

January 9, 2020

North Carolina Department of Transportation Geotechnical Unit Mail Service Center 1592 Raleigh, North Carolina 27699-1592

Attention: Mr. Craig Haden email: cehaden@ncdot.gov

Reference: Preliminary Site Assessment Report

NCDOT Project I-5878, WBS Element 53078.1.1

Parcel 83-Throw Back Thursday

608 Spring Branch Road

Dunn, Harnett County, North Carolina

S&ME Project 4305-19-161

Dear Mr. Haden:

S&ME, Inc. (S&ME) is submitting this Preliminary Site Assessment (PSA) Report to the North Carolina Department of Transportation (NCDOT). This report presents the background/project information, field activities, findings, conclusions, and recommendations. These services were performed in general accordance with S&ME Proposal No. 43-1900576 REV-01 dated August 9, 2019, and Contract Number 7000018853 dated April 12, 2018 between NCDOT and S&ME, Inc., authorized by NCDOT in its September 5,2019 Notice to Proceed Letter.

♦ Background/Project Information

Based on NCDOT's July 24, 2019, Request for Technical and Cost Proposal, the PSA was conducted within the NCDOT right-of-way (ROW) and/or easement as indicated on the preliminary plan sheets provided by NCDOT at the following property:

NCDOT Parcel No.	Property Owner	Site Address
83	Randy Emanuel Sr	(Throw Back Thursday)
		608 Spring Branch (aka Pope Road), Dunn, NC



The property is developed with a commercial building that is currently occupied by a bar named Throw Back Thursday. The building is reported to have previously been occupied by a gasoline/convenience store. According to a 1993 aerial photograph of the site, what appears to be a pump island canopy is shown in front of the building. Information regarding the location of USTs were not provided or observed from the aerial photograph. The property is not listed with registered petroleum underground storage tanks (USTs) (active or closed). The property is not listed with North Carolina Department of Environmental Quality (NCDEQ) Incidents associated with petroleum releases from USTs or aboveground storage tanks.

The PSA included a geophysical survey, subsequent limited soil sampling (five soil borings up to 10 feet below ground surface (ft.-bgs)) and limited groundwater sampling (one groundwater sample), within accessible areas of the proposed ROW/easement in preparation for construction activities. **Figure 1** shows the vicinity and site location, and **Figure 2** shows the site and boring locations. Soil and groundwater sampling results are shown on **Figure 3**.

Field Services

Prior to field activities, a site specific Health and Safety Plan was prepared as required by the Occupational Health and Safety Act (OSHA). Underground utilities were located and marked by the North Carolina One-Call Service. A private utility locator (East Coast Underground, LLC) was also used to locate and mark underground utilities.

Geophysical Survey

On July 25, 2019, S&ME completed Time Domain Electromagnetic (TDEM) and Ground Penetrating Radar (GPR) surveys within accessible areas of the proposed ROW/easement at Parcel 83. Brief descriptions of these complementary geophysical techniques are presented in the following paragraphs.

Time Domain Electromagnetics (TDEM)

TDEM measures the electrical conductivity of subsurface materials and discriminates between moderately conductive earth materials and very conductive metallic targets within the shallow subsurface. The conductivity is determined by transmitting a time-varying magnetic pulse into the subsurface and measuring the amplitude and phase shift of the secondary magnetic field. The secondary magnetic field is created when the conductive materials become an inductor as the primary magnetic field is passed through them. TDEM data are acquired continuously at a walking pace typically along a series of parallel or perpendicular lines. The system generates audible and visual indications when metallic targets are encountered. These measurements can also be supported with a global positioning system (GPS) which is output directly into the TDEM data file.

We used a Geonics Limited EM-61 MK2 TDEM system in general accordance with ASTM D6820 "Standard Guide for Use of the Time Domain Electromagnetic Method for Subsurface Investigation." Data was collected along lines spaced at approximately five feet using a Juniper® Systems GeodeTM sub-meter GPS as positioning support. The approximate TDEM data collection paths are presented in **Figure 4.** Golden Software's Surfer® program was used to grid and plot the data (**Figures 5 and 6**). The TDEM data has been presented as Plots A and B in order to provide both opaque and semi-transparent views, respectively.



Ground Penetrating Radar (GPR)

GPR transmits electromagnetic waves into the subsurface from an antenna at a specific frequency and measures the time for wave reflections to be received by interfaces between materials with differing material properties (e.g. soil/metal, etc.). The intensity of the reflected GPR wave is a function of the contrast in the material properties (i.e. dielectric permittivity) at the interface, the conductivity of the material that the wave is traveling through, and the frequency of the signal.

We used a Geophysical Survey Systems, Inc. (GSSI) SIR® 4000 GPR system equipped with a 350 MHz antenna in general accordance with ASTM D6432 "Standard Guide for Using the Surface Ground Penetrating Radar Method for Subsurface Investigation" to further characterize anomalies/features identified during the TDEM survey.

A total of seven GPR profiles (Lines 1 through 7) were collected for documentation (**Figure 7**). The data was post-processed using the GSSI Radan® 7 GPR software program for additional analysis.

Geophysical Findings

Responses indicative of a potential UST were not identified in the geophysical data sets collected at the site. Two anomalous features unrelated to known surficial targets were identified in the geophysical data sets (Anomalies A and B; **Figures 6 and 7**). Anomaly A is characterized by high amplitude GPR responses located within the upper one foot below ground surface (bgs) and may be related to an isolated buried metallic object (**Figure 8**). Anomaly B was characterized by a low TDEM response that was unable to be identified in the GPR data sets, which suggests the potential metallic target is either relatively small or deeper than the maximum penetration depth of the GPR signal. The anomalies were marked in the field using white spray paint. Example GPR profiles are presented in **Figures 8 and 9**.

Soil Sampling

On October 29, 2019, Troxler Geologic, Inc. (Troxler's) drill crew utilized a track mounted Geoprobe® rig to advance five soil borings (B-1 through B-5) and to collect soil samples within accessible areas of the proposed ROW/easement at Parcel 83. The approximate location of the soil borings are shown in **Figure 2**. A photographic log is included in **Appendix I.** Troxler's drill crew advanced the Geoprobe® borings up to a depth of approximately 10 ft.-bgs. During the advancement of the soil borings, groundwater was encountered at depths ranging from approximately six to 9.3 ft.-bgs. Soil samples were continuously collected in four-foot long disposable acetate-plastic sleeves that line the hollow stainless-steel sample probes. Soil recovered from the sleeves was classified on-site by S&ME personnel and screened with a Photoionization Detector (PID) at approximately two foot depth intervals to measure relative headspace concentrations of volatile organic compounds (VOCs).

VOC headspace readings were obtained from an aliquot of each soil sample that was placed in a re-sealable bag. Another portion of the sample was placed in a separate re-sealable bag and stored in an insulated container with ice for possible laboratory analyses. After waiting approximately 15 minutes to allow the sample to reach ambient temperature and headspace equilibrium, the PID probe was inserted into the bag to obtain a headspace reading. A summary of the PID readings and logs of the soil borings are included in **Appendix II.**



Petroleum odors and elevated PID readings were not noted in the collected soil samples. Therefore, a soil sample was selected from each boring at varying depth intervals. The soil samples were placed into laboratory supplied containers and transported to RED Lab, LLC (Red Lab) in an insulated cooler with ice for analysis. A total of five soil samples (one soil sample per boring) were analyzed by RED Lab for TPH-GRO and TPH-DRO using ultra-violet fluorescence (UVF) spectroscopy with product (fuel) identification.

Soil Analytical Results

Based upon analytical results of soil samples analyzed by RED Lab using UVP spectroscopy, TPH-GRO and TPH-DRO were not reported at concentrations exceeding their respective North Carolina TPH Action Levels. TPH-DRO was reported in borings B-1, B-3 and B-4 at the six to eight foot depth interval, at concentrations ranging from one milligram per kilograms (mg/kg) to 1.3 mg/kg, which is slightly above the laboratory method reporting limits but well below its North Carolina TPH Action Level of 100 mg/kg. TPH-GRO and TPH-DRO were not reported at concentrations exceeding the laboratory method reporting limits at the remaining soil samples. A summary of the soil analytical results is presented in **Table 1** and shown on **Figure 3**. A copy of the laboratory analytical report provided by RED Lab is presented in **Appendix III**.

Groundwater Sampling

During the advancement of the soil borings, groundwater was encountered at depths ranging from approximately six to 9.3 ft.-bgs. Therefore, the Geoprobe® was used to advance one of the soil borings into the groundwater table for the collection of a groundwater sample. Due to the lack of petroleum odors or elevated PID readings, boring B-2 was selected for collection of a groundwater sample, due to its reported location at the former pump island. A temporary monitor well (TW-1) was installed at boring B-2 to a depth of approximately 10.5 ft.-bgs using a five foot section of one-inch diameter, Schedule 40 PVC well riser attached to a ten foot section of 0.01-inch slotted screen that intersected the groundwater table. Groundwater within the temporary monitor well at boring B-2 was measured at six ft.-bgs. Groundwater was purged from the temporary well until relatively clear using disposable tubing attached to a peristaltic pump. The flow rate was reduced and laboratory supplied containers were filled directly from the tubing, labeled as TW-1 and placed in an insulated cooler with ice for transport to Con-Test Laboratories (Con-Test) for analysis of VOCs by EPA Method 8260 and polycyclic aromatic compounds (PAHs) by EPA Method 8270.

Upon completion of the soil and groundwater sampling, the well materials were removed and the soil borings backfilled with bentonite pellets and soil cuttings. Investigative derived wastes (IDW), such as soil cuttings generated during the soil boring advancement and decontamination water, were spread on the ground in accordance with the procedures specified by NCDEQ. Used gloves and tubing were bagged and disposed offsite.

Groundwater Analytical Results

Based upon analytical results of the groundwater sample analyzed by Con-Test, several petroleum related target constituents were reported at concentrations above the laboratory method reporting limits but below their 15A NCAC 2L Groundwater Quality Standards (2L Standards). A summary of the groundwater analytical results is presented in **Table 2** and shown on **Figure 3**. A copy of the laboratory analytical report provided by Con-Test is presented in **Appendix III**.



Conclusion and Recommendations

The geophysical survey identified two anomalies (Anomalies A and B) which may be related to isolated buried metallic objects. Responses indicative of a potential UST were not identified in the geophysical data sets collected at the site.

S&ME advanced five soil borings (B-1 through B-5) to a depth of up to approximately 10 ft.-bgs at the site. Petroleum odors and elevated PID readings were not noted in soil samples collected at the borings. Selected soil samples from the soil borings were analyzed for TPH-GRO and TPH-DRO using UVF spectroscopy.

TPH-GRO and TPH-DRO were not reported at concentrations exceeding their respective North Carolina TPH Action Levels. TPH-DRO was reported in borings B-1, B-3 and B-4 at the six to eight foot depth interval at concentrations slightly above the laboratory method reporting limits but well below its North Carolina TPH Action Level. TPH-GRO and TPH-DRO were not reported at concentrations exceeding the laboratory method reporting limits at the remaining soil samples.

During the soil boring advancement, groundwater was encountered at depths ranging from six to 9.3 ft.-bgs. One temporary well (TW-1) was installed at soil boring B-2 located at the former pump island. Groundwater at TW-1 was measured at six ft.-bgs and analyzed by Con-Test for VOCs by EPA Method 8260 and PAHs by EPA Method 8270. Several petroleum related target constituents were reported at concentrations above the laboratory reporting limits but below their respective 2L Standards.

Based on the findings of the geophysical survey and analytical results of soil and groundwater samples, it is likely that during construction, NCDOT may encounter marginally impacted soil (below TPH Action Levels) and groundwater marginally impacted with petroleum (below 2L Standards) at the site.

It should also be assumed that saturated petroleum impacted soil will be encountered if construction excavations extend deeper than six ft.-bgs on the site. If construction dewatering is required, petroleum impacted groundwater must be properly disposed or treated at a licensed facility. If petroleum stained or odorous soils are encountered during construction, these soils should be properly handled and disposed at a licensed facility.

S&ME recommends maintaining an awareness level for the presence of marginally impacted petroleum in the soil (below TPH Action Levels) and marginally impacted groundwater (below 2L Standards) at the site for the safety of workers and the public.

Limitations

The results of this preliminary investigation are limited to the boring locations presented herein. The results of this Preliminary Site Assessment are not all inclusive and may not represent existing conditions across the entire property. These results only reflect the current conditions at the locations sampled on the date this Preliminary Site Assessment was performed. This report has been prepared in accordance with generally accepted environmental engineering and geophysical practice for specific application to this project. The conclusions and recommendations contained in this report are based upon applicable standards of our practice in this geographic area at the time this report was prepared. No other warranty, expressed or implied, is made.

January 9, 2020



The geophysical methods used for this survey have inherent limitations. Site metallic features (e.g., surficial debris, reinforced concrete, utilities, etc.) and overhead transmission lines can produce a false electromagnetic response and may mask subsurface features. The depth of exploration of the GPR signal is highly site specific and is greatly limited by signal attenuation (absorption) of the subsurface materials. Signal attenuation is dependent upon the electrical conductivity of the subsurface materials. Signal attenuation is greatest in materials with relatively high electrical conductivities such as clay soils, and lowest in relatively low conductivity materials such as unsaturated sand. For this project location, the GPR data sets appear to have a maximum depth of penetration of about five ft.-bgs.

Regardless of the thoroughness of a geophysical study, there is always a possibility that actual conditions may not match the interpretations. The results should be considered accurate only to the degree implied by the methods used and the method's limitations and data coverage. Accordingly, the possibility exists that not all features at a project site will be located due to either subsurface soil conditions or the occurrence of features outside the lateral limits and below the depth of penetration of the methods used. As with most surface geophysical methods, resolution of the subsurface will also decrease with depth. As such, the size and/or contrast of features compared to the imaged subsurface media must be significant enough to produce the anticipated response. The location and/or determination (or the lack thereof) of potential buried features is based on our review of the provided information and of the geophysical survey. Under no circumstances does S&ME assume any responsibility for damages resulting from the presence of subsurface features that may exist but were not identified by our survey.

This Preliminary Site Assessment was performed solely for NCDOT regarding the above-referenced site and assessment area. This report is provided for the sole use of NCDOT. Use of this report by any other parties will be at such party's sole risk. S&ME disclaims liability for any such use or reliance by third parties. The observations presented in this report are indicative of conditions during the time of the assessment and of the specific areas referenced.



Closing

S&ME appreciates the opportunity to provide these services to you. If you have any questions or comments regarding this report, please contact us at your convenience.

Sincerely,

S&ME, Inc.

DocuSigned by:

Jamie Honeratt

-4C890EAEC25F488...

Jamie T Honeycutt Environmental Professional jhoneycutt@smeinc.com



Thomas P. Raymond, P.E., P.M.P. Senior Consultant traymond@smeinc.com

Attachments:

Table 1: Summary of Soil Sampling Results

Table 2: Summary of Groundwater Sampling Results

Figure 1: Vicinity Map Figure 2: Site Map

Figure 3: Soil and Groundwater Constituent Map

Figure 4: TDEM Path Location Plan

Figure 5: TDEM Data Plot A **Figure 6:** TDEM Data Plot B

Figure 7: Geophysical Anomaly Location Plan **Figure 8:** Example GPR Data – Lines 2 and 3 **Figure 9:** Example GPR Data – Lines 4 and 7

Appendix I: Photographs **Appendix II:** Boring Logs

Appendix III: Laboratory Analytical Reports and Chain of Custody

Docusigned by:
Midual Pfrifer
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Michael W. Pfeifer Senior Project Manager mpfeifer@smeinc.com

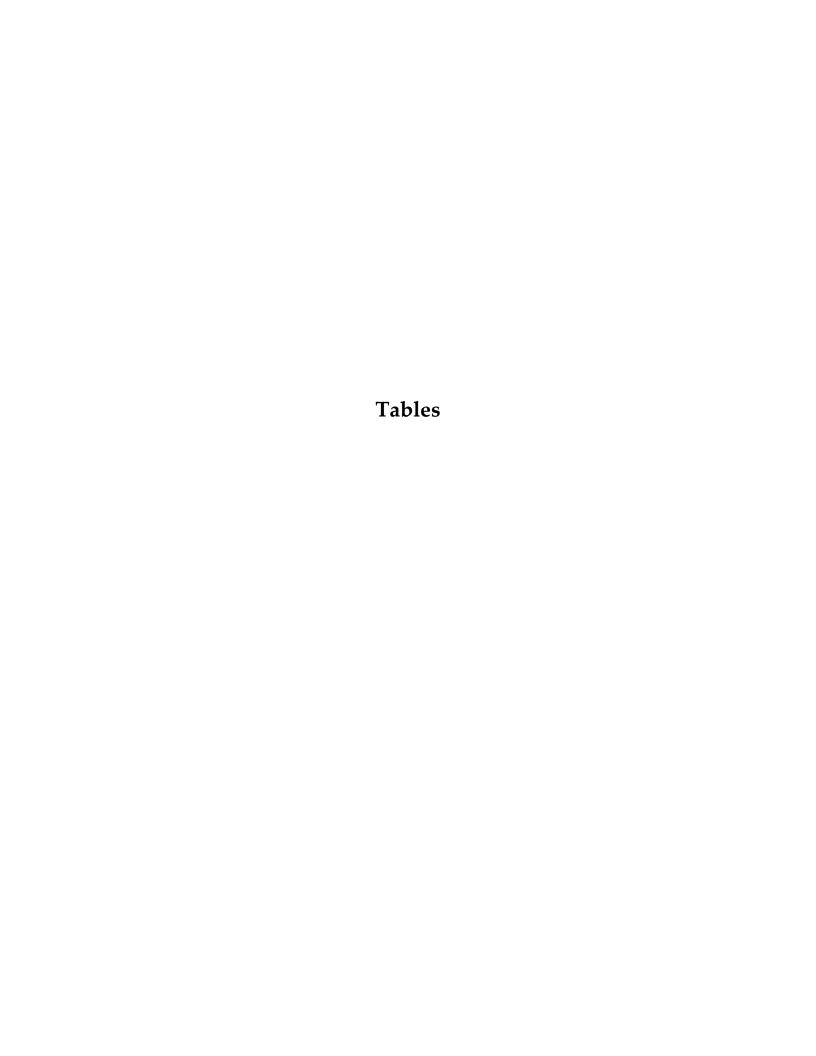


TABLE 1 SUMMARY OF SOIL SAMPLING RESULTS NCDOT Project I-5878



Parcel 83 - (Throw Back Thursday) 608 Spring Branch Road Dunn, Harnett County, North Carolina S&ME Project No. 4305-19-161

Ar	nalytical Metho	d→	Total Petroleum Hydrocarbons (TPH) Gasoline Range Organics (GRO) and Diesel Range Organics (DRO) by Ultraviolet Fluorescence (UVF) Spectrometry						
Sample ID	Date	Contaminant of Concern→	TPH-GRO	TPH-DRO					
		(ftbgs)							
B-1	10/29/2019	6 to 8	<0.45	1.1					
B-2	10/29/2019	2 to 4	<0.29	<0.29					
B-3	10/29/2019	6 to 8	<0.5	1					
B-4	10/29/2019	6 to 8	<0.51	1.3					
B-5	10/29/2019	6 to 8	<0.52	<0.52					
No	orth Carolina T	PH Action Levels	50	100					

Notes:

- 1. UVF analysis performed by RED Lab, LLC
- 2. Concentrations are reported in milligrams per kilogram (mg/Kg).
- 3. ft.-bgs:- feet below ground surface.
- 4. Concentrations exceeding the laboratory's reporting limits are shown in **BOLD** fields.
- 5. Concentrations exceeding the North Carolina TPH Action Levels are shown in Shaded and **BOLD** fields.



TABLE 2 SUMMARY OF GROUNDWATER SAMPLING RESULTS

NCDOT Project I-5878

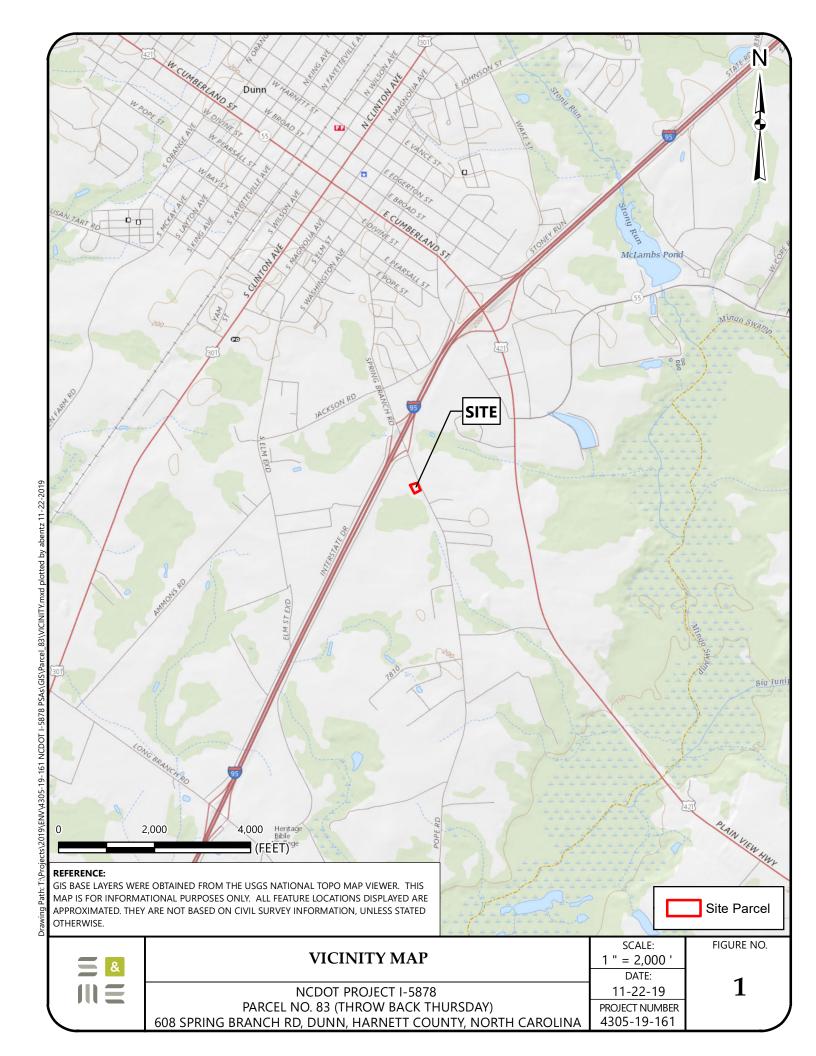
Parcel 83 - (Throw Back Thursday) 608 Spring Branch Road Dunn, Harnett County, North Carolina S&ME Project No. 4305-19-161

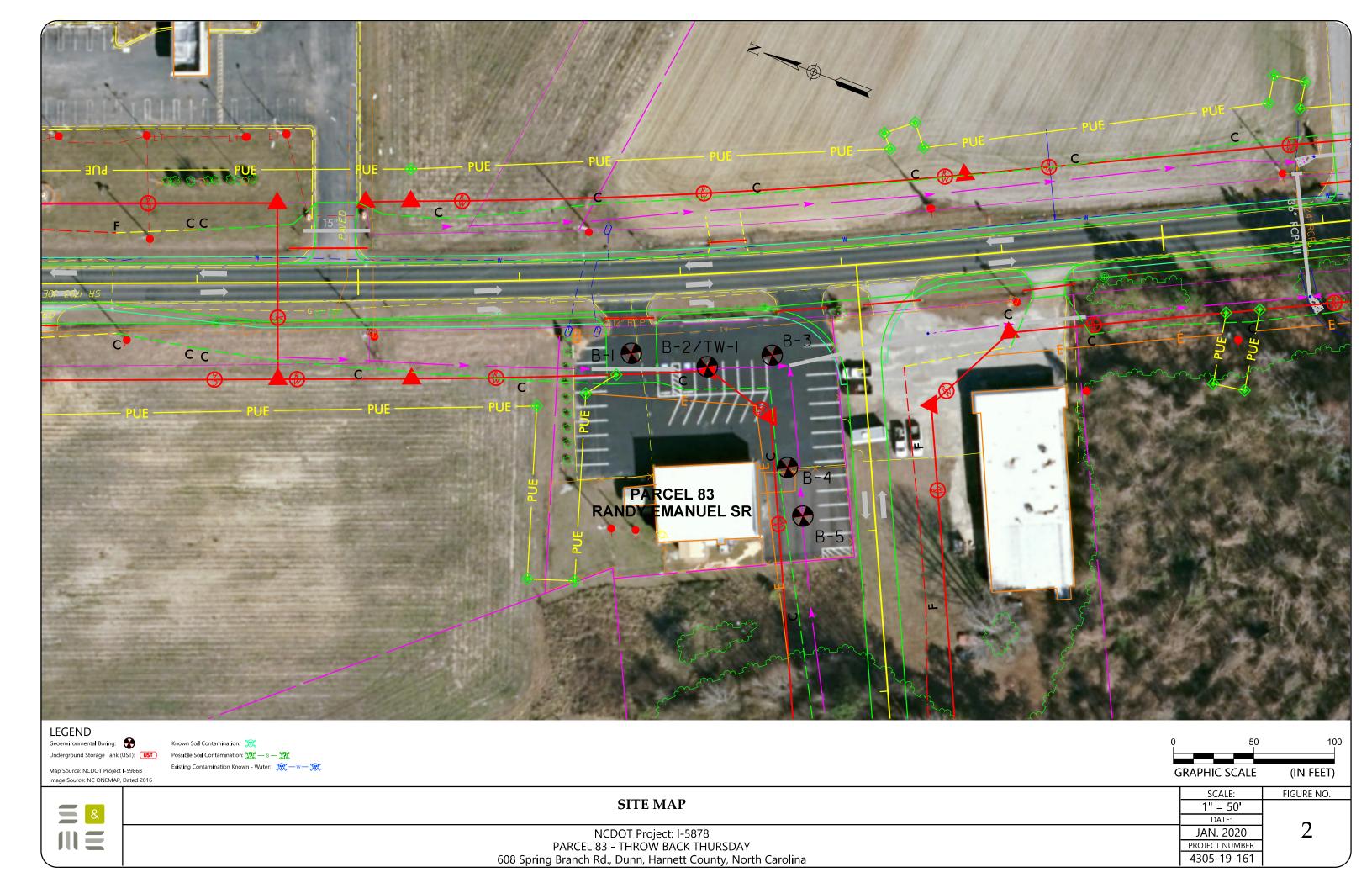
Analytical	Method→		Volatile Organic Compounds by EPA Method 8260 Polycycl									clic Aromatic Compounds (PAHs) by EPA Method 8270											
Sample ID	Contaminant of Concern→	Benzene	Ethylbenzene	Isopropylbenzene	Naphthalene	n-Butylbenzene	sec-Butylbenzene	Styrene	tert-Butyl Alcohol	n-Propylbenzene	p-Isopropyltoluene	Toluene	1,2,4- Trimethylbenzene	1,3,5- Trimethylbenzene	Total Xylenes	Acenaphthene	Acenaphthylene	Fluoranthene	Fluorene	Naphthalene	Phenanthrene	Pyrene	2-Methylnaphthalene
B-2/TW-1	10/29/2019	0.66 J	4.9	43	3.9	3.5	15	0.28 J	0.24 J	1.8	0.92 J	0.25 J	4.4	2.2	1.9	0.30	0.095 J	0.027 J	0.38 J	3.2	0.42	0.025J	4.9
2L S	Standard (µg/L)	1	600	70	6	70	70	70	10	70	25	600	400	400	500	80	200	300	300	6	200	200	30
	GCL (µg/L)	5,000	84,500	25,000	6,000	6,900	8,500	70,000	10,000	30,000	11,700	260,000	28,500	25,000	85,500	2,120	1,965	300	990	6,000	410	200	12,500

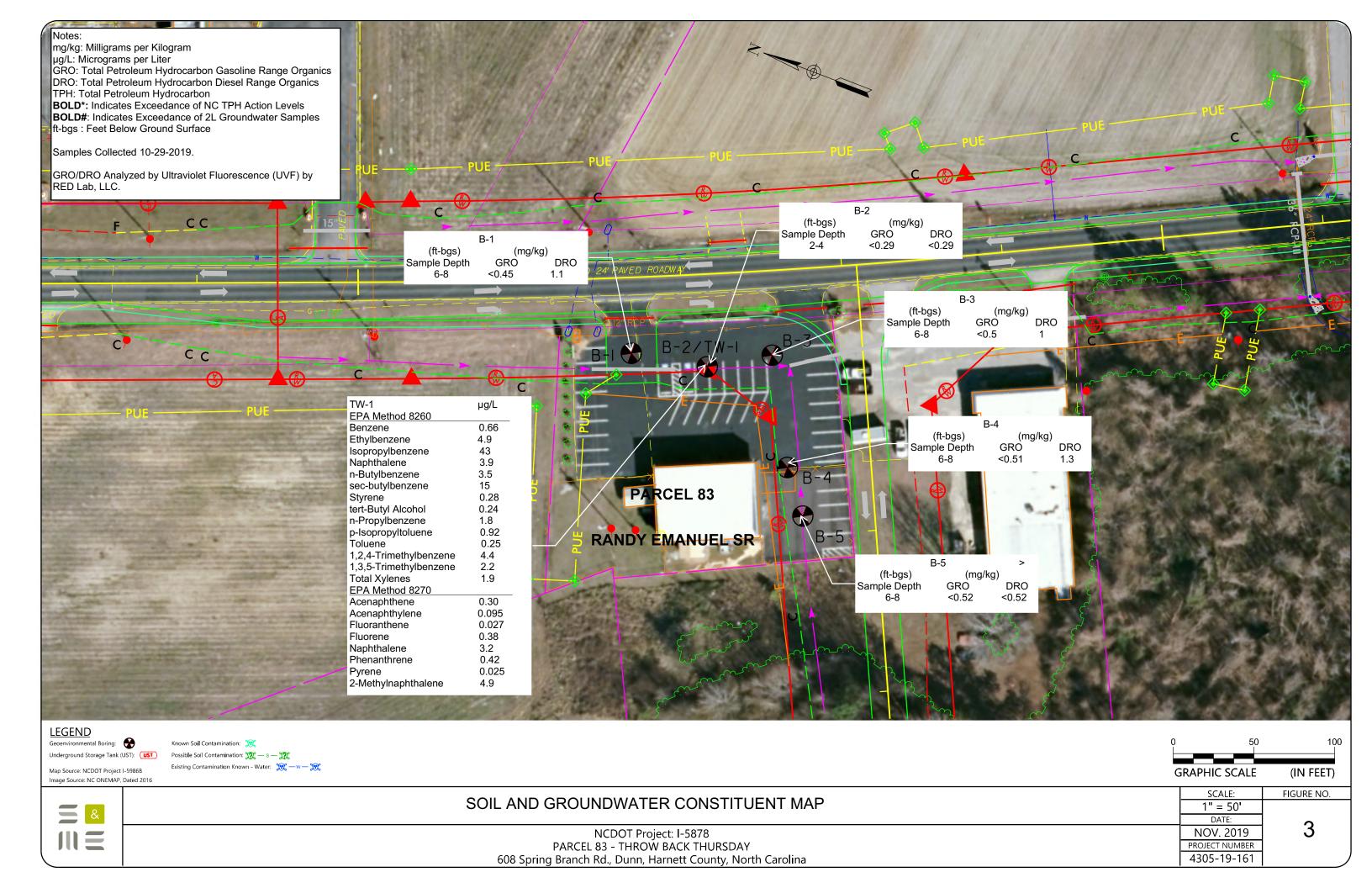
Notes:

- 1. Analytes that are not shown for the method were not detected.
- 2. Concentrations are reported in micrograms per liter (µg/L).
- 3. 2L Standard: North Carolina Groundwater Quality Standards: 15A NCAC 2L.0202
- 4. Concentrations exceeding the laboratory's reporting limits are shown in **BOLD** fields.
- 5. Concentrations exceeding the 2L Standards are shown in Shaded and BOLD fields.
- 6. GCL: Gross Contamination Level.
- 7. J: Estimated concentration detected below the reporting limit.









REFERENCE:

GOOGLE EARTH PRO AERIAL PHOTOGRAPH (DATED, MARCH 4, 2018)





LEGEND

---- Approximate TDEM Path

Approximate Requested Survey Area

SCALE: AS SHOWN

NCDOT PROJECT: 1-5878 PARCEL #83 - (THROW BACK THURSDAY) 608 SPRING BRANCH ROAD, DUNN, HARNETT COUNTY, NORTH CAROLINA

TDEM PATH LOCATION PLAN

DATE: 11/26/2019

PROJECT NUMBER 4305-19-161

FIGURE NO.

4





GOOGLE EARTH PRO AERIAL PHOTOGRAPH (DATED, MARCH 4, 2018)





TDEM DATA PLOT A

NCDOT PROJECT: 1-5878 PARCEL #83 - (THROW BACK THURSDAY) 608 SPRING BRANCH ROAD, DUNN, HARNETT COUNTY, NORTH CAROLINA

SCALE: AS SHOWN

DATE: 11/26/2019

PROJECT NUMBER 4305-19-161

FIGURE NO.

5



REFERENCE:

GOOGLE EARTH PRO AERIAL PHOTOGRAPH (DATED, MARCH 4, 2018)





LEGEND

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Approximate Location of Geophysical Anomaly

SCALE: AS SHOWN

NCDOT PROJECT: 1-5878 PARCEL #83 - (THROW BACK THURSDAY) 608 SPRING BRANCH ROAD, DUNN, HARNETT COUNTY, NORTH CAROLINA

TDEM DATA PLOT B

DATE: 11/26/2019

PROJECT NUMBER 4305-19-161

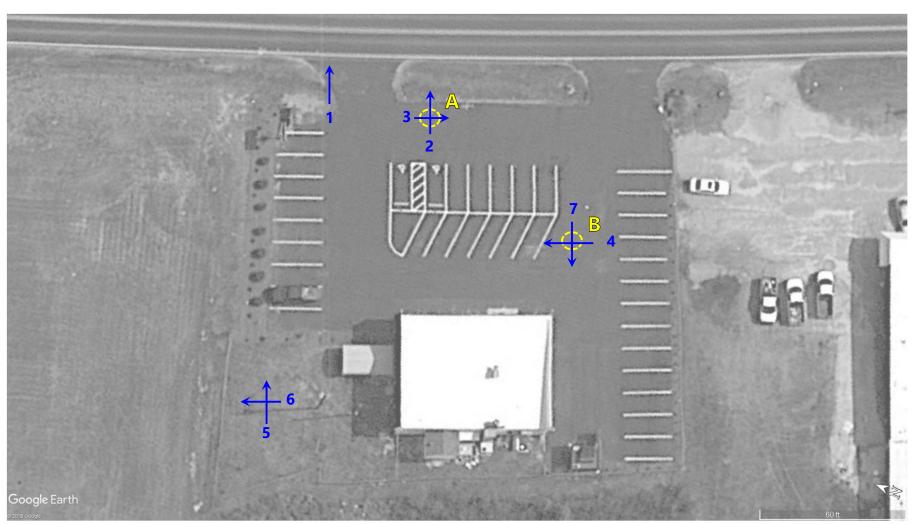
FIGURE NO.



REFERENCE:

GOOGLE EARTH PRO AERIAL PHOTOGRAPH (DATED, MARCH 4, 2018)





LEGEND

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Approximate Location of Geophysical Anomaly



Approximate Location of GPR Profile

GEOPHYSICAL ANOMALY LOCATION PLAN

NCDOT PROJECT: I-5878 PARCEL #83 - (THROW BACK THURSDAY) 608 SPRING BRANCH ROAD, DUNN, HARNETT COUNTY, NORTH CAROLINA

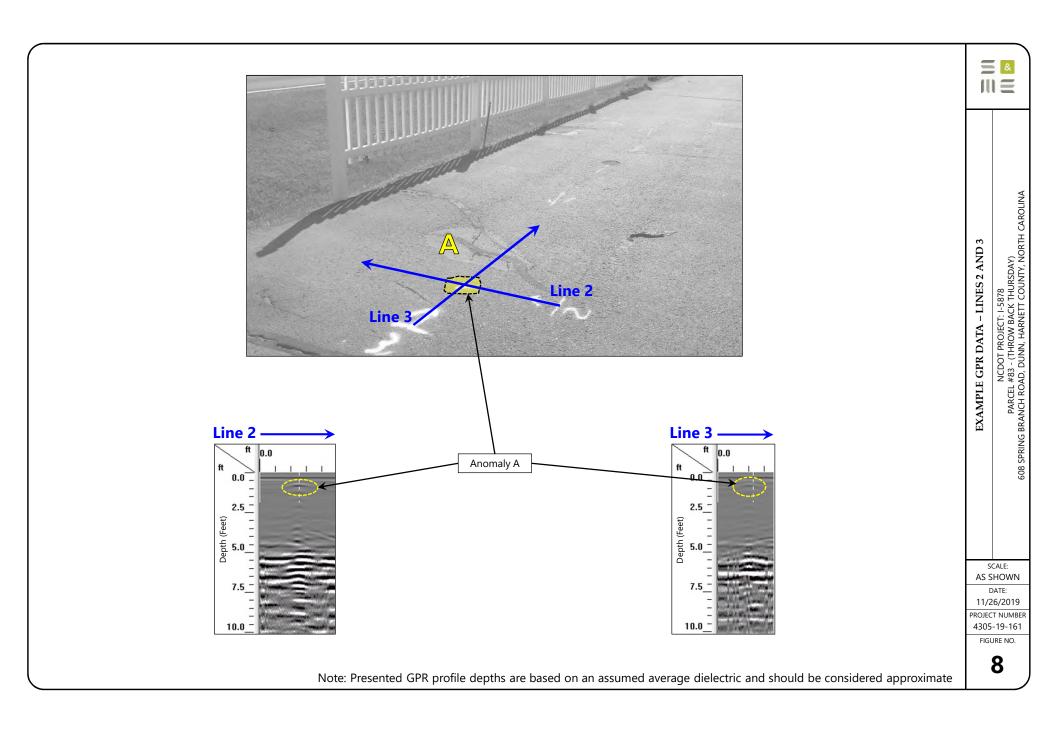
SCALE: AS SHOWN

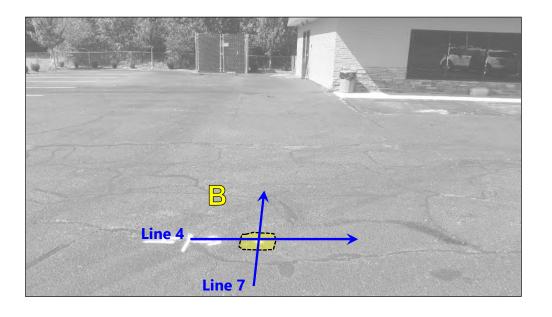
DATE: 11/26/2019

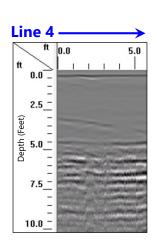
PROJECT NUMBER 4305-19-161

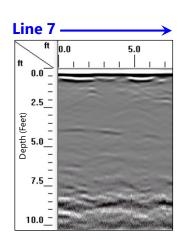
FIGURE NO.

7





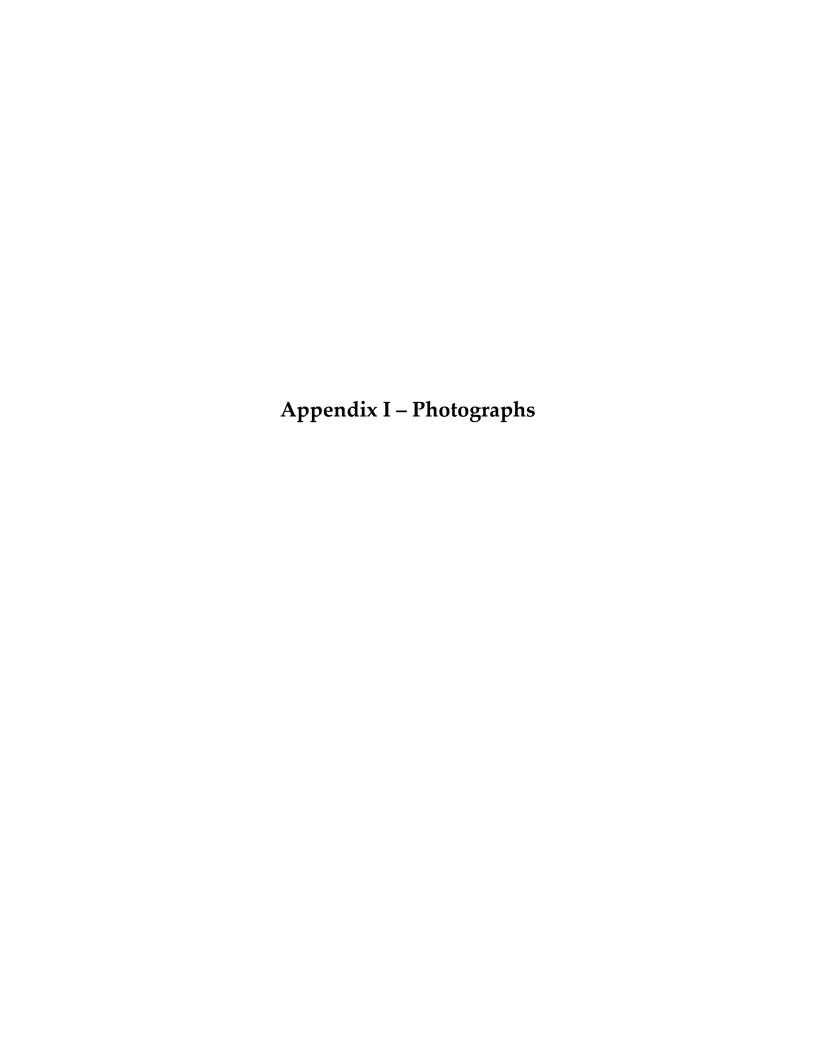




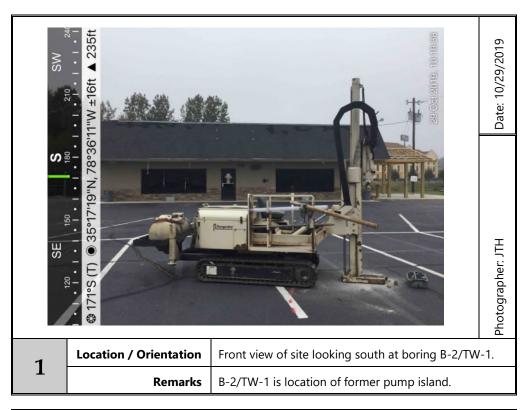
NCDOT PROJECT: I-5878 PARCEL #83 - (THROW BACK THURSDAY) 608 SPRING BRANCH ROAD, DUNN, HARNETT COUNTY, NORTH CAROLINA EXAMPLE GPR DATA – LINES 4 AND 7 SCALE: AS SHOWN DATE: 11/26/2019 PROJECT NUMBER 4305-19-161 FIGURE NO.

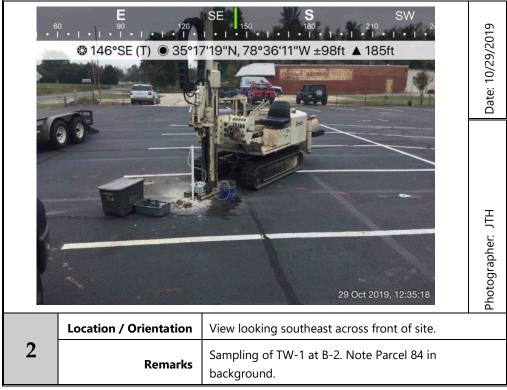
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Note: Presented GPR profile depths are based on an assumed average dielectric and should be considered approximate











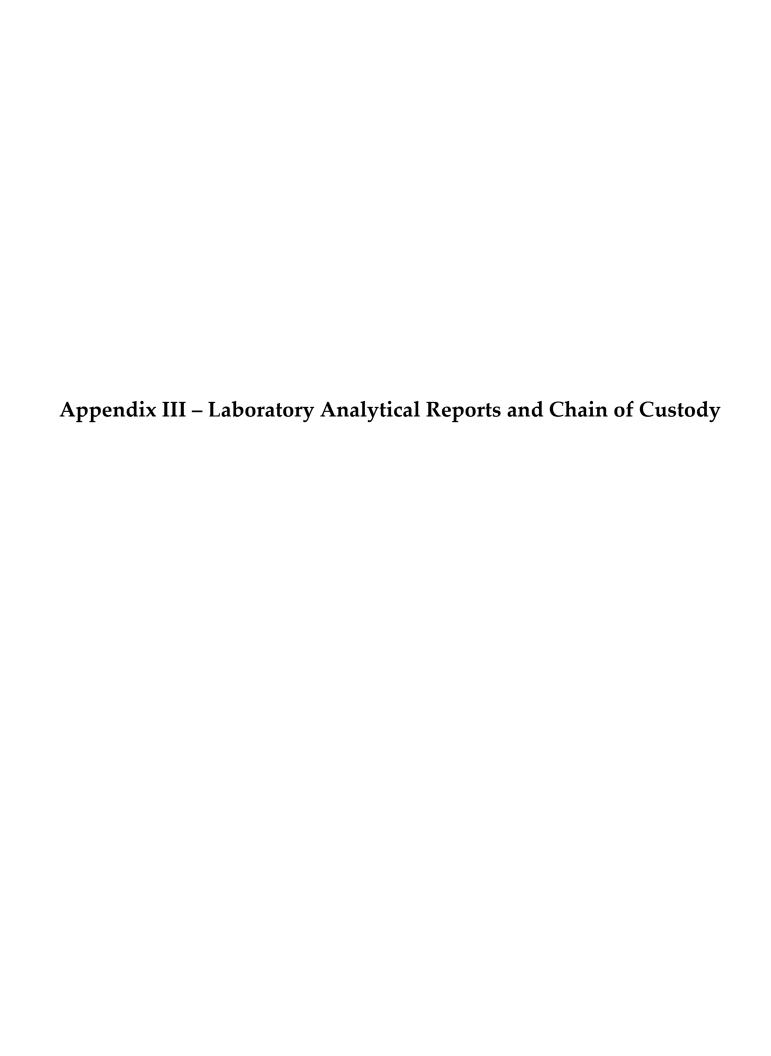
PROJECT:	NCDOT I-5878 Parcel 83-608 Spring Branch Rd (Throw Back Thurs	sday) Dunn NC			PODIA	NG LOG:	D 1			
	S&ME Project No. 4305-19-161	suay), Dunn, NC			DUKII	NG LOG	D-1			
DATE DRILLED:	Tuesday, October 29, 2019	BORING DEPTH (FT):	10							
DRILL RIG:	Geoprobe 54DT	WATER LEVEL:								
DRILLER:	Troxler Geologic, Inc.	CAVE-IN DEPTH:	Not App	licable						
HAMMER TYPE:	Not Applicable	LOGGED BY:								
SAMPLING METHOD:	Macro-Core Sampler	NORTHING:								
DRILLING METHOD:	Macro-Core Sampler (3-in. OD)	EASTING:								
DIVIDENTO INICITIOD.	macro core sumpler (5 m. 65)	LASTING.		1			1			1
DEPTH (feet) GRAPHIC LOG	MATERIAL DESCRIPTION		WATER LEVEL	SAMPLE	PID READING (PPM)	LABORATORY ANALYSES	Sample Time / 1st 6in	2nd 6in	3rd 6in	N VALUE
Si Si	sphalt, Gravel, andy Clay, Tan, Orange, Red,				2.3	No No				
					1.3 4.2	No Yes	1000			
	layey Sand, Tan, White,			H	4.2	163	1000			
10	layey Sand, Tan, Writte,			816	3.5	No				
15 — 15 — 20 — 25 — — — — — — — — — — — — — — — — —										
30										

PROJECT:	NCDOT I-5878 Parcel 83-608 Spring Branch Rd (Throw Back Thurso			BORIN	NG LOG:	B-2/	TW-1			
0.475.0001.50	S&ME Project No. 4305-19-161		4.2							
DATE DRILLED:	Tuesday, October 29, 2019	BORING DEPTH (FT):								
DRILL RIG:	Geoprobe 54DT	WATER LEVEL:								
DRILLER:	Troxler Geologic, Inc.	CAVE-IN DEPTH:								
HAMMER TYPE:	Not Applicable		J. Honeyc	utt						
Sampling method:	Macro-Core Sampler	NORTHING:								
DRILLING METHOD:	Macro-Core Sampler (3-in. OD)	EASTING:								
DEPTH (feet) GRAPHIC LOG	MATERIAL DESCRIPTION		WATER LEVEL	SAMPLE	PID READING (PPM)	LABORATORY ANALYSES	Sample Time / 1st 6in	2nd 6in	3rd 6in	N VALUE
Conc	rete, Gravel,									
Sand	y Clay, Tan, Orange, Red,			ŧ	2.9	No Yes	1020			
5 —				ł	4.2	No	1020			
			•	ł	5.2	No				
Clave	ey Sand, Tan, White,			ж						
L ₁₀ ////	Sand, Red,			ш	11.5	No				
	ng Terminated at 13 Ft-BGS									
15 —										
 _										
20 —										
25 —										
30										

PROJECT:	NCDOT I-5878 Parcel 83-608 Spring Branch Rd (Throw Back Thursd	lay), Dunn, NC	BORING LOG: B-3												
	S&ME Project No. 4305-19-161	- , -													
DATE DRILLED:	Tuesday, October 29, 2019	BORING DEPTH (FT):	10												
DRILL RIG:	Geoprobe 54DT	WATER LEVEL:													
DRILLER:	Troxler Geologic, Inc.	CAVE-IN DEPTH:	Not Appl	icable											
HAMMER TYPE:	Not Applicable	LOGGED BY:	J. Honey	cutt											
SAMPLING METHOD:	Macro-Core Sampler	NORTHING:													
DRILLING METHOD:	Macro-Core Sampler (3-in. OD)	EASTING:													
DEPTH (feet) GRAPHIC LOG	MATERIAL DESCRIPTION		WATER LEVEL	SAMPLE	PID READING (PPM)	LABORATORY ANALYSES	Sample Time / 1st 6in	2nd 6in	3rd 6in	N VALUE					
San	dy Clay, Tan, Orange, Red,				1.5 4.5	No No									
5 - Cla	yey Sand, Tan, Orange, Red,				4.9	No									
				I	4.7	Yes	1020								
10 Bor	ing Terminated at 10 Ft-BGS		•	Н	4.1	No									
15 —															
20 —															
25 —															
30															

PROJECT:	:	NCDOT I-5878 Parcel 83-608 Spring Branch Rd (Throw Back Th	BORING LOG: B-4												
DATE DRUG		S&ME Project No. 4305-19-16		10											
DATE DRILL	.ED:	Tuesday, October 29, 2019	BORING DEPTH (FT):												
DRILL RIG:		Geoprobe 54DT	WATER LEVEL:												
DRILLER:		Troxler Geologic, Inc.	CAVE-IN DEPTH:												
HAMMER T		Not Applicable		J. Honey	cutt										
SAMPLING		Macro-Core Sampler	NORTHING:												
DRILLING M	ИЕТНОD:	Macro-Core Sampler (3-in. OD)	EASTING:												
DEPTH (feet)	GRAPHIC LOG	MATERIAL DESCRIPTION		WATER LEVEL	SAMPLE	PID READING (PPM)	LABORATORY ANALYSES	Sample Time / 1st 6in	2nd 6in	3rd 6in	N VALUE				
5 — 10 — 15 — —		Asphalt, Gravel, Sandy Clay, Tan, Orange, Red, Boring Terminated at 10 Ft-BGS			SA	2.8 4.3 4.1 5.2 3.2	No No Yes No	1130		υ <u>ε</u>	N				
20 —															
<u> </u>															
25 —															
_															
30 —															
55															

PROJECT:	NCDOT I-5878 Parcel 83-608 Spring Branch Rd (Throw Back Thurs	dav). Dunn. NC			BORIN	NG LOG:	B-5			
	S&ME Project No. 4305-19-161	aay,, 2 a, 110			50 1411					
DATE DRILLED:	Tuesday, October 29, 2019	BORING DEPTH (FT):	10							
DRILL RIG:	Geoprobe 54DT	WATER LEVEL:								
DRILLER:	Troxler Geologic, Inc.	CAVE-IN DEPTH:		icable						
HAMMER TYPE:	Not Applicable									
SAMPLING METHOD:	Macro-Core Sampler	NORTHING:								
DRILLING METHOD:	Macro-Core Sampler (3-in. OD)	EASTING:								
		ı	Æ		9	RY S	/ ei			
DEPTH (feet) GRAPHIC LOG	MATERIAL DESCRIPTION		WATER LEVEL	SAMPLE	PID READING (PPM)	LABORATORY ANALYSES	Sample Time / 1st 6in	2nd 6in	3rd 6in	N VALUE
- Aspt	nalt, Gravel, dy Clay, Tan, Orange, Red,			ł	1.7	No				
5 —				ŧ	1.9	No				
				ł	1.8	No				
			•	ŧ	3.0 4.2	Yes No	1150			
10 — Borii	ng Terminated at 10 Ft-BGS				4.6	140				
15 —										
20 —										
25 —										
30										









Hydrocarbon Analysis Results

Client: S&ME

Address: 3201 SPRING FOREST ROAD

Samples taken
Samples extracted

Tuesday, October 29, 2019 Tuesday, October 29, 2019

Samples analysed Friday, November 1, 2019

Contact: JAMIE HONEYCUTT Operator MAX MOYER

Project: NCDOT I-5878 PARCEL 83

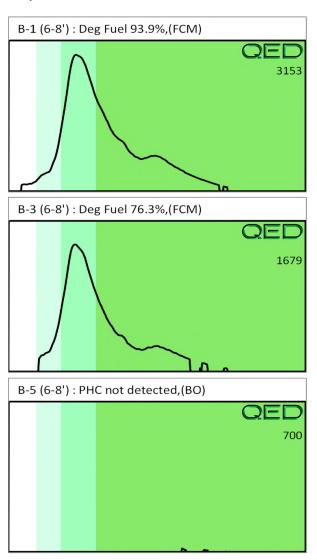
													F03640								
Matrix	Sample ID	Dilution used	BTEX (C6 - C9)	GRO (C5 - C10)	DRO (C10 - C35)	TPH (C5 - C35)	Total Aromatics (C10-C35)	16 EPA PAHs	ВаР	Ratio		Ratios		Ratios		Ratios		Ratios			HC Fingerprint Match
										% light % mid % heavy											
S	B-1 (6-8')	18.2	<0.45	<0.45	1.1	1.1	0.75	<0.15	<0.018	0	84.9	15.1	Deg Fuel 93.9%,(FCM)								
S	B-2 (2-4')	11.8	<0.29	<0.29	< 0.29	< 0.29	< 0.06	< 0.09	<0.012	0	100	0	,(FCM)								
S	B-3 (6-8')	20.2	<0.5	<0.5	1	1	0.44	<0.16	< 0.02	0	89	11	Deg Fuel 76.3%,(FCM)								
S	B-4 (6-8')	20.5	<0.51	<0.51	1.3	1.3	0.82	<0.16	< 0.02	0	81.3	18.7	Deg Fuel 76.7%,(FCM)								
S	B-5 (6-8')	20.6	< 0.52	< 0.52	< 0.52	< 0.52	<0.1	<0.17	<0.021	0	0	0	PHC not detected,(BO)								
	Initial C	QC check	OK					Final F	CM QC	Check	OK	99.5 %									

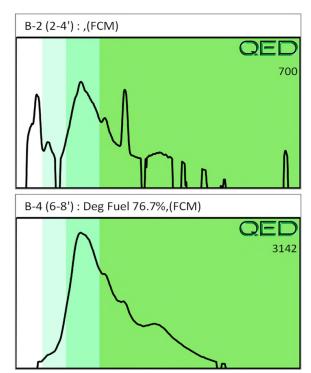
Results generated by a QED HC-1 analyser. Concentration values in mg/kg for soil samples and mg/L for water samples. Soil values are not corrected for moisture or stone content

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

Project: NCDOT I-5878 PARCEL 83





RAPID ENVIRONMENTAL DIAGNOSTICS CHAIN OF CUSTODY AND ANALYTICAL ∞ **REQUEST FORM** NCOOT I-5878 Money Is a smeller con Forcet 125 COST I-5878 DENN Jamie T Harar D Jamie Client Name: Collected by: Project Ref .: Address: Phone #: Contact: Email:

5598 Marvin K Moss Lane RED Lab, LLC

MARBIONC Bldg, Suite 2003 Wilmington, NC 28409

B143

Each sample will be analyzed for

BTEX, GRO, DRO, TPH, PAH total aromatics and BaP

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

November 7, 2019

Jamie Honeycutt S&ME, Inc - Raleigh, NC 3201 Spring Forest Rd. Raleigh, NC 27616

Project Location: Dunn, NC

Client Job Number:

Project Number: 4305-19-161

Laboratory Work Order Number: 19K0021

Keny K. Mille

Enclosed are results of analyses for samples received by the laboratory on October 31, 2019. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Kerry K. McGee Project Manager

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

S&ME, Inc - Raleigh, NC 3201 Spring Forest Rd. Raleigh, NC 27616 ATTN: Jamie Honeycutt

REPORT DATE: 11/7/2019

PURCHASE ORDER NUMBER:

PROJECT NUMBER: 4305-19-161

ANALYTICAL SUMMARY

WORK ORDER NUMBER: 19K0021

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

PROJECT LOCATION: Dunn, NC

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

TW-1 19K0021-01 Ground Water SW-846 8260D

SW-846 8270E



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

EXECUTIVE SUMMARY

Client ID: TW-1 Lab ID: 19K0021-01

Analyte	Results/Qua	l	DL	RL	Units	Method
1,2,4-Trimethylbenzene	4.4		0.18	1.0	$\mu g/L$	SW-846 8260D
1,3,5-Trimethylbenzene	2.2		0.14	1.0	$\mu g/L$	SW-846 8260D
Benzene	0.66	J	0.18	1.0	$\mu g/L$	SW-846 8260D
Ethylbenzene	4.9		0.13	1.0	$\mu g/L$	SW-846 8260D
Isopropylbenzene (Cumene)	43		0.17	1.0	$\mu g/L$	SW-846 8260D
m+p Xylene	1.5	J	0.30	2.0	$\mu g/L$	SW-846 8260D
Naphthalene	3.9		0.31	2.0	$\mu g/L$	SW-846 8260D
n-Butylbenzene	3.5		0.21	1.0	$\mu g/L$	SW-846 8260D
n-Propylbenzene	1.8		0.13	1.0	$\mu g/L$	SW-846 8260D
o-Xylene	0.40	J	0.17	1.0	$\mu g/L$	SW-846 8260D
p-Isopropyltoluene (p-Cymene)	0.92	J	0.20	1.0	$\mu g/L$	SW-846 8260D
sec-Butylbenzene	15		0.16	1.0	$\mu g/L$	SW-846 8260D
Styrene	0.28	J	0.11	1.0	$\mu g/L$	SW-846 8260D
tert-Butylbenzene	0.24	J	0.17	1.0	$\mu g/L$	SW-846 8260D
Toluene	0.25	J	0.14	1.0	$\mu g/L$	SW-846 8260D
2-Methylnaphthalene (SIM)	4.9		0.062	1.0	$\mu g/L$	SW-846 8270E
Acenaphthene (SIM)	0.30		0.033	0.30	$\mu g/L$	SW-846 8270E
Acenaphthylene (SIM)	0.095	J	0.035	0.20	$\mu g/L$	SW-846 8270E
Fluoranthene (SIM)	0.027	J	0.025	0.50	$\mu g/L$	SW-846 8270E
Fluorene (SIM)	0.38	J	0.034	1.0	$\mu g/L$	SW-846 8270E
Naphthalene (SIM)	3.2		0.26	1.0	$\mu g/L$	SW-846 8270E
Phenanthrene (SIM)	0.42		0.030	0.050	$\mu g/L$	SW-846 8270E
Pyrene (SIM)	0.025	J	0.023	1.0	$\mu g/L$	SW-846 8270E

Con-Test does not accept liability for the consequences of any actions taken solely on the basis of the information provided in the Executive Summary section of this report. Users must review this report in its entirety to determine data usability and assessment.



CASE NARRATIVE SUMMARY

All reported results are within defined laboratory quality control objectives unless listed below or otherwise qualified in this report.

For method 8270, only PAHs were requested and reported.

SW-846 8260D

Qualifications:

V-20

Continuing calibration verification (CCV) did not meet method specifications and was biased on the high side. Data validation is not affected since sample result was "not detected" for this compound. Analyte & Samples(s) Qualified:

Bromomethane

B245122-BS1, B245122-BSD1, S042311-CCV1

Chloromethane

B245122-BS1, B245122-BSD1, S042311-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.

Lua Watslengton Technical Representative



Project Location: Dunn, NC Sample Description: Work Order: 19K0021

Date Received: 10/31/2019
Field Sample #: TW-1

Sampled: 10/29/2019 12:30

Sample ID: 19K0021-01
Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

Volatile Organic Compounds by GC/MS											
Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst	
Acetone	ND	50	3.8	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Acrylonitrile	ND	5.0	0.52	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
tert-Amyl Methyl Ether (TAME)	ND	0.50	0.14	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Benzene	0.66	1.0	0.18	μg/L	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Bromobenzene	ND	1.0	0.15	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Bromochloromethane	ND	1.0	0.32	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Bromodichloromethane	ND	0.50	0.16	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Bromoform	ND	1.0	0.46	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Bromomethane	ND	2.0	0.78	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
2-Butanone (MEK)	ND	20	1.9	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
tert-Butyl Alcohol (TBA)	ND	20	4.2	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
n-Butylbenzene	3.5	1.0	0.21	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
sec-Butylbenzene	15	1.0	0.16	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
tert-Butylbenzene	0.24	1.0	0.17	μg/L	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
tert-Butyl Ethyl Ether (TBEE)	ND	0.50	0.16	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Carbon Disulfide	ND	5.0	4.4	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Carbon Tetrachloride	ND	1.0	0.11	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Chlorobenzene	ND	1.0	0.15	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Chlorodibromomethane	ND	0.50	0.21	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Chloroethane	ND	2.0	0.35	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Chloroform	ND	2.0	0.17	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Chloromethane	ND	2.0	0.45	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
2-Chlorotoluene	ND	1.0	0.12	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
4-Chlorotoluene	ND	1.0	0.14	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,2-Dibromo-3-chloropropane (DBCP)	ND	5.0	0.53	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,2-Dibromoethane (EDB)	ND	0.50	0.19	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Dibromomethane	ND	1.0	0.37	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,2-Dichlorobenzene	ND	1.0	0.16	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,3-Dichlorobenzene	ND	1.0	0.12	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,4-Dichlorobenzene	ND	1.0	0.13	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
trans-1,4-Dichloro-2-butene	ND	2.0	0.31	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Dichlorodifluoromethane (Freon 12)	ND	2.0	0.26	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,1-Dichloroethane	ND	1.0	0.16	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,2-Dichloroethane	ND	1.0	0.41	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,1-Dichloroethylene	ND	1.0	0.32	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
cis-1,2-Dichloroethylene	ND	1.0	0.13	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
trans-1,2-Dichloroethylene	ND	1.0	0.31	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,2-Dichloropropane	ND	1.0	0.20	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,3-Dichloropropane	ND	0.50	0.11	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
2,2-Dichloropropane	ND	1.0	0.20	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
1,1-Dichloropropene	ND	2.0	0.16	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
cis-1,3-Dichloropropene	ND	0.50	0.13	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
trans-1,3-Dichloropropene	ND	0.50	0.23	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	
Diethyl Ether	ND	2.0	0.34	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH	

Page 6 of 23



Project Location: Dunn, NC Sample Description: Work Order: 19K0021

Date Received: 10/31/2019 Field Sample #: TW-1

Sampled: 10/29/2019 12:30

Sample ID: 19K0021-01 Sample Matrix: Ground Water

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	0.50	0.17	μg/L	1	•	SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,4-Dioxane	ND	50	22	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Ethylbenzene	4.9	1.0	0.13	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Hexachlorobutadiene	ND	0.60	0.47	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
2-Hexanone (MBK)	ND	10	1.5	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Isopropylbenzene (Cumene)	43	1.0	0.17	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
p-Isopropyltoluene (p-Cymene)	0.92	1.0	0.20	μg/L	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Methyl tert-Butyl Ether (MTBE)	ND	1.0	0.25	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Methylene Chloride	ND	5.0	0.34	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
4-Methyl-2-pentanone (MIBK)	ND	10	1.7	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Naphthalene	3.9	2.0	0.31	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
n-Propylbenzene	1.8	1.0	0.13	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Styrene	0.28	1.0	0.11	μg/L	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,1,1,2-Tetrachloroethane	ND	1.0	0.27	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,1,2,2-Tetrachloroethane	ND	0.50	0.22	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Tetrachloroethylene	ND	1.0	0.18	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Tetrahydrofuran	ND	10	0.51	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Toluene	0.25	1.0	0.14	μg/L	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,2,3-Trichlorobenzene	ND	5.0	0.57	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,2,4-Trichlorobenzene	ND	1.0	0.40	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,3,5-Trichlorobenzene	ND	1.0	0.30	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,1,1-Trichloroethane	ND	1.0	0.20	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,1,2-Trichloroethane	ND	1.0	0.16	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Trichloroethylene	ND	1.0	0.24	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Trichlorofluoromethane (Freon 11)	ND	2.0	0.33	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,2,3-Trichloropropane	ND	2.0	0.25	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	ND	1.0	0.32	μg/L	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,2,4-Trimethylbenzene	4.4	1.0	0.18	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
1,3,5-Trimethylbenzene	2.2	1.0	0.14	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Vinyl Chloride	ND	2.0	0.45	$\mu g/L$	1		SW-846 8260D	11/5/19	11/6/19 10:45	EEH
m+p Xylene	1.5	2.0	0.30	$\mu g/L$	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH
o-Xylene	0.40	1.0	0.17	$\mu g/L$	1	J	SW-846 8260D	11/5/19	11/6/19 10:45	EEH
Surrogates		% Reco	very	Recovery Limits	S	Flag/Qual				
1,2-Dichloroethane-d4		94.7		70-130					11/6/19 10:45	
Toluene-d8		99.0		70-130					11/6/19 10:45	



Work Order: 19K0021 Sample Description:

Project Location: Dunn, NC Date Received: 10/31/2019 Field Sample #: TW-1

Sampled: 10/29/2019 12:30

Sample ID: 19K0021-01 Sample Matrix: Ground Water

Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst
Acenaphthene (SIM)	0.30	0.30	0.033		Dilution	riag/Quai	SW-846 8270E	11/5/19	11/6/19 18:18	CLA
1 ,				μg/L	1					
Acenaphthylene (SIM)	0.095	0.20	0.035	μg/L	1	J	SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Anthracene (SIM)	ND	0.20	0.032	μg/L	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Benzo(a)anthracene (SIM)	ND	0.050	0.016	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Benzo(a)pyrene (SIM)	ND	0.10	0.012	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Benzo(b)fluoranthene (SIM)	ND	0.050	0.015	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Benzo(g,h,i)perylene (SIM)	ND	0.50	0.018	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Benzo(k)fluoranthene (SIM)	ND	0.20	0.012	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Chrysene (SIM)	ND	0.20	0.015	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Dibenz(a,h)anthracene (SIM)	ND	0.10	0.017	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Fluoranthene (SIM)	0.027	0.50	0.025	$\mu g/L$	1	J	SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Fluorene (SIM)	0.38	1.0	0.034	$\mu g/L$	1	J	SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Indeno(1,2,3-cd)pyrene (SIM)	ND	0.10	0.018	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
2-Methylnaphthalene (SIM)	4.9	1.0	0.062	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Naphthalene (SIM)	3.2	1.0	0.26	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Phenanthrene (SIM)	0.42	0.050	0.030	$\mu g/L$	1		SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Pyrene (SIM)	0.025	1.0	0.023	$\mu g/L$	1	J	SW-846 8270E	11/5/19	11/6/19 18:18	CLA
Surrogates		% Reco	very	Recovery Limits	s	Flag/Qual				
Nitrobenzene-d5		65.4		30-130					11/6/19 18:18	
2-Fluorobiphenyl		48.5		30-130					11/6/19 18:18	
p-Terphenyl-d14		63.0		30-130					11/6/19 18:18	



Sample Extraction Data

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

Lab Number [Field ID]	Batch	Initial [mL]	Final [mL]	Date
19K0021-01 [TW-1]	B245122	5	5.00	11/05/19

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

Lab Number [Field ID]	Batch	Initial [mL]	Final [mL]	Date
19K0021-01 [TW-1]	B245268	1000	1.00	11/05/19



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
Batch B245122 - SW-846 5030B										

Record (RAMS)	Batch B245122 - SW-846 5030B				
Accyolatifich ND 5.0 ggL Bettene ND 1.0 ggL Bettene ND 1.0 ggL Bettene ND 1.0 ggL Bettene ND 1.0 ggL Bettenenthane ND 1.0 ggL Bronochementhane ND 1.0 ggL Bronochem ND 1.0 ggL Bronochem ND 2.0 ggL Pattaneon ND 2.0 ggL Pattaneon ND 1.0 ggL Pattaneon ND 1.0 ggL ButtyBeriche ND 1.0<	Blank (B245122-BLK1)				Prepared: 11/05/19 Analyzed: 11/06/19
text-Amy Modyl Eller (TAME) ND 0.50 ug/L Bromone Bromene ND 1.0 ug/L Bromoneldoromethane ND 1.0 ug/L Bromoneldoromethane ND 0.50 ug/L Bromoneldoromethane ND 0.50 ug/L Bromoneldoromethane ND 2.0 ug/L Bromoneldoromethane ND 2.0 ug/L Bromoneldoromethane ND 2.0 ug/L Palmatines (MIK) ND 2.0 ug/L Intermonelloridoromethane ND 1.0 ug/L Intermonelloridoromethane ND 1.0 ug/L Carbon Danifide ND 5.0 ug/L Chlorodrinoromethane ND 1.0 ug/L Chlorodrinoromethane ND 2.0 ug/L Chlorodrinoromethane ND 2.0 ug/L Chlorodrinoromethane ND 1.0 ug/L Chlorodrinoromethane ND 1.0 ug/L		ND			
Blacenee ND 1.0 upf. Bromondectromechane ND 1.0 upg. Bromondectromechane ND 0.50 upg. Bromondectromechane ND 0.50 upg. Bromondectromechane ND 2.0 upg. Bromondectromechane ND 2.0 upg. Bromondectromechane ND 2.0 upg. Betablishumme ND 2.0 upg. betablishumme ND 1.0 upg. see-Burgheneme ND 1.0 upg. see-Burgheneme ND 1.0 upg. cert-Burgheneme ND 1.0 upg. cert-Burgheneme ND 1.0 upg. Carbon Tearithoride ND 1.0 upg. Chloromechane ND 1.0 upg. Chloromechane ND 2.0 upg. Chloromechane ND 2.0 upg. Chloromechane ND 2.0 <td>-</td> <td>ND</td> <td></td> <td></td> <td></td>	-	ND			
Bomonabanama ND 1.0 ug/L Bomonchioromatiane ND 1.0 ug/L Bomonderboromatiane ND 1.0 ug/L Bomonderboromatiane ND 2.0 ug/L Bomonderboromatiane ND 2.0 ug/L Politations (MEK) ND 2.0 ug/L Publication ND 1.0 ug/L Publication ND 1.0 ug/L See-Butylbaceane ND 1.0 ug/L tear-Basy Bay Benezine ND 1.0 ug/L carbon Strandinde ND 1.0 ug/L Carbon Daniafde ND 1.0 ug/L Chloroderace ND 1.0 ug/L Chloroderace ND 1.0 ug/L Chloroderace ND 2.0 ug/L Chloroderace ND 2.0 ug/L Chloroderace ND 2.0 ug/L Chloroderace ND 2.0	tert-Amyl Methyl Ether (TAME)	ND	0.50		
Bomodelbromethane ND 0.50 agE Bromoderbromethane ND 0.50 agE Bromoderbromethane ND 0.50 agE Bromoderbromethane ND 2.00 agE Bromoderbromethane ND 2.00 agE be-Buryl Accord (TBA) ND 2.00 agE be-Buryl Excerce ND 1.00 agE be-Buryl-Burzene ND 1.00 agE bert-Hugel Experiment ND 1.00 agE Carbon Errarbierde ND 1.00 agE Chlorofrom ND 1.00 agE Chlorofrom ND 2.00 agE Chlorofrom ND 2.00 agE Chlorofrom	Benzene	ND	1.0	$\mu g/L$	
Bemondsfeme ND 0.50 µgL Benomentame ND 0.20 µgL Benomentame ND 0.20 µgL Jubiamene (MIK) ND 2.00 µgL vert-Butyl Acchol (TBA) ND 2.00 µgL sebhylbenzene ND 0.10 µgL tert-Butyl Bible (TBEE) ND 0.00 µgL Carban Desalfide ND 5.00 µgL Carban Bealfide ND 5.00 µgL Chloredimonamen ND 0.10 µgL Chloredimonenthane ND 0.10 µgL Chloredimonenthane ND 2.0 µgL L'Albeitonbertene ND	Bromobenzene	ND	1.0		
None	Bromochloromethane	ND	1.0	$\mu \text{g/L}$	
Brownethane		ND	0.50	μg/L	
2-Buttone (MEK) ND 20 μgL ter-Butyl Achol (TBA) ND 20 μgL ter-Butyl Achol (TBA) ND 10 μgL sce-Butylbenzene ND 10 μgL ter-Butyl Ebbyr Ebbr (TBEE) ND 0.50 μgL Carbon Distiffer ND 10 μgL Carbon Carteraloride ND 10 μgL Carbon Carteraloride ND 10 μgL Chloroditromomethane ND 0.50 μgL Chloroditromomethane ND 2.0 μgL Chloroditromomethane ND 1.0 μgL 2-Chloroditromomethane ND 1.0 μgL 1_2-Di	Bromoform	ND	1.0	$\mu \text{g/L}$	
ter-Bary Alcohed (TBA) ND 20 gst. neurylbrozone ND 1.0 pgt. sec-Budylbenzone ND 1.0 gst. ter-Bary Bibly fiber (TBEE) ND 0.50 gst. Carbon Datalifide ND 5.0 gst. Carbon Datalifide ND 1.0 gst. Chlorocharene ND 1.0 gst. Chlorocharene ND 1.0 gst. Chlorocharene ND 0.50 gst. Chlorocharene ND 2.0 gst. Chlorocharene ND 2.0 gst. Chlorocharene ND 1.0 gst. Chlorocharene ND 1.0 gst. 1.2-Drbomo-3-chloropropane (DBCP) ND 5.0 ggt. 1.2-Drbomo-denne (EDB) ND 0.50 ggt. 1.2-Drbomo-denne (EDB) ND 0.50 ggt. 1.2-Drbomo-denne (EDB) ND 0.50 ggt. 1.2-Drbomo-denne	Bromomethane	ND	2.0	$\mu \text{g/L}$	
sa-Bary-fluenzene	2-Butanone (MEK)	ND	20	$\mu g\!/\!L$	
sex-BartyNethoraces ND 1.0 pgL tert-Buryl Ethyl Ether (TBEF) ND 0.50 pgL Carbon Dissulife ND 5.0 pgL Chloro-Shurife ND 1.0 pgL Chloro-Chromethane ND 0.50 pgL Chloro-Ghromethane ND 2.0 pgL Chloro-Ghromethane ND 2.0 pgL Chloro-Ghromethane ND 2.0 pgL Chloro-Ghromethane ND 1.0 pgL 2-Chloro-Ghromethane ND 1.0 pgL 2-Chloro-Ghromethane ND 1.0 pgL 1.2-Dirbomos-Achtoropropane (DBCP) ND 5.0 pgL 1.2-Dirbomos-Achtoropropane ND 1.0 pgL 1.2-Dirbomos-Achtoropropane ND 1.0 pgL 1.2-Dirbomos-Achtoropropane ND 1.0 pgL 1.2-Dirbomos-Achtoropropane ND 1.0 pgL 1.4-Dirbomos-Achtoropropane ND 1.0 p	tert-Butyl Alcohol (TBA)	ND	20	$\mu g\!/\!L$	
tor-Barly Ethyl Ethyr (TBEE) ND 0.50 pgL Carbon Disalifide ND 5.0 pgL Carbon Carbon Certachiorde ND 1.0 pgL Chlorocharea ND 0.50 pgL Chlorocharea ND 0.50 pgL Chlorocharea ND 2.0 pgL Chlorocharea ND 2.0 pgL Chlorocharea ND 2.0 pgL Chlorocharea ND 1.0 pgL 1,2-Dirborocharea ND 1.0 pgL 1,2-Dirborocharea ND 1.0 pgL 1,3-Dirborocharea ND 1.0 pgL 1,3-Dirborocharea ND 1.0 pgL 1,3-Dirborocharea ND 2.0 pg	n-Butylbenzene	ND	1.0	$\mu g\!/\!L$	
ten-Buy Liby Libry	sec-Butylbenzene	ND	1.0	$\mu g\!/\!L$	
Carbon Disulfide ND 5.0 µgL Carbon Tetrachloride ND 1.0 µgL Chlorochanzene ND 0.50 µgL Chlorochanzene ND 2.0 µgL Chlorochane ND 2.0 µgL Chloroform ND 2.0 µgL Chlorofoluene ND 1.0 µgL 4 Chlorofoluene ND 1.0 µgL 4 Chlorofoluene ND 5.0 µgL 1.2-Dibromo-schloropropane (DBCP) ND 5.0 µgL 1.2-Dibromo-chloropropane (DBCP) ND 5.0 µgL 1.2-Dichlorochane ND 1.0 µgL 1.2-Dichlorochane ND 1.0 µgL 1.2-Dichlorochane ND 1.0 µgL 1.3-Dichlorochane ND 1.0 µgL 1.3-Dichlorochane ND 1.0 µgL 1.3-Dichlorochane ND 1.0 µgL 1.3-Dichlorochane ND	tert-Butylbenzene	ND	1.0	$\mu g\!/\!L$	
Carbon Tetrachloride ND 1.0 µg/L Chlorobazone ND 1.0 µg/L Chlorodhrane ND 0.5 µg/L Chlorodhrane ND 2.0 µg/L Chlorodhrane ND 2.0 µg/L Chlorodhane ND 1.0 µg/L Chlorodhone ND 1.0 µg/L Chloromethane ND 1.0 µg/L 1.2-Dibromo-Schloropopane (DBCP) ND 0.5 µg/L 1.2-Dibromo-Schloropopane (DBCP) ND 0.5 µg/L 1.2-Dibromo-Schloropopane (EDB) ND 0.5 µg/L 1.2-Dibromo-Schloropopane (EDCP) ND 0.5 µg/L 1.2-Dibromo-Schloropopane (EDCP) ND 0.0 µg/L 1.2-Dibromo-Schloropopane (EDCP) ND 1.0 µg/L 1.2-Dibromo-Schloropopane (EDCP) ND 1.0 µg/L 1.4-Dichloropopane (EDCP) ND 1.0 µg/L 1.1-Dichloropopane (EDCP) ND 0.0	tert-Butyl Ethyl Ether (TBEE)	ND	0.50	$\mu \text{g/L}$	
Chlorodbenzene ND 1.0 µg/L Chlorodbromomethane ND 0.50 µg/L Chlorodbromomethane ND 2.0 µg/L Chloroform ND 2.0 µg/L Chlorodbrome ND 1.0 µg/L 4-Chlorodbene ND 1.0 µg/L 4-Chlorodbrome ND 5.0 µg/L 1,2-Dibromos-beloropopane (DBCP) ND 5.0 µg/L 1,2-Dibromoschane (FDB) ND 0.50 µg/L 1,2-Dichlorodenzene ND 1.0 µg/L 1,3-Dichlorodenzene ND 1.0 µg/L 1,4-Dichlorodenzene ND 1.0 µg/L 1,4-Dichlorodenzene ND 2.0 µg/L 1,4-Dichlorodenzene ND 2.0 µg/L 1,4-Dichlorodenzene ND 2.0 µg/L 1,4-Dichlorodenzene ND 2.0 µg/L 1,4-Dichlorodenzene ND 1.0 µg/L 1,4-Dichlorod	Carbon Disulfide	ND	5.0	$\mu \text{g/L}$	
Chlorochane ND 0.50 µg.L Chlorochane ND 2.0 µg.L Chloromehane ND 2.0 µg.L Chlorocholune ND 1.0 µg.L 4.Chlorotolune ND 1.0 µg.L 1.2-Dibromo-3-chloropropane (DBCP) ND 5.0 µg.L 1.2-Dibromo-4-chloropropane (DBCP) ND 5.0 µg.L 1.2-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L 1.3-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L 1.4-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L 1.1-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L	Carbon Tetrachloride	ND	1.0	$\mu \text{g/L}$	
Chlorochane ND 0.50 µg.L Chlorochane ND 2.0 µg.L Chloromehane ND 2.0 µg.L Chlorocholune ND 1.0 µg.L 4.Chlorotolune ND 1.0 µg.L 1.2-Dibromo-3-chloropropane (DBCP) ND 5.0 µg.L 1.2-Dibromo-4-chloropropane (DBCP) ND 5.0 µg.L 1.2-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L 1.3-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L 1.4-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L 1.1-Dibromo-4-chloropropane (DBCP) ND 1.0 µg.L	Chlorobenzene		1.0		
Chlorothame ND 2,0 µg/L Chloroform ND 2,0 µg/L Chlorotoluene ND 2,0 µg/L 2-Chlorotoluene ND 1,0 µg/L 1,2-Dibromo-2-biloropropane (DBCP) ND 5,0 µg/L 1,2-Dibromo-chlane (EDB) ND 0,50 µg/L 1,2-Dichlorobenzene ND 1,0 µg/L 1,3-Dichlorobenzene ND 1,0 µg/L 1,3-Di	Chlorodibromomethane		0.50		
Chlorome ND 2.0 µg/L Chloromethane ND 1.0 µg/L 4.Chlorotoluene ND 1.0 µg/L 4.Chlorotoluene ND 1.0 µg/L 1.2-Dibromos-3-chloropropane (DBCP) ND 5.0 µg/L 1.2-Dibromoshane (EDB) ND 1.0 µg/L 1.2-Dichlorobenzene ND 1.0 µg/L 1,2-Dichlorobenzene ND 1.0 µg/L 1,4-Dichloro-2-butene ND 2.0 µg/L 1,4-Dichloro-2-butene ND 2.0 µg/L 1,1-Dichloro-2-butene ND 1.0 µg/L 1,1-Dichloro-2-butene ND 1.0 µg/L	Chloroethane		2.0		
Chlorotoluene ND 2.0 µg/L 4-Chlorotoluene ND 1.0 µg/L 4-Chlorotoluene ND 5.0 µg/L 1,2-Dibromoe-3-chloropropane (DBCP) ND 5.0 µg/L 1,2-Dibromoethane (EDB) ND 0.0 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 1,4-Dichlorothylene ND 1.0 µg/L 1,1-Dichlorothylene ND 1.0 µg/L 1,2-Dichlorothylene ND 1.0 µg/L 1,2-Dichlorothylene ND 0.0 µg/L 1,2-Dichlorothylene ND 0.0 µg/L	Chloroform		2.0		
2-Chlorotoluene ND 1,0 pg/L 4-Chlorotoluene ND 1,0 pg/L 1,2-Dibromoethane (EDBY) ND 0,50 pg/L 1,2-Dibromoethane (EDB) ND 0,50 pg/L 1,2-Dichlorobenzene ND 1,0 pg/L 1,3-Dichlorobenzene ND 1,0 pg/L 1,4-Dichlorobenzene ND 1,0 pg/L 1,4-Dichlorobenzene ND 2,0 pg/L 1,4-Dichlorobenzene ND 2,0 pg/L 1,1-Dichlorobenzene ND 2,0 pg/L 1,1-Dichlorobenzene ND 2,0 pg/L 1,1-Dichlorobenzene ND 1,0 pg/L <	Chloromethane				
4-Chlorotoluene ND 1.0 µg/L 1,2-Dibromo-S-chloropropane (DBCP) ND 5.0 µg/L 1,2-Dibromo-Shoropropane (DBCP) ND 0.50 µg/L 1,2-Dibromo-Shoropropane (DBCP) ND 0.50 µg/L 1,2-Dibromo-Shoropropane (DBCP) ND 1.0 µg/L 1,2-Dibromo-Shoropropane ND 1.0 µg/L 1,3-Dichlorobenzene ND 1.0 µg/L 1,4-Dichlorobenzene ND 1.0 µg/L 1,4-Dichloro-S-butene ND 1.0 µg/L 1,4-Dichloro-S-butene ND 1.0 µg/L 1,1-Dichloro-Shoropropane (Fren 12) ND 2.0 µg/L 1,1-Dichloro-Shoropropane ND 1.0 µg/L 1,1-Dichloro-Shoropropane ND 1.0 µg/L 1,1-Dichloro-Shoropropane ND 1.0 µg/L 1,1-Dichloro-Shoropropane ND 1.0 µg/L 1,2-Dichloro-Shoropropane ND 1.0 µg/L 1,2-Dichloropropane ND 1.0 µg/L 1,2-Dichloropropane ND 1.0 µg/L 1,1-Dichloropropane ND 0.50 µg/L 1,1-Dichloropropa	2-Chlorotoluene				
1,2-Dibromo-s-holropropane (DBCP) ND 5.0 µgL 1,2-Dibromoethane (EDB) ND 0.50 µgL Dibromomethane ND 1.0 µgL 1,2-Dichlorobenzene ND 1.0 µgL 1,3-Dichlorobenzene ND 1.0 µgL 1,4-Dichloro-butene ND 1.0 µgL Dichlorodifluoromethane (Fren 12) ND 2.0 µgL 1,1-Dichlorodethane ND 1.0 µgL 1,1-Dichlorothylene ND 1.0 µgL 1,1-Dichlorothylene ND 1.0 µgL 1,2-Dichlorothylene ND 1.0 µgL 1,3-Dichlorothylene ND 1.0 µgL 1,3-Dichloropropane ND 1.0 µgL 1,3-Dichloropropane ND 0.5 µgL 1,1-Dichloropropane ND 0.5 µgL 1,1-Dichloropropane ND 0.5 µgL 1,1-Dichloropropane ND 0.5 µgL					
1,2-Dibromoethane (EDB) ND 0.50 µg/L Dibromoethane ND 1.0 µg/L 1,2-Dichlorobenzene ND 1.0 µg/L 1,3-Dichlorobenzene ND 1.0 µg/L 1,4-Dichloro-2-butene ND 2.0 µg/L Dichlorodifluoromethane (Fren 12) ND 2.0 µg/L 1,1-Dichloroethane ND 1.0 µg/L 1,2-Dichloroethylene ND 1.0 µg/L 1,1-Dichloroethylene ND 1.0 µg/L 1,1-Dichloroethylene ND 1.0 µg/L 1,1-Dichloroethylene ND 1.0 µg/L 1,1-Dichloroethylene ND 1.0 µg/L 1,2-Dichloroethylene ND 1.0 µg/L 1,2-Dichloropropane ND 1.0 µg/L 1,2-Dichloropropane ND 0.0 µg/L 1,1-Dichloropropane ND 0.5 µg/L 1,1-Dichloropropane ND 0.5 µg/L					
Dibromomethane ND 1.0 µg/L 1,2-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 1,1-Dichloroethane ND 1.0 µg/L 1,2-Dichloroethane ND 1.0 µg/L 1,1-Dichloroethylene ND 1.0 µg/L 1,1-Dichloroethylene ND 1.0 µg/L 1,2-Dichloroptopane ND 1.0 µg/L 1,2-Dichloroptopane ND 1.0 µg/L 1,3-Dichloroptopane ND 0.5 µg/L 1,3-Dichloroptopane ND 0.5 µg/L cis-1,3-Dichloroptopene ND 0.5 µg/L cis-1,3-Dichloroptopene ND 0.5 µg/L Dichtyl Ether ND 0.5 µg/L Dichtyl Ether (DIPE) ND 0.0 µg/L					
1.2-Dichlorobenzene ND 1.0 μg/L 1.3-Dichlorobenzene ND 1.0 μg/L 1.4-Dichloroc-butene ND 1.0 μg/L Dichlorodifluoromethane (Freon 12) ND 2.0 μg/L 1,1-Dichlorocethane ND 1.0 μg/L 1,1-Dichlorocethylene ND 1.0 μg/L 1,1-Dichlorocethylene ND 1.0 μg/L cis-1,2-Dichlorocethylene ND 1.0 μg/L trans-1,2-Dichlorocethylene ND 1.0 μg/L 1,3-Dichloropropane ND 1.0 μg/L 1,3-Dichloropropane ND 0.5 μg/L 1,1-Dichloropropane ND 0.5 μg/L 1,1-Dichloropropane ND 0.5 μg/L 1,1-Dichloropropene ND 0.5 μg/L trans-1,3-Dichloropropene ND 0.5 μg/L trans-1,3-Dichloropropene ND 0.5 μg/L trans-1,4-Dichloropropene ND 0.5					
1,3-Dichlorobenzene ND 1,0 µg/L 1,4-Dichloro-2-butene ND 1,0 µg/L trans-1,4-Dichloro-2-butene ND 2,0 µg/L 1,1-Dichloroethane (Freon 12) 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 1,2-Dichloropropane ND 0.50 µg/L 1,3-Dichloropropane ND 0.50 µg/L 1,1-Dichloropropene ND 0.50 µg/L cis-1,3-Dichloropropene ND 0.50 µg/L trans-1,3-Dichloropropene ND 0.50 µg/L Dichyl Ether ND 0.50 µg/L Disopropyl Ether (DIPE) ND 0.50 µg/L Hexachlorobutadiene ND 0.60					
1,4-Dichlorobenzene ND 1.0 µg/L bichlorodifluoromethane (Freon 12) ND 2.0 µg/L 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 cis-1,2-Dichloroethylene ND 1.0 µg/L trans-1,2-Dichloropropane ND 1.0 µg/L 1,3-Dichloropropane ND 1.0 µg/L 1,3-Dichloropropane ND 1.0 µg/L 2,2-Dichloropropane ND 1.0 µg/L 1,1-Dichloropropene ND 0.50 µg/L 1,1-Dichloropropene ND 0.50 µg/L Dichly1 Ether ND 0.50 µg/L Dichly1 Ether (DIPE) ND 0.50 µg/L Disporpyl Ether (DIPE) ND 0.50 µg/L Ethylbenzene ND 0.60 µg/L					
trans-1,4-Dichloro-2-butene ND 2.0 µg/L Dichlorodifluoromethane (Freon 12) ND 2.0 µg/L 1,1-Dichloroethane ND 1.0 µg/L 1,2-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-Dichloroethylene ND 1.0 µg/L 1,2-Dichloroethylene ND 1.0 µg/L 1,2-Dichloropropane ND 1.0 µg/L 1,2-Dichloropropane ND 1.0 µg/L 2,2-Dichloropropane ND 1.0 µg/L 1,1-Dichloropropene ND 0.50 µg/L 1,1-Dichloropropene ND 0.50 µg/L Dichly1 Ether ND 0.50 µg/L Dichly1 Ether ND 0.50 µg/L Dichly1 Ether ND 0.50 µg/L Hexachlorobutadiene ND 0.60 µg/L </td <td></td> <td></td> <td></td> <td></td> <td></td>					
Dichlorodifluoromethane (Freon 12) ND 2.0 µg/L 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 1.0 µg/L 1,1-Dichloropropane ND 1.0 µg/L 1,1-Dichloropropene ND 0.50 µg/L 1,1-Dichloropropene ND 0.50 µg/L 1trans-1,3-Dichloropropene ND 0.50 µg/L Dichlyl Ether ND 0.50 µg/L Dichlyl Ether (DIPE) ND 0.50 µg/L Disporpyl Ether (DIPE) ND 0.50 µg/L Ethylbenzene ND 0.60 µg/L Ethylbenzene ND 0.60 µg/L </td <td></td> <td></td> <td></td> <td></td> <td></td>					
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 0.50 μg/L cis-1,3-Dichloropropene ND 0.50 μg/L trans-1,3-Dichloropropene ND 0.50 μg/L Dicthyl Ether ND 0.50 μg/L Diisopropyl Ether (DIPE) ND 0.50 μg/L 1,4-Dioxane ND 1.0 μg/L Hexachlorobutadiene ND 0.60 μg/L Hexachlorobutadiene ND 0.60 μg/L 2-Hexanone (MBK) ND 1.0 μg/L					
1,2-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-Dichloroprophylene ND 1.0 µg/L 1,2-Dichloropropane ND 1.0 µg/L 1,3-Dichloropropane ND 1.0 µg/L 2,2-Dichloropropane ND 2.0 µg/L 1,1-Dichloropropene ND 0.5 µg/L cis-1,3-Dichloropropene ND 0.5 µg/L trans-1,3-Dichloropropene ND 0.5 µg/L Dictyl Ether ND 0.5 µg/L Dictyl Ether (DIPE) ND 0.5 µg/L 1,4-Dioxane ND 5 µg/L Ethylbenzene ND 0.6 µg/L Hexachlorobutadiene ND 0.6 µg/L 2-Hexanone (MBK) ND 0.0 µg/L 1-sopropyllounee (p-Cymene) ND 0.0 µg/L					
1,1-Dichloroethylene ND 1.0 µg/L cis-1,2-Dichloroethylene ND 1.0 µg/L 1,2-Dichloroptopane ND 1.0 µg/L 1,3-Dichloroptopane ND 0.50 µg/L 2,2-Dichloroptopane ND 1.0 µg/L 1,1-Dichloroptopene ND 2.0 µg/L 1,1-Dichloroptopene ND 0.50 µg/L cis-1,3-Dichloroptopene ND 0.50 µg/L Dicthyl Ether ND 0.50 µg/L Disopropyl Ether (DIPE) ND 0.50 µg/L 1,4-Dioxane ND 0.50 µg/L Ethylbenzene ND 0.60 µg/L Hexachlorobutadiene ND 0.60 µg/L 1-Lexanone (MBK) ND 0.60 µg/L 1-sopropylbourene (Cumene) ND 0.0 µg/L 1-sopropylbourene (Cumene) ND 0.0 µg/L					
cis-1,2-Dichloroethylene ND 1.0 μg/L 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 Dicityl Ether ND 0.50 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 1.0 μg/L Isopropylbourene (Cumene) ND 1.0 μg/L p-Isopropyltoluene (p-Cymene) ND 1.0 μg/L					
trans-1,2-Dichloroethylene ND 1.0 µg/L 1,2-Dichloropropane ND 0.50 µg/L 2,2-Dichloropropane ND 1.0 µg/L 1,1-Dichloropropane ND 1.0 µg/L 1,1-Dichloropropane ND 0.50 µg/L 1,1-Dichloropropene ND 0.50 µg/L trans-1,3-Dichloropropene ND 0.50 µg/L Dicthyl Ether ND 0.50 µg/L Disopropyl Ether (DIPE) ND 0.50 µg/L 1,4-Dioxane ND 0.50 µg/L Ethylbenzene ND 0.60 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 Diethyl Ether ND 0.50 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 Isopropylboluene (p-Cymene) 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					
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 0.50 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					
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 1.0 μg/L Isopropylbenzene (Cumene) ND 1.0 μg/L p-Isopropyltoluene (p-Cymene) ND 1.0 μg/L					
cis-1,3-Dichloropropene ND 0.50 $\mu g/L$ trans-1,3-Dichloropropene ND 0.50 $\mu g/L$ Diethyl Ether ND 2.0 $\mu g/L$ Diisopropyl Ether (DIPE) ND 0.50 $\mu g/L$ 1,4-Dioxane ND 50 $\mu g/L$ Ethylbenzene ND 1.0 $\mu g/L$ Hexachlorobutadiene ND 0.60 $\mu g/L$ 2-Hexanone (MBK) ND 1.0 $\mu g/L$ Isopropylbenzene (Cumene) ND 1.0 $\mu g/L$ p-Isopropyltoluene (p-Cymene) ND 1.0 $\mu g/L$					
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 1.0 μg/L Isopropylbenzene (Cumene) ND 1.0 μg/L p-Isopropyltoluene (p-Cymene) ND 1.0 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					
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 1.0 μg/L Isopropylbenzene (Cumene) ND 1.0 μg/L p-Isopropyltoluene (p-Cymene) ND 1.0 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					
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					
HexachlorobutadieneND0.60μg/L2-Hexanone (MBK)ND10μg/LIsopropylbenzene (Cumene)ND1.0μg/Lp-Isopropyltoluene (p-Cymene)ND1.0μg/L					
2-Hexanone (MBK) ND 10 μ g/L Isopropylbenzene (Cumene) ND 1.0 μ g/L p-Isopropyltoluene (p-Cymene) ND 1.0 μ g/L	-				
Isopropylbenzene (Cumene) ND 1.0 $\mu g/L$ p-Isopropyltoluene (p-Cymene) ND 1.0 $\mu g/L$					
p-Isopropyltoluene (p-Cymene) ND $1.0 \mu g/L$					
Methyl tert-Butyl Ether (MTBE) ND $1.0 \mu g/L$					
	Methyl tert-Butyl Ether (MTBE)	ND	1.0	μg/L	



QUALITY CONTROL

Spike

Source

%REC

RPD

Volatile Organic Compounds by GC/MS - Quality Control

Reporting

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Satch B245122 - SW-846 5030B										
lank (B245122-BLK1)				Prepared: 11	/05/19 Analy	/zed: 11/06/1	9			
lethylene Chloride	ND	5.0	μg/L							
Methyl-2-pentanone (MIBK)	ND	10	μg/L							
aphthalene	ND	2.0	μg/L							
Propylbenzene	ND	1.0	$\mu g/L$							
yrene	ND	1.0	μg/L							
1,1,2-Tetrachloroethane	ND	1.0	μg/L							
1,2,2-Tetrachloroethane	ND	0.50	μg/L							
trachloroethylene	ND	1.0	μg/L							
trahydrofuran	ND	10	μg/L							
luene	ND	1.0	μg/L							
2,3-Trichlorobenzene		5.0	μg/L							
2,4-Trichlorobenzene	ND	1.0	μg/L μg/L							
3,5-Trichlorobenzene	ND									
•	ND	1.0	μg/L							
1,1-Trichloroethane	ND	1.0	μg/L							
1,2-Trichloroethane	ND	1.0	μg/L							
ichloroethylene	ND	1.0	μg/L							
ichlorofluoromethane (Freon 11)	ND	2.0	μg/L							
2,3-Trichloropropane	ND	2.0	μg/L							
1,2-Trichloro-1,2,2-trifluoroethane (Freon	ND	1.0	μg/L							
2,4-Trimethylbenzene	ND	1.0	μg/L							
3,5-Trimethylbenzene	ND	1.0	μg/L							
nyl Chloride	ND	2.0	μg/L							
+p Xylene	ND	2.0	$\mu g/L$							
Xylene	ND	1.0	$\mu g \! / \! L$							
rrogate: 1,2-Dichloroethane-d4	23.6		μg/L	25.0		94.4	70-130			
rrogate: Toluene-d8	24.4		$\mu g/L$	25.0		97.4	70-130			
rrogate: 4-Bromofluorobenzene	24.4		$\mu g/L$	25.0		97.4	70-130			
CS (B245122-BS1)				Prepared: 11	/05/19 Analy	zed: 11/06/1	9			
	88.7	50	μg/L	100		88.7	70-160			
eetone										
		5.0	μg/L	10.0		83.0	70-130			
erylonitrile	8.30		$\mu g/L$	10.0 10.0		83.0 94.8	70-130 70-130			
erylonitrile t-Amyl Methyl Ether (TAME)	8.30 9.48	5.0	μg/L μg/L			94.8	70-130			
erylonitrile rt-Amyl Methyl Ether (TAME) enzene	8.30 9.48 10.7	5.0 0.50	μg/L μg/L μg/L	10.0						
erylonitrile rt-Amyl Methyl Ether (TAME) enzene omobenzene	8.30 9.48 10.7 10.4	5.0 0.50 1.0 1.0	μg/L μg/L μg/L μg/L	10.0 10.0 10.0		94.8 107 104	70-130 70-130 70-130			
erylonitrile et-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane	8.30 9.48 10.7 10.4	5.0 0.50 1.0 1.0	μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0		94.8 107 104 104	70-130 70-130 70-130 70-130			
erylonitrile et-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane	8.30 9.48 10.7 10.4 10.4	5.0 0.50 1.0 1.0 1.0 0.50	μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105	70-130 70-130 70-130 70-130 70-130			
erylonitrile t-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform	8.30 9.48 10.7 10.4 10.4 10.5 9.77	5.0 0.50 1.0 1.0 1.0 0.50	μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7	70-130 70-130 70-130 70-130 70-130 70-130			V-20
erylonitrile t-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane	8.30 9.48 10.7 10.4 10.5 9.77 9.88	5.0 0.50 1.0 1.0 1.0 0.50 1.0	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8	70-130 70-130 70-130 70-130 70-130 70-130 40-160			V-20
erylonitrile tt-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK)	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5	5.0 0.50 1.0 1.0 0.50 1.0 2.0	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160			V-20
erylonitrile rt-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) rt-Butyl Alcohol (TBA)	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3	5.0 0.50 1.0 1.0 0.50 1.0 2.0 20	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160			V-20
erylonitrile tt-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) tt-Butyl Alcohol (TBA) Butylbenzene	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 20 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130			V-20
rylonitrile t-Amyl Methyl Ether (TAME) nzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) t-Butyl Alcohol (TBA) Butylbenzene c-Butylbenzene	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1	5.0 0.50 1.0 1.0 0.50 1.0 2.0 20 20 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 100 10		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130			V-20
rylonitrile t-Amyl Methyl Ether (TAME) nzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) t-Butyl Alcohol (TBA) Butylbenzene t-Butylbenzene t-Butylbenzene	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130			V-20
erylonitrile t-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) t-Butyl Alcohol (TBA) Butylbenzene e-Butylbenzene t-Butylbenzene t-Butylbenzene t-Butyl Ether (TBEE)	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-130 70-130 70-130 70-130			V-20
erylonitrile t-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) t-Butyl Alcohol (TBA) Butylbenzene e-Butylbenzene t-Butylbenzene t-Butyl Ethyl Ether (TBEE) rbon Disulfide	8.30 9.48 10.7 10.4 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24 11.8	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 5.0 5.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4 118	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130			V-20
erylonitrile tt-Amyl Methyl Ether (TAME) enzene omobenzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) tt-Butyl Alcohol (TBA) Butylbenzene tt-Butylbenzene tt-Butyl Ethyl Ether (TBEE) rbon Disulfide rbon Tetrachloride	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 20 1.0 1.0 0.50 5.0 1.0	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130			V-20
erylonitrile tt-Amyl Methyl Ether (TAME) enzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) tt-Butyl Alcohol (TBA) Butylbenzene tt-Butylbenzene tt-Butyl Ethyl Ether (TBEE) rbon Disulfide rbon Tetrachloride allorobenzene	8.30 9.48 10.7 10.4 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24 11.8	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 20 1.0 1.0 0.50 5.0 1.0	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4 118	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130			V-20
erylonitrile tt-Amyl Methyl Ether (TAME) enzene omochloromethane omodichloromethane omoform omomethane Butanone (MEK) tt-Butyl Alcohol (TBA) Butylbenzene tt-Butylbenzene tt-Butyl Ethyl Ether (TBEE) rbon Disulfide rbon Tetrachloride allorobenzene	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24 11.8 10.3	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 20 1.0 1.0 0.50 5.0 1.0	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4 118 103	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130			V-20
erylonitrile tt-Amyl Methyl Ether (TAME) enzene comobenzene comochloromethane comodichloromethane comodichloromethane comomethane Butanone (MEK) tt-Butyl Alcohol (TBA) Butylbenzene ct-Butylbenzene tt-Butyl Ethyl Ether (TBEE) arbon Disulfide arbon Tetrachloride allorobenzene thlorodibromomethane	8.30 9.48 10.7 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24 11.8 10.3 11.4	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 20 1.0 1.0 0.50 5.0 1.0	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4 118 103 114	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130			V-20
cetone crylonitrile rt-Amyl Methyl Ether (TAME) enzene romobenzene romochloromethane romodichloromethane romomethane Butanone (MEK) rt-Butyl Alcohol (TBA) Butylbenzene rc-Butylbenzene rt-Butyl Ethyl Ether (TBEE) arbon Disulfide arbon Tetrachloride hlorodibromomethane hloroethane	8.30 9.48 10.7 10.4 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24 11.8 10.3 11.4 10.0	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 20 1.0 1.0 0.50 5.0 1.0 0.50	Hg/L Hg/L Hg/L Hg/L Hg/L Hg/L Hg/L Hg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4 118 103 114	70-130 70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130 70-130			V-20
crylonitrile rt-Amyl Methyl Ether (TAME) enzene romobenzene romochloromethane romodichloromethane romomethane Butanone (MEK) rt-Butyl Alcohol (TBA) Butylbenzene ec-Butylbenzene rt-Butyl Ethyl Ether (TBEE) arbon Disulfide arbon Tetrachloride hlorobenzene hlorodibromomethane	8.30 9.48 10.7 10.4 10.4 10.5 9.77 9.88 87.5 80.3 9.81 11.1 10.9 9.24 11.8 10.3 11.4 10.0 11.6	5.0 0.50 1.0 1.0 1.0 0.50 1.0 2.0 20 1.0 1.0 1.0 0.50 5.0 1.0 0.50	Hg/L Hg/L Hg/L Hg/L Hg/L Hg/L Hg/L Hg/L	10.0 10.0 10.0 10.0 10.0 10.0 10.0 10.0		94.8 107 104 104 105 97.7 98.8 87.5 80.3 98.1 111 109 92.4 118 103 114 100 116	70-130 70-130 70-130 70-130 70-130 40-160 40-160 40-160 70-130 70-130 70-130 70-130 70-130 70-130 70-130 70-130			V-20 V-20



QUALITY CONTROL

Spike

Source

%REC

RPD

Volatile Organic Compounds by GC/MS - Quality Control

Reporting

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result %REC	%REC Limits	RPD	RPD Limit	Notes
Batch B245122 - SW-846 5030B									
LCS (B245122-BS1)				Prepared: 11	/05/19 Analyzed: 11/06/	19			
4-Chlorotoluene	10.5	1.0	μg/L	10.0	105	70-130			
1,2-Dibromo-3-chloropropane (DBCP)	9.08	5.0	μg/L	10.0	90.8	70-130			
1,2-Dibromoethane (EDB)	10.8	0.50	μg/L	10.0	108	70-130			
Dibromomethane	10.4	1.0	μg/L	10.0	104	70-130			
1,2-Dichlorobenzene	11.0	1.0	μg/L	10.0	110	70-130			
1,3-Dichlorobenzene	11.5	1.0	$\mu g/L$	10.0	115	70-130			
1,4-Dichlorobenzene	11.2	1.0	μg/L	10.0	112	70-130			
trans-1,4-Dichloro-2-butene	7.67	2.0	$\mu g/L$	10.0	76.7	70-130			
Dichlorodifluoromethane (Freon 12)	8.34	2.0	$\mu g/L$	10.0	83.4	40-160			
1,1-Dichloroethane	10.7	1.0	$\mu g/L$	10.0	107	70-130			
1,2-Dichloroethane	9.98	1.0	$\mu g/L$	10.0	99.8	70-130			
1,1-Dichloroethylene	11.0	1.0	$\mu g/L$	10.0	110	70-130			
cis-1,2-Dichloroethylene	10.6	1.0	$\mu g/L$	10.0	106	70-130			
trans-1,2-Dichloroethylene	10.5	1.0	$\mu g/L$	10.0	105	70-130			
1,2-Dichloropropane	10.8	1.0	$\mu g/L$	10.0	108	70-130			
1,3-Dichloropropane	10.3	0.50	$\mu g/L$	10.0	103	70-130			
2,2-Dichloropropane	8.49	1.0	$\mu g/L$	10.0	84.9	40-130			
1,1-Dichloropropene	10.5	2.0	μg/L	10.0	105	70-130			
cis-1,3-Dichloropropene	10.0	0.50	μg/L	10.0	100	70-130			
trans-1,3-Dichloropropene	9.41	0.50	μg/L	10.0	94.1	70-130			
Diethyl Ether	10.8	2.0	μg/L	10.0	108	70-130			
Diisopropyl Ether (DIPE)	10.1	0.50	μg/L	10.0	101	70-130			
1,4-Dioxane	86.6	50	μg/L	100	86.6	40-130			
Ethylbenzene	11.0	1.0	μg/L	10.0	110	70-130			
Hexachlorobutadiene	9.83	0.60	μg/L	10.0	98.3	70-130			
2-Hexanone (MBK)	84.3	10	μg/L	100	84.3	70-160			
Isopropylbenzene (Cumene)	11.1	1.0	μg/L	10.0	111	70-130			
p-Isopropyltoluene (p-Cymene)	10.7	1.0	μg/L	10.0	107	70-130			
Methyl tert-Butyl Ether (MTBE)	10.3	1.0	μg/L	10.0	103	70-130			
Methylene Chloride	11.0	5.0	μg/L	10.0	110	70-130			
4-Methyl-2-pentanone (MIBK)	86.6	10	μg/L	100	86.6	70-160			
Naphthalene	7.64	2.0	μg/L	10.0	76.4	40-130			
n-Propylbenzene	10.6	1.0	μg/L	10.0	106	70-130			
Styrene	10.9	1.0	μg/L	10.0	109	70-130			
1,1,1,2-Tetrachloroethane	11.3	1.0	μg/L	10.0	113	70-130			
1,1,2,2-Tetrachloroethane	10.8	0.50	μg/L	10.0	108	70-130			
Tetrachloroethylene	11.5	1.0	μg/L μg/L	10.0	115	70-130			
Tetrahydrofuran	9.60	10	μg/L	10.0	96.0	70-130			J
Toluene	10.9	1.0	μg/L	10.0	109	70-130			3
1,2,3-Trichlorobenzene	8.06	5.0	μg/L μg/L	10.0	80.6	70-130			
1,2,4-Trichlorobenzene	8.70	1.0	μg/L μg/L	10.0	87.0	70-130			
1,3,5-Trichlorobenzene	9.18	1.0	μg/L μg/L	10.0	91.8	70-130			
1,1,1-Trichloroethane		1.0	μg/L μg/L	10.0	104	70-130			
1,1,2-Trichloroethane	10.4	1.0	μg/L μg/L	10.0	110	70-130			
Trichloroethylene	11.0 11.1	1.0	μg/L μg/L	10.0	111	70-130			
Trichlorofluoromethane (Freon 11)		2.0	μg/L μg/L	10.0	96.1	70-130			
1,2,3-Trichloropropane	9.61	2.0	μg/L μg/L	10.0	103	70-130			
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	10.3 11.5	1.0	μg/L μg/L	10.0	103	70-130 70-130			
1,2,4-Trimethylbenzene	10.8	1.0	μg/L	10.0	108	70-130			
1,3,5-Trimethylbenzene	10.4	1.0	μg/L	10.0	104	70-130			
Vinyl Chloride	9.58	2.0	μg/L	10.0	95.8	40-160			



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	
Batch B245122 - SW-846 5030B											
LCS (B245122-BS1)				Prepared: 11	/05/19 Anal	yzed: 11/06/1	9				
m+p Xylene	21.3	2.0	μg/L	20.0		106	70-130				
o-Xylene	11.1	1.0	$\mu g/L$	10.0		111	70-130				
Surrogate: 1,2-Dichloroethane-d4	23.4		μg/L	25.0		93.6	70-130				
Surrogate: Toluene-d8	25.0		μg/L	25.0		99.9	70-130				
Surrogate: 4-Bromofluorobenzene	24.6		μg/L	25.0		98.4	70-130				
LCS Dup (B245122-BSD1)			10		/05/19 Analy						
Acetone	01.2	50	μg/L	100	703/17 / Mai			2.72	25		
Acrylonitrile	91.2	5.0	μg/L μg/L	10.0		91.2 90.1	70-160 70-130	2.72 8.20	25		
tert-Amyl Methyl Ether (TAME)	9.01	0.50	μg/L μg/L								
Benzene	9.41	1.0		10.0 10.0		94.1 104	70-130 70-130	0.741 3.23	25 25		
Bromobenzene	10.4	1.0	μg/L μg/L								
Bromochloromethane	10.0			10.0		100	70-130	4.00	25		
Bromochloromethane Bromodichloromethane	10.4	1.0	μg/L	10.0		104	70-130	0.865	25		
	10.2	0.50	μg/L	10.0		102	70-130	3.68	25		
Bromoform	10.2	1.0	μg/L	10.0		102	70-130	3.91	25	***	
Bromomethane	9.76	2.0	μg/L	10.0		97.6	40-160	1.22	25	V-20	
2-Butanone (MEK)	89.8	20	μg/L	100		89.8	40-160	2.63	25		
tert-Butyl Alcohol (TBA)	83.6	20	μg/L	100		83.6	40-160	4.00	25		
n-Butylbenzene	9.65	1.0	μg/L	10.0		96.5	70-130	1.64	25		
sec-Butylbenzene	11.0	1.0	μg/L	10.0		110	70-130	1.18	25		
tert-Butylbenzene	10.8	1.0	μg/L	10.0		108	70-130	1.11	25		
tert-Butyl Ethyl Ether (TBEE)	9.17	0.50	μg/L	10.0		91.7	70-130	0.760	25		
Carbon Disulfide	10.8	5.0	μg/L	10.0		108	70-130	8.59	25		
Carbon Tetrachloride	10.2	1.0	μg/L	10.0		102	70-130	1.46	25		
Chlorobenzene	11.5	1.0	μg/L	10.0		115	70-130	0.610	25		
Chlorodibromomethane	9.73	0.50	μg/L	10.0		97.3	70-130	2.84	25		
Chloroethane	10.8	2.0	μg/L	10.0		108	70-130	7.24	25		
Chloroform	10.1	2.0	μg/L	10.0		101	70-130	2.53	25		
Chloromethane	8.77	2.0	μg/L	10.0		87.7	40-160	2.81	25	V-20	
2-Chlorotoluene	10.7	1.0	$\mu g/L$	10.0		107	70-130	0.845	25		
4-Chlorotoluene	10.4	1.0	$\mu g/L$	10.0		104	70-130	1.34	25		
1,2-Dibromo-3-chloropropane (DBCP)	8.90	5.0	$\mu g/L$	10.0		89.0	70-130	2.00	25		
1,2-Dibromoethane (EDB)	10.7	0.50	$\mu g/L$	10.0		107	70-130	1.58	25		
Dibromomethane	10.2	1.0	$\mu g/L$	10.0		102	70-130	2.33	25		
1,2-Dichlorobenzene	11.0	1.0	μg/L	10.0		110	70-130	0.181	25		
1,3-Dichlorobenzene	11.4	1.0	μg/L	10.0		114	70-130	0.872	25		
1,4-Dichlorobenzene	10.9	1.0	μg/L	10.0		109	70-130	3.07	25		
trans-1,4-Dichloro-2-butene	8.71	2.0	$\mu g/L$	10.0		87.1	70-130	12.7	25		
Dichlorodifluoromethane (Freon 12)	8.23	2.0	μg/L	10.0		82.3	40-160	1.33	25		
1,1-Dichloroethane	10.2	1.0	μg/L	10.0		102	70-130	4.40	25		
1,2-Dichloroethane	9.91	1.0	μg/L	10.0		99.1	70-130	0.704	25		
1,1-Dichloroethylene	10.6	1.0	μg/L	10.0		106	70-130	3.90	25		
cis-1,2-Dichloroethylene	10.1	1.0	μg/L	10.0		101	70-130	5.01	25		
trans-1,2-Dichloroethylene	10.3	1.0	μg/L	10.0		103	70-130	2.50	25		
1,2-Dichloropropane	10.8	1.0	μg/L	10.0		108	70-130	0.649	25		
1,3-Dichloropropane	10.4	0.50	μg/L	10.0		104	70-130	0.870	25		
2,2-Dichloropropane	8.15	1.0	μg/L	10.0		81.5	40-130	4.09	25		
1,1-Dichloropropene	10.4	2.0	μg/L	10.0		104	70-130	1.06	25		
cis-1,3-Dichloropropene	9.73	0.50	μg/L μg/L	10.0		97.3	70-130	3.14	25		
trans-1,3-Dichloropropene	9.73	0.50	μg/L μg/L	10.0		96.6	70-130	2.62	25		
Diethyl Ether	10.3	2.0	μg/L μg/L	10.0		103	70-130	4.65	25		
Diisopropyl Ether (DIPE)	9.94	0.50	μg/L μg/L	10.0		99.4	70-130	1.99	25		



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	
Batch B245122 - SW-846 5030B											
LCS Dup (B245122-BSD1)				Prepared: 11	/05/19 Analy	yzed: 11/06/1	19				
1,4-Dioxane	92.6	50	μg/L	100		92.6	40-130	6.72	50		 † ‡
Ethylbenzene	10.7	1.0	μg/L	10.0		107	70-130	2.12	25		
Hexachlorobutadiene	10.1	0.60	μg/L	10.0		101	70-130	2.41	25		
2-Hexanone (MBK)	84.4	10	μg/L	100		84.4	70-160	0.0712	25		†
Isopropylbenzene (Cumene)	10.6	1.0	μg/L	10.0		106	70-130	3.96	25		
p-Isopropyltoluene (p-Cymene)	10.5	1.0	μg/L	10.0		105	70-130	2.36	25		
Methyl tert-Butyl Ether (MTBE)	10.2	1.0	μg/L	10.0		102	70-130	0.585	25		
Methylene Chloride	10.6	5.0	μg/L	10.0		106	70-130	3.60	25		
4-Methyl-2-pentanone (MIBK)	90.1	10	μg/L	100		90.1	70-160	4.02	25		†
Naphthalene	8.02	2.0	μg/L	10.0		80.2	40-130	4.85	25		†
n-Propylbenzene	10.3	1.0	μg/L	10.0		103	70-130	2.11	25		
Styrene	10.4	1.0	μg/L	10.0		104	70-130	4.90	25		
1,1,1,2-Tetrachloroethane	11.3	1.0	$\mu g/L$	10.0		113	70-130	0.265	25		
1,1,2,2-Tetrachloroethane	11.1	0.50	μg/L	10.0		111	70-130	3.01	25		
Tetrachloroethylene	11.5	1.0	μg/L	10.0		115	70-130	0.348	25		
Tetrahydrofuran	10.0	10	μg/L	10.0		100	70-130	4.28	25		
Toluene	10.5	1.0	μg/L	10.0		105	70-130	3.82	25		
1,2,3-Trichlorobenzene	7.86	5.0	μg/L	10.0		78.6	70-130	2.51	25		
1,2,4-Trichlorobenzene	8.34	1.0	$\mu g/L$	10.0		83.4	70-130	4.23	25		
1,3,5-Trichlorobenzene	8.42	1.0	$\mu g/L$	10.0		84.2	70-130	8.64	25		
1,1,1-Trichloroethane	9.84	1.0	$\mu g/L$	10.0		98.4	70-130	5.24	25		
1,1,2-Trichloroethane	11.0	1.0	μg/L	10.0		110	70-130	0.364	25		
Trichloroethylene	11.1	1.0	$\mu g/L$	10.0		111	70-130	0.180	25		
Trichlorofluoromethane (Freon 11)	9.30	2.0	$\mu g/L$	10.0		93.0	70-130	3.28	25		
1,2,3-Trichloropropane	10.0	2.0	$\mu g/L$	10.0		100	70-130	2.85	25		
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	11.1	1.0	μg/L	10.0		111	70-130	4.07	25		
1,2,4-Trimethylbenzene	10.7	1.0	$\mu \text{g/L}$	10.0		107	70-130	0.650	25		
1,3,5-Trimethylbenzene	9.97	1.0	$\mu g\!/\!L$	10.0		99.7	70-130	3.93	25		
Vinyl Chloride	9.25	2.0	$\mu g/L$	10.0		92.5	40-160	3.51	25		†
m+p Xylene	21.3	2.0	$\mu g/L$	20.0		106	70-130	0.0470	25		
o-Xylene	11.1	1.0	$\mu g/L$	10.0		111	70-130	0.541	25		
Surrogate: 1,2-Dichloroethane-d4	22.9		μg/L	25.0		91.5	70-130				
Surrogate: Toluene-d8	25.2		μg/L	25.0		101	70-130				
Surrogate: 4-Bromofluorobenzene	24.6		μg/L	25.0		98.5	70-130				



QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch B245268 - SW-846 3510C										
Blank (B245268-BLK1)				Prepared & A	Analyzed: 11	/06/19				
Acenaphthene (SIM)	ND	0.30	μg/L							
Acenaphthylene (SIM)	ND	0.20	$\mu g/L$							
Anthracene (SIM)	ND	0.20	$\mu g/L$							
Benzo(a)anthracene (SIM)	ND	0.050	$\mu g/L$							
Benzo(a)pyrene (SIM)	ND	0.10	$\mu g/L$							
Benzo(b)fluoranthene (SIM)	ND	0.050	$\mu g/L$							
Benzo(g,h,i)perylene (SIM)	ND	0.50	$\mu g/L$							
Benzo(k)fluoranthene (SIM)	ND	0.20	$\mu g/L$							
Chrysene (SIM)	ND	0.20	$\mu g/L$							
Dibenz(a,h)anthracene (SIM)	ND	0.10	$\mu g/L$							
Fluoranthene (SIM)	ND	0.50	$\mu g \! / \! L$							
Fluorene (SIM)	ND	1.0	$\mu g/L$							
ndeno(1,2,3-cd)pyrene (SIM)	ND	0.10	$\mu g/L$							
-Methylnaphthalene (SIM)	ND	1.0	$\mu g/L$							
Taphthalene (SIM)	ND	1.0	$\mu g/L$							
Phenanthrene (SIM)	ND	0.050	$\mu g/L$							
yrene (SIM)	ND	1.0	$\mu \text{g/L}$							
Surrogate: Nitrobenzene-d5	73.7		μg/L	100		73.7	30-130			
urrogate: 2-Fluorobiphenyl	51.1		μg/L	100		51.1	30-130			
urrogate: p-Terphenyl-d14	65.3		$\mu g/L$	100		65.3	30-130			
LCS (B245268-BS1)				Prepared & A	Analyzed: 11	/06/19				
Acenaphthene (SIM)	35.2	6.0	μg/L	50.0		70.4	40-140			
Acenaphthylene (SIM)	35.7	4.0	$\mu g/L$	50.0		71.4	40-140			
Anthracene (SIM)	38.2	4.0	$\mu g/L$	50.0		76.3	40-140			
Benzo(a)anthracene (SIM)	37.0	1.0	$\mu g/L$	50.0		74.1	40-140			
Benzo(a)pyrene (SIM)	37.4	2.0	$\mu g/L$	50.0		74.8	40-140			
Benzo(b)fluoranthene (SIM)	39.6	1.0	$\mu g/L$	50.0		79.2	40-140			
Benzo(g,h,i)perylene (SIM)	40.1	10	μg/L	50.0		80.2	40-140			
Benzo(k)fluoranthene (SIM)	39.8	4.0	$\mu g/L$	50.0		79.6	40-140			
Chrysene (SIM)	35.8	4.0	$\mu g/L$	50.0		71.7	40-140			
Dibenz(a,h)anthracene (SIM)	42.6	2.0	$\mu g/L$	50.0		85.2	40-140			
luoranthene (SIM)	37.3	10	$\mu g/L$	50.0		74.7	40-140			
Fluorene (SIM)	36.3	20	$\mu g/L$	50.0		72.6	40-140			
ndeno(1,2,3-cd)pyrene (SIM)	42.8	2.0	$\mu g/L$	50.0		85.6	40-140			
-Methylnaphthalene (SIM)	34.3	20	$\mu g/L$	50.0		68.7	40-140			
Naphthalene (SIM)	32.7	20	$\mu g/L$	50.0		65.4	40-140			
Phenanthrene (SIM)	35.9	1.0	μg/L	50.0		71.8	40-140			
Pyrene (SIM)	36.4	20	$\mu g/L$	50.0		72.7	40-140			
Surrogate: Nitrobenzene-d5	74.9		μg/L	100		74.9	30-130			
Surrogate: 2-Fluorobiphenyl	56.8		μg/L	100		56.8	30-130			
Surrogate: p-Terphenyl-d14	59.6		μg/L	100		59.6	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
Batch B245268 - SW-846 3510C										
LCS Dup (B245268-BSD1)				Prepared &	Analyzed: 11	/06/19				
Acenaphthene (SIM)	36.1	6.0	μg/L	50.0		72.2	40-140	2.47	20	
Acenaphthylene (SIM)	36.5	4.0	$\mu g/L$	50.0		73.0	40-140	2.22	20	
Anthracene (SIM)	39.5	4.0	$\mu g/L$	50.0		79.0	40-140	3.40	20	
Benzo(a)anthracene (SIM)	38.1	1.0	$\mu g/L$	50.0		76.2	40-140	2.82	20	
Benzo(a)pyrene (SIM)	38.7	2.0	$\mu g/L$	50.0		77.4	40-140	3.47	20	
Benzo(b)fluoranthene (SIM)	41.0	1.0	$\mu g/L$	50.0		82.0	40-140	3.43	20	
Benzo(g,h,i)perylene (SIM)	41.6	10	$\mu g/L$	50.0		83.3	40-140	3.82	20	
Benzo(k)fluoranthene (SIM)	41.8	4.0	$\mu g/L$	50.0		83.6	40-140	4.80	20	
Chrysene (SIM)	37.1	4.0	$\mu g/L$	50.0		74.1	40-140	3.35	20	
Dibenz(a,h)anthracene (SIM)	44.3	2.0	$\mu g/L$	50.0		88.5	40-140	3.78	20	
Fluoranthene (SIM)	38.4	10	$\mu g/L$	50.0		76.8	40-140	2.80	20	
Fluorene (SIM)	37.1	20	$\mu g/L$	50.0		74.2	40-140	2.23	20	
Indeno(1,2,3-cd)pyrene (SIM)	44.5	2.0	$\mu g/L$	50.0		88.9	40-140	3.76	20	
2-Methylnaphthalene (SIM)	35.8	20	$\mu g/L$	50.0		71.7	40-140	4.27	20	
Naphthalene (SIM)	33.0	20	$\mu g/L$	50.0		66.0	40-140	1.04	20	
Phenanthrene (SIM)	37.1	1.0	$\mu g/L$	50.0		74.2	40-140	3.18	20	
Pyrene (SIM)	37.4	20	$\mu g/L$	50.0		74.7	40-140	2.71	20	
Surrogate: Nitrobenzene-d5	77.8		μg/L	100		77.8	30-130			
Surrogate: 2-Fluorobiphenyl	59.4		$\mu g/L$	100		59.4	30-130			
Surrogate: p-Terphenyl-d14	60.7		μg/L	100		60.7	30-130			



FLAG/QUALIFIER SUMMARY

- QC result is outside of established limits.
- † Wide recovery limits established for difficult compound.
- ‡ Wide RPD limits established for difficult compound.
- # Data exceeded client recommended or regulatory level

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).

V-20 Continuing calibration verification (CCV) did not meet method specifications and was biased on the high side.

Data validation is not affected since sample result was "not detected" for this compound.



CERTIFICATIONS

Certified Analyses included in this Report

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



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

Code	Description	Number	Expires
AIHA	AIHA-LAP, LLC - ISO17025:2017	100033	03/1/2020
MA	Massachusetts DEP	M-MA100	06/30/2020
CT	Connecticut Department of Publilc Health	PH-0567	09/30/2021
NY	New York State Department of Health	10899 NELAP	04/1/2020
NH-S	New Hampshire Environmental Lab	2516 NELAP	02/5/2020
RI	Rhode Island Department of Health	LAO00112	12/30/2019
NC	North Carolina Div. of Water Quality	652	12/31/2019
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	07/30/2020
ME	State of Maine	2011028	06/9/2021
VA	Commonwealth of Virginia	460217	12/14/2019
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

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Page		# of Containers	Preservation Code	container code	UksalVaatMateiks vamplas			Selentist of Englastification	Field Filtered	Lab to Filter		1 Matrix Codes:	GW = Ground Water WW = Waste Water	DW = Drinking Water	A = Air N = Soil	SL = Sludge	O = Other (please	define)	² Preservation Codes:	l = lced	H = HCL M = Methanol	N = Nitric Acid	B = Sodium Bisulfate	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)		PCB ONLY	Soxhlet Non Soxhlet	
242017 39 Spruce Street East Longmeadow, MA 01028				7	ANALYSIS REQUESTED		***************************************																			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		Program Information NCA	andfill	IHSB Orphaned Landfill	State Lead		NELAC and AINA-LAP, LLC Accredited	Other	Chromatogram AIHA-LAP, LLC	
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IMPORTANT!

The wildfires are causing hazardous conditions in California. Learn More



411359783302



Delivered Friday 11/01/2019 at 9:02 am



DELIVERED

Signed for by: R.PETRIAS

GET STATUS UPDATES OBTAIN PROOF OF DELIVERY

FROM

Autryville, NC US

TO

East Longmeadow, MA US

Shipment Facts

TRACKING NUMBER

411359783302

SERVICE

FedEx Priority Overnight

WEIGHT

53.4 lbs / 24.22 kgs

DELIVERED TO

Shipping/Receiving

TOTAL PIECES

1

TOTAL SHIPMENT WEIGHT

53.4 lbs / 24.22 kgs

RETURN REASON

TERMS

Third Party

PACKAGING

Your Packaging

SPECIAL HANDLING SECTION

Deliver Weekday

STANDARD TRANSIT

DARD TRANSII

5\

11/01/2019 by 10:30 am

SHIP DATE

3) -

Thu 10/31/2019

ACTUAL DELIVERY

Fri 11/01/2019 9:02 am

Travel History

Local Scan Time

V

Friday, 11/01/2019

9:02 am

East Longmeadow, MA

Delivered

7:45 am

WINDSOR LOCKS, CT

On FedEx vehicle for delivery

6:27 am

EAST GRANBY, CT

At destination sort facility

Page 22 of 23

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



						Doc# 27	7 Rev 5 2017	.	
Login S	•	ceipt Checklist -	• -		-			ny False	
	Statem	ent will be broug	ght to the at	ttention of	the Client	- State True	or False		
Client	StMI	2			1 \				
Receiv	ed By	INP		Date	191	۱۹	Time	902	
How were th		In Cooler		No Cooler	\ '	On Ice	7	No Ice	
receiv	/ed?	Direct from Samp	oling			Ambient		Melted Ice	
Were samp	oles within		By Gun #	2		Actual Tem	p- 4.1		
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ls COC in in	k/ Legible?	1	·	Were san	nples rece	ived within h	olding time?	T	_
Did COC is	nclude all	Client		Analysis	1	Sampl	er Name	7	<u> </u>
pertinent Inf	formation?	Project		ID's		Collection	Dates/Times		_
Are Sample	labels filled	l out and legible?		•					
Are there La	b to Filters?	1	<u> É </u>		Who wa	s notified?			_
Are there Ru	ishes?			•	Who wa	s notified?			<u>.</u>
Are there Sh	ort Holds?		<u> </u>		Who wa	s notified?			
s there enou	_			•		r			
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Proper Media						samples red	quired?	t	•
Mere trip bla					On COC?		_		
Do all sampl	es have the	proper pH?	NA	Acid		•	Base .		•
/ials	#	Containers:	#			#			#
Jnp-		1 Liter Amb.	2	1 Liter			16 oz		
HCL-	3,	500 mL Amb.		500 mL	····		8oz Am		
Meoh-		250 mL Amb.		250 mL			4oz Am		
Bisulfate-		Flashpoint		Col./Ba			2oz Am		
DI-		Other Glass		Other	*******		 	ore	
Thiosulfate- Sulfuric-		SOC Kit Perchlorate		Plasti			Frozen:		
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222				Unused I	Viedia				
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Jnp-		1 Liter Amb.			Plastic		16 oz		
HCL-		500 mL Amb.			Plastic		8oz Am	~~~~~	
Vleoh-		250 mL Amb.			Plastic	 	4oz Am		
Bisulfate-		Col./Bacteria		Flash			2oz Am		
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