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June 18, 2020

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

Reference: Preliminary Site Assessment for the Judith L. Moss Property (Parcel 53) 1710 W Wilson Street Tarboro, Edgecombe County, North Carolina State Project: U-4424 WBS Element 39062.1.2 DAA Project No. 20080204-010103

Dear Mr. Haden:

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

Location and Description

The Judith L. Moss Property (Parcel #53) is located at 1710 W. Wilson Street in Tarboro, Edgecombe County, North Carolina. The property is situated on the west side of W. Wilson Street approximately 500 feet north of the intersection with Simpson Drive (**Figure 1**). The property was a former automotive fueling station, but as of the date of the fieldwork, it was vacant and unused (**Figure 2**).

The structure on site is a single-story brick building with a concrete apron in front of the building. The concrete apron may have been part of a dispenser island, but no visual evidence at ground surface of a UST was noted. The NCDOT requested a Preliminary Site Assessment for the proposed ROW/easement because the property is a former automotive fueling station. The scope of work as defined in the Request for Technical and Cost Proposal was to evaluate the proposed ROW/easement with respect to the presence of known and unknown USTs and assess whether contamination existed in the subsurface within the proposed ROW/easement. An estimate of the quantity of impacted soil was to be provided if impacted soils were encountered.

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

Geophysical Survey

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

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

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

DAA detected several anomalies in the study area. The survey attributed the all but one of the anomalies to visible cultural features or underground utilities. The geophysical data indicated two anomalies that required further investigation with GPR. A magnetic anomaly (Anomaly 1) was located at the northeast corner of the right-of-way/easement, and one anomaly (Anomaly 2) was located about 60 feet to the southwest of Anomaly 1. GPR signatures suggest USTs. Based on the lack of additional information or visual evidence of a tank, DAA classified the anomaly as a possible UST. **Attachment A** presents DAA's detailed report of findings and interpretations.

Site Assessment Activities

On May 18, 2020, DAA mobilized to the site to conduct a Geoprobe[®] direct-push investigation to evaluate subsurface soil and groundwater conditions within the proposed ROW/easement to a depth of 10 to 12 feet below ground surface (ft bgs). DAA advanced three direct-push probes (SB-1 through SB-3) at select locations (**Figure 2**). The soil boring logs are included as **Attachment B**. The borings were located to evaluate the subsurface conditions in the study area (**Attachment C**).

The lithology encountered by the direct-push samples was generally consistent throughout the site. The ground surface was covered by approximately 6 inches of gravel or topsoil. Below this surface cover was a brown to light brown medium- to fine-grained sand, which appeared to be fill to a depth of about 2 to 3 ft bgs. Below this sand was a sandy clay grading into a clay. Neither bedrock nor groundwater was encountered in any of the borings. Each boring was backfilled with a mix of bentonite (swelling clay to seal the boring) and drill cuttings to the surface after completion.

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

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

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

Following completion of the soil sampling, the boring with the highest recorded PID field screen results or indications of potential contamination (odors, staining, etc.) was converted to a temporary groundwater monitoring well using the direct push screen point. No groundwater was noted in the boring during drilling; however, the Geoprobe[®] screen was advanced to 16 feet and a groundwater sample was collected using low-flow techniques.

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

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

Analytical Results

Table 1 summarizes the soil laboratory results for the three soil samples for TPH DRO/GRO, and **Table 2** summarizes the groundwater laboratory results. **Attachment D** presents the complete laboratory reports.

One soil sample, SB-1, contained a detectable DRO concentration of 3.6 milligrams per kilogram (mg/kg). No other sample contained detectable DRO or GRO concentrations. The action levels are 50 mg/kg for GRO and 100 mg/kg for DRO¹. None of the soil samples analyzed for this site contained DRO or GRO concentrations above their respective action levels.

The groundwater analytical results (**Table 2**) indicate the detection of several petroleum compounds. DAA compared these concentrations to the Groundwater Quality Standards established in 15A NCAC 2L (2L Standards). None of the VOC or SVOC compounds detected were present in concentrations above the applicable 2L Standards.

Contaminated Soil Volume Estimate

The UVF analytical results (**Table 1, Figure 3**) of the soil samples collected on May 18, 2020, indicate that none of the soil samples contained DRO or GRO concentrations above applicable action level. Therefore, no estimate of the volume of soil necessitating possible remediation was required for this report.

Conclusions and Recommendations

DAA conducted a Preliminary Site Assessment to evaluate the NCDOT proposed ROW/easement on the Judith L, Moss Property (Parcel #53) located at 1710 W. Wilson Street in Tarboro, Edgecombe County, North Carolina. A geophysical survey conducted at the site indicated the presence of two possible USTs within or near the proposed ROW/easement; however, no visual signs of a UST were noted at ground surface.

Three soil borings and one temporary well screen point were advanced to evaluate the subsurface soil and groundwater conditions within the proposed ROW/easement. None of the soil samples analyzed contained a GRO or DRO concentration above the action level. Groundwater contained several compounds, but none above the 2L Standards.

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

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

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

Sincerely,

DocuSigned by: **Draper Aden Associates** Mike Branson 942B7ACDE09841E... Micha 6/25/2020

Michael W. Branson, P.G. Project Manager

Attachments



Willen D. Never

William D. Newcomb, P.G. Senior Hydrogeologist

TABLES

TABLE 1 SOIL FIELD SCREENING AND ANALYTICAL RESULTS JUDITH L MOSS PROPERTY (PARCEL 46) TARBORO, EDGECOMBE COUNTY, NORTH CAROLINA STATE PROJECT: U-4424 WBS ELEMENT 39062.1.2 DAA PROJECT NO. 20080204-010103

ANALYTICAL RESULTS PID READING SAMPLE ID DEPTH (ft) SAMPLE ID (mg/kg) (ppm) UVF GRO UVF DRO 50 100 Action Level (mg/kg) <0.74 2.5 3.6 0 - 2 SB-1 2 - 4 2.2 SB-1 4 - 6 1.3 6 - 8 1 8 - 10 1.4 0 - 2 4 4.2 2 - 4 3.9 4 - 6 SB-2 4.9 6 - 8 SB-2 < 0.76 < 0.76 8 - 10 2.1 10 - 12 1.8 3.9 0 - 2 2 - 4 4.1 SB-3 4 - 6 4.8 SB-3 < 0.68 <0.68 4.2 6 - 8 8 - 10 4.2

1) ft - feet

2) ppm - parts per million

3) PID - photoionization detector

4) mg/kg - milligrams per kilogram

5) UVF DRO - Diesel range organics by ultraviolet fluorescence (UVF)

6) UVF GRO - Gasoline range organics by UVF

7) Action level for TPH based upon NCDEQ memo *Guidelines for North Carolina Action Limits for Total Petroleum Hydrocarbons* - July 29, 2016. VOC action levels based on Maximum Soil Contaminant Concentrations.

8) Soil samples were collected on May 18, 2020.

GROUND JUDITH L TARBORO, EDGE S ^T W DAA PR	TABLE 2 WATER ANALYTICAL MOSS PROPERTY (P/ COMBE COUNTY, NC FATE PROJECT: U-442 BS ELEMENT 39062.1 OJECT NO. 20080204	RESULTS ARCEL 46) DRTH CAROLINA 24 1.2 -010103
SAMPLE ID		SB-2
Analyte	15A NCAC 2L Standard (µq/L)	Concentration (µg/L)
Vola	atile Organic Compou	inds
Ethylbenzene	600	1.2
n-Propylbenzene	70	1.0
Toluene	600	1.6
1,2,4-Trimethylbenzene	400	1.2
Xylenes (Total)	500	5.5
Semiv	olatile Organic Comp	ounds
2-Methylnaphthalene	30	0.12 J
Naphthalene	6	0.36 J

1) µg/L - micrograms per liter

г

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

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

4) IMAC - Interim Maximum Allowable Concentration

FIGURES







ATTACHMENT A

Geophysical Study For Possible USTs 1710 West Wilson Street Tarboro, North Carolina



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

May 28, 2020

DAA Project Number: 20080204-010203





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

May 28, 2020

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

RE: Geophysical Study for Possible USTs 1710 West Wilson Street, Tarboro, North Carolina Draper Aden Associates Project No. 20080204-010203

Dear Mr. Pilipchuck:

Draper Aden Associates has completed the geophysical study at 1710 West Wilson Street near Tarboro, North Carolina. The objective of this study was to assist in determining if any underground storage tanks (USTs) may be present beneath the study area. To meet this objective, a combination of ground penetrating radar (GPR) and electromagnetic induction (EM) techniques were utilized. The following report documents our methodologies and findings.

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

Sincerely, Draper Aden Associates

Johanna Vaughan, P.G. Geologist

7.02

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



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- Figure 8. Summary of GPR and EM61 Results

1.0 EXECUTIVE SUMMARY

Draper Aden Associates (DAA) was retained by N.C. Department of Transportation to conduct a geophysical study at 1710 West Wilson Street in Tarboro, North Carolina. The objective of this study was to assist in determining if any underground storage tanks (USTs) may be present beneath the study area. To meet this objective, a combination of ground penetrating radar (GPR) and EM61 electromagnetic induction techniques were utilized.

The EM61 data were collected on April 29 and May 4, 2020 in grid fashion with sub-parallel traverses spaced approximately four feet apart, tracked by a GPS unit capable of sub-foot accuracy. The contoured results from the EM61 data reveal the locations of two anomalies which may represent possible USTs, identified as Anomaly 1 and Anomaly 2. Anomaly 1 is located approximately 22 feet southwest of the southwest corner of the former store building. Anomaly 2 is located approximately 23 feet east of the southeast corner of the former store building.

GPR data were collected at the site on April 29, 2020 utilizing the same grid as the EM data collection, with sub-parallel traverses spaced approximately two feet apart, a GPS unit capable of sub-foot accuracy. Analysis of the GPR data in cross-section revealed a limited number of hyperbolic reflectors consistent with possible USTs, located approximately 1.5 feet below the ground surface. These features correlate to Anomalies 1 and 2 identified in the EM61 data. Planview analysis of the GPR data reveals an anomaly in the depth slices between 2.5 and 3.5 feet depth which correlate well with EM61 Anomaly 1. The depth slices from 4.5 to 5.5 feet depth reveal an anomaly that is coincident with EM61 Anomaly 2.

The combined analysis of the EM61 and GPR data reveals generally good correlation between the two methods, with two coincident anomalies observed in the data from each method (Anomalies 1 and 2). It is uncertain if Anomalies 1 and 2 represent USTs, but of the collected geophysical data, these features are the most likely to represent potential USTs.

2.0 INTRODUCTION

Draper Aden Associates (DAA) was retained by N.C. Department of Transportation to conduct a geophysical study at 1710 West Wilson Street in Tarboro, North Carolina (Figure 1). The objective of this study was to assist in determining if any underground storage tanks (USTs) may be present beneath the study area. To meet this objective, a combination of ground penetrating radar (GPR) and electromagnetic induction (EM) techniques were utilized. The following report documents our methodologies and findings.

The tasks involved in this study included:

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

3.0 ELECTROMAGNETIC INDUCTION (EM) STUDY

3.1 EM Field Methods

The instrument used for this investigation was the EM61 manufactured by Geonics, LTD. The EM61 data were collected on April 29 and May 4, 2020 in grid fashion with sub-parallel traverses spaced approximately four feet apart (Figure 2). The distribution of the EM61 data was tracked by a global positioning system (GPS) unit capable of sub-foot accuracy. The collected data were subsequently contoured laterally and analyzed for evidence of any possible USTs.

3.2 EM61 Results

The contoured results from the EM61 data are presented in Figure 3, overlain onto Google Earth aerial imagery. The data reveals a pair of anomalies which may represent possible USTs, identified as Anomaly 1 and Anomaly 2. Anomaly 1 is located approximately 22 feet southwest of the southwest corner of the former store building. Anomaly 2 is located approximately 23 feet east of the southeast corner of the former store building. It is uncertain if these two anomalies represent USTs, but of the collected data, these features are the most likely to possibly represent USTs. Since the EM61 instrument is particularly sensitive to buried metallic materials or objects, we consider it likely that these anomalies are metallic in composition.

4.0 GPR STUDY

4.1 GPR Field Methods

The instrument used for this investigation was the Noggin 250 manufactured by Sensors and Software, Inc. in Ontario, Canada, which utilizes a 250 MHz antenna mounted on a moveable cart. GPR data were collected on April 29, 2020 utilizing the same grid as the EM61 data collection, with sub-parallel traverses spaced approximately two feet apart, tracked by a GPS unit capable of sub-foot accuracy (Figure 4).

4.2 GPR Results

The GPR data were analyzed as vertical cross-sections and as depth slices, or plan-view maps of the GPR response from various depth intervals for evidence of possible USTs. Figure 5 depicts three sample GPR cross-sections from the collected data which contain broad hyperbolic reflectors consistent with possible USTs, located approximately 1.5 feet below the ground surface. These features correlate to Anomalies 1 and 2 identified in the EM61 data.

Figures 6 and 7 illustrate GPR depth slices in 6-inch-thick depth intervals spanning from 1.0 to 6.0 feet depth. An anomaly seen in the depth slices from 2.5 to 3.0 feet depth and 3.0 to 3.5 feet depth correlates well with EM Anomaly 1. The depth slices from 4.5 to 5.0 feet depth and 5.5 to 5.5 feet depth reveal an anomaly that is coincident with EM Anomaly 2. Numerous other areas of elevated GPR response observed throughout the depth slices may represent miscellaneous buried objects, materials, or conditions, such as intermittent clay layers or zones of wet soil. However, due to a general lack of elevated EM response in other areas (beyond the vicinities of Anomalies 1 and 2), these other areas of anomalous GPR response are considered likely to not be metallic in composition.

5.0 CONCLUSIONS

The combined analysis of the EM61 and GPR data reveals generally good correlation between the two methods, with two coincident anomalies observed in the data from each method (Anomalies 1 and 2).

It is uncertain if Anomalies 1 and 2 represent USTs, but of the collected geophysical data, these features are the most likely to represent potential USTs. Since the EM61 instrument is particularly sensitive to buried metallic materials or objects, we consider it likely that Anomalies 1 and 2 are metallic in composition. Furthermore, the locations of Anomalies 1 and 2 were characterized in profile view of the GPR data as broad hyperbolic reflectors, which is consistent with possible USTs. The results of the study are summarized in Figure 8, depicting the locations of Anomalies 1 and 2.

6.0 LIMITATIONS

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

7.0 FIGURES



Draper Aden Associates
Engineering • Surveying • Environmental Services

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

Richmond, VA Raleigh Charlottesville, VA Fayette Hampton Roads, VA Norther Virginia

Raleigh, NC Fayetteville, NC Northern Virginia Virginia Beach, VA DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE: 05/15/2020 FIGURE 1











Google Earth 2020 Coogle	R C C C C C C C C C C C C C C C C C C C	Earth 20 coogle	50 ft
GPR Depth Slices from 1.0 Foot to 4.0 Feet De	Geophysical Study for Possible USTs 1710 West Wilson Street, Tarboro, NC Pth	PROJECT: 200802	204-010203
Draper Aden Associates Engineering • Surveying • Environmental Services 2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291	DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE: 05/15/2020		FIGURE



GPR Depth Slices from 4.0 Feet to 6.0 Feet De	Geophysical Study for Possible USTs 1710 West Wilson Street, Tarboro, NC	PROJECT: 200802	204-010203
Draper Aden Associates Engineering • Surveying • Environmental Services 2206 South Main Street Blacksburg, VA 24060 540-552-0444 Fax: 540-552-0291	DESIGNED: JMV DRAWN: JMV CHECKED: CMP DATE: 05/15/2020		figure 7



ATTACHMENT B



		aper .	Ade	en Ass g • Envirom	Sociate mental Service	es es		BORING I	D: SB-2
PRO		IAME:	Tarbo	oro PSA			PROJECT NUMBER: 200802	04-010103	
CLIE	NT: NC	DOT-	John	Pilipchul	k, PE		DATE: 5/18/2020		
SITE	LOCA	TION: 1	1710	W. Wilso	on St., Ta	arboro, NC	TOTAL DEPTH (ft bgs): 8		
DRIL	LING C	ONTR	ACT	OR: DAA	A- Sean J	larvah	BORING COORDINATES:		
DRIL	LING N	/IETHO	D: Di	irect Pus	h		BOREHOLE DIAMETER (inch	es): 2	
DRIL	LING E	QUIPE	MEN	IT: Geop	robe		DEPTH TO WATER (ft bgs): N	NE	
LOG	GED B	Y: Brar	ndy B	arnes			PROJECT MANAGER: Mike E	Branson, PG	
	SAM	IPLES	(m						
DEPTH (ft bgs)	Sample ID	Recovery %	PID Reading (pl	US	CS	LITI	HOLOGIC DESCRIPTION:		Notes:
		50	4	SP		POORLY GRADED) SAND, Light brown to brown with d lium grained, upper one foot to two	orange, feet is fill	
2— 		50	4	- SC		CLAYEY SAND, Da medium grained	ark brown to black, with gray, fine-g	rained to	
5	- - - - - -		4			Wet, with some ora	ange and staining		
6— 7—	SB-2	70	5						
				-					
9		100	2	CH/CL		CLAY, Gray and or course-grained clay	ange, with seams of medium-graine yey sand	ed to	
11			2						
12-					//////	Dark reduish brown	rand orange		
						End of Borehol	e at 12 feet		
									Page 1 of 1

Draper Aden Associates BORING ID: SB-3 Engineering • Surveying • Environmental Services PROJECT NAME: Tarboro PSA PROJECT NUMBER: 20080204-010103 CLIENT: NCDOT- John Pilipchuk, PE DATE: 5/18/2020 SITE LOCATION: 1710 W. Wilson St., Tarboro, NC TOTAL DEPTH (ft bgs): 8 DRILLING CONTRACTOR: DAA- Sean Jarvah BORING COORDINATES: **DRILLING METHOD: Direct Push** BOREHOLE DIAMETER (inches): 2 DRILLING EQUIPEMENT: Geoprobe DEPTH TO WATER (ft bgs): NE PROJECT MANAGER: Mike Branson, PG LOGGED BY: Brandy Barnes SAMPLES PID Reading (ppm) DEPTH (ft bgs) Notes: Sample ID Recovery USCS LITHOLOGIC DESCRIPTION: 0 **GRAVEL FILL** 1 4 POORLY GRADED SAND, Light brown, fine-grained to medium-SP grained 2 40 CLAYEY SAND, Light brown to brown with light gray gravel and 3 4 SC burned debris 4 5 SB-3 5 70 6 CLAY, Light gray, tan, and orange, with dark brown, fine-grained CH/CL 7 4 to medium-grained, mottled 8 100 9 4 10 End of Borehole at 10 feet

ATTACHMENT C



PHOTOGRAPH 1. Location of SB-1.



PHOTOGRAPH 2. Location of SB-2, no photograph of SB-3..



Tarboro Preliminary Site Assessment Parcel 53 1710 W. Wilson Street, Tarboro, NC DAA PN: 20080204-010103 ATTACHMENT D

Q	ED		E				B					\int	<u>QROS</u>		
				Hydroca	arbon An	alysis Ro	esults								
Client: Address:	DAA 114 Edinburgh S. Dr. Cary, NC 27511								Sa Sample Sampl	mples es extr es ana	taken acted lysed		Monday, May 18, 2020 Monday, May 18, 2020 Wednesday, May 20, 2020		
Contact:	Michael Branson									One	erator		Harry Wooten		
Project:	Tarboro PSA 20080204-010103														
													F03640		
Matrix	Sample ID	Dilution used	BTEX (C6 - C9)	GRO (C5 - C10)	DRO (C10 - C35)	TPH (C5 - C35)	Total Aromatics	16 EPA PAHs	BaP	Ratios		Ratios			HC Fingerprint Match
							(010-035)			% light	% mid	% heavy			
S	SB-1	29.7	<0.74	<0.74	3.6	3.6	1.6	<0.24	<0.03	0	84.6	15.4	Deg.Fuel 88.8%,(FCM)		
s	SB-2	30.6	<0.76	<0.76	<0.76	<0.76	<0.15	<0.24	<0.031	0	0	0	PHC not detected,(BO)		
S	SB-3	27.2	<0.68	<0.68	<0.68	<0.68	<0.14	<0.22	<0.027	0	0	0	PHC not detected,(BO)		
Deculto gon	Initial C		QC check	OK	e and mg/l fo	r water comp		oo oro pot d	Final FC	CM QC	Check	OK	100.9 %		
Fingerprints (SBS) or (LE	provide a tentative hydrocarbon identificatio 3S) = Site Specific or Library Background Su	n. The abbr	eviations are pplied to resi	e:- FCM = Re ult : (PFM) = F	esults calculat Poor Fingerpri	ed using Fund int Match : (T)	damental Calib = Turbid : (P)	ration Mode Particulat	e : % = confi e present	dence fo	r sample	e conter e fingerp	print match to library		

QED Hydrocarbon Fingerprints





		Relinquished by	And nee	Relinquished by	COMMENTS/REQUESTS:							- B.	1 				1325 X	5116172 1350 X	Date/Time 24 Hour 48 Hou	Sample Collection TAT Requested	Collected by: Browy	Phone #:	Email: whoren shed	Project Ref .: Tabaro PS,	Contact: M. Brons C.	Audress: Covy, NC 3	Ily Edinburg	Client Name: IDAA
			021813	1200											>	<u></u>	X	×	IT UVF	Analysis	S CHAIN			VINGOR V		E	ans.or	
P		Accepted by	Okan ar	Accepted by	TARGET GC/UVF			~						- 1.5 		222 02-2	123 SA	RR CR-	GC Initials	s Type	OF CUSTODY AND ANALY	RAPID ENVIRONMENTAL D						Durilaz
		Date/Time	7. 14. 10.	Date/Time	ANALYTES:														Sample ID		TICAL REQUEST FORM	NAGNOSTICS				TM	Ū	
	Ref. No HOV			RED Lab USE ONLY	-3°	1.1										E2 2 40 1 2 . 5	5 A 42.8 7 2 2	NET UTA NIS	Total Wt. Tare Wt. Sample Wt.		Solvents: VC, 1,1 DCE, 1,2 cis DCE, 1,2 trans DCE, TCE, and PCE. Specify target analytes in the space provided below.	Analyses are for BTEX and Chlorinated	total BTEX, GRO, DRO, TPH, PAH total	Each UVF sample will be analyzed for	Wilmington, NC 28409	MARBIONC Bldg, Suite 2003	5598 Marvin K Moss Lane	RED Lab, LLC

*

10 × 1

5



May 28, 2020

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

Project Location: Tarboro, NC - Parcel 53 Client Job Number: Project Number: 20080204-010103 Laboratory Work Order Number: 20E0838

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

Sincerely,

Beny K. Millee

Kerry K. McGee Project Manager

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B258487	12
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Draper Aden Associates 114 Edinburgh Drive South, Suite 200 Cary, NC 27519 ATTN: Mike Branson

REPORT DATE: 5/28/2020

SUB LAB

PURCHASE ORDER NUMBER:

OKCHASE OKDER NOMBER.

PROJECT NUMBER: 20080204-010103

ANALYTICAL SUMMARY

WORK ORDER NUMBER: 20E0838

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

MATRIX

PROJECT LOCATION: Tarboro, NC - Parcel 53

FIELD SAMPLE #

SB-2

20E0838-01 Ground W

LAB ID:

SAMPLE DESCRIPTION

TEST SW-846 8260D

Ground Water

SW-846 8200D

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CASE NARRATIVE SUMMARY

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



SW-846 8260D

Qualifications:

L-04

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

Trichlorofluoromethane (Freon 11)

20E0838-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1

V-05

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

Analyte & Samples(s) Qualified:

1.1.1.2-Tetrachloroethane

20E0838-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

Bromoform

20E0838-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

Dichlorodifluoromethane (Freon 12

20E0838-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

tert-Butyl Alcohol (TBA)

20E0838-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

Trichlorofluoromethane (Freon 11)

20E0838-01[SB-2], B258487-BLK1, B258487-BS1, B258487-BSD1, S048715-CCV1

SW-846 8270E

Qualifications:

V-04

Initial calibration did not meet method specifications. Compound was calibrated using a response factor where %RSD is outside of method specified criteria. Reported result is estimated. Analyte & Samples(s) Qualified:

Benzidine

20E0838-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

V-05

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

Analyte & Samples(s) Qualified:

Aniline

20E0838-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

Benzidine 20E0838-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

V-06

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

Analyte & Samples(s) Qualified:

Benzoic Acid

20E0838-01[SB-2], S048793-CCV1

V-34

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

estimated. Analyte & Samples(s) Qualified:

4-Chloroaniline

20E0838-01[SB-2], B258532-BLK1, B258532-BS1, B258532-BSD1, S048748-CCV1, S048793-CCV1

Aniline

20E0838-01[SB-2], S048793-CCV1



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

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

Kappenne

Tod E. Kopyscinski Laboratory Director



Volatile Organic Compounds by GC/MS

Sample Description:

Sampled: 5/18/2020 14:40

Project Location: Tarboro, NC - Parcel 53 Date Received: 5/19/2020

Field Sample #: SB-2

Sample ID: 20E0838-01

Sample Matrix: Ground Water

								Date	Date/Time	
Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Prepared	Analyzed	Analyst
Acetone	ND	50	3.8	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Acrylonitrile	ND	5.0	0.52	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
tert-Amyl Methyl Ether (TAME)	ND	0.50	0.14	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Benzene	ND	1.0	0.18	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Bromobenzene	ND	1.0	0.15	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Bromochloromethane	ND	1.0	0.32	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Bromodichloromethane	ND	0.50	0.16	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Bromoform	ND	1.0	0.46	$\mu g/L$	1	V-05	SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Bromomethane	ND	2.0	1.4	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
2-Butanone (MEK)	ND	20	1.9	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
tert-Butyl Alcohol (TBA)	ND	20	4.2	$\mu g/L$	1	V-05	SW-846 8260D	5/21/20	5/21/20 12:19	MFF
n-Butylbenzene	ND	1.0	0.21	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
sec-Butylbenzene	ND	1.0	0.16	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
tert-Butylbenzene	ND	1.0	0.17	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
tert-Butyl Ethyl Ether (TBEE)	ND	0.50	0.16	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Carbon Disulfide	ND	5.0	4.4	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Carbon Tetrachloride	ND	1.0	0.11	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Chlorobenzene	ND	1.0	0.15	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Chlorodibromomethane	ND	0.50	0.21	$\mu g/L$	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Chloroethane	ND	2.0	0.36	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Chloroform	ND	2.0	0.17	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Chloromethane	ND	2.0	0.45	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
2-Chlorotoluene	ND	1.0	0.12	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
4-Chlorotoluene	ND	1.0	0.14	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,2-Dibromo-3-chloropropane (DBCP)	ND	5.0	0.53	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,2-Dibromoethane (EDB)	ND	0.50	0.19	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Dibromomethane	ND	1.0	0.37	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,2-Dichlorobenzene	ND	1.0	0.16	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,3-Dichlorobenzene	ND	1.0	0.12	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,4-Dichlorobenzene	ND	1.0	0.13	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
trans-1,4-Dichloro-2-butene	ND	2.0	0.31	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Dichlorodifluoromethane (Freon 12)	ND	2.0	0.26	μg/L	1	V-05	SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,1-Dichloroethane	ND	1.0	0.16	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,2-Dichloroethane	ND	1.0	0.41	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,1-Dichloroethylene	ND	1.0	0.32	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
cis-1,2-Dichloroethylene	ND	1.0	0.13	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
trans-1,2-Dichloroethylene	ND	1.0	0.31	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,2-Dichloropropane	ND	1.0	0.20	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,3-Dichloropropane	ND	0.50	0.11	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
2,2-Dichloropropane	ND	1.0	0.20	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,1-Dichloropropene	ND	2.0	0.16	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
cis-1,3-Dichloropropene	ND	0.50	0.13	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
trans-1,3-Dichloropropene	ND	0.50	0.23	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Diethyl Ether	ND	2.0	0.34	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF

Work Order: 20E0838

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Volatile Organic Compounds by GC/MS

Work Order: 20E0838

Project Location: Tarboro, NC - Parcel 53 Date Received: 5/19/2020 Field Sample #: SB-2

Sample ID: 20E0838-01 Sample Matrix: Ground Water Sampled: 5/18/2020 14:40

Sample Description:

Analyta	Doculto	DI	ы	Units	Dilution	Flog/Quol	Method	Date Prepared	Date/Time	Analyst
Diisopropyl Ether (DIPE)	ND	0.50	0.17		1	Tiag/Quai	SW 846 8260D	5/21/20	5/21/20 12:10	MEE
1 4-Dioxane	ND	50	22	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Ethylbenzene	1.2	1.0	0.13	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Hexachlorobutadiene	ND	0.60	0.15	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
2-Hexanone (MBK)	ND	10	1.5	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Isopropylbenzene (Cumene)	ND	2.0	0.17	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
p-Isopropyltoluene (p-Cymene)	ND	2.0	0.20	μ <u>σ</u> /L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Methyl tert-Butyl Ether (MTBE)	ND	1.0	0.25	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Methylene Chloride	ND	5.0	0.23	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
4-Methyl-2-pentanone (MIBK)	ND	10	1.7	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Nanhthalene	ND	5.0	0.31	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
n-Pronylhenzene	1.0	1.0	0.13	µg/L	1		SW 846 8260D	5/21/20	5/21/20 12:19	MEE
Styrene	ND	2.0	0.15	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1 1 1 2-Tetrachloroethane	ND	1.0	0.27	μg/L	1	V 05	SW 846 8260D	5/21/20	5/21/20 12:19	MEE
1 1 2 2-Tetrachloroethane	ND	0.50	0.27	μg/L μg/I	1	V-05	SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Tetrachloroethylene	ND	1.0	0.22	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Tetrahydrofuran	ND	10	0.10	μg/L	1		SW 846 8260D	5/21/20	5/21/20 12:19	MEE
Toluene	16	10	0.14	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1 2 3-Trichlorobenzene	ND	5.0	0.57	μ <u>σ</u> /L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1 2 4-Trichlorobenzene	ND	1.0	0.40	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1 3 5-Trichlorobenzene	ND	1.0	0.40	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1 1 1-Trichloroethane	ND	1.0	0.20	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1 1 2-Trichloroethane	ND	1.0	0.16	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Trichloroethylene	ND	1.0	0.10	μg/L μg/I	1		SW-846 8260D	5/21/20	5/21/20 12:19	MEE
Trichlorofluoromethane (Freon 11)	ND	2.0	0.24	μg/L μg/I	1	V-05 I-04	SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1.2.3-Trichloropropane	ND	2.0	0.25	μg/L μg/I	1	V-05, E-04	SW-846 8260D	5/21/20	5/21/20 12:19	MEE
1,1,2-Trichloro-1,2,2-trifluoroethane (Freen 113)	ND	1.0	0.32	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,2,4-Trimethylbenzene	1.2	1.0	0.18	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
1,3,5-Trimethylbenzene	ND	1.0	0.14	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Vinyl Chloride	ND	2.0	0.45	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
m+p Xylene	4.0	2.0	0.30	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
o-Xylene	1.5	1.0	0.17	μg/L	1		SW-846 8260D	5/21/20	5/21/20 12:19	MFF
Surrogates		% Reco	overy	Recovery Limits	5	Flag/Qual				
1,2-Dichloroethane-d4		115		70-130					5/21/20 12:19	
Toluene-d8		97.0		70-130					5/21/20 12:19	
4-Bromotluorobenzene		98.6		/0-130					5/21/20 12:19	



Work Order: 20E0838

Project Location: Tarboro, NC - Parcel 53 Date Received: 5/19/2020

Field Sample #: SB-2

Sample ID: 20E0838-01

Sample Matrix: Ground Water

Sampled: 5/18/2020 14:40

Sample Description:

Semivolatile Organic Compounds by GC/MS													
Analyte	Results	RI.	DL	Units	Dilution	Flag/Qual	Method	Date Prenared	Date/Time Analyzed	Analyst			
Acenaphthene (SIM)	ND	0.31	0.034	ug/L	1	Ting/Quin	SW-846 8270F	5/21/20	5/28/20 2:00	IMR			
Acenaphthylene (SIM)	ND	0.21	0.036	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Acetophenone	ND	10	0.41	ug/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Aniline	ND	5.2	0.77	не/L	1	V-05, V-34	SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Anthracene (SIM)	ND	0.21	0.033	не/L	1	,	SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Benzidine	ND	21	17	μg/L	1	V-04, V-05	SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Benzo(a)anthracene (SIM)	ND	0.052	0.016	μg/L	1	,	SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Benzo(a)pyrene (SIM)	ND	0.10	0.012	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Benzo(b)fluoranthene (SIM)	ND	0.052	0.015	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Benzo(g,h,i)perylene (SIM)	ND	0.52	0.019	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Benzo(k)fluoranthene (SIM)	ND	0.21	0.012	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Benzoic Acid	ND	10	5.5	μg/L	1	V-06	SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Bis(2-chloroethoxy)methane	ND	10	0.49	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Bis(2-chloroethyl)ether	ND	10	0.53	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Bis(2-chloroisopropyl)ether	ND	10	0.75	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Bis(2-Ethylhexyl)phthalate	ND	10	0.54	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
4-Bromophenylphenylether	ND	10	0.31	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Butylbenzylphthalate	ND	10	0.30	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Carbazole	ND	10	0.29	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
4-Chloroaniline	ND	10	0.35	μg/L	1	V-34	SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
4-Chloro-3-methylphenol	ND	10	0.50	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2-Chloronaphthalene	ND	10	0.47	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2-Chlorophenol	ND	10	0.39	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
4-Chlorophenylphenylether	ND	10	0.32	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Chrysene (SIM)	ND	0.21	0.015	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Dibenz(a,h)anthracene (SIM)	ND	0.10	0.018	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Dibenzofuran	ND	5.2	0.27	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Di-n-butylphthalate	ND	10	0.47	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
1,2-Dichlorobenzene	ND	5.2	0.47	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
1,3-Dichlorobenzene	ND	5.2	0.48	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
1,4-Dichlorobenzene	ND	5.2	0.39	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
3,3-Dichlorobenzidine	ND	10	0.37	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2,4-Dichlorophenol	ND	10	0.31	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Diethylphthalate	ND	10	0.23	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2,4-Dimethylphenol	ND	10	0.82	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Dimethylphthalate	ND	10	0.32	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
4,6-Dinitro-2-methylphenol	ND	10	2.0	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2,4-Dinitrophenol	ND	10	1.7	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2,4-Dinitrotoluene	ND	10	0.34	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
2,6-Dinitrotoluene	ND	10	0.36	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Di-n-octylphthalate	ND	10	0.54	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
1,2-Diphenylhydrazine/Azobenzene	ND	10	0.39	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL			
Fluoranthene (SIM)	ND	0.52	0.026	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			
Fluorene (SIM)	ND	1.0	0.035	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR			

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Work Order: 20E0838

Project Location: Tarboro, NC - Parcel 53 Date Received: 5/19/2020

Sample Description:

Field Sample #: SB-2

Sample ID: 20E0838-01

Sampled: 5/18/2020 14:40

			Semivo	latile Organic Co	mpounds by	GC/MS				
Analyte	Results	RL	DL	Units	Dilution	Flag/Qual	Method	Date Prepared	Date/Time Analyzed	Analyst
Hexachlorobenzene	ND	10	0.45	μg/L	1	-	SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Hexachlorobutadiene	ND	10	0.61	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Hexachlorocyclopentadiene	ND	10	4.9	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Hexachloroethane	ND	10	0.55	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Indeno(1,2,3-cd)pyrene (SIM)	ND	0.10	0.019	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR
Isophorone	ND	10	0.31	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
1-Methylnaphthalene	ND	5.2	0.29	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
2-Methylnaphthalene (SIM)	0.12	1.0	0.064	μg/L	1	J	SW-846 8270E	5/21/20	5/28/20 2:00	IMR
2-Methylphenol	ND	10	0.47	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
3/4-Methylphenol	ND	10	0.21	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Naphthalene (SIM)	0.36	1.0	0.26	μg/L	1	J	SW-846 8270E	5/21/20	5/28/20 2:00	IMR
2-Nitroaniline	ND	10	0.41	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
3-Nitroaniline	ND	10	0.43	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
4-Nitroaniline	ND	10	0.51	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Nitrobenzene	ND	10	0.42	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
2-Nitrophenol	ND	10	0.43	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
4-Nitrophenol	ND	10	0.65	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
N-Nitrosodimethylamine	ND	10	1.9	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
N-Nitrosodiphenylamine/Diphenylamine	ND	10	0.30	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
N-Nitrosodi-n-propylamine	ND	10	0.53	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Pentachloronitrobenzene	ND	10	1.5	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Pentachlorophenol	ND	10	0.34	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Phenanthrene (SIM)	ND	0.052	0.031	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR
Phenol	ND	10	0.20	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Pyrene (SIM)	ND	1.0	0.024	μg/L	1		SW-846 8270E	5/21/20	5/28/20 2:00	IMR
Pyridine	ND	5.2	2.9	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
1,2,4,5-Tetrachlorobenzene	ND	10	0.35	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
1,2,4-Trichlorobenzene	ND	5.2	0.58	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
2,4,5-Trichlorophenol	ND	10	0.49	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
2,4,6-Trichlorophenol	ND	10	0.34	μg/L	1		SW-846 8270E	5/21/20	5/27/20 13:38	BGL
Surrogates		% Reco	very	Recovery Limits	;	Flag/Qual				

2-Fluorophenol	38.2	15-110	5/27/20 13:38
Phenol-d6	28.1	15-110	5/27/20 13:38
Nitrobenzene-d5	70.2	30-130	5/28/20 2:00
Nitrobenzene-d5	62.3	30-130	5/27/20 13:38
2-Fluorobiphenyl	62.7	30-130	5/27/20 13:38
2-Fluorobiphenyl	66.1	30-130	5/28/20 2:00
2,4,6-Tribromophenol	68.5	15-110	5/27/20 13:38
p-Terphenyl-d14	62.9	30-130	5/28/20 2:00
p-Terphenyl-d14	71.0	30-130	5/27/20 13:38



Sample Extraction Data

Prep Method: SW-846 5030B	Analytical Method: SW-846 8260)D			
Lab Number [Field ID]	Ba	atch Initi	ial [mL]	Final [mL]	Date
20E0838-01 [SB-2]	B2	58487 5	5	5.00	05/21/20
Prep Method: SW-846 3510C	Analytical Method: SW-846 827(DE			
Lab Number [Field ID]	Ba	atch Initi	ial [mL]	Final [mL]	Date
20E0838-01 [SB-2]	B2	58532 97	70	1.00	05/21/20
Prep Method: SW-846 3510C	Analytical Method: SW-846 827(DE			
Lab Number [Field ID]	Ba	atch Initi	ial [mL]	Final [mL]	Date
20E0838-01 [SB-2]	B2	58763 97	70	1.00	05/21/20



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 B258487 - SW-846 5030B										
Blank (B258487-BLK1)				Prepared &	Analyzed: 05	/21/20				
Acetone	ND	50	μg/L							
Acrylonitrile	ND	5.0	μg/L							
tert-Amyl Methyl Ether (TAME)	ND	0.50	μg/L							
Benzene	ND	1.0	μg/L							
Bromobenzene	ND	1.0	μg/L							
Bromochloromethane	ND	1.0	μg/L							
Bromodichloromethane	ND	0.50	μg/L							
Bromoform	ND	1.0	μg/L							V-05
Bromomethane	ND	2.0	μg/L							
2-Butanone (MEK)	ND	20	μg/L							
tert-Butyl Alcohol (TBA)	ND	20	μg/L							V-05
n-Butylbenzene	ND	1.0	μg/L							
sec-Butylbenzene	ND	1.0	μg/L							
tert-Butylbenzene	ND	1.0	μg/L							
tert-Butyl Ethyl Ether (TBEE)	ND	0.50	μg/L							
Carbon Disulfide	ND	5.0	μg/L							
Carbon Tetrachloride	ND	1.0	μg/L							
Chlorobenzene	ND	1.0	μg/L							
Chlorodibromomethane	ND	0.50	μg/L							
Chloroethane	ND	2.0	μg/L							
Chloroform	ND	2.0	μg/L							
Chloromethane	ND	2.0	μg/L							
2-Chlorotoluene	ND	1.0	μg/L							
4-Chlorotoluene	ND	1.0	μg/L							
1,2-Dibromo-3-chloropropane (DBCP)	ND	5.0	μg/L							
1,2-Dibromoethane (EDB)	ND	0.50	μg/L							
Dibromomethane	ND	1.0	μg/L							
1,2-Dichlorobenzene	ND	1.0	μg/L							
1,3-Dichlorobenzene	ND	1.0	μg/L							
1,4-Dichlorobenzene	ND	1.0	μg/L							
trans-1,4-Dichloro-2-butene	ND	2.0	μg/L							
Dichlorodifluoromethane (Freon 12)	ND	2.0	μg/L							V-05
1,1-Dichloroethane	ND	1.0	μg/L							
1,2-Dichloroethane	ND	1.0	μg/L							
I,I-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							
2.2 Dichloropropane	ND	0.50	µg/L							
1.1 Disklaranranana	ND	1.0	µg/L							
ais 1.2 Dichleropropene	ND	2.0	µg/L							
trong 1.2 Dichloropropene	ND	0.50	µg/L							
Diethyl Ether	ND	2.0	μg/L μg/I							
Diisonronyl Ether (DIPF)	ND	0.50	μg/L 11σ/Ι							
1 4-Dioxane	ND	50	μg/L μα/Ι							
Fthylhenzene	ND	1.0	μg/L μα/Ι							
Heyachlorobutadiene	ND	0.60	μg/L μα/Ι							
2-Hevanone (MBK)	ND	10	μg/L α/I							
2-HOAMUNE (MDK)	ND	10	μg/L α/I							
n-Isonronyltoluene (n-Cymene)	ND	1.0	μg/L μα/Ι							
Methyl tert-Butyl Ether (MTRE)	ND	1.0	μg/L μσ/Γ							
monification (MTDE)	ND	1.0	μg/ L							



QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	Notes
Batch B258487 - SW-846 5030B										
Blank (B258487-BLK1)				Prepared &	Analyzed: 05	/21/20				
Methylene Chloride	ND	5.0	μg/L							
4-Methyl-2-pentanone (MIBK)	ND	10	$\mu g/L$							
Naphthalene	ND	2.0	μg/L							
n-Propylbenzene	ND	1.0	μg/L							
Styrene	ND	1.0	μg/L							
1,1,1,2-Tetrachloroethane	ND	1.0	μg/L							V-05
1,1,2,2-Tetrachloroethane	ND	0.50	μg/L							
Tetrachloroethylene	ND	1.0	μg/L							
Tetrahydrofuran	ND	10	μg/L							
Toluene	ND	1.0	μg/L							
1,2,3-Trichlorobenzene	ND	5.0	μg/L							
1,2,4-Trichlorobenzene	ND	1.0	μg/L							
1,3,5-Trichlorobenzene	ND	1.0	μg/L							
1,1,1-Trichloroethane	ND	1.0	μg/L							
1,1,2-Trichloroethane	ND	1.0	μg/L							
Trichloroethylene	ND	1.0	μg/L							
Trichlorofluoromethane (Freon 11)	ND	2.0	μg/L							L-04, V-05
1,2,3-Trichloropropane	ND	2.0	μg/L							
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	ND	1.0	μg/L							
1,2,4-Trimethylbenzene	ND	1.0	μg/L							
1,3,5-Trimethylbenzene	ND	1.0	μg/L							
Vinyl Chloride	ND	2.0	μg/L							
m+p Xylene	ND	2.0	μg/L							
o-Xylene	ND	1.0	μg/L							
Surrogate: 1 2-Dichloroethane-d4	27.5		ug/L	25.0		110	70-130			
Surrogate: Toluene-d8	23.6		ug/L	25.0		94.5	70-130			
Surrogate: 4-Bromofluorobenzene	25.2		μg/L	25.0		101	70-130			
LCS (R 258487-RS1)				Prenared &	Analyzed: 05	/21/20				
Acetone	89.2	50	μg/L	100	rinaryzea. 05	89.2	70-160			
Acrylonitrile	11.2	5.0	ug/L	10.0		112	70-130			
tert-Amyl Methyl Ether (TAME)	8 36	0.50	μg/L	10.0		83.6	70-130			
Benzene	10.5	1.0	μ <u>α</u> /L	10.0		105	70-130			
Bromobenzene	10.3	1.0	μ <u>α</u> /L	10.0		103	70-130			
Bromochloromethane	10.5	1.0	μ <u>α</u> /L	10.0		108	70-130			
Bromodichloromethane	8 78	0.50	μ <u>α</u> /L	10.0		87.8	70-130			
Bromoform	8.70	1.0	це/Г.	10.0		88.7	70-130			V-05
Bromomethane	5 25	2.0	μg/L	10.0		52.5	40-160			
2-Butanone (MEK)	110	20	μ <u>α</u> /L	100		110	40-160			
tert-Butyl Alcohol (TBA)	68.8	20	μ <u>α</u> /L	100		68.8	40-160			V-05
n-Butylbenzene	11 7	1.0	μ <u>α</u> /L	10.0		117	70-130			. 00
sec-Butylbenzene	12.0	1.0	це/Г.	10.0		120	70-130			
tert-Butylbenzene	11.0	1.0	не/Г.	10.0		115	70-130			
tert-Butyl Ethyl Ether (TBEE)	0 22	0.50	цø/L	10.0		93.2	70-130			
Carbon Disulfide	9.34	5.0	цø/L	10.0		82.4	70-130			
Carbon Tetrachloride	0.24 8.42	1.0	кв/Е це/Г.	10.0		84 2	70-130			
Chlorobenzene	0.42	1.0	не/Г.	10.0		977	70-130			
Chlorodibromomethane	7.// 0.01	0.50	нø/Г.	10.0		82.1	70-130			
Chloroethane	0.21	2.0	но/L	10.0		74 3	70-130			
Chloroform	10.40	2.0	но/L	10.0		106	70-130			
Chloromethane	10.0	2.0	нь/L 110/L	10.0		61.9	40-160			
2-Chlorotoluene	10.0	2.0	но/L	10.0		109	70-130			



QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	Notes
Batch B258487 - SW-846 5030B										
LCS (B258487-BS1)				Prepared &	Analyzed: 05/	21/20				
4-Chlorotoluene	10.9	1.0	μg/L	10.0		109	70-130			
1,2-Dibromo-3-chloropropane (DBCP)	9.15	5.0	μg/L	10.0		91.5	70-130			
1,2-Dibromoethane (EDB)	9.22	0.50	μg/L	10.0		92.2	70-130			
Dibromomethane	8.99	1.0	μg/L	10.0		89.9	70-130			
1,2-Dichlorobenzene	10.2	1.0	μg/L	10.0		102	70-130			
1,3-Dichlorobenzene	10.8	1.0	μg/L	10.0		108	70-130			
1,4-Dichlorobenzene	10.1	1.0	μg/L	10.0		101	70-130			
trans-1,4-Dichloro-2-butene	10.5	2.0	μg/L	10.0		105	70-130			
Dichlorodifluoromethane (Freon 12)	6.26	2.0	μg/L	10.0		62.6	40-160			V-05
1,1-Dichloroethane	10.8	1.0	μg/L	10.0		108	70-130			
1,2-Dichloroethane	8.15	1.0	μg/L	10.0		81.5	70-130			
1,1-Dichloroethylene	8.71	1.0	$\mu g/L$	10.0		87.1	70-130			
cis-1,2-Dichloroethylene	11.3	1.0	μg/L	10.0		113	70-130			
trans-1,2-Dichloroethylene	9.51	1.0	μg/L	10.0		95.1	70-130			
1,2-Dichloropropane	10.1	1.0	μg/L	10.0		101	70-130			
1,3-Dichloropropane	9.46	0.50	μg/L	10.0		94.6	70-130			
2,2-Dichloropropane	9.00	1.0	μg/L	10.0		90.0	40-130			
1,1-Dichloropropene	9.29	2.0	μg/L	10.0		92.9	70-130			
cis-1,3-Dichloropropene	8.99	0.50	μg/L	10.0		89.9	70-130			
trans-1,3-Dichloropropene	8.97	0.50	μg/L	10.0		89.7	70-130			
Diethyl Ether	9.14	2.0	μg/L	10.0		91.4	70-130			
Diisopropyl Ether (DIPE)	10.5	0.50	μg/L	10.0		105	70-130			
1,4-Dioxane	125	50	μg/L	100		125	40-130			
Ethylbenzene	10.6	1.0	μg/L	10.0		106	70-130			
Hexachlorobutadiene	9.47	0.60	μg/L	10.0		94.7	70-130			
2-Hexanone (MBK)	96.2	10	μg/L	100		96.2	70-160			
Isopropylbenzene (Cumene)	10.3	1.0	μg/L	10.0		103	70-130			
p-Isopropyltoluene (p-Cymene)	10.8	1.0	μg/L	10.0		108	70-130			
Methyl tert-Butyl Ether (MTBE)	8.07	1.0	μg/L	10.0		80.7	70-130			
Methylene Chloride	10.0	5.0	μg/L	10.0		100	70-130			
4-Methyl-2-pentanone (MIBK)	99.2	10	μg/L	100		99.2	70-160			
Naphthalene	10.8	2.0	μg/L	10.0		108	40-130			
n-Propylbenzene	10.4	1.0	μg/L	10.0		104	70-130			
Styrene	10.4	1.0	μg/L	10.0		104	70-130			
1,1,2,2 T (,)	8.45	1.0	μg/L	10.0		84.5	70-130			V-05
1,1,2,2-1etrachloroethane	9.22	0.50	μg/L	10.0		92.2	70-130			
Tetrakudra furan	9.18	1.0	μg/L	10.0		91.8	70-130			
Teluana	11.2	10	μg/L	10.0		112	70-130			
1.2.2 Tricklorohonzene	9.73	1.0	μg/L	10.0		97.3	70-130			
1,2,3- ITICHIOIODENZENE	9.65	5.0	μg/L	10.0		96.5	/0-130			
1,2,4- IIICHIOFODENZENE	10.4	1.0	μg/L	10.0		104	70-130			
1, J. J. Trichloroothana	9.75	1.0	μg/L	10.0		97.5	/0-130			
1,1,1-111CHIOFORTHAINE	9.20	1.0	μg/L	10.0		92.0	/0-130			
r, 1, 2- memoroculane	9.18	1.0	µg/L	10.0		91.8	70-130			
Trichlorofluoromethons (From 11)	8.72	1.0	µg/L	10.0		81.2 CAE **	70-130			V 05 T 04
1 2 3-Trichloropropage	6.45	2.0	μg/L	10.0		04.5 *	70-130			v-05, L-04
1.2.5- memoropropane	9.56	2.0	µg/L	10.0		95.0 01.2	70-130			
113)	9.13	1.0	μg/L	10.0		91.5	/0-130			
1,2,4-Trimethylbenzene	10.6	1.0	μg/L	10.0		106	70-130			
1,3,5-Trimethylbenzene	10.0	1.0	μg/L	10.0		100	70-130			
Vinyl Chloride	6.50	2.0	μg/L	10.0		65.0	40-160			

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QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

		Reporting		Spike	Source		%REC		RPD		
Analyte	Result	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	Notes	
Batch B258487 - SW-846 5030B											
LCS (B258487-BS1)				Prepared &	Analyzed: 05/	/21/20					
m+p Xylene	21.4	2.0	μg/L	20.0		107	70-130				
o-Xylene	10.7	1.0	μg/L	10.0		107	70-130				
Surrogate: 1,2-Dichloroethane-d4	24.7		μg/L	25.0		98.8	70-130				
Surrogate: Toluene-d8	24.6		μg/L	25.0		98.5	70-130				
Surrogate: 4-Bromofluorobenzene	26.9		$\mu g/L$	25.0		107	70-130				
LCS Dup (B258487-BSD1)				Prepared &	Analyzed: 05/	/21/20					
Acetone	91.1	50	μg/L	100		91.1	70-160	2.08	25		
Acrylonitrile	10.4	5.0	μg/L	10.0		104	70-130	7.52	25		
tert-Amyl Methyl Ether (TAME)	8.33	0.50	μg/L	10.0		83.3	70-130	0.359	25		
Benzene	10.4	1.0	μg/L	10.0		104	70-130	0.960	25		
Bromobenzene	10.2	1.0	μg/L	10.0		102	70-130	0.195	25		
Bromochloromethane	10.5	1.0	μg/L	10.0		105	70-130	2.92	25		
Bromodichloromethane	8.94	0.50	μg/L	10.0		89.4	70-130	1.81	25		
Bromoform	8.36	1.0	μg/L	10.0		83.6	70-130	5.92	25	V-05	
Bromomethane	5.82	2.0	μg/L	10.0		58.2	40-160	10.3	25		
2-Butanone (MEK)	111	20	μg/L	100		111	40-160	0.987	25		
tert-Butyl Alcohol (TBA)	64.0	20	μg/L	100		64.0	40-160	7.29	25	V-05	
n-Butylbenzene	11.7	1.0	μg/L	10.0		117	70-130	0.171	25		
sec-Butylbenzene	12.1	1.0	μg/L	10.0		121	70-130	1.16	25		
tert-Butylbenzene	11.6	1.0	μg/L	10.0		116	70-130	0.782	25		
tert-Butyl Ethyl Ether (TBEE)	9.64	0.50	μg/L	10.0		96.4	70-130	3.38	25		
Carbon Disulfide	7.55	5.0	μg/L	10.0		75.5	70-130	8.74	25		
Carbon Tetrachloride	8.48	1.0	μg/L	10.0		84.8	70-130	0.710	25		
Chlorobenzene	9.61	1.0	μg/L	10.0		96.1	70-130	1.65	25		
Chlorodibromomethane	8.24	0.50	μg/L	10.0		82.4	70-130	0.365	25		
Chloroethane	7.52	2.0	μg/L	10.0		75.2	70-130	1.20	25		
Chloroform	10.8	2.0	μg/L	10.0		108	70-130	1.22	25		
Chloromethane	6.08	2.0	μg/L	10.0		60.8	40-160	1.79	25		
2-Chlorotoluene	10.8	1.0	μg/L	10.0		108	70-130	0.739	25		
4-Chlorotoluene	10.9	1.0	μg/L	10.0		109	70-130	0.276	25		
1,2-Dibromo-3-chloropropane (DBCP)	9.06	5.0	μg/L	10.0		90.6	70-130	0.988	25		
1,2-Dibromoethane (EDB)	9.22	0.50	μg/L	10.0		92.2	70-130	0.00	25		
Dibromomethane	8.93	1.0	μg/L	10.0		89.3	70-130	0.670	25		
1,2-Dichlorobenzene	10.3	1.0	μg/L	10.0		103	70-130	1.46	25		
1,3-Dichlorobenzene	10.7	1.0	μg/L	10.0		107	70-130	1.21	25		
1,4-Dichlorobenzene	10.3	1.0	μg/L	10.0		103	70-130	1.67	25		
trans-1,4-Dichloro-2-butene	9.18	2.0	μg/L	10.0		91.8	70-130	13.1	25		
Dichlorodifluoromethane (Freon 12)	6.23	2.0	μg/L	10.0		62.3	40-160	0.480	25	V-05	
1,1-Dichloroethane	11.1	1.0	μg/L	10.0		111	70-130	2.65	25		
1,2-Dichloroethane	8.22	1.0	μg/L	10.0		82.2	70-130	0.855	25		
1,1-Dichloroethylene	8.87	1.0	μg/L	10.0		88.7	70-130	1.82	25		
cis-1,2-Dichloroethylene	11.2	1.0	μg/L	10.0		112	70-130	1.42	25		
trans-1,2-Dichloroethylene	9.45	1.0	μg/L	10.0		94.5	70-130	0.633	25		
1,2-Dichloropropane	9.71	1.0	μg/L	10.0		97.1	70-130	4.04	25		
1,3-Dichloropropane	9.45	0.50	μg/L	10.0		94.5	70-130	0.106	25		
2,2-Dichloropropane	9.06	1.0	μg/L	10.0		90.6	40-130	0.664	25		
1,1-Dichloropropene	9.00	2.0	μg/L	10.0		90.0	70-130	3.17	25		
cis-1,3-Dichloropropene	9.34	0.50	μg/L	10.0		93.4	70-130	3.82	25		
trans-1,3-Dichloropropene	8.87	0.50	μg/L	10.0		88.7	70-130	1.12	25		
Diethyl Ether	9.63	2.0	μg/L	10.0		96.3	70-130	5.22	25		
Diisopropyl Ether (DIPE)	10.7	0.50	μg/L	10.0		107	70-130	1.60	25		
	10.,										



QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

Avelat	Dervit	Reporting	I Inite	Spike	Source	0/DEC	%REC	DDD	RPD	N-4	
Anaryte	Kesuit	Limit	Units	Level	Result	%KEC	Limits	KPD	Limit	Notes	
Batch B258487 - SW-846 5030B											_
LCS Dup (B258487-BSD1)				Prepared &	Analyzed: 05	5/21/20					
1,4-Dioxane	127	50	μg/L	100		127	40-130	1.84	50		†‡
Ethylbenzene	10.4	1.0	μg/L	10.0		104	70-130	1.05	25		
Hexachlorobutadiene	9.30	0.60	μg/L	10.0		93.0	70-130	1.81	25		
2-Hexanone (MBK)	97.7	10	μg/L	100		97.7	70-160	1.60	25		t
Isopropylbenzene (Cumene)	10.2	1.0	μg/L	10.0		102	70-130	1.37	25		
p-Isopropyltoluene (p-Cymene)	10.7	1.0	μg/L	10.0		107	70-130	0.928	25		
Methyl tert-Butyl Ether (MTBE)	8.13	1.0	μg/L	10.0		81.3	70-130	0.741	25		
Methylene Chloride	9.70	5.0	μg/L	10.0		97.0	70-130	3.15	25		
4-Methyl-2-pentanone (MIBK)	102	10	μg/L	100		102	70-160	2.60	25		t
Naphthalene	11.1	2.0	μg/L	10.0		111	40-130	2.92	25		†
n-Propylbenzene	10.4	1.0	μg/L	10.0		104	70-130	0.673	25		
Styrene	10.4	1.0	μg/L	10.0		104	70-130	0.481	25		
1,1,1,2-Tetrachloroethane	8.44	1.0	μg/L	10.0		84.4	70-130	0.118	25	V-05	
1,1,2,2-Tetrachloroethane	9.22	0.50	μg/L	10.0		92.2	70-130	0.00	25		
Tetrachloroethylene	8.98	1.0	μg/L	10.0		89.8	70-130	2.20	25		
Tetrahydrofuran	10.6	10	μg/L	10.0		106	70-130	5.32	25		
Toluene	9.56	1.0	μg/L	10.0		95.6	70-130	1.76	25		
1,2,3-Trichlorobenzene	9.77	5.0	μg/L	10.0		97.7	70-130	1.24	25		
1,2,4-Trichlorobenzene	10.6	1.0	μg/L	10.0		106	70-130	1.62	25		
1,3,5-Trichlorobenzene	9.98	1.0	μg/L	10.0		99.8	70-130	2.33	25		
1,1,1-Trichloroethane	8.89	1.0	μg/L	10.0		88.9	70-130	3.43	25		
1,1,2-Trichloroethane	9.44	1.0	μg/L	10.0		94.4	70-130	2.79	25		
Trichloroethylene	9.17	1.0	μg/L	10.0		91.7	70-130	5.03	25		
Trichlorofluoromethane (Freon 11)	6.68	2.0	μg/L	10.0		66.8 *	70-130	3.50	25	L-04, V-05	
1,2,3-Trichloropropane	9.81	2.0	μg/L	10.0		98.1	70-130	2.58	25		
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon	8.91	1.0	μg/L	10.0		89.1	70-130	2.44	25		
113)											
1,2,4-Trimethylbenzene	10.7	1.0	μg/L	10.0		107	70-130	0.282	25		
1,3,5-Trimethylbenzene	9.97	1.0	μg/L	10.0		99.7	70-130	0.500	25		
Vinyl Chloride	6.51	2.0	μg/L	10.0		65.1	40-160	0.154	25		†
m+p Xylene	21.3	2.0	μg/L	20.0		106	70-130	0.749	25		
o-Xylene	10.6	1.0	μg/L	10.0		106	70-130	1.41	25		
Surrogate: 1,2-Dichloroethane-d4	25.2		μg/L	25.0		101	70-130				_
Surrogate: Toluene-d8	24.6		μg/L	25.0		98.4	70-130				
Surrogate: 4-Bromofluorobenzene	26.3		μg/L	25.0		105	70-130				



QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

	_	Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	Notes
Batch B258532 - SW-846 3510C										
Blank (B258532-BLK1)				Prepared: 05	5/21/20 Anal	yzed: 05/26/2	20			
Acetophenone	ND	10	μg/L							
Aniline	ND	5.0	μg/L							V-05
Benzidine	ND	20	μg/L							V-04, V-05
Benzoic Acid	ND	10	μg/L							
Bis(2-chloroethoxy)methane	ND	10	μg/L							
Bis(2-chloroethyl)ether	ND	10	μg/L							
Bis(2-chloroisopropyl)ether	ND	10	μg/L							
Bis(2-Ethylhexyl)phthalate	ND	10	μg/L							
4-Bromophenylphenylether	ND	10	μg/L							
Butylbenzylphthalate	ND	10	μg/L							
Carbazole	ND	10	μg/L							
4-Chloroaniline	ND	10	μg/L							V-34
4-Chloro-3-methylphenol	ND	10	μg/L							
2-Chloronaphthalene	ND	10	μg/L							
2-Chlorophenol	ND	10	μg/L							
4-Chlorophenylphenylether	ND	10	μg/L							
Dibenzofuran	ND	5.0	μg/L							
Di-n-butylphthalate	ND	10	μg/L							
1,2-Dichlorobenzene	ND	5.0	μg/L							
1,3-Dichlorobenzene	ND	5.0	μg/L							
1,4-Dichlorobenzene	ND	5.0	μg/L							
3,3-Dichlorobenzidine	ND	10	μg/L							
2,4-Dichlorophenol	ND	10	μg/L							
Diethylphthalate	ND	10	μg/L							
2,4-Dimethylphenol	ND	10	μg/L							
Dimethylphthalate	ND	10	μg/L							
4,6-Dinitro-2-methylphenol	ND	10	μg/L							
2,4-Dinitrophenol	ND	10	μg/L							
2,4-Dinitrotoluene	ND	10	μg/L							
2,6-Dinitrotoluene	ND	10	μg/L							
Di-n-octylphthalate	ND	10	μg/L							
1,2-Diphenylhydrazine/Azobenzene	ND	10	μg/L							
Hexachlorobenzene	ND	10	μg/L							
Hexachlorobutadiene	ND	10	μg/L							
Hexachlorocyclopentadiene	ND	10	μg/L							
Hexachloroethane	ND	10	μg/L							
Isophorone	ND	10	μg/L							
1-Methylnaphthalene	ND	5.0	μg/L							
2-Methylphenol	ND	10	μg/L							
3/4-Methylphenol	ND	10	μg/L							
2-Nitroaniline	ND	10	μg/L							
3-Nitroaniline	ND	10	μg/L							
4-Nitroaniline	ND	10	μg/L							
Nitrobenzene	ND	10	μg/L							
2-Nitrophenol	ND	10	μg/L							
4-Nitrophenol	ND	10	μg/L							
N-Nitrosodimethylamine	ND	10	μg/L							
N-Nitrosodiphenylamine/Diphenylamine	ND	10	μg/L							
N-Nitrosodi-n-propylamine	ND	10	μg/L							
Pentachloronitrobenzene	ND	10	μg/L							
Pentachlorophenol	ND	10	μg/L							
Phenol	ND	10	μg/L							

QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch B258532 - SW-846 3510C										
Blank (B258532-BLK1)				Prepared: 05	5/21/20 Analy	zed: 05/26/2	20			
Pyridine	ND	5.0	μg/L							
1,2,4,5-Tetrachlorobenzene	ND	10	μg/L							
1,2,4-Trichlorobenzene	ND	5.0	μg/L							
2,4,5-Trichlorophenol	ND	10	μg/L							
2,4,6-Trichlorophenol	ND	10	μg/L							
Surrogate: 2-Fluorophenol	101		ug/L	200		50.3	15-110			
Surrogate: Phenol-d6	70.3		μg/L μg/L	200		35.2	15-110			
Surrogate: Nitrobenzene-d5	80.6		μg/L μg/L	100		80.6	30-130			
Surrogate: 2-Fluorobinhenyl	80.5		μg/L μg/L	100		80.5	30-130			
Surrogate: 2.4.6-Tribromophenol	169		иg/L	200		84.5	15-110			
Surrogate: p-Terphenyl-d14	93.6		μg/L	100		93.6	30-130			
L CS (D259532 DS1)			10	Prepared: 05	5/21/20 Anals	rzed: 05/26/	20			
Acetophenone	20.4	10	ug/L	50.0	//21/20 Analy	78 7	40-140			
Aniline	21.7	5.0	119/L	50.0		63.4	40-140			V-05
Benzidine	31.7	20	µg/L	50.0		60.9	40-140			V-04 V-05
Benzoic Acid	50.5	10	µg/L	50.0		35.2	10 130			v-04, v-05
Bis(2-chloroethoxy)methane	17.0	10	μg/L μg/Ι	50.0		93.2 81.8	40 140			
Bis(2-chloroethyl)ether	40.9	10	μg/L μg/I	50.0		77.5	40-140			
Bis(2 chloroisopropul)ether	38.8	10	μg/L	50.0		96.0	40-140			
Bis(2 Ethylboxyl)phthelate	43.4	10	µg/L	50.0		80.9	40-140			
4. Dromonhanvlahanvlathar	44.4	10	µg/L	50.0		88.7	40-140			
4-Bromophenyiphenyiether	40.1	10	µg/L	50.0		80.2	40-140			
Carbanala	40.6	10	µg/L	50.0		81.2	40-140			
	40.7	10	μg/L	50.0		81.4	40-140			
4-Chioroaniline	35.8	10	μg/L	50.0		71.7	40-140			V-34
4-Chloro-3-methylphenol	38.5	10	μg/L	50.0		77.1	30-130			
2-Chloronaphthalene	34.0	10	μg/L	50.0		68.0	40-140			
2-Chlorophenol	35.1	10	μg/L	50.0		70.2	30-130			
4-Chlorophenylphenylether	39.1	10	μg/L	50.0		78.2	40-140			
Dibenzoturan	40.6	5.0	μg/L	50.0		81.1	40-140			
Di-n-butylphthalate	43.6	10	μg/L	50.0		87.3	40-140			
1,2-Dichlorobenzene	28.9	5.0	μg/L	50.0		57.9	40-140			
1,3-Dichlorobenzene	27.2	5.0	μg/L	50.0		54.3	40-140			
1,4-Dichlorobenzene	27.5	5.0	μg/L	50.0		55.0	40-140			
3,3-Dichlorobenzidine	44.5	10	μg/L	50.0		88.9	40-140			
2,4-Dichlorophenol	38.3	10	μg/L	50.0		76.6	30-130			
Diethylphthalate	40.4	10	μg/L	50.0		80.7	40-140			
2,4-Dimethylphenol	32.7	10	μg/L	50.0		65.4	30-130			
Dimethylphthalate	40.0	10	μg/L	50.0		80.1	40-140			
4,6-Dinitro-2-methylphenol	39.5	10	μg/L	50.0		79.1	30-130			
2,4-Dinitrophenol	36.2	10	μg/L	50.0		72.3	30-130			
2,4-Dinitrotoluene	38.5	10	μg/L	50.0		77.0	40-140			
2,6-Dinitrotoluene	40.6	10	μg/L	50.0		81.2	40-140			
Di-n-octylphthalate	43.3	10	μg/L	50.0		86.6	40-140			
1,2-Diphenylhydrazine/Azobenzene	45.1	10	μg/L	50.0		90.2	40-140			
Hexachlorobenzene	39.5	10	μg/L	50.0		79.0	40-140			
Hexachlorobutadiene	28.0	10	μg/L	50.0		56.1	40-140			
Hexachlorocyclopentadiene	24.6	10	μg/L	50.0		49.3	30-140			
Hexachloroethane	26.3	10	μg/L	50.0		52.6	40-140			
Isophorone	39.2	10	μg/L	50.0		78.3	40-140			
1-Methylnaphthalene	34.4	5.0	μg/L	50.0		68.8	40-140			
2-Methylphenol	34.9	10	μg/L	50.0		69.8	30-130			



QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

International and the set of the	Analyte	Result	Reporting	Unite	Spike Level	Source Result	%REC	%REC	RPD	RPD Limit	Notes	
Name decision of the standard s	Datab D359522 SW 946 2510C	Result	Linn	emo	Lever	resure	/utile	Linits	IU D	Linit	rotes	
IA NLEWSKA 163 (97.47919400 (97.47994.4794.0 </td <td>Balcii B230332 - 5 W-040 3510C</td> <td></td> <td></td> <td></td> <td>D 1.05</td> <td></td> <td>1.05/26/</td> <td>20</td> <td></td> <td></td> <td></td> <td>_</td>	Balcii B230332 - 5 W-040 3510C				D 1.05		1.05/26/	20				_
44. Metry optimize14.2000001.480.004.140.105-Mitonaline9.80.0 </td <td>LCS (B258532-BS1)</td> <td></td> <td></td> <td>17</td> <td>Prepared: 05</td> <td>5/21/20 Anal</td> <td>yzed: 05/26/</td> <td>20</td> <td></td> <td></td> <td></td> <td></td>	LCS (B258532-BS1)			17	Prepared: 05	5/21/20 Anal	yzed: 05/26/	20				
2-Abiomanine50.6ininin50.010140.1	3/4-Methylphenol	34.2	10	μg/L	50.0		68.3	30-130				
s-Advanamine 39.8 00 µD 1 500 717 4040 Ninoskaname 42 00 µD 2 500 72.3 4040 Shringshend 360 00 µD 1 500 72.3 4040 Shringshend 20.4 00 µD 500 72.3 4040 Shringshend 20.4 00 µD 500 40.9 01.9 US 100 40.9 10.9 US 100 Shringshend 20.4 00 µD 1 500 40.9 10.9 US 100 100 Ninoskaname program 42.4 00 µD 1 500 40.9 10.9 US 100 Shringshend 20.4 00 µD 1 500 40.9 10.9 US 100 Shringshend 20.4 00 µD 1 500 40.9 10.9 US 100 Shringshend 20.4 00 µD 1 500 40.9 10.9 US 100 Shringshend 20.0 00 Shringshend 20.0 00 Shrings	2-Nitroaniline	50.6	10	μg/L	50.0		101	40-140				
α-Antonalina 60.4 00 μpL 50.0 72.0 90.1 90.0	3-Nitroaniline	39.8	10	μg/L	50.0		79.7	40-140				
Nationaliania 3.6.2 00 µE 50.0 7.3 00.100 4-Kinophend 2.4 00 µE 50.0 40.0 90.100 10.000 10.000 4-Kinophend 2.4 00 µE 50.0 40.0 90.100 60.10 10.000 <	4-Nitroaniline	40.4	10	μg/L	50.0		80.9	40-140				
2-Autophead3a00μ2500(n)09101010N-Ninoscinetylamine2.60μ50049.240.140N-Ninoscinetylamine4.60μ50048.840.140N-Ninoscinetylamine4.200μ50070.040.140N-Ninoscinetylamine3.00μ50070.060.130 </td <td>Nitrobenzene</td> <td>36.2</td> <td>10</td> <td>μg/L</td> <td>50.0</td> <td></td> <td>72.3</td> <td>40-140</td> <td></td> <td></td> <td></td> <td></td>	Nitrobenzene	36.2	10	μg/L	50.0		72.3	40-140				
+-Autoprint20400μµL50040.900.1-101.1.11	2-Nitrophenol	38.0	10	μg/L	50.0		76.0	30-130				
NAMIONAL NA	4-Nitrophenol	20.4	10	μg/L	50.0		40.9	10-130				Ť
A-Ninosodi-perjunance Lipschanne44.610µgL50.083.840.140Petadhironophano63.010µgL50.087.880.070.080.10070.0Petadhironophano18.310µgL50.070	N-Nitrosodimethylamine	24.6	10	μg/L	50.0		49.2	40-140				
N-Natrosci-3-propriamne42.410µgL50.087.840-14Pentadhovaphenal50.010µg.750.070.030.1311Pentadhovaphenal50.010µg.750.070.010.140111Pridac20.150.0µg.750.070.140.14011111.2.4.5 1retachiovaphenal37.610µg.750.072.230.130111 </td <td>N-Nitrosodiphenylamine/Diphenylamine</td> <td>44.6</td> <td>10</td> <td>μg/L</td> <td>50.0</td> <td></td> <td>89.3</td> <td>40-140</td> <td></td> <td></td> <td></td> <td></td>	N-Nitrosodiphenylamine/Diphenylamine	44.6	10	μg/L	50.0		89.3	40-140				
Partial/conjunction00000000.0 </td <td>N-Nitrosodi-n-propylamine</td> <td>42.4</td> <td>10</td> <td>μg/L</td> <td>50.0</td> <td></td> <td>84.8</td> <td>40-140</td> <td></td> <td></td> <td></td> <td></td>	N-Nitrosodi-n-propylamine	42.4	10	μg/L	50.0		84.8	40-140				
Pendenkoppkenol 3.5 10 µµL 50.0 70.0 30-130 Pjrdine 21.1 5.0 µµL 50.0 40.1 10-140 1 1.2.4 <trictholkorberzene< td=""> 36.3 5.0 µµL 50.0 71.3 40-140 1 1.2.4<trictholkorberzene< td=""> 36.8 5.0 µµL 50.0 77.3 30-130 1<!--</td--><td>Pentachloronitrobenzene</td><td>40.3</td><td>10</td><td>μg/L</td><td>50.0</td><td></td><td>80.7</td><td>40-140</td><td></td><td></td><td></td><td></td></trictholkorberzene<></trictholkorberzene<>	Pentachloronitrobenzene	40.3	10	μg/L	50.0		80.7	40-140				
Phende 18.3 10 μp1 50.0 36.6 20.130 14.0 14.1 12,4,5-Trichlorobenzene 36.3 10 μp1 50.0 7.6 40.140 1 2,4,5-Trichlorobenzene 36.8 10 µp1 50.0 7.3 30.130 1 1 2,4,5-Trichlorophenol 37.6 10 µp1 20.0 7.2 31.10 1 1 Surrogate 2-Flucorophenol 37.6 10 µp1 200 37.6 15.10 1	Pentachlorophenol	35.0	10	μg/L	50.0		70.0	30-130				
Prindine 20.1 5.0 µp.1 50.0 40.1 10-140 ···· 12.4-Trickhorbenzene 36.3 10 µp.1 50.0 72.6 40-140 · 2.4.5-Trickhorbenzene 36.6 10 µp.1 50.0 77.2 30-130 · · 2.4.5-Trickhorbenzene 36.6 10 µp.1 200 50.6 15.10 ·<	Phenol	18.3	10	μg/L	50.0		36.6	20-130				Ť
12,45 12,45 50.0 72.6 40.140 2,4.5 10 µg/L 50.0 67.7 40.140 2,4.5 100 µg/L 50.0 77.2 30.130 2,4.6 Tichlorophenol 37.6 10 µg/L 20.0 57.3 30.130 Surrogate: Phenologhenol 7.4 µg/L 200 50.6 15.110 Surrogate: Phenologhenol 7.4 µg/L 200 37.7 15.110 Surrogate: 2,4.6 100 R2.5 30.130 Surrogate: 2,4.6 100 R2.5 30.130 Surrogate: 2,4.6 100 R2.1 100 R2.5 30.130 Surrogate: 2,4.6 100 µg/L 100 R2.5 30.130	Pyridine	20.1	5.0	μg/L	50.0		40.1	10-140				Ť
1.2.4. Trichlorobenzene 30.8 5.0 µµL 50.0 6.7. 40.40 2.4.6. Trichlorobenzon 37.6 10 µµL 50.0 77.2 30.130 2.4.6. Trichlorobenzon 10/1 µµL 200 50.6 15.110 Surrogate: Z-Fluorobenol 10/1 µµL 100 78.2 30.130 Surrogate: Z-Honorobhenol 76.4 µµL 100 82.5 30.130 Surrogate: Z-AG-Trithcomophenol 76.3 µµL 100 86.1 30.130 Surrogate: Z-AG-Trithcomophenol 76.3 µµL 100 86.1 30.130 Surogate: Z-AG-Trithcomophenol 76.3 µµL 50.0 72.7 40.140 80.3 20 Autiline 26.4 5.0 µµL 50.0 72.7 40.140 91.6 20 V-0.4 14 Benzichnorethorethynethen 36.7 10 µµL 50.0 72.4 40.140 11.8 20 - Benzichnorethynethen 36.7 10 µµL 50.	1,2,4,5-Tetrachlorobenzene	36.3	10	μg/L	50.0		72.6	40-140				
2.4.5 Trichlorophenol 37.6 10 µgL 50.0 72.3 30.130 2.4.6 Trichlorophenol 38.6 10 µgL 50.0 72.2 30.130 Surrogate: 2-Fluorophenol 75.4 µgL 200 50.6 15.110 Surrogate: Schoorophenol 75.2 µgL 100 75.2 30.130 Surrogate: 2-AC-Trichromophenol 76.3 µgL 100 75.2 30.130 Surrogate: 2-AC-Trichromophenol 76.3 µgL 200 81.3 15.110	1,2,4-Trichlorobenzene	30.8	5.0	μg/L	50.0		61.7	40-140				
24.6 Triblorophenol 38.6 10 μgL 50.0 7.2 30.130 Surrogate: Penolodic 101 μgL 200 37.7 15.110 Surrogate: Nitrobenzone dS 78.2 μgL 100 82.5 30.130 Surrogate: Nitrobenzone dS 78.2 μgL 100 82.5 30.130 Surrogate: 2,4.6 Tribromophenol 163 μgL 200 81.3 15.110 Surrogate: 2,4.6 Tribromophenol 163 μgL 50.0 72.7 40.140 80.3 20 Actiphenone 66.7 μgL 50.0 56.7 40.140 81.3 20 Aniline 28.4 50 μgL 50.0 56.7 40.140 81.3 20 Aniline 28.4 50 μgL 50.0 73.5 40.140 81.3 20 Barzidine 27.7 80.140 9.16 20 $V.04, V.05$ 4 Barzidine 28.4 50 μgL 50.0 73.5 40.140 11.3 20 Big2-chloroethylphether 36.7 10 μgL 50.0 73.5 40.140 18.3 20 Big2-chloroethylphether 36.7 10 μgL 50.0 73.5 40.140 18.3 20 Big2-chloroethylphether 36.7 10 μgL 50.0 73.5 40.140 18.3 20 Big2-chloroethylphether 36.7 10 μgL 50.0 76.4 40.14	2,4,5-Trichlorophenol	37.6	10	μg/L	50.0		75.3	30-130				
Sarragae: 2-Encophenol101µgL20050.615.10Sarragae: 2-Encophenol75.7µgL20077.715.10Sarrogae: 2-Encophenol62.5µgL10078.230-13Sarrogae: 2-Encophenol63.3µgL10081.315.10Sarrogae: 2-Encophenol63.4µgL10081.430-13Sarrogae: 2-Encophenol63.7µgL50.072.740.1081.420Actophenon76.3µgL50.072.740.1081.450\$0.0\$2.5Benzoinchi71.710µgL50.055.640.14011.350\$0.4\$1.6Benzoinchi71.710µgL50.055.640.14011.320\$1.6\$1	2,4,6-Trichlorophenol	38.6	10	μg/L	50.0		77.2	30-130				
Surrogate: Phono-Info 75.4 µg/L 200 37.7 15.10 Surrogate: Phono-Information 78.2 µg/L 100 78.2 30-130 Surrogate: 2-Incorrolphenol 16.3 µg/L 200 81.3 15-10 Surrogate: 2-Incorrolphenol 16.3 µg/L 200 81.3 15-10 CS Dup (B2S8S3-BSD1) PErpert: 05/21/20 multicette 80.3 20 Actiophenon 36.3 10 µg/L 50.0 72.7 40.140 8.03 20 Aniline 28.4 5.0 µg/L 50.0 73.5 40.140 9.16 20 V-04, V-0.7 Benzic Acid 17.1 10 µg/L 50.0 73.5 40.140 13.3 20 Bis(2-chlorothoxy)methane 36.7 10 µg/L 50.0 73.5 40.140 14.3 20 Bis(2-chlorothoxy)methane 36.7 10 µg/L 50.0 73.4 40.140 4.32 20 Bis(2	Surrogate: 2-Fluorophenol	101		μg/L	200		50.6	15-110				
Surogate: Nitrobenzene-d5 78.2 µg/L 100 78.2 30-130 Surogate: 2-Fluotobjhenyl 82.5 µg/L 100 82.5 30-130 Surogate: 2-Fluotobjhenyl 86.7 µg/L 200 81.3 15-10 Surogate: p-Terphenyl-d14 86.7 µg/L 100 86.1 30-130 Actophenone 36.3 10 µg/L 50.0 72.7 40-140 11.1 50 \$\nu_05.7 Antinine 28.4 50 µg/L 50.0 55.6 40-140 9.16 20 \$\nu_04.1 50 \$\nu_05.6 40-140 9.16 20 \$\nu_04.1 50 \$\nu_05.6 40-140 9.16 20 \$\nu_04.1 50 \$\nu_05.6 40-140 11.3 20 \$\nu_04.1 10 \$\nu_06.1 \$\nu_06.1 \$\nu_05.1 \$\nu_06.1 \$\nu_06.1<	Surrogate: Phenol-d6	75.4		μg/L	200		37.7	15-110				
Surogate: 2-Huorobiphenyl §2.5 µg/L 100 82.5 30-130 Surogate: 2-A.6-Trihoroophenol 163 µg/L 100 81.3 15-110 Surogate: 2-A.6-Trihoroophenol 86.1 192/L 100 81.3 15-110 Surogate: 2-A.6-Trihoroophenol 36.3 10 µg/L 50.0 7.7 40-40 8.03 20 Aniline 28.4 50 µg/L 50.0 56.6 40-140 11.1 50 V04, V05 7 Benzidine 27.8 20 µg/L 50.0 55.6 40-140 11.8 20 V04, V05 7 Bic2-chloroethypithem 36.7 10 µg/L 50.0 7.2 40-140 1.8 20 Bic2-chloroethypithem 36.7 10 µg/L 50.0 7.2 40-140 1.4 20 Bic2-chloroethypithem 36.7 10 µg/L 50.0 7.4 40.4 49.2 20 Bic2-chloroethypithens/them	Surrogate: Nitrobenzene-d5	78.2		μg/L	100		78.2	30-130				
Surogais: 2,4,6-Tribromphenol/63µg/L20081.315-110Surogais: p-Terphenyl-01486.1np/L10086.130-10Conspan=100 (2000)Conspan=100 (2000)Actiophenone30.310µg/L50.072.740-1408.0320Aniine28.45.0µg/L50.056.740-14011.150V-04, V-05\$Benzidine27.820µg/L50.063.240-1401.0820\$Benzidine36.7100µg/L50.073.540-1401.0820Bis(2-chlorothoxy)methane36.610µg/L50.073.240-1401.320Bis(2-chlorothoxy)methane34.610µg/L50.075.440-1404.8320Bis(2-chlorothoxy)methane34.210µg/L50.076.440-1404.8320Bis(2-chlorothy)tehref38.210µg/L50.076.440-1404.8320Carbazole32.210µg/L50.076.440-1404.8320Carbazole32.210µg/L50.076.440-1404.8320Carbazole32.210µg/L50.076.440-1404.8320Carbazole32.410µg/L50.076.440-1404.8320Carbazole33.4 <t< td=""><td>Surrogate: 2-Fluorobiphenyl</td><td>82.5</td><td></td><td>μg/L</td><td>100</td><td></td><td>82.5</td><td>30-130</td><td></td><td></td><td></td><td></td></t<>	Surrogate: 2-Fluorobiphenyl	82.5		μg/L	100		82.5	30-130				
Surogate: p-Tephenyl-Id186./µg/L10086.196.196.196.196.1Departe: 05/21/0 Autro: 05/200Actop (025832-35D)Departe: 05/21/0 Autro: 05/2007.740.408.030Aniline28.4509/L50.07.740.408.030Vol. (Vol. (Surrogate: 2,4,6-Tribromophenol	163		μg/L	200		81.3	15-110				
LCS Dap (B25852-BSD1) Prepared: 05/21/20 Analyzed: 05/26/20 Acetophenone 36.3 10 µg/L 50.0 72.7 40.140 8.03 20 Aniline 28.4 5.0 µg/L 50.0 56.7 40.140 11.1 50 V-05.5 4 Benzidine 27.8 20 µg/L 50.0 34.2 10.130 2.77 50 V-04, V-05 Benzoic Acid 17.1 10 µg/L 50.0 73.5 40.140 10.8 20 Bis(2-chlorosthy)methane 36.7 10 µg/L 50.0 75.2 40.140 1.4 20 Bis(2-chlorosthy)methane 37.6 10 µg/L 50.0 75.4 40.140 4.33 20 Bis(2-chlorosthy)methane 38.2 10 µg/L 50.0 76.4 40.140 4.52 20 Carbazole 39.2 10 µg/L 50.0 76.4 40.140 5.21 20 V-34 4	Surrogate: p-Terphenyl-d14	86.1		$\mu g/L$	100		86.1	30-130				
Accophenone 36.3 10 $\mu g/L$ 50.0 72.7 40.140 8.03 20 Antline 28.4 50 $\mu g/L$ 50.0 56.7 40.140 11.1 50 $V.05$ \ddagger Berzidine 27.8 20 $\mu g/L$ 50.0 55.6 40.140 9.16 20 $V.04$ $V.05$ \ddagger Berzoic Acid 17.1 10 $\mu g/L$ 50.0 53.6 40.140 10.8 20 $V.04$ $V.05$ \ddagger Bis(2-chlorochy)methane 36.7 10 $\mu g/L$ 50.0 73.5 40.140 11.3 20 Bis(2-chlorochy)methane 36.7 10 $\mu g/L$ 50.0 75.2 40.140 14.4 20 Bis(2-chlorochy)methane 37.6 10 $\mu g/L$ 50.0 76.4 40.140 4.83 20 Bis(2-chlorochy)methane 38.7 10 $\mu g/L$ 50.0 76.4 40.140 4.83 20 Butylbenzylphthalate 38.7 10 $\mu g/L$ 50.0 76.4 40.140 4.83 20 Carbazole 39.2 10 $\mu g/L$ 50.0 76.4 40.140 4.83 20 4-Chloro-shmethylphenol 37.0 10 $\mu g/L$ 50.0 76.4 40.140 3.75 20 4-Chloro-shmethylphenol 37.0 10 $\mu g/L$ 50.0 76.8 40.140 3.75 20 2-Chlorophenylphthalate 31.7 10 $\mu g/L$ <	LCS Dup (B258532-BSD1)				Prepared: 05	5/21/20 Anal	yzed: 05/26/2	20				
Aniline28.45.0µg/L50.056.740-14011.150V.05‡Benzidine27.820µg/L50.055.640-1409.1620V.04, V.05Benzoic Acid17.110µg/L50.034.210-1302.7501* ±Bis(2-chloroethxy)methane36.710µg/L50.069.240-14011.3201.41.420Bis(2-chloroethxy)methar37.610µg/L50.072.240-1401.4201.4	Acetophenone	36.3	10	μg/L	50.0		72.7	40-140	8.03	20		
Benzidine 27,8 20 µg/L 50.0 55.6 40-140 9.16 20 V-04, V-05 Benzoic Acid 17,1 10 µg/L 50.0 34.2 10-130 2.77 50 1* Bis(2-chloroethxy)methane 36.6 10 µg/L 50.0 73.5 40-140 10.8 20 Bis(2-chloroethy)lether 37.6 10 µg/L 50.0 75.2 40-140 11.3 20 Bis(2-chloroisopropy)lether 37.6 10 µg/L 50.0 75.2 40-140 14.4 20 Bis(2-chloroisopropy)lether 38.2 10 µg/L 50.0 76.4 40-140 4.83 20 Burylbenz/ylhhalate 41.2 10 µg/L 50.0 76.4 40-140 3.75 20 Carbazole 39.2 10 µg/L 50.0 78.4 40-140 5.21 20 V-34 4-Chloro-3-methylphenol 37.0 10 µg/L 50.0 78.1 40-140 1.51 20 2-Chlorophenylphenylether <td< td=""><td>Aniline</td><td>28.4</td><td>5.0</td><td>μg/L</td><td>50.0</td><td></td><td>56.7</td><td>40-140</td><td>11.1</td><td>50</td><td>V-05</td><td>‡</td></td<>	Aniline	28.4	5.0	μg/L	50.0		56.7	40-140	11.1	50	V-05	‡
Benzoic Acid17.110µg/L50.034.210-1302.7750\$1Bis(2-chloroethoxy)methane36.710µg/L50.073.540-14010.820Bis(2-chloroethoy)methane37.610µg/L50.069.240-14011.320Bis(2-chloroethy)methane37.610µg/L50.075.240-1407.2920Bis(2-chloroethy)methane38.210µg/L50.076.440-1404.8320Butylbenzylphthalate38.710µg/L50.078.440-1404.8320Carbazole39.210µg/L50.078.440-1404.92204-Chloroethylphenol37.010µg/L50.078.440-1404.92204-Chloroethylphenol37.010µg/L50.078.440-1404.92204-Chloroethylphenol37.010µg/L50.078.440-1404.92204-Chloroethylphenol37.010µg/L50.068.140-1407.5204-Chloroethylphenol38.410µg/L50.078.440-1407.5204-Chloroethylphenol38.410µg/L50.077.740-1407.62010-houtylphenol38.410µg/L50.078.840-1401.62011-houtphenolymether88.95.0µg/L50.0 <td>Benzidine</td> <td>27.8</td> <td>20</td> <td>μg/L</td> <td>50.0</td> <td></td> <td>55.6</td> <td>40-140</td> <td>9.16</td> <td>20</td> <td>V-04, V-05</td> <td></td>	Benzidine	27.8	20	μg/L	50.0		55.6	40-140	9.16	20	V-04, V-05	
Bis(2-chloroethoxy)methane 36.7 10 µg/L 50.0 73.5 40.140 10.8 20 Bis(2-chloroethy)lether 34.6 10 µg/L 50.0 62.2 40.140 11.3 20 Bis(2-chloroethy)lether 37.6 10 µg/L 50.0 75.2 40.140 7.29 20 4-Bromophenylphenylether 38.2 10 µg/L 50.0 76.4 40.140 4.83 20 A-Bromophenylphenylether 38.2 10 µg/L 50.0 76.4 40.140 4.83 20 Carbazole 39.2 10 µg/L 50.0 76.4 40.140 3.75 20 4-Chloro-amethylphenol 37.0 10 µg/L 50.0 76.4 40.140 5.21 20 V-34 4-Chloro-amethylphenol 37.0 10 µg/L 50.0 76.8 40.140 7.5 20 2-Chloroaphthalene 31.7 10 µg/L 50.0 76.8 40.140 1.6 20 2-Chlorophenol 33.4 10 µ	Benzoic Acid	17.1	10	μg/L	50.0		34.2	10-130	2.77	50		†‡
Bis(2-chlorosehyl)ether 34.6 10 µg/L 50.0 69.2 40-140 11.3 20 Bis(2-chlorosioporpyl)ether 37.6 10 µg/L 50.0 75.2 40-140 14.4 20 Bis(2-chlorosioporpyl)ether 38.2 10 µg/L 50.0 76.4 40-140 4.33 20 4-Bromophenylphenylether 38.2 10 µg/L 50.0 77.3 40-140 4.32 20 Carbazole 39.2 10 µg/L 50.0 77.4 40-140 5.21 20 V-34 4-Chloro-3-methylphenol 37.0 10 µg/L 50.0 74.0 30-130 4.11 20 2-Chlorophenol 31.7 10 µg/L 50.0 74.0 30-130 4.11 20 2-Chlorophenol 33.4 10 µg/L 50.0 66.9 30-130 4.84 20 2-Chlorophenol 38.4 10 µg/L 50.0 77.7 40-140 1.66 20 1,3-Dichlorobenzene 29.4 5.0 µg/L	Bis(2-chloroethoxy)methane	36.7	10	μg/L	50.0		73.5	40-140	10.8	20		
Bis(2-chloroisopropyl)ether 37,6 10 µg/L 50.0 75.2 40-140 7.4 20 Bis(2-Ethylhexyl)phthalate 41.2 10 µg/L 50.0 76.4 40-140 4.83 20 4-Bromophenylphenylether 38.2 10 µg/L 50.0 76.4 40-140 4.83 20 Carbazole 38.7 10 µg/L 50.0 77.3 40-140 4.92 20 Carbazole 39.2 10 µg/L 50.0 78.4 40-140 5.21 20 V-34 4-Chloroaniline 34.0 10 µg/L 50.0 68.1 40-140 5.21 20 V-34 2-Chloroaphthalene 37.0 10 µg/L 50.0 63.3 40-140 7.5 20 2-Chlorophenol 33.4 10 µg/L 50.0 66.9 30-130 4.84 20 1-2-bichorophenylphenylether 38.4 10 µg/L 50.0 76.8 40-140 1.61 20 1-2-bichorophenzine 38.4 10 <	Bis(2-chloroethyl)ether	34.6	10	μg/L	50.0		69.2	40-140	11.3	20		
Bis(2-Ethylhexyl)phthalate 41.2 10 $\mu g/L$ 50.0 82.5 40.140 7.29 20 4-Bromophenylphenylether 38.2 10 $\mu g/L$ 50.0 76.4 40.140 4.83 20 Butylbenzylphthalate 38.7 10 $\mu g/L$ 50.0 77.3 40.140 4.92 20 Carbazole 39.2 10 $\mu g/L$ 50.0 78.4 40.140 5.7 20 $V.34$ 4-Chloroaniline 34.0 10 $\mu g/L$ 50.0 74.0 30.130 4.11 20 $V.34$ 2-Chloroaphthalene 31.7 10 $\mu g/L$ 50.0 66.9 30.130 4.84 20 2-Chlorophenol 33.4 10 $\mu g/L$ 50.0 66.9 30.130 4.84 20 2-Chlorophenylphenylether 38.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 2-Chlorophenylphenylether 38.4 10 $\mu g/L$ 50.0 76.8 40.140 1.76 20 2-Lhorophenylphenylether 38.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 1.2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 88.8 40.140 1.61 20 1.3-Dichlorobenzene 28.1 5.0 $\mu g/L$ 50.0 56.9 40.140 3.26 20 1.4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 75.9 40.140 <td>Bis(2-chloroisopropyl)ether</td> <td>37.6</td> <td>10</td> <td>μg/L</td> <td>50.0</td> <td></td> <td>75.2</td> <td>40-140</td> <td>14.4</td> <td>20</td> <td></td> <td></td>	Bis(2-chloroisopropyl)ether	37.6	10	μg/L	50.0		75.2	40-140	14.4	20		
4-Bromophenylphenylether 38.2 10 $\mu g/L$ 50.0 76.4 40.140 4.83 20 Butylbenzylphthalate 38.7 10 $\mu g/L$ 50.0 77.3 40.140 4.92 20 Carbazole 39.2 10 $\mu g/L$ 50.0 78.4 40.140 5.21 20 $V-34$ 4-Chloroaniline 34.0 10 $\mu g/L$ 50.0 68.1 40.140 5.21 20 $V-34$ 4-Chloro-3-methylphenol 37.0 10 $\mu g/L$ 50.0 63.3 40.140 7.5 20 2-Chlorophenol 31.7 10 $\mu g/L$ 50.0 63.3 40.140 7.5 20 2-Chlorophenylphenylether 38.4 10 $\mu g/L$ 50.0 63.3 40.140 7.5 20 Dibenzofuran 38.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 77.7 40.140 4.28 20 1/2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 58.8 40.140 1.61 20 1/3-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.32 20 1/4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.32 20 2/4-Dichlorophenol 36.3 10 $\mu g/L$ 50.0 72.5 30.130 50.7 20	Bis(2-Ethylhexyl)phthalate	41.2	10	μg/L	50.0		82.5	40-140	7.29	20		
Butylbenzylphhalate 38.7 10 $\mu g/L$ 50.0 77.3 40.140 4.92 20 Carbazole 39.2 10 $\mu g/L$ 50.0 78.4 40.140 3.75 20 4-Chloroaniline 34.0 10 $\mu g/L$ 50.0 68.1 40.140 5.21 20 $V.34$ 4-Chloro-3-methylphenol 37.0 10 $\mu g/L$ 50.0 63.3 40.140 7.15 20 $V.34$ 2-Chloronaphthalene 31.7 10 $\mu g/L$ 50.0 63.3 40.140 7.15 20 2-Chlorophenol 33.4 10 $\mu g/L$ 50.0 66.9 30.130 4.84 20 2-Chlorophenol 38.4 10 $\mu g/L$ 50.0 77.7 40.140 4.28 20 Dibenzoftran 88.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 77.7 40.140 4.28 20 1,2-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 58.8 40.140 1.61 20 1,3-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.32 20 1,4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 72.5 30.130 5.0 20 1,4-Dichlorobenzene 36.3 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 <tr< td=""><td>4-Bromophenylphenylether</td><td>38.2</td><td>10</td><td>μg/L</td><td>50.0</td><td></td><td>76.4</td><td>40-140</td><td>4.83</td><td>20</td><td></td><td></td></tr<>	4-Bromophenylphenylether	38.2	10	μg/L	50.0		76.4	40-140	4.83	20		
Carbazole 39.2 10 $\mu g/L$ 50.0 78.4 40.140 3.75 20 4-Chloroaniline 34.0 10 $\mu g/L$ 50.0 68.1 40.140 5.21 20 $V-34$ 4-Chloroa-3-methylphenol 37.0 10 $\mu g/L$ 50.0 63.3 40.140 7.15 20 2-Chloronaphthalene 31.7 10 $\mu g/L$ 50.0 63.3 40.140 7.15 20 2-Chlorophenol 33.4 10 $\mu g/L$ 50.0 66.9 30.130 4.84 20 4-Chlorophenylphenylether 38.4 10 $\mu g/L$ 50.0 76.8 40.140 1.76 20 Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 77.7 40.140 4.28 20 1.2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 80.0 40.140 4.28 20 1.3-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.1 40.140 4.28 20 1.4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.1 40.140 3.26 20 1.4-Dichlorobenzene 41.1 10 $\mu g/L$ 50.0 82.2 40.140 7.83 20 2.4-Dichlorobenzene 36.3 10 $\mu g/L$ 50.0 72.5 30.130 5.50 20 2.4-Dichlorophenol 36.3 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20	Butylbenzylphthalate	38.7	10	μg/L	50.0		77.3	40-140	4.92	20		
4-Chloroaniline 34.0 10 $\mu g/L$ 50.0 68.1 40.140 5.21 20 $V.34$ 4-Chloroa-3-methylphenol 37.0 10 $\mu g/L$ 50.0 74.0 30.130 4.11 20 2-Chloronaphthalene 31.7 10 $\mu g/L$ 50.0 63.3 40.140 7.15 20 2-Chlorophenol 33.4 10 $\mu g/L$ 50.0 66.9 30.130 4.84 20 4-Chlorophenylphenylether 38.4 10 $\mu g/L$ 50.0 76.8 40.140 1.76 20 Dienz-butylphthalate 40.0 10 $\mu g/L$ 50.0 77.7 40.140 4.28 20 1.2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 80.0 40.140 4.65 20 1.3-Dichlorobenzene 28.1 5.0 $\mu g/L$ 50.0 56.1 40.140 3.26 20 1.4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.26 20 1.4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 82.2 40.140 3.32 20 2.4-Dichlorobenzene 36.3 10 $\mu g/L$ 50.0 72.5 30.130 5.50 20 2.4-Dichlorobenzene 30.1 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 2.4-Dichlorobenzene 30.1 10 $\mu g/L$ 50.0 78.1 40.140 3.27	Carbazole	39.2	10	μg/L	50.0		78.4	40-140	3.75	20		
4-Chloro-3-methylphenol 37,0 10 µg/L 50.0 74.0 30-130 4.11 20 2-Chloronaphthalene 31.7 10 µg/L 50.0 63.3 40-140 7.15 20 2-Chlorophenol 33.4 10 µg/L 50.0 66.9 30-130 4.84 20 4-Chlorophenylphenylether 38.4 10 µg/L 50.0 76.8 40-140 1.76 20 Dibenzofuran 38.9 5.0 µg/L 50.0 77.7 40-140 4.28 20 1.2-Dichlorobenzene 29.4 5.0 µg/L 50.0 80.0 40-140 8.65 20 1.3-Dichlorobenzene 29.4 5.0 µg/L 50.0 58.8 40-140 1.61 20 1.3-Dichlorobenzene 28.1 5.0 µg/L 50.0 56.9 40-140 3.26 20 1.4-Dichlorobenzene 28.4 5.0 µg/L 50.0 72.5 30-130 5.50 20 2.4-Dichlorobenzidine 41.1 10 µg/L 50.0	4-Chloroaniline	34.0	10	μg/L	50.0		68.1	40-140	5.21	20	V-34	
2-Chloronaphthalene 31.7 10 $\mu g/L$ 50.0 63.3 40.140 7.15 20 2-Chlorophenol 33.4 10 $\mu g/L$ 50.0 66.9 30.130 4.84 20 4-Chlorophenylphenylether 38.4 10 $\mu g/L$ 50.0 76.8 40.140 1.76 20 Dibenzofuran 38.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 80.0 40.140 8.65 20 1,2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 56.1 40.140 1.61 20 1,3-Dichlorobenzene 28.1 5.0 $\mu g/L$ 50.0 56.1 40.140 3.26 20 1,4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.32 20 2,4-Dichlorobenzidine 41.1 10 $\mu g/L$ 50.0 82.2 40.140 7.83 20 2,4-Dichlorophenol 36.3 10 $\mu g/L$ 50.0 72.5 30.130 5.50 20 Diethylphthalate 39.1 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 2,4-Dimethylphenol 30.4 10 $\mu g/L$ 50.0 78.1 40.140 2.58 50 50	4-Chloro-3-methylphenol	37.0	10	μg/L	50.0		74.0	30-130	4.11	20		
2-Chlorophenol 33.4 10 $\mu g/L$ 50.0 66.9 30.130 4.84 204-Chlorophenylphenylether 38.4 10 $\mu g/L$ 50.0 76.8 40.140 1.76 20 Dibenzofuran 38.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 80.0 40.140 8.65 20 1,2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 56.1 40.140 3.26 20 1,3-Dichlorobenzene 28.1 5.0 $\mu g/L$ 50.0 56.9 40.140 3.26 20 1,4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 82.2 40.140 3.32 20 2,4-Dichlorobenzidine 41.1 10 $\mu g/L$ 50.0 72.5 $30-130$ 5.50 20 2,4-Dichlorobenzidine 39.1 10 $\mu g/L$ 50.0 72.5 $30-130$ 5.50 20 2,4-Dichlorobenzidine 39.1 10 $\mu g/L$ 50.0 78.1 $40-140$ 3.27 20 2,4-Dichlorobenzidine 30.4 10 $\mu g/L$ 50.0 78.1 $40-140$ 3.27 20 2,4-Direthylphenol 30.4 10 $\mu g/L$ 50.0 78.1 $40-140$ 2.58 50 50 2,4-Direthylphenol 30.4 10 $\mu g/L$ 50.0 78.1 $40-140$ 2.58 50 <td< td=""><td>2-Chloronaphthalene</td><td>31.7</td><td>10</td><td>μg/L</td><td>50.0</td><td></td><td>63.3</td><td>40-140</td><td>7.15</td><td>20</td><td></td><td></td></td<>	2-Chloronaphthalene	31.7	10	μg/L	50.0		63.3	40-140	7.15	20		
4-Chlorophenylpheny	2-Chlorophenol	33.4	10	μg/L	50.0		66.9	30-130	4.84	20		
Dibenzofuran 38.9 5.0 $\mu g/L$ 50.0 77.7 40.140 4.28 20 Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 80.0 40.140 8.65 20 1,2-Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 58.8 40.140 1.61 20 1,3-Dichlorobenzene 28.1 5.0 $\mu g/L$ 50.0 56.1 40.140 3.26 20 1,4-Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.32 20 3,3-Dichlorobenzidine 41.1 10 $\mu g/L$ 50.0 82.2 40.140 7.83 20 2,4-Dichlorophenol 36.3 10 $\mu g/L$ 50.0 72.5 30.130 5.50 20 Diethylphthalate 39.1 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 2,4-Dimethylphenol 30.4 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 Dimethylphthalate 39.0 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20	4-Chlorophenylphenylether	38.4	10	μg/L	50.0		76.8	40-140	1.76	20		
Di-n-butylphthalate 40.0 10 $\mu g/L$ 50.0 80.0 40.140 8.65 20 $1,2$ -Dichlorobenzene 29.4 5.0 $\mu g/L$ 50.0 58.8 40.140 1.61 20 $1,3$ -Dichlorobenzene 28.1 5.0 $\mu g/L$ 50.0 56.1 40.140 3.26 20 $1,4$ -Dichlorobenzene 28.4 5.0 $\mu g/L$ 50.0 56.9 40.140 3.32 20 $3,3$ -Dichlorobenzidine 41.1 10 $\mu g/L$ 50.0 82.2 40.140 7.83 20 $2,4$ -Dichlorophenol 36.3 10 $\mu g/L$ 50.0 72.5 30.130 5.50 20 Diethylphthalate 39.1 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 $2,4$ -Dimethylphenol 30.4 10 $\mu g/L$ 50.0 60.8 30.130 7.29 20 Dimethylphthalate 39.0 10 $\mu g/L$ 50.0 78.1 40.140 2.58 50 $\frac{1}{2}$	Dibenzofuran	38.9	5.0	μg/L	50.0		77.7	40-140	4.28	20		
1,2-Dichlorobenzene29,45.0µg/L50.058.840-1401.61201,3-Dichlorobenzene28.15.0µg/L50.056.140-1403.26201,4-Dichlorobenzene28.45.0µg/L50.056.940-1403.32203,3-Dichlorobenzidine41.110µg/L50.082.240-1407.83202,4-Dichlorobenzidine36.310µg/L50.072.530-1305.5020Diethylphthalate39.110µg/L50.078.140-1403.27202,4-Dimethylphenol30.410µg/L50.060.830-1307.2920Dimethylphthalate39.010µg/L50.078.140-1402.5850\$	Di-n-butylphthalate	40.0	10	μg/L	50.0		80.0	40-140	8.65	20		
1,3-Dichlorobenzene28.15.0µg/L50.056.140-1403.26201,4-Dichlorobenzene28.45.0µg/L50.056.940-1403.32203,3-Dichlorobenzidine41.110µg/L50.082.240-1407.83202,4-Dichlorobenzidine36.310µg/L50.072.530-1305.5020Diethylphthalate39.110µg/L50.078.140-1403.27202,4-Dimethylphthalate30.410µg/L50.060.830-1307.2920Dimethylphthalate39.010µg/L50.078.140-1402.5850‡	1,2-Dichlorobenzene	29.4	5.0	μg/L	50.0		58.8	40-140	1.61	20		
1,4-Dichlorobenzene 28.4 5.0 µg/L 50.0 56.9 40.140 3.32 20 3,3-Dichlorobenzidine 41.1 10 µg/L 50.0 82.2 40.140 7.83 20 2,4-Dichlorobenzidine 36.3 10 µg/L 50.0 72.5 30-130 5.50 20 Diethylphthalate 39.1 10 µg/L 50.0 78.1 40-140 3.27 20 2,4-Dimethylphenol 30.4 10 µg/L 50.0 60.8 30-130 7.29 20 Dimethylphthalate 39.0 10 µg/L 50.0 78.1 40-140 2.58 50 \$	1,3-Dichlorobenzene	28.1	5.0	μg/L	50.0		56.1	40-140	3.26	20		
3,3-Dichlorobenzidine 41.1 10 µg/L 50.0 82.2 40-140 7.83 20 2,4-Dichlorobenzidine 36.3 10 µg/L 50.0 72.5 30-130 5.50 20 Diethylphthalate 39.1 10 µg/L 50.0 78.1 40-140 3.27 20 2,4-Dimethylphenol 30.4 10 µg/L 50.0 60.8 30-130 7.29 20 Dimethylphthalate 39.0 10 µg/L 50.0 78.1 40-140 2.58 50 \$\$	1,4-Dichlorobenzene	20.1	5.0	. υ μg/L	50.0		56.9	40-140	3.32	20		
2,4-Dichlorophenol36.310 $\mu g/L$ 50.072.530-1305.5020Diethylphthalate39.110 $\mu g/L$ 50.078.140-1403.27202,4-Dimethylphenol30.410 $\mu g/L$ 50.060.830-1307.2920Dimethylphthalate39.010 $\mu g/L$ 50.078.140-1402.5850 \ddagger	3,3-Dichlorobenzidine	41 1	10	μg/L	50.0		82.2	40-140	7.83	20		
Diethylphthalate 39.1 10 $\mu g/L$ 50.0 78.1 40.140 3.27 20 $2,4$ -Dimethylphthalate 30.4 10 $\mu g/L$ 50.0 60.8 30.130 7.29 20 Dimethylphthalate 39.0 10 $\mu g/L$ 50.0 78.1 40.140 2.58 50 \ddagger	2,4-Dichlorophenol	36.3	10	μg/L	50.0		72.5	30-130	5.50	20		
2,4-Dimethylphenol 30.4 10 $\mu g/L$ 50.0 60.8 $30-130$ 7.29 20 Dimethylphthalate 39.0 10 $\mu g/L$ 50.0 78.1 $40-140$ 2.58 50 \ddagger	Diethylphthalate	39.1	10	μg/L	50.0		78.1	40-140	3.27	20		
Dimethylphthalate 39.0 $10 \mu g/L$ 50.0 78.1 40.140 2.58 50 \ddagger	2,4-Dimethylphenol	30.4	10	. υ μg/L	50.0		60.8	30-130	7.29	20		
	Dimethylphthalate	39.0	10	. ς μg/L	50.0		78.1	40-140	2.58	50		t

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QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes	
Batch B258532 - SW-846 3510C	Tesur		onto	20101	result	, vite e	2	10.0			
LCS Dup (B258532-BSD1)				Prepared: 0:	5/21/20 Anal	yzed: 05/26/	20				_
4,6-Dinitro-2-methylphenol	37.7	10	μg/L	50.0		75.4	30-130	4.77	50		— ‡
2,4-Dinitrophenol	35.7	10	μg/L	50.0		71.4	30-130	1.31	50		t
2,4-Dinitrotoluene	38.0	10	μg/L	50.0		76.1	40-140	1.20	20		•
2,6-Dinitrotoluene	39.2	10	μg/L	50.0		78.4	40-140	3.56	20		
Di-n-octylphthalate	40.2	10	μg/L	50.0		80.3	40-140	7.55	20		
1,2-Diphenylhydrazine/Azobenzene	37.8	10	μg/L	50.0		75.6	40-140	17.7	20		
Hexachlorobenzene	37.6	10	μg/L	50.0		75.1	40-140	5.03	20		
Hexachlorobutadiene	28.1	10	μg/L	50.0		56.3	40-140	0.356	20		
Hexachlorocyclopentadiene	22.6	10	μg/L	50.0		45.3	30-140	8.42	50		†‡
Hexachloroethane	27.0	10	μg/L	50.0		54.0	40-140	2.70	50		\$
Isophorone	34.9	10	μg/L	50.0		69.7	40-140	11.6	20		
1-Methylnaphthalene	32.9	5.0	μg/L	50.0		65.7	40-140	4.61	20		
2-Methylphenol	34.0	10	μg/L	50.0		67.9	30-130	2.76	20		
3/4-Methylphenol	31.3	10	μg/L	50.0		62.5	30-130	8.87	20		
2-Nitroaniline	44.4	10	μg/L	50.0		88.7	40-140	13.2	20		
3-Nitroaniline	39.1	10	μg/L	50.0		78.2	40-140	1.93	20		
4-Nitroaniline	41.1	10	μg/L	50.0		82.1	40-140	1.55	20		
Nitrobenzene	32.7	10	μg/L	50.0		65.4	40-140	10.0	20		
2-Nitrophenol	35.6	10	μg/L	50.0		71.1	30-130	6.63	20		
4-Nitrophenol	20.0	10	μg/L	50.0		39.9	10-130	2.38	50		†‡
N-Nitrosodimethylamine	23.9	10	μg/L	50.0		47.8	40-140	3.01	20		
N-Nitrosodiphenylamine/Diphenylamine	40.9	10	μg/L	50.0		81.8	40-140	8.74	20		
N-Nitrosodi-n-propylamine	37.1	10	μg/L	50.0		74.2	40-140	13.3	20		
Pentachloronitrobenzene	40.0	10	μg/L	50.0		79.9	40-140	0.971	20		
Pentachlorophenol	34.6	10	μg/L	50.0		69.2	30-130	1.15	50		\$
Phenol	16.7	10	μg/L	50.0		33.4	20-130	9.26	20		Ť
Pyridine	19.5	5.0	μg/L	50.0		39.0	10-140	2.73	50		†‡
1,2,4,5-Tetrachlorobenzene	33.6	10	μg/L	50.0		67.2	40-140	7.73	20		
1,2,4-Trichlorobenzene	29.9	5.0	μg/L	50.0		59.9	40-140	2.93	20		
2,4,5-Trichlorophenol	36.4	10	μg/L	50.0		72.9	30-130	3.29	20		
2,4,6-Trichlorophenol	36.8	10	μg/L	50.0		73.7	30-130	4.61	50		\$
Surrogate: 2-Fluorophenol	101		μg/L	200		50.3	15-110				—
Surrogate: Phenol-d6	69.5		μg/L	200		34.8	15-110				
Surrogate: Nitrobenzene-d5	71.6		μg/L	100		71.6	30-130				
Surrogate: 2-Fluorobiphenyl	76.8		μg/L	100		76.8	30-130				
Surrogate: 2,4,6-Tribromophenol	170		μg/L	200		84.8	15-110				
Surrogate: p-Terphenyl-d14	82.9		μg/L	100		82.9	30-130				

Batch B258763 - SW-846 3510C

			Prepared: 05/21/20 Analyzed: 05/27/20
ND	0.30	μg/L	
ND	0.20	μg/L	
ND	0.20	μg/L	
ND	0.050	μg/L	
ND	0.10	μg/L	
ND	0.050	μg/L	
ND	0.50	μg/L	
ND	0.20	μg/L	
ND	0.20	μg/L	
ND	0.10	μg/L	
ND	0.50	μg/L	
ND	1.0	μg/L	
	ND ND ND ND ND ND ND ND ND ND ND	ND 0.30 ND 0.20 ND 0.20 ND 0.050 ND 0.10 ND 0.50 ND 0.50 ND 0.20 ND 0.50 ND 0.20 ND 0.20 ND 0.20 ND 0.10 ND 0.50 ND 0.50 ND 1.0	ND 0.30 μg/L ND 0.20 μg/L ND 0.20 μg/L ND 0.050 μg/L ND 0.20 μg/L ND 0.20 μg/L ND 0.20 μg/L ND 0.10 μg/L ND 0.10 μg/L ND 0.50 μg/L ND 0.50 μg/L ND 0.50 μg/L ND 1.0 μg/L



QUALITY CONTROL

Semivolatile Organic Compounds by GC/MS - Quality Control

		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	Notes
Batch B258763 - SW-846 3510C										
Blank (B258763-BLK1)				Prepared: 05	5/21/20 Analy	yzed: 05/27/2	20			
Indeno(1,2,3-cd)pyrene (SIM)	ND	0.10	μg/L							
2-Methylnaphthalene (SIM)	ND	1.0	μg/L							
Naphthalene (SIM)	ND	1.0	μg/L							
Phenanthrene (SIM)	ND	0.050	μg/L							
Pyrene (SIM)	ND	1.0	μg/L							
Surrogate: Nitrobenzene-d5	86.0		μg/L	100		86.0	30-130			
Surrogate: 2-Fluorobiphenyl	76.3		μg/L	100		76.3	30-130			
Surrogate: p-Terphenyl-d14	74.9		μg/L	100		74.9	30-130			
LCS (B258763-BS1)				Prepared: 05	5/21/20 Analy	vzed: 05/27/2	20			
Acenaphthene (SIM)	45.4	6.0	ug/L	50.0		90.8	40-140			
Acenaphthylene (SIM)	44.2	4.0	μg/L	50.0		88.4	40-140			
Anthracene (SIM)	49.1	4.0	μg/L	50.0		98.3	40-140			
Benzo(a)anthracene (SIM)	48.5	1.0	μg/L	50.0		97.1	40-140			
Benzo(a)pyrene (SIM)	51.6	2.0	μg/L	50.0		103	40-140			
Benzo(b)fluoranthene (SIM)	53.7	1.0	μg/L	50.0		107	40-140			
Benzo(g,h,i)perylene (SIM)	54.6	10	μg/L	50.0		109	40-140			
Benzo(k)fluoranthene (SIM)	52.5	4.0	μg/L	50.0		105	40-140			
Chrysene (SIM)	47.3	4.0	μg/L	50.0		94.6	40-140			
Dibenz(a,h)anthracene (SIM)	57.9	2.0	μg/L	50.0		116	40-140			
Fluoranthene (SIM)	51.2	10	μg/L	50.0		102	40-140			
Fluorene (SIM)	48.4	20	μg/L	50.0		96.9	40-140			
Indeno(1,2,3-cd)pyrene (SIM)	60.7	2.0	μg/L	50.0		121	40-140			
2-Methylnaphthalene (SIM)	43.4	20	μg/L	50.0		86.8	40-140			
Naphthalene (SIM)	39.6	20	μg/L	50.0		79.2	40-140			
Phenanthrene (SIM)	46.6	1.0	μg/L	50.0		93.2	40-140			
Pyrene (SIM)	45.4	20	μg/L	50.0		90.9	40-140			
Surrogate: Nitrobenzene-d5	87.1		$\mu g/L$	100		87.1	30-130			
Surrogate: 2-Fluorobiphenyl	90.5		$\mu g/L$	100		90.5	30-130			
Surrogate: p-Terphenyl-d14	73.7		$\mu g/L$	100		73.7	30-130			
LCS Dup (B258763-BSD1)				Prepared: 05	5/21/20 Anal	yzed: 05/27/2	20			
Acenaphthene (SIM)	43.9	6.0	μg/L	50.0		87.7	40-140	3.49	20	
Acenaphthylene (SIM)	42.5	4.0	μg/L	50.0		85.1	40-140	3.83	20	
Anthracene (SIM)	47.5	4.0	μg/L	50.0		95.0	40-140	3.39	20	
Benzo(a)anthracene (SIM)	47.2	1.0	μg/L	50.0		94.5	40-140	2.71	20	
Benzo(a)pyrene (SIM)	50.2	2.0	μg/L	50.0		100	40-140	2.75	20	
Benzo(b)fluoranthene (SIM)	52.7	1.0	μg/L	50.0		105	40-140	1.92	20	
Benzo(g,h,i)perylene (SIM)	53.0	10	μg/L	50.0		106	40-140	2.90	20	
Benzo(k)fluoranthene (SIM)	51.4	4.0	μg/L	50.0		103	40-140	2.12	20	
Chrysene (SIM)	46.0	4.0	μg/L	50.0		92.0	40-140	2.79	20	
Dibenz(a,h)anthracene (SIM)	56.8	2.0	μg/L	50.0		114	40-140	1.99	20	
Fluoranthene (SIM)	49.4	10	μg/L	50.0		98.8	40-140	3.58	20	
Fluorene (SIM)	47.3	20	μg/L	50.0		94.6	40-140	2.34	20	
Indeno(1,2,3-cd)pyrene (SIM)	59.0	2.0	μg/L	50.0		118	40-140	2.94	20	
2-Methylnaphthalene (SIM)	42.2	20	μg/L	50.0		84.5	40-140	2.66	20	
INAPIRIALIZZA (SIM)	39.2	20	μg/L	50.0		78.4	40-140	1.07	20	
Purene (SIM)	44.9	1.0	μg/L	50.0		89.9	40-140	3.67	20	
	44.5	20	µg/L	50.0		89.0	40-140	2.14	20	
Surrogate: Nitrobenzene-d5	83.6		μg/L	100		83.6	30-130			
Surrogate: 2-Fluorobiphenyl	88.7		μg/L	100		88.7	30-130			
Surrogate: p-Terphenyl-d14	72.7		μg/L	100		72.7	30-130			

‡



FLAG/QUALIFIER SUMMARY

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

compound. Reported result is estimated.



CERTIFICATIONS

Certified Analyses included in this Report

Analyte	Certifications
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-Dichloro 2 hutana	NC
Dicklorediflueremethane (Freen 12)	
1.1 Dishlaraathana	NC
1.2 Dichloroethane	NC
1.1 Dichloroethylene	NC
cis-1 2-Dichloroethylene	NC
trans-1 2-Dichloroethylene	NC
1 2-Dichloronronane	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
SW-846 8270E in Soil	
Acetophenone	NY,NH,ME,NC,VA
Aniline	NY,NH,ME,NC,VA
Benzidine	CT,NY,NH,ME,NC,VA
Benzoic Acid	NY,NH,ME,NC,VA
Bis(2-chloroethoxy)methane	CT,NY,NH,ME,NC,VA
Bis(2-chloroethyl)ether	CT,NY,NH,ME,NC,VA
Bis(2-chloroisopropyl)ether	CT,NY,NH,ME,NC,VA
Bis(2-Ethylhexyl)phthalate	CT,NY,NH,ME,NC,VA
4-Bromophenylphenylether	CT,NY,NH,ME,NC,VA
Butylbenzylphthalate	CT,NY,NH,ME,NC,VA
Carbazole	NC
4-Chloroaniline	CT,NY,NH,ME,NC,VA
4-Chloro-3-methylphenol	CT,NY,NH,ME,NC,VA
2-Chloronaphthalene	CT,NY,NH,NC,VA
2-Chlorophenol	CT,NY,NH,ME,NC,VA
4-Chlorophenylphenylether	CT,NY,NH,ME,NC,VA



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



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Certified Analyses included in this Report

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



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Certified Analyses included in this Report

Analyte	Certifications	
SW-846 8270E in Water		
Pentachlorophenol	CT,NY,NC,ME,NH,VA	
Phenol	CT,NY,NC,ME,NH,VA	
Pyridine	CT,NY,NC,ME,NH,VA	
1,2,4,5-Tetrachlorobenzene	NY,NC	
1,2,4-Trichlorobenzene	CT,NY,NC,ME,NH,VA	
2,4,5-Trichlorophenol	CT,NY,NC,ME,NH,VA	
2,4,6-Trichlorophenol	CT,NY,NC,ME,NH,VA	
2-Fluorophenol	NC	

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

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

		[]									_																		Tat	ole c	of C	Conte	ents
	Page of	# of Containers	² Preservation Code	³ Container Code	and a strate literation Sources	Field Filtered		Orthophics and the Sommer		Lab to Filter		Matrix Codes: GW = Ground Water	WW = Waste Water DW = Drinking Water	A = Air S = Soil	SL = Studge	0 = Other (please	define)	² <u>Preservation Codes</u> : I = loed	H = HCL M = Methanol	N = Nitric Acid S = Sulfuric Acid	B = Sodium Bisulfate Y = Sodium Hydrovide	Thissifiate	0 = Other (please define)		³ <u>Container Codes</u> : A = Amher Glace	G = Glass D = Plastic	ST = Sterile	V = Vial S = Summa Canister	T = Tedlar Bag 0 = Other (please	denne)		PCB ONLY Soxhiet	Non Soxhlet
03242017	39 Spruce Street East Longmeadow, MA 01028				ANALYSIS REQUESTED																		s to indicate possible sample concentration	Conc Code column above:	L- LOW, C - CICAI, U - CINIOWI	Program Information		HSB Orphaned Landfill	State Lead Other:		NELAC and AlHA-LAP, LLC Accredited	Other	AIHA-LAP, LLC
pavel 53 http://www.contestlabs.com boc # 379 Rev 1	CHAIN OF CUSTODY RECORD (North Carolina)	7-Day 10-Day 10-Day 22	Due Date:	Rush Approval Regulted				Other:	CLP Like Data Pkg Required:		Endino 10 *:	Date (ing Composite Grab Code Code		XX XCM XX									Please use the following code:	within the C		North Carolina Detection Limit Requirements	ZT	Swst	HISB MSCC			tity Government Municipality	Federal Brownfield City School
NEOS38	-LESE Accounter Fax: 413-525-2332 Fax: 413-525-6405	Email: info@contestlabs.com	raper Holen Hssociates	dinburgh S. Dr., Cory, NC. 27511		avoore YJA D'53	70050704-010102	M. Bronson	me/Number:			der# Client Sample ID / Description		1 5R-2 5/16/20												ignateme) Date/Time:			ignatured Date/Time: (6 30	athre) (A A Date/Time: 011=	5 C 8 2 16 10	ignature) Date/Time: Project En	ature) Date/Time:
	COD-	5.4 South resources for the second of the second se Second second sec		Address: 1/4 E	Phone:		Project Location: Project Number:	Project Manager:	Con-Test Quote Nar	Invoice Recipient:		Work On					- Address -						Comments:			Bernjourshed by: (s	Decoved by (com	LOCA	Re Dy: (s	and by: (signe	28	Re D lished by: (s	Re ed by: sign

IMPORTANT!

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



Your Packaging



(?)

5/20/2020 by 10:30 am

Packaging

Deliver Weekday, Non Standard

Have Not Confirm	ned Sample Container		con-test	, (B)
Numbers With Lab S	taff Before Relinquishing		ANALYTICAL LABORATORY	
Over Sample	S		Doc# 277 Rev 5 2017	
Login Sample Rec	eipt Checklist - (Rejectio	n Criteria Listing - Usi	ng Acceptance Policy) Any False	
Statem	ent will be brought to the	e attention of the Clien	t - State I rue or Paise	
Client Draper	Ader		Time	
Received By	Par	Date <u>5 μ</u>		
How were the samples	In Cooler T	No Cooler	On Ice No Ice	
received?	Direct from Sampling		Ambient Melted Ice	
	By Gun	# 2	Actual Temp - 28	
Were samples within	Dy Our		Actual Temp -	
Temperature? 2-6°C	By Blan	K #	es Tampered with?	
Was Custody Se	al Intact? <u>(A</u>	Does Chain A	oree With Samples?	
Was COC Relin	quisned (amples?	<u> </u>	
Are there broken/le	eaking/loose caps on any s	Were samples rec	eived within holding time?	
Is COC in ink/ Legible ?		Analysis	Sampler Name	
DIG COC Include an	Project	ID's T	Collection Dates/Times	
Are Semple labels filled	Lout and legible?			
Are sample labels filled	F	Who w	as notified?	
Are there Rushes?	MKF	> Who w	as notified?	
Are there Short Holds?	F	Who w	as notified?	
is there enough Volume	? 1		T	
is there Headspace whe	ere applicable?	MS/MSD	? <u> </u>	
Proper Media/Container	s Used?	Is splittin	g samples required?	· · · · · · · · · · · · · · · · · · ·
Were trip blanks receive	əd?E	On COC	Passa	and a state of the second s Second second
Do all samples have the	proper pH?	A Acid		
	Containers:			
Unp-	1 Liter Amb. 2	1 Liter Plastic	16 oz Amb.	
HCL- 3	500 mL Amb.	500 mL Plastic	4oz Amb/Clear	
Meoh-	250 mL Amb.	250 mL Plastic	207 Amb/Clear	
Bisulfate-	Flashpoint	Other Plastic	Encore	
		Plastic Bag	Frozen:	
Thiosulfate-	Perchlorate	Ziplock		
Sulturic-				
		UNUSEO MEDIA		
Viale F		1 Liter Plastic	16 oz Amb.	
Unp-	500 ml Amb	500 mL Plastic	8oz Amb/Clear	
HUL-	250 ml Amb	250 mL Plastic	4oz Amb/Clear	
Neon-	Col /Bacteria	Flashpoint	2oz Amb/Clear	
	Other Plastic	Other Glass	Encore	
Thiosulfate-	SOC Kit	Plastic Bag	Frozen:	
Culturio	Perchlorate	Ziplock		
ISUBURG* F				