8.24	5.0	µg/L	10.0		82.4	70-130			
Carbon Tetrachloride	8.42	1.0	µg/L	10.0		84.2	70-130			
Chlorobenzene	9.77	1.0	µg/L	10.0		97.7	70-130			
Chlorodibromomethane	8.21	0.50	µg/L	10.0		82.1	70-130			
Chloroethane	7.43	2.0	µg/L	10.0		74.3	70-130			
Chloroform	10.6	2.0	µg/L	10.0		106	70-130			
Chloromethane	6.19	2.0	µg/L	10.0		61.9	40-160			†
2-Chlorotoluene	10.9	1.0	µg/L	10.0		109	70-130			

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
<b>Batch B258487 - SW-846 5030B</b>										
<b>LCS (B258487-BS1)</b>										
Prepared & Analyzed: 05/21/20										
4-Chlorotoluene	10.9	1.0	µg/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	µg/L	10.0		92.2	70-130			
Dibromomethane	8.99	1.0	µg/L	10.0		89.9	70-130			
1,2-Dichlorobenzene	10.2	1.0	µg/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	µg/L	10.0		101	70-130			
trans-1,4-Dichloro-2-butene	10.5	2.0	µg/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	µg/L	10.0		108	70-130			
1,2-Dichloroethane	8.15	1.0	µg/L	10.0		81.5	70-130			
1,1-Dichloroethylene	8.71	1.0	µg/L	10.0		87.1	70-130			
cis-1,2-Dichloroethylene	11.3	1.0	µg/L	10.0		113	70-130			
trans-1,2-Dichloroethylene	9.51	1.0	µg/L	10.0		95.1	70-130			
1,2-Dichloropropane	10.1	1.0	µg/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	µg/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	70-160			†
Isopropylbenzene (Cumene)	10.3	1.0	µg/L	10.0		103	70-130			
p-Isopropyltoluene (p-Cymene)	10.8	1.0	µg/L	10.0		108	70-130			
Methyl tert-Butyl Ether (MTBE)	8.07	1.0	µg/L	10.0		80.7	70-130			
Methylene Chloride	10.0	5.0	µg/L	10.0		100	70-130			
4-Methyl-2-pentanone (MIBK)	99.2	10	µg/L	100		99.2	70-160			†
Naphthalene	10.8	2.0	µg/L	10.0		108	40-130			†
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			
1,1,1,2-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			
<b>Trichlorofluoromethane (Freon 11)</b>	6.45	2.0	µg/L	10.0		<b>64.5</b> *	70-130			V-05, L-04
1,2,3-Trichloropropane	9.56	2.0	µg/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	µg/L	10.0		100	70-130			
Vinyl Chloride	6.50	2.0	µg/L	10.0		65.0	40-160			†

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

**QUALITY CONTROL**

**Volatile Organic Compounds by GC/MS - Quality Control**

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
<b>Batch B258487 - SW-846 5030B</b>										
<b>LCS (B258487-BS1)</b>										
Prepared & Analyzed: 05/21/20										
m+p Xylene	21.4	2.0	µg/L	20.0		107	70-130			
o-Xylene	10.7	1.0	µg/L	10.0		107	70-130			
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			
<b>LCS Dup (B258487-BSD1)</b>										
Prepared & Analyzed: 05/21/20										
Acetone	91.1	50	µg/L	100		91.1	70-160	2.08	25	†
Acrylonitrile	10.4	5.0	µg/L	10.0		104	70-130	7.52	25	
tert-Amyl Methyl Ether (TAME)	8.33	0.50	µg/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	µg/L	10.0		105	70-130	2.92	25	
Bromodichloromethane	8.94	0.50	µg/L	10.0		89.4	70-130	1.81	25	
Bromoform	8.36	1.0	µg/L	10.0		83.6	70-130	5.92	25	V-05
Bromomethane	5.82	2.0	µg/L	10.0		58.2	40-160	10.3	25	†
2-Butanone (MEK)	111	20	µg/L	100		111	40-160	0.987	25	†
tert-Butyl Alcohol (TBA)	64.0	20	µg/L	100		64.0	40-160	7.29	25	V-05 †
n-Butylbenzene	11.7	1.0	µg/L	10.0		117	70-130	0.171	25	
sec-Butylbenzene	12.1	1.0	µg/L	10.0		121	70-130	1.16	25	
tert-Butylbenzene	11.6	1.0	µg/L	10.0		116	70-130	0.782	25	
tert-Butyl Ethyl Ether (TBEE)	9.64	0.50	µg/L	10.0		96.4	70-130	3.38	25	
Carbon Disulfide	7.55	5.0	µg/L	10.0		75.5	70-130	8.74	25	
Carbon Tetrachloride	8.48	1.0	µg/L	10.0		84.8	70-130	0.710	25	
Chlorobenzene	9.61	1.0	µg/L	10.0		96.1	70-130	1.65	25	
Chlorodibromomethane	8.24	0.50	µg/L	10.0		82.4	70-130	0.365	25	
Chloroethane	7.52	2.0	µg/L	10.0		75.2	70-130	1.20	25	
Chloroform	10.8	2.0	µg/L	10.0		108	70-130	1.22	25	
Chloromethane	6.08	2.0	µg/L	10.0		60.8	40-160	1.79	25	†
2-Chlorotoluene	10.8	1.0	µg/L	10.0		108	70-130	0.739	25	
4-Chlorotoluene	10.9	1.0	µg/L	10.0		109	70-130	0.276	25	
1,2-Dibromo-3-chloropropane (DBCP)	9.06	5.0	µg/L	10.0		90.6	70-130	0.988	25	
1,2-Dibromoethane (EDB)	9.22	0.50	µg/L	10.0		92.2	70-130	0.00	25	
Dibromomethane	8.93	1.0	µg/L	10.0		89.3	70-130	0.670	25	
1,2-Dichlorobenzene	10.3	1.0	µg/L	10.0		103	70-130	1.46	25	
1,3-Dichlorobenzene	10.7	1.0	µg/L	10.0		107	70-130	1.21	25	
1,4-Dichlorobenzene	10.3	1.0	µg/L	10.0		103	70-130	1.67	25	
trans-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 †
1,1-Dichloroethane	11.1	1.0	µg/L	10.0		111	70-130	2.65	25	
1,2-Dichloroethane	8.22	1.0	µg/L	10.0		82.2	70-130	0.855	25	
1,1-Dichloroethylene	8.87	1.0	µg/L	10.0		88.7	70-130	1.82	25	
cis-1,2-Dichloroethylene	11.2	1.0	µg/L	10.0		112	70-130	1.42	25	
trans-1,2-Dichloroethylene	9.45	1.0	µg/L	10.0		94.5	70-130	0.633	25	
1,2-Dichloropropane	9.71	1.0	µg/L	10.0		97.1	70-130	4.04	25	
1,3-Dichloropropane	9.45	0.50	µg/L	10.0		94.5	70-130	0.106	25	
2,2-Dichloropropane	9.06	1.0	µg/L	10.0		90.6	40-130	0.664	25	†
1,1-Dichloropropene	9.00	2.0	µg/L	10.0		90.0	70-130	3.17	25	
cis-1,3-Dichloropropene	9.34	0.50	µg/L	10.0		93.4	70-130	3.82	25	
trans-1,3-Dichloropropene	8.87	0.50	µg/L	10.0		88.7	70-130	1.12	25	
Diethyl Ether	9.63	2.0	µg/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	

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

**QUALITY CONTROL**

**Volatile Organic Compounds by GC/MS - Quality Control**

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

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**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
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**Batch B258532 - SW-846 3510C**

**Blank (B258532-BLK1)**

Prepared: 05/21/20 Analyzed: 05/26/20

Acetophenone	ND	10	µg/L							
Aniline	ND	5.0	µg/L							V-05
Benzidine	ND	20	µg/L							V-04, V-05
Benzoic Acid	ND	10	µg/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							
4-Bromophenylphenylether	ND	10	µg/L							
Butylbenzylphthalate	ND	10	µg/L							
Carbazole	ND	10	µg/L							
4-Chloroaniline	ND	10	µg/L							V-34
4-Chloro-3-methylphenol	ND	10	µg/L							
2-Chloronaphthalene	ND	10	µg/L							
2-Chlorophenol	ND	10	µg/L							
4-Chlorophenylphenylether	ND	10	µg/L							
Dibenzofuran	ND	5.0	µg/L							
Di-n-butylphthalate	ND	10	µg/L							
1,2-Dichlorobenzene	ND	5.0	µg/L							
1,3-Dichlorobenzene	ND	5.0	µg/L							
1,4-Dichlorobenzene	ND	5.0	µg/L							
3,3-Dichlorobenzidine	ND	10	µg/L							
2,4-Dichlorophenol	ND	10	µg/L							
Diethylphthalate	ND	10	µg/L							
2,4-Dimethylphenol	ND	10	µg/L							
Dimethylphthalate	ND	10	µg/L							
4,6-Dinitro-2-methylphenol	ND	10	µg/L							
2,4-Dinitrophenol	ND	10	µg/L							
2,4-Dinitrotoluene	ND	10	µg/L							
2,6-Dinitrotoluene	ND	10	µg/L							
Di-n-octylphthalate	ND	10	µg/L							
1,2-Diphenylhydrazine/Azobenzene	ND	10	µg/L							
Hexachlorobenzene	ND	10	µg/L							
Hexachlorobutadiene	ND	10	µg/L							
Hexachlorocyclopentadiene	ND	10	µg/L							
Hexachloroethane	ND	10	µg/L							
Isophorone	ND	10	µg/L							
1-Methylnaphthalene	ND	5.0	µg/L							
2-Methylphenol	ND	10	µg/L							
3/4-Methylphenol	ND	10	µg/L							
2-Nitroaniline	ND	10	µg/L							
3-Nitroaniline	ND	10	µg/L							
4-Nitroaniline	ND	10	µg/L							
Nitrobenzene	ND	10	µg/L							
2-Nitrophenol	ND	10	µg/L							
4-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	µg/L							
Phenol	ND	10	µg/L							

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
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Batch B258532 - SW-846 3510C

Blank (B258532-BLK1)

Prepared: 05/21/20 Analyzed: 05/26/20

Pyridine	ND	5.0	µg/L							
1,2,4,5-Tetrachlorobenzene	ND	10	µg/L							
1,2,4-Trichlorobenzene	ND	5.0	µg/L							
2,4,5-Trichlorophenol	ND	10	µg/L							
2,4,6-Trichlorophenol	ND	10	µg/L							
Surrogate: 2-Fluorophenol	101		µg/L	200		50.3	15-110			
Surrogate: Phenol-d6	70.3		µg/L	200		35.2	15-110			
Surrogate: Nitrobenzene-d5	80.6		µg/L	100		80.6	30-130			
Surrogate: 2-Fluorobiphenyl	80.5		µg/L	100		80.5	30-130			
Surrogate: 2,4,6-Tribromophenol	169		µg/L	200		84.5	15-110			
Surrogate: p-Terphenyl-d14	93.6		µg/L	100		93.6	30-130			

LCS (B258532-BS1)

Prepared: 05/21/20 Analyzed: 05/26/20

Acetophenone	39.4	10	µg/L	50.0		78.7	40-140			
Aniline	31.7	5.0	µg/L	50.0		63.4	40-140			V-05
Benzidine	30.5	20	µg/L	50.0		60.9	40-140			V-04, V-05
Benzoic Acid	17.6	10	µg/L	50.0		35.2	10-130			†
Bis(2-chloroethoxy)methane	40.9	10	µg/L	50.0		81.8	40-140			
Bis(2-chloroethyl)ether	38.8	10	µg/L	50.0		77.5	40-140			
Bis(2-chloroisopropyl)ether	43.4	10	µg/L	50.0		86.9	40-140			
Bis(2-Ethylhexyl)phthalate	44.4	10	µg/L	50.0		88.7	40-140			
4-Bromophenylphenylether	40.1	10	µg/L	50.0		80.2	40-140			
Butylbenzylphthalate	40.6	10	µg/L	50.0		81.2	40-140			
Carbazole	40.7	10	µg/L	50.0		81.4	40-140			
4-Chloroaniline	35.8	10	µg/L	50.0		71.7	40-140			V-34
4-Chloro-3-methylphenol	38.5	10	µg/L	50.0		77.1	30-130			
2-Chloronaphthalene	34.0	10	µg/L	50.0		68.0	40-140			
2-Chlorophenol	35.1	10	µg/L	50.0		70.2	30-130			
4-Chlorophenylphenylether	39.1	10	µg/L	50.0		78.2	40-140			
Dibenzofuran	40.6	5.0	µg/L	50.0		81.1	40-140			
Di-n-butylphthalate	43.6	10	µg/L	50.0		87.3	40-140			
1,2-Dichlorobenzene	28.9	5.0	µg/L	50.0		57.9	40-140			
1,3-Dichlorobenzene	27.2	5.0	µg/L	50.0		54.3	40-140			
1,4-Dichlorobenzene	27.5	5.0	µg/L	50.0		55.0	40-140			
3,3-Dichlorobenzidine	44.5	10	µg/L	50.0		88.9	40-140			
2,4-Dichlorophenol	38.3	10	µg/L	50.0		76.6	30-130			
Diethylphthalate	40.4	10	µg/L	50.0		80.7	40-140			
2,4-Dimethylphenol	32.7	10	µg/L	50.0		65.4	30-130			
Dimethylphthalate	40.0	10	µg/L	50.0		80.1	40-140			
4,6-Dinitro-2-methylphenol	39.5	10	µg/L	50.0		79.1	30-130			
2,4-Dinitrophenol	36.2	10	µg/L	50.0		72.3	30-130			
2,4-Dinitrotoluene	38.5	10	µg/L	50.0		77.0	40-140			
2,6-Dinitrotoluene	40.6	10	µg/L	50.0		81.2	40-140			
Di-n-octylphthalate	43.3	10	µg/L	50.0		86.6	40-140			
1,2-Diphenylhydrazine/Azobenzene	45.1	10	µg/L	50.0		90.2	40-140			
Hexachlorobenzene	39.5	10	µg/L	50.0		79.0	40-140			
Hexachlorobutadiene	28.0	10	µg/L	50.0		56.1	40-140			
Hexachlorocyclopentadiene	24.6	10	µg/L	50.0		49.3	30-140			†
Hexachloroethane	26.3	10	µg/L	50.0		52.6	40-140			
Isophorone	39.2	10	µg/L	50.0		78.3	40-140			
1-Methylnaphthalene	34.4	5.0	µg/L	50.0		68.8	40-140			
2-Methylphenol	34.9	10	µg/L	50.0		69.8	30-130			

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
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Batch B258532 - SW-846 3510C

LCS (B258532-BS1)

Prepared: 05/21/20 Analyzed: 05/26/20

3/4-Methylphenol	34.2	10	µg/L	50.0		68.3	30-130			
2-Nitroaniline	50.6	10	µg/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	50.0		70.0	30-130			
Phenol	18.3	10	µg/L	50.0		36.6	20-130			†
Pyridine	20.1	5.0	µg/L	50.0		40.1	10-140			†
1,2,4,5-Tetrachlorobenzene	36.3	10	µg/L	50.0		72.6	40-140			
1,2,4-Trichlorobenzene	30.8	5.0	µg/L	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		µg/L	200		50.6	15-110			
Surrogate: Phenol-d6	75.4		µg/L	200		37.7	15-110			
Surrogate: Nitrobenzene-d5	78.2		µg/L	100		78.2	30-130			
Surrogate: 2-Fluorobiphenyl	82.5		µg/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		µg/L	100		86.1	30-130			

LCS Dup (B258532-BS1)

Prepared: 05/21/20 Analyzed: 05/26/20

Acetophenone	36.3	10	µg/L	50.0		72.7	40-140	8.03	20	
Aniline	28.4	5.0	µg/L	50.0		56.7	40-140	11.1	50	V-05 ‡
Benzidine	27.8	20	µg/L	50.0		55.6	40-140	9.16	20	V-04, V-05
Benzoic Acid	17.1	10	µg/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	50.0		82.5	40-140	7.29	20	
4-Bromophenylphenylether	38.2	10	µg/L	50.0		76.4	40-140	4.83	20	
Butylbenzylphthalate	38.7	10	µg/L	50.0		77.3	40-140	4.92	20	
Carbazole	39.2	10	µg/L	50.0		78.4	40-140	3.75	20	
4-Chloroaniline	34.0	10	µg/L	50.0		68.1	40-140	5.21	20	V-34
4-Chloro-3-methylphenol	37.0	10	µg/L	50.0		74.0	30-130	4.11	20	
2-Chloronaphthalene	31.7	10	µg/L	50.0		63.3	40-140	7.15	20	
2-Chlorophenol	33.4	10	µg/L	50.0		66.9	30-130	4.84	20	
4-Chlorophenylphenylether	38.4	10	µg/L	50.0		76.8	40-140	1.76	20	
Dibenzofuran	38.9	5.0	µg/L	50.0		77.7	40-140	4.28	20	
Di-n-butylphthalate	40.0	10	µg/L	50.0		80.0	40-140	8.65	20	
1,2-Dichlorobenzene	29.4	5.0	µg/L	50.0		58.8	40-140	1.61	20	
1,3-Dichlorobenzene	28.1	5.0	µg/L	50.0		56.1	40-140	3.26	20	
1,4-Dichlorobenzene	28.4	5.0	µg/L	50.0		56.9	40-140	3.32	20	
3,3-Dichlorobenzidine	41.1	10	µg/L	50.0		82.2	40-140	7.83	20	
2,4-Dichlorophenol	36.3	10	µg/L	50.0		72.5	30-130	5.50	20	
Diethylphthalate	39.1	10	µg/L	50.0		78.1	40-140	3.27	20	
2,4-Dimethylphenol	30.4	10	µg/L	50.0		60.8	30-130	7.29	20	
Dimethylphthalate	39.0	10	µg/L	50.0		78.1	40-140	2.58	50	‡



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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
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Batch B258532 - SW-846 3510C

LCS Dup (B258532-BSD1)

Prepared: 05/21/20 Analyzed: 05/26/20

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	µg/L	50.0		71.4	30-130	1.31	50	‡
2,4-Dinitrotoluene	38.0	10	µg/L	50.0		76.1	40-140	1.20	20	
2,6-Dinitrotoluene	39.2	10	µg/L	50.0		78.4	40-140	3.56	20	
Di-n-octylphthalate	40.2	10	µg/L	50.0		80.3	40-140	7.55	20	
1,2-Diphenylhydrazine/Azobenzene	37.8	10	µg/L	50.0		75.6	40-140	17.7	20	
Hexachlorobenzene	37.6	10	µg/L	50.0		75.1	40-140	5.03	20	
Hexachlorobutadiene	28.1	10	µg/L	50.0		56.3	40-140	0.356	20	
Hexachlorocyclopentadiene	22.6	10	µg/L	50.0		45.3	30-140	8.42	50	† ‡
Hexachloroethane	27.0	10	µg/L	50.0		54.0	40-140	2.70	50	‡
Isophorone	34.9	10	µg/L	50.0		69.7	40-140	11.6	20	
1-Methylnaphthalene	32.9	5.0	µg/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	µg/L	50.0		62.5	30-130	8.87	20	
2-Nitroaniline	44.4	10	µg/L	50.0		88.7	40-140	13.2	20	
3-Nitroaniline	39.1	10	µg/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	200		34.8	15-110			
Surrogate: Nitrobenzene-d5	71.6		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	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 Analyzed: 05/27/20

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							
Fluoranthene (SIM)	ND	0.50	µg/L							
Fluorene (SIM)	ND	1.0	µg/L							

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
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Batch B258763 - SW-846 3510C

Blank (B258763-BLK1)

Prepared: 05/21/20 Analyzed: 05/27/20

Indeno(1,2,3-cd)pyrene (SIM)	ND	0.10	µg/L							
2-Methylnaphthalene (SIM)	ND	1.0	µg/L							
Naphthalene (SIM)	ND	1.0	µg/L							
Phenanthrene (SIM)	ND	0.050	µg/L							
Pyrene (SIM)	ND	1.0	µg/L							
Surrogate: Nitrobenzene-d5	86.0		µg/L	100		86.0	30-130			
Surrogate: 2-Fluorobiphenyl	76.3		µg/L	100		76.3	30-130			
Surrogate: p-Terphenyl-d14	74.9		µg/L	100		74.9	30-130			

LCS (B258763-BS1)

Prepared: 05/21/20 Analyzed: 05/27/20

Acenaphthene (SIM)	45.4	6.0	µg/L	50.0		90.8	40-140			
Acenaphthylene (SIM)	44.2	4.0	µg/L	50.0		88.4	40-140			
Anthracene (SIM)	49.1	4.0	µg/L	50.0		98.3	40-140			
Benzo(a)anthracene (SIM)	48.5	1.0	µg/L	50.0		97.1	40-140			
Benzo(a)pyrene (SIM)	51.6	2.0	µg/L	50.0		103	40-140			
Benzo(b)fluoranthene (SIM)	53.7	1.0	µg/L	50.0		107	40-140			
Benzo(g,h,i)perylene (SIM)	54.6	10	µg/L	50.0		109	40-140			
Benzo(k)fluoranthene (SIM)	52.5	4.0	µg/L	50.0		105	40-140			
Chrysene (SIM)	47.3	4.0	µg/L	50.0		94.6	40-140			
Dibenz(a,h)anthracene (SIM)	57.9	2.0	µg/L	50.0		116	40-140			
Fluoranthene (SIM)	51.2	10	µg/L	50.0		102	40-140			
Fluorene (SIM)	48.4	20	µg/L	50.0		96.9	40-140			
Indeno(1,2,3-cd)pyrene (SIM)	60.7	2.0	µg/L	50.0		121	40-140			
2-Methylnaphthalene (SIM)	43.4	20	µg/L	50.0		86.8	40-140			
Naphthalene (SIM)	39.6	20	µg/L	50.0		79.2	40-140			
Phenanthrene (SIM)	46.6	1.0	µg/L	50.0		93.2	40-140			
Pyrene (SIM)	45.4	20	µg/L	50.0		90.9	40-140			
Surrogate: Nitrobenzene-d5	87.1		µg/L	100		87.1	30-130			
Surrogate: 2-Fluorobiphenyl	90.5		µg/L	100		90.5	30-130			
Surrogate: p-Terphenyl-d14	73.7		µg/L	100		73.7	30-130			

LCS Dup (B258763-BSD1)

Prepared: 05/21/20 Analyzed: 05/27/20

Acenaphthene (SIM)	43.9	6.0	µg/L	50.0		87.7	40-140	3.49	20	
Acenaphthylene (SIM)	42.5	4.0	µg/L	50.0		85.1	40-140	3.83	20	
Anthracene (SIM)	47.5	4.0	µg/L	50.0		95.0	40-140	3.39	20	
Benzo(a)anthracene (SIM)	47.2	1.0	µg/L	50.0		94.5	40-140	2.71	20	
Benzo(a)pyrene (SIM)	50.2	2.0	µg/L	50.0		100	40-140	2.75	20	
Benzo(b)fluoranthene (SIM)	52.7	1.0	µg/L	50.0		105	40-140	1.92	20	
Benzo(g,h,i)perylene (SIM)	53.0	10	µg/L	50.0		106	40-140	2.90	20	
Benzo(k)fluoranthene (SIM)	51.4	4.0	µg/L	50.0		103	40-140	2.12	20	
Chrysene (SIM)	46.0	4.0	µg/L	50.0		92.0	40-140	2.79	20	
Dibenz(a,h)anthracene (SIM)	56.8	2.0	µg/L	50.0		114	40-140	1.99	20	
Fluoranthene (SIM)	49.4	10	µg/L	50.0		98.8	40-140	3.58	20	
Fluorene (SIM)	47.3	20	µg/L	50.0		94.6	40-140	2.34	20	
Indeno(1,2,3-cd)pyrene (SIM)	59.0	2.0	µg/L	50.0		118	40-140	2.94	20	‡
2-Methylnaphthalene (SIM)	42.2	20	µg/L	50.0		84.5	40-140	2.66	20	
Naphthalene (SIM)	39.2	20	µg/L	50.0		78.4	40-140	1.07	20	
Phenanthrene (SIM)	44.9	1.0	µg/L	50.0		89.9	40-140	3.67	20	
Pyrene (SIM)	44.5	20	µg/L	50.0		89.0	40-140	2.14	20	
Surrogate: Nitrobenzene-d5	83.6		µg/L	100		83.6	30-130			
Surrogate: 2-Fluorobiphenyl	88.7		µg/L	100		88.7	30-130			
Surrogate: p-Terphenyl-d14	72.7		µg/L	100		72.7	30-130			

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

**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

## Certified Analyses included in this Report

Analyte	Certifications
<i>SW-846 8260D in Water</i>	
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	NC
cis-1,2-Dichloroethylene	NC
trans-1,2-Dichloroethylene	NC
1,2-Dichloropropane	NC
1,3-Dichloropropane	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

**CERTIFICATIONS**

**Certified Analyses included in this Report**

Analyte	Certifications
<b><i>SW-846 8260D in Water</i></b>	
Hexachlorobutadiene	NC
2-Hexanone (MBK)	NC
Isopropylbenzene (Cumene)	NC
p-Isopropyltoluene (p-Cymene)	NC
Methyl tert-Butyl Ether (MTBE)	NC
Methylene Chloride	NC
4-Methyl-2-pentanone (MIBK)	NC
Naphthalene	NC
n-Propylbenzene	NC
Styrene	NC
1,1,1,2-Tetrachloroethane	NC
1,1,2,2-Tetrachloroethane	NC
Tetrachloroethylene	NC
Tetrahydrofuran	NC
Toluene	NC
1,2,3-Trichlorobenzene	NC
1,2,4-Trichlorobenzene	NC
1,3,5-Trichlorobenzene	NC
1,1,1-Trichloroethane	NC
1,1,2-Trichloroethane	NC
Trichloroethylene	NC
Trichlorofluoromethane (Freon 11)	NC
1,2,3-Trichloropropane	NC
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	NC
1,2,4-Trimethylbenzene	NC
1,3,5-Trimethylbenzene	NC
Vinyl Chloride	NC
m+p Xylene	NC
o-Xylene	NC
<b><i>SW-846 8270E in Soil</i></b>	
Acetophenone	NY,NH,ME,NC,VA
Aniline	NY,NH,ME,NC,VA
Benzidine	CT,NY,NH,ME,NC,VA
Benzoic Acid	NY,NH,ME,NC,VA
Bis(2-chloroethoxy)methane	CT,NY,NH,ME,NC,VA
Bis(2-chloroethyl)ether	CT,NY,NH,ME,NC,VA
Bis(2-chloroisopropyl)ether	CT,NY,NH,ME,NC,VA
Bis(2-Ethylhexyl)phthalate	CT,NY,NH,ME,NC,VA
4-Bromophenylphenylether	CT,NY,NH,ME,NC,VA
Butylbenzylphthalate	CT,NY,NH,ME,NC,VA
Carbazole	NC
4-Chloroaniline	CT,NY,NH,ME,NC,VA
4-Chloro-3-methylphenol	CT,NY,NH,ME,NC,VA
2-Chloronaphthalene	CT,NY,NH,NC,VA
2-Chlorophenol	CT,NY,NH,ME,NC,VA
4-Chlorophenylphenylether	CT,NY,NH,ME,NC,VA

**CERTIFICATIONS**

**Certified Analyses included in this Report**

Analyte	Certifications
<b>SW-846 8270E in Soil</b>	
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
<b>SW-846 8270E in Water</b>	
Acetophenone	NY,NC
Aniline	CT,NY,NC,ME,VA

**CERTIFICATIONS**

**Certified Analyses included in this Report**

Analyte	Certifications
<i>SW-846 8270E in Water</i>	
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**

Analyte	Certifications
<i>SW-846 8270E in Water</i>	
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
2-Fluorophenol	NC

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

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 Public 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



20E0837

Parcel 46

Page 1 of 1

Doc # 379 Rev 1\_03242017

39 Spruce Street  
East Longmeadow, MA 01028

http://www.contestlabs.com



Phone: 413-525-2332  
Fax: 413-525-6405

Email: info@contestlabs.com

Company Name: Drape Aelon Associates  
Address: 114 Edinburgh S. Dr., Cary, NC 275  
Project Name: Tarboro PSA Parcel 46  
Project Location: Tarboro, NC Parcel 46  
Project Number: 20080204-0103  
Project Manager: M. Branson  
Con-Test Quote Name/Number:

Invoice Recipient:  
Sampled By: Brady Barnes

Con-Test Work Order #	Client Sample ID / Description	Beginning Date/Time	Ending Date/Time	Composite	Grab	Matrix Code	Conc Code
	<u>SB-2</u>	<u>5/18/20</u>	<u>1135</u>		<u>X</u>	<u>GW</u>	

Requested Turnaround Time:  
 7-Day  10-Day   
 Due Date:  
 Rush-Approval Required:  
 1-Day  3-Day   
 2-Day  4-Day   
 Data Delivery:  
 Format:  PDF  EXCEL   
 Other:  
 CLP Like Data Pkg Required:   
 Email To:  
 Fax To #:

# of Containers	Preservation Code	Container Code
<u>32</u>	<u>H</u>	<u>A</u>

ANALYSIS REQUESTED

<u>20VS</u>	<u>0628</u>	<u>X</u>
<u>2270</u>	<u>SVEC</u>	<u>X</u>
<u>8260</u>	<u>VOC</u>	<u>X</u>

**1 Matrix Codes:**  
 GW = Ground Water  
 WW = Waste Water  
 DW = Drinking Water  
 A = Air  
 S = Soil  
 SL = Sludge  
 SOL = Solid  
 O = Other (please define)

**2 Preservation Codes:**  
 I = Iced  
 H = HCL  
 M = Methanol  
 N = Nitric Acid  
 S = Sulfuric Acid  
 B = Sodium Bisulfate  
 X = Sodium Hydroxide  
 T = Sodium Thiosulfate  
 O = Other (please define)

**3 Container Codes:**  
 A = Amber Glass  
 G = Glass  
 P = Plastic  
 ST = Sterile  
 V = Vial  
 S = Summa Canister  
 T = Tedlar Bag  
 O = Other (please define)

Requested Turnaround Time	Due Date	1-Day	2-Day	3-Day	4-Day	Format	Other	CLP Like Data Pkg Required	Email To	Fax To #	Ending Date/Time	Composite	Grab	Matrix Code	Conc Code
<u>32</u>	<u>H</u>	<u>A</u>	<u>20VS</u>	<u>0628</u>	<u>X</u>	<u>2270</u>	<u>SVEC</u>	<u>X</u>	<u>8260</u>	<u>VOC</u>	<u>X</u>	<u>GW</u>			

Please use the following codes to indicate possible sample concentration within the Conc Code column above:  
 H - High; M - Medium; L - Low; C - Clean; U - Unknown

North Carolina Detection Limit Requirements

2L  
 GWPC  
 SWSL  
 IHSB  
 MSCC

Date/Time: 5/13/20 115  
 Date/Time: 5/19/20 1155  
 Date/Time: 5/19/20 6030  
 Date/Time: 5/26 1014

Project Entity:  
 Government  
 Federal  
 City  
 Municipality  
 Brownfield  
 School

Program Information:  
 DSCA  
 SWS Landfill  
 IHSB Orphaned Landfill  
 State Lead  
 Other:  
 US/Trust Fund  
 REC

PCB ONLY  
 Soxhlet  
 Non Soxhlet

NEIAC and AHA-LAP, LLC Accredited

Other:  
 Chromatogram  
 AHA-LAP, LLC

**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  
Wednesday 5/20/2020 at 10:14 am



**DELIVERED**

Signed for by: R.PIETRIAS

**GET STATUS UPDATES**  
**OBTAIN PROOF OF DELIVERY**

**FROM**  
Raleigh, NC US

**TO**  
EAST LONGMEADOW, MA US

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		5/20/2020	East Longmeadow, MA
149454201810	RALEIGH, NC	5/19/2020		5/20/2020	East Longmeadow, MA
149454201820	RALEIGH, NC	5/19/2020		5/20/2020	East Longmeadow, MA

Shipment Facts

**TRACKING NUMBER**  
149454201820

**SERVICE**  
FedEx Priority Overnight

**MASTER TRACKING NUMBER**  
149454201809

**WEIGHT**  
10 lbs / 4.54 kgs

**DIMENSIONS**  
25x14x14 in.

**DELIVERED TO**  
Receptionist/Front Desk

**TOTAL PIECES**  
3

**TOTAL SHIPMENT WEIGHT**  
10 lbs / 4.54 kgs

**TERMS**  
Third Party

**PACKAGING**  
Your Packaging

**SPECIAL HANDLING SECTION**  
Deliver Weekday, Non Standard Packaging

**STANDARD TRANSIT**  
 5/20/2020 by 10:30 am

I Have Not Confirmed Sample Container Numbers With Lab Staff Before Relinquishing Over Samples \_\_\_\_\_



**con-test**<sup>®</sup>  
ANALYTICAL LABORATORY

Doc# 277 Rev 5 2017

**Login Sample Receipt Checklist - (Rejection Criteria Listing - Using Acceptance Policy) Any False Statement will be brought to the attention of the Client - State True or False**

Client Dropper Ader  
 Received By DAF Date 5/26 Time 10:14

How were the samples received?  
 In Cooler T No Cooler \_\_\_\_\_ On Ice T No Ice \_\_\_\_\_  
 Direct from Sampling \_\_\_\_\_ Ambient \_\_\_\_\_ Melted Ice \_\_\_\_\_

Were samples within Temperature? 2-6°C T By Gun # 2 Actual Temp - 28  
 By Blank # \_\_\_\_\_ Actual Temp - \_\_\_\_\_

Was Custody Seal Intact? NA Were Samples Tampered with? NA  
 Was COC Relinquished? T Does Chain Agree With Samples? T

Are there broken/leaking/loose caps on any samples? F

Is COC in ink/ Legible? T Were samples received within holding time? T  
 Did COC include all pertinent Information? Client T Analysis T Sampler Name \_\_\_\_\_  
 Project T ID's T Collection Dates/Times T

Are Sample labels filled out and legible? T  
 Are there Lab to Filters? F Who was notified? \_\_\_\_\_  
 Are there Rushes? WBS F Who was notified? \_\_\_\_\_  
 Are there Short Holds? F Who was notified? \_\_\_\_\_  
 Is there enough Volume? T  
 Is there Headspace where applicable? F MS/MSD? F  
 Proper Media/Containers Used? T Is splitting samples required? f  
 Were trip blanks received? F On COC? f  
 Do all samples have the proper pH? NA Acid \_\_\_\_\_ Base \_\_\_\_\_

Vials	#	Containers:	#	#	#
Unp-		1 Liter Amb.	2	1 Liter Plastic	16 oz Amb.
HCL-	3	500 mL Amb.		500 mL Plastic	8oz Amb/Clear
Meoh-		250 mL Amb.		250 mL Plastic	4oz Amb/Clear
Bisulfate-		Flashpoint		Col./Bacteria	2oz Amb/Clear
DI-		Other Glass		Other Plastic	Encore
Thiosulfate-		SOC Kit		Plastic Bag	Frozen:
Sulfuric-		Perchlorate		Ziplock	

**Unused Media**

Vials	#	Containers:	#	#	#
Unp-		1 Liter Amb.		1 Liter Plastic	16 oz Amb.
HCL-		500 mL Amb.		500 mL Plastic	8oz Amb/Clear
Meoh-		250 mL Amb.		250 mL Plastic	4oz Amb/Clear
Bisulfate-		Col./Bacteria		Flashpoint	2oz Amb/Clear
DI-		Other Plastic		Other Glass	Encore
Thiosulfate-		SOC Kit		Plastic Bag	Frozen:
Sulfuric-		Perchlorate		Ziplock	

Comments: