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August 5, 2020

ATO 449

*Via Email

Mr. Joshua Haugh Project Manager New York State Department of Environmental Conservation Division of Environmental Remediation 1130 North Westcott Road Schenectady, NY 12306 joshua.haugh@dec.ny.gov

RE: Post Remedial Action Groundwater Evaluation Hamilton Hill II – Target Area 1 Site City of Schenectady, Schenectady County BCP Site ID No.: C447052 C.T. Male Project No. 16.6334

Dear Mr. Haugh:

This letter report presents the findings of a post remedial action groundwater evaluation conducted at the Hamilton Hill – Target Area 1 Brownfield Cleanup Program (BCP) Site (BCP Site No. C447052) located at 830 & 834 Albany Street in the City of Schenectady, Schenectady County, New York (see Site Location Map in Attachment A). The groundwater evaluation was conducted as a component of the Site's remedy contained in the November 2019 Decision Document for the Site.

The remedial action at the Site involved the remediation of source areas of contamination including impacted fill/soil and petroleum bulk storage tanks. The impacted fill/soil was remediated to Unrestricted Use Soil Cleanup Objectives (SCOs) via excavation and off-site disposal as documented by endpoint soil sampling analyses. The excavations were backfilled with imported fill that was tested to meet Unrestricted Use SCOs criteria. Two (2) aboveground and two (2) underground petroleum bulk storage tanks were closed by removal and disposed off-site.

The post remedial action groundwater evaluation involved the collection of groundwater samples from six (6) strategically placed monitoring wells installed after completion of the above referenced remedial actions. The means and methods for the groundwater evaluation were presented in the Department approved Post Remedial Action Groundwater Sampling Plan, dated April 27, 2020. The Sampling Plan and Department approval letter are presented as Exhibit 1.

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<u>Methods</u>

The newly installed monitoring wells are depicted as RAMW1 to RAMW6 (highlighted in yellow) on Figures 2 and 3 in Attachment A. Figure 2 depicts the monitoring well locations in relation to sampling locations completed as part of the Remedial Investigation (RI) and previous Phase II Environmental Site Assessments (ESAs) of the Site. Figure 3 depicts the monitoring well locations in relation to buildings currently being developed on the Site.

On June 25, 2020, Precision Environmental Services, Inc. (PES) advanced six (6) soil borings across the Site to facilitate the installation of the monitoring wells. The borings were completed via direct-push drilling methods employing a track mounted Geoprobe unit.

At each boring location, a two-inch diameter macro-core sampler was advanced at continuous four (4) foot depth intervals to the termination depths of the borings. The recovered soil samples were visually classified by a C.T. Male field technician and recorded on individual Direct-Push Exploration Logs, which are presented in Attachment B.

Soil samples were collected from the soil borings at continuous intervals by a C.T. Male field geologist for field screening for organic vapors with a photoionization detector (PID) meter and for visual/olfactory evidence of contamination. The Organic Vapor Headspace Analysis Logs are presented in Attachment C.

Each of the borings were converted to groundwater monitoring wells for the purpose of collecting groundwater samples for laboratory analysis. The monitoring wells were constructed of one (1)-inch diameter PVC slotted screen and riser pipe. The monitoring wells were protected at the ground surface with curb box enclosures set in concrete pads. Monitoring Well Construction Logs are presented in Attachment D.

To prevent the potential for cross contamination between the test boring locations, drilling tools and sampling equipment that came into contact with the Site soils and groundwater were decontaminated prior to the start of the drilling activities and between boring locations utilizing a detergent/water wash and tap water rinse. Soil and groundwater samples were handled with a new pair of gloves to deter cross contamination of the soil and groundwater samples collected for screening and/or laboratory analysis.

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Drill cuttings of soil located beneath the clean fill that was imported onto the Site as part of the remedial action, and groundwater development/purge water, were transferred into designated 55-gallon drums (2 total) with covers. The drum contents will be profiled and disposed of off-site at a treatment, storage and/or disposal facility (TSDF) at a later date. The methods for profiling/disposal of the drum contents will be presented to the Department Project Manager for approval.

The monitoring wells were developed on June 29, 2020 by a C.T. Male field engineer. The wells were surged, and approximately five (5) well volumes were evacuated from each monitoring well. Temperature, pH, specific conductivity and turbidity field parameters were recorded for each well volume evacuated from the monitoring wells.

Groundwater samples were collected from the monitoring wells on June 30, 2020. Prior to sampling, the water levels were recorded in each well by measuring the static water levels from the top of the PVC casing utilizing a water level meter. The groundwater samples were collected employing low flow purging/sampling methods utilizing a peristaltic pump with new tubing dedicated to each well. The groundwater samples were transferred into new laboratory supplied sampling containers and delivered to Alpha Analytical, Inc.'s (Alpha) Albany, New York service center. Temperature, pH, specific conductivity and turbidity field parameters had stabilized prior to collection of the groundwater samples. The groundwater samples were submitted to Alpha for laboratory analysis for volatile organic compounds (VOCs) plus 10 tentatively identified compounds (TICs) by EPA Method 8260 and semi volatile organic compounds (SVOCs) plus 20 TICs by EPA Method 8270.

Findings

The soil borings were advanced to depths ranging from 18 to 20 feet below the ground surface (bgs). Approved fill imported onto the Site to backfill the remedial excavations was observed in all of the borings from the ground surface to depths ranging from six (6) to eight (8) feet bgs. The fill was composed primarily of fine to coarse sands with occurrences of silt and gravel. Underlying the fill was native soil which generally consisted of sands with varying percentages of silt and gravel grading into silt and clay with varying percentages of sand and gravel. The soils became wet at depths that ranged from 11 feet bgs at RAMW4 to 15 feet bgs at RAMW6. See Direct Push Logs in Attachment B.

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As presented on the Organic Vapor Headspace Analysis Logs in Attachment C, the PID readings for the soil samples collected from the borings were less than one (1) part per million (ppm) above background with the exception of a saturated native soil sample collected from the 12 to 14-foot sampling depth interval at RAMW3, which is located outside of the eastern boundary of the 830 Albany Street parcel defined under the Brownfield Cleanup Program (BCP). Soils collected at this depth interval were saturated, exhibited a PID reading of 69.7 ppm above background, and exhibited black staining and a petroleum-type odor. In consultation with the Department's Project Manager for this site within a Brownfield Program, the NYSDEC Spills Hotline was contacted and spill #2002279 was assigned to the Site relative to the subjectively impacted soil.

Groundwater conditions were assessed during collection of groundwater samples from monitoring wells RAMW1 to RAMW6 on June 30, 2020. The following table presents static groundwater levels and the field parameters (temperature, pH, specific conductivity, turbidity, olfactory observations) recorded at each monitoring well during the groundwater sampling event.

Monitoring	Water	Temp.	pН	Specific	Turbidity	Olfactory Observation
Well	Level			Conductivity		
RAMW1	13.64	14.7	8.02	765 μS	> 50	No Odor/No Sheen
RAMW2	12.90	16.2	7.65	676 μS	> 50	No Odor/No Sheen
RAMW3	11.54	15.8	7.53	1,811 µS	6.67	Very Slight Petroleum-
						Type Odor/No Sheen
RAMW4	10.55	15.6	7.68	1,792 μS	3.88	No Odor/No Sheen
RAMW5	11.96	14.6	7.51	2,420 µS	4.61	No Odor/No Sheen
RAMW6	12.83	14.5	7.57	1,904 µS	4.31	No Odor/No Sheen

Water Level measured in feet below the top of the monitoring well PVC riser casing.

Temperature recorded in degrees Celsius (°C).

Specific Conductivity measured in microsiemens (μ S).

Turbidity recorded in nephelometric turbidity units (NTU).

Static groundwater levels in the six (6) monitoring wells on the date of sampling ranged from approximately 11.96 feet bgs at RAMW5 to 13.64 feet bgs at RAMW1. The horizontal location and vertical elevation of each monitoring well was surveyed by a C.T. Male survey crew on July 24, 2020. Based on the measured depth to groundwater and the monitoring well elevations, groundwater contours were developed. As depicted on Figure 2 in Attachment A, the localized groundwater flow direction is from southwest to northeast, which is consistent with the groundwater flow observed during the RI.

The groundwater samples were analyzed for VOCs plus 10 TICS and SVOCs plus 20 TICs. The analytical results summary table is presented in Appendix E. The laboratory

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report is presented in Appendix F. The laboratory data was independently validated by Environmental Data Services, Inc. (EDS) of Virginia Beach, Virginia. The Data Usability Summary Report (DUSR) is presented in Appendix G. The analytical results disclosed the following.

- Petroleum related compounds and TICs were not detected in groundwater sampled from monitoring well RAMW3. Saturated soil at RAMW3 had exhibited elevated PID readings, staining and petroleum-type odors during advancement of the soil boring and had exhibited a slight petroleum-type odor during groundwater sample collection. RAMW3 is outside of the eastern boundary of the 830 Albany Street parcel defined under the BCP.
- Tetrachloroethene was the only compound detected above its New York State (NYS) Ambient Water Quality Standard (AWQS). Tetrachloroethene was detected slightly above its NYS AWQS of 5 parts per billion (ppb) in three (3) monitoring wells within the 830 Albany Street parcel only (not 834 Albany Street). These wells included RAMW1 (7.3 ppb), RAMW3 (6.8 ppb) and RAMW4 (8.7 ppb) only. Tetrachloroethene was not detected above its NYS AWQS in the two (2) monitoring wells (RAMW5, RAMW6) within the 834 Albany Street parcel.
- Chloroform was recently detected at a concentration range of 1.7 ppm to 2.8 ppm in two (2) of the six (6) monitoring wells (RAMW2, RAMW3), which is below its NYS AWQS of 7 ppb. These wells are each located within the 830 Albany Street parcel. Chloroform, which was detected slightly above its NYS AWQS in groundwater samples collected during the past RI was not detected above NYS AWQS in the post remedial action groundwater samples. Chloroform was detected at a concentration range of 8 ppm to 10 ppm in seven (7) monitoring wells (MW1, MW2, MW5, RIGP1, RIGP2, RIMW2, RIMW4) sampled during the past RI. The wells were each located within the 830 Albany Street parcel.
- SVOCs were not detected above the laboratory's method detection limits in any of the post remedial action groundwater samples. SVOCS were detected above NYS AWQS at varying frequencies in past RI sampled monitoring well MW2 (7 SVOCS) bordering the eastern boundary of the 830 Albany Street parcel, and past RI sampled monitoring wells 834-MW2 (1 SVOC), RIMW5 (3 SVOCs) and RIGP5 (5 SVOCs) within the 834 Albany Street parcel.

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Conclusions and Recommendations

The remedial action was successful at eliminating known and potential source areas of contamination from the Site as demonstrated by the remedial excavation endpoint sampling results and observations. These sources included impacted fill/soil and petroleum bulk storage tanks. Fill imported onto the Site after remediation of the source areas met criteria for unrestricted use of the Site.

Petroleum-type compounds were not detected in groundwater sampled from monitoring well RAMW3, which was the location that exhibited subjective impacts (elevated PID readings, staining, petroleum-type odor) in a single isolated sample of saturated soil collected during advancement of the soil boring for installation of the monitoring well. Based on these facts, and that RAMW3 is located outside of the BCP Site boundary, it is recommended that associated spill #2002279 be closed by the Department requiring no further action.

The tetrachloroethene detections above NYS AWQS are consistent in persistence with tetrachloroethene detections from nearby monitoring wells sampled as part of the RI (see Figure 2 for comparisons). For RAMW1 (tetrachloroethene = 7.3 ppb), tetrachloroethene was detected at a concentration of 6.1 ppb in nearby RI monitoring well RIMW2. For RAMW3 (tetrachloroethene = 6.8 ppb), tetrachloroethene was detected at a concentration range of 6.9 to 8.6 ppb in nearby RI monitoring wells RIGP2 (6.9 ppb), RIMW3 (7.2 ppb) and RIGP3 (8.6 ppb). For RAMW4 (tetrachloroethene = 8.7 ppb), tetrachloroethene was not detected above its NYS AWQS at RI monitoring well MW2.

Chloroform, which was detected slightly above its NYS AWQS of 7 ppb in groundwater samples collected from seven (7) RI monitoring wells was recently detected below its NYS AWQS in groundwater samples collected from two (2) of the six (6) post remedial action monitoring wells within the 830 Albany Street parcel.

The post remedial action groundwater sampling analytical results were also compared to analytical results of groundwater samples collected from the Site during Phase II ESA investigations conducted in 2016 and 2018 that exceeded NYS AWQS (see Figure 4 in Attachment A for analytes that were detected above NYS AWQS). As depicted on Figure 4, detections of cis-1,2-dichloroethene above NYS AWQS in Phase II ESA monitoring wells CTM-MW-02 and CTM-MW-03, and acetone and toluene above NYS AWQS in Phase II ESA monitoring wells CTM-MW-01, CTM-MW-02, CTM-MW-03 and CTM-MW-05 were non-detect in groundwater samples collected from all of the post remedial action

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monitoring wells. The aforementioned Phase II ESA monitoring wells were each located within the 830 Albany Street parcel. Detections of tetrachloroethene above NYS AWQS in the Phase II ESA monitoring wells ranged in concentration from 6.2 ppb to 14 ppb and were confined to Phase II ESA monitoring wells CTM-MW-01, CTM-MW-02, CTM-MW-04 and CTM-MW-05 within the 830 Albany Street parcel. Detections of tetrachloroethene above NYS AWQS in the nearby post remedial action monitoring wells ranged in concentration from 6.8 ppb to 8.7 ppb.

The Site was successfully remediated to unrestricted use and potential/known source areas of contamination have been removed. As tetrachloroethene was detected slightly above its NYS AWQS and is consistent with detections of tetrachloroethene in past RI and 2016/2018 Phase II ESA groundwater samples, and chloroform and SVOCs were detected below their NYS AWQS, a bulk reduction of chloroform and SVOCs in groundwater has been attained by the remedial action. Concentrations of tetrachloroethene have also attained asymptotic conditions, and that potential/known source areas that could contribute to future groundwater contamination have been eliminated.

Based on the remedial actions completed, and analytical data developed prior to and as a function of the BCP, the following conclusions and recommendations are made.

- The groundwater evaluation presented in this report satisfies the groundwater evaluation criteria of the Decision Document since groundwater is not used as a potable water supply and the remedial action removed all on-site sources that would contribute to groundwater contamination. Further groundwater monitoring is not required as Track 1 Unrestricted Use criteria has been substantially attained for the groundwater component of the remedy.
- The concentrations of compounds remaining in groundwater are below or approaching their regulatory standards at asymptotic levels. This, coupled with the depth to groundwater being greater than 11 feet below the finished floor of the new building, and the installation of a new vapor barrier and new radon reduction system as part of building construction, potential vapor intrusion issues for future occupants of the buildings is considered negligible.

Please do not hesitate to contact me should you require further information.

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Respectfully submitted, C.T. MALE ASSOCIATES

Jeffry A. Marp

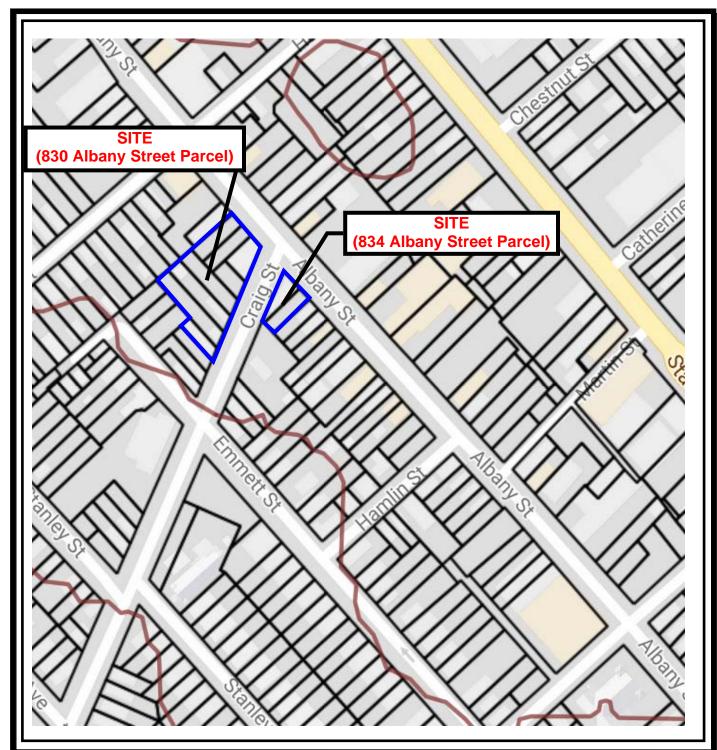
Jeffrey A. Marx, P.E. Sr. Environmental Engineer

Attachment A:	Figures
Attachment B:	Direct Push Logs
Attachment C:	Organic Vapor Headspace Analysis Logs
Attachment D:	Monitoring Well Construction Logs
Attachment E:	Analytical Results Summary Table
Attachment F:	Analytical Report
Attachment G:	Data Usability Summary Report
Exhibit 1:	Department Approved Post Remedial Action Groundwater
	Sampling Plan

ec: Sue McCann, Hamilton Hill II Limited Partnership Kelly Melaragno, Hamilton Hill II Limited Partnership Janis Stewart, Hamilton Hill II Limited Partnership Christine Vooris, NYSDOH Kirk Moline, P.G., C.T. Male Associates Steve Bieber, CHMM, C.T. Male Associates

ATTACHMENT A

FIGURES



MAP REFERENCE

USGS 7.5 Minute Topographical Map Schenectady, New York Quadrangle Year 2013



FIGURE 1: SITE LOCATION MAP HAMILTON HILL II - TARGET AREA 1 SITE

CITY OF SCHENECTADY

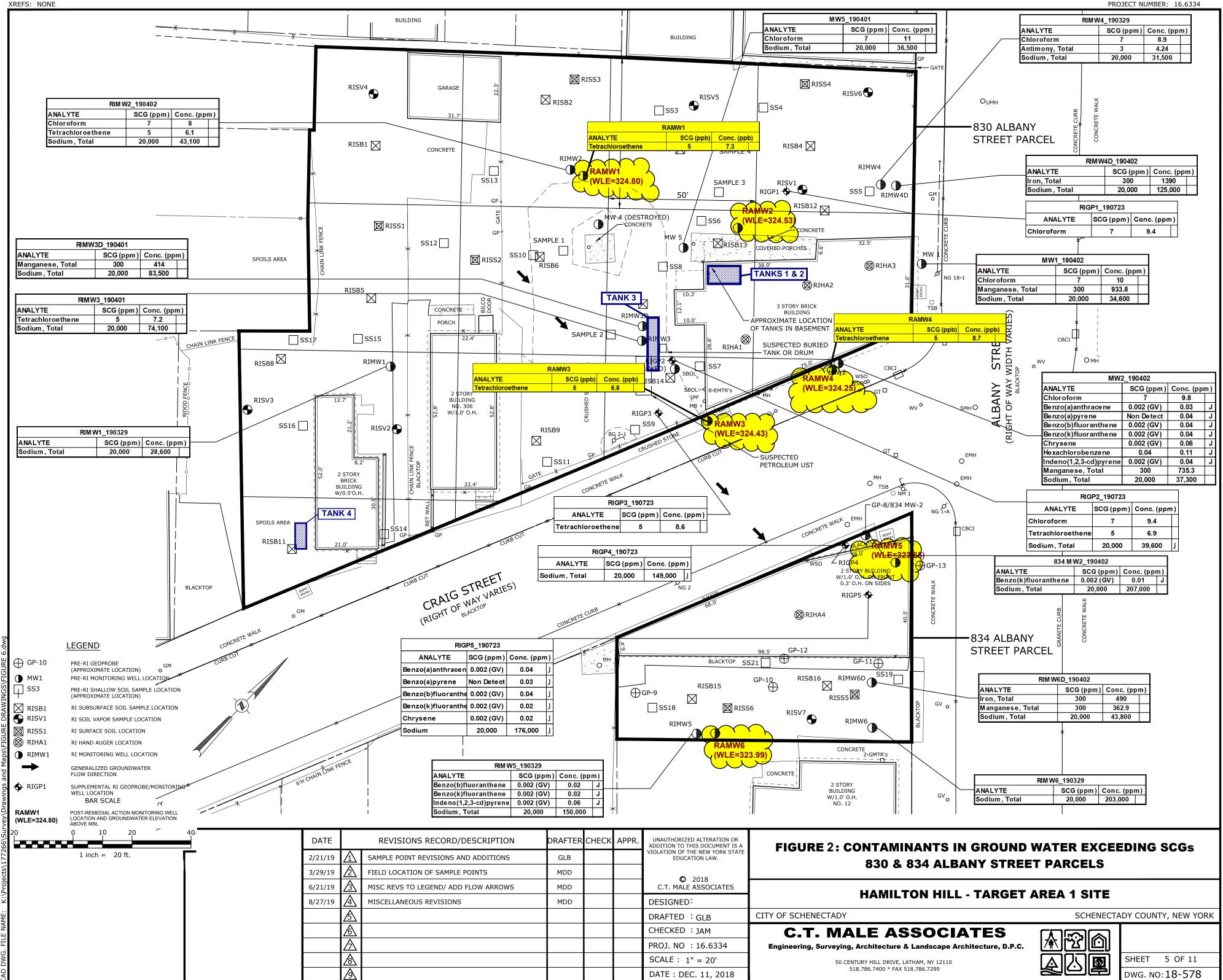
SCHENECTADY COUNTY, NY

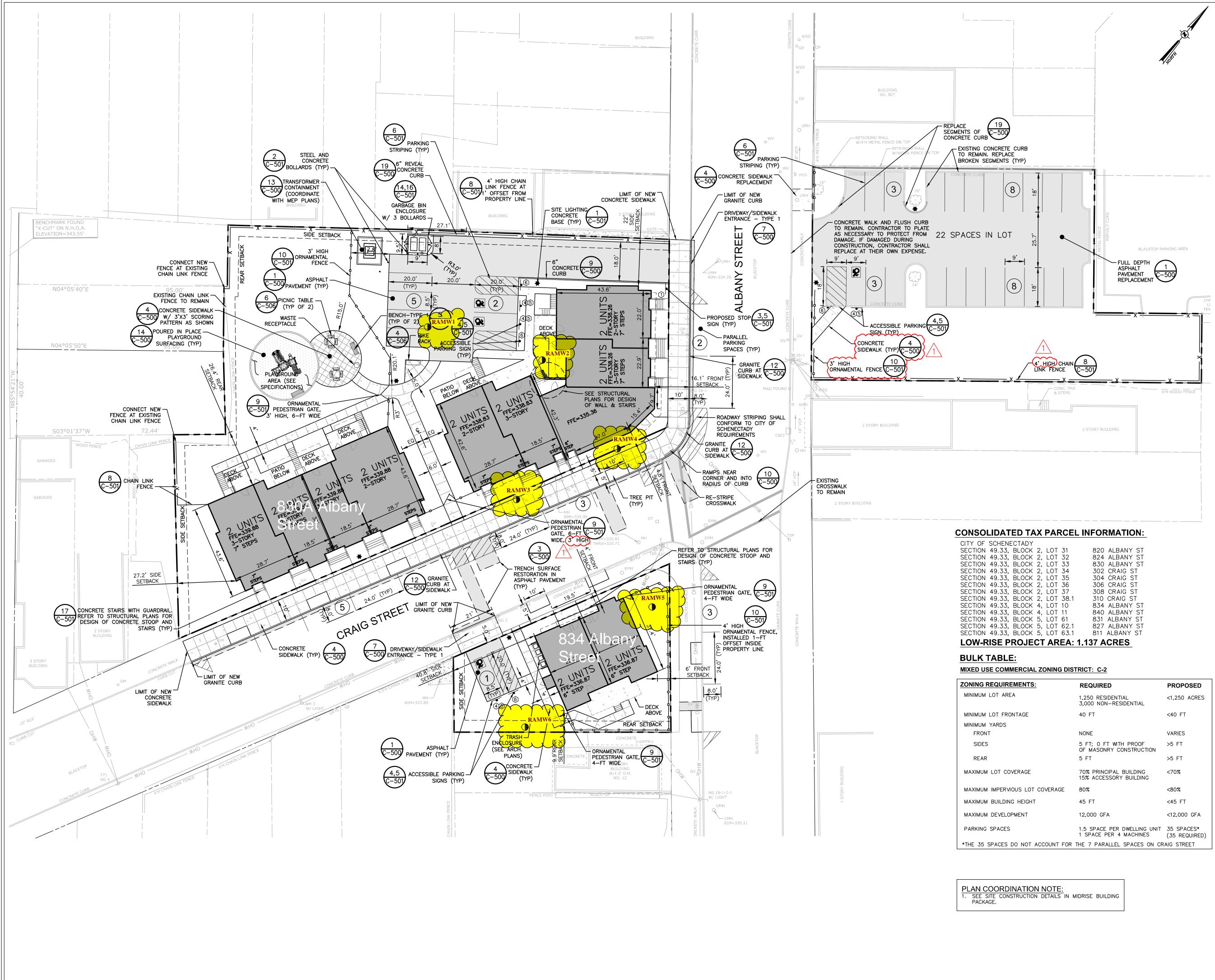
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PROJECT No: 16.6334

The locations and features depicted on this map are approximate and do not represent an actual survey.







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	830 /	ALBANY ST
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	304 (CRAIG ST
	306 (CRAIG ST
	308 (CRAIG ST
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	831 A	ALBANY ST
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	REQUIRED	PROPOSED
	1,250 RESIDENTIAL 3,000 NON-RESIDENTIAL	<1,250 ACRES
	40 FT	<40 FT
	NONE	VARIES
	5 FT; 0 FT WITH PROOF OF MASONRY CONSTRUCTION	>5 FT
	5 FT	>5 FT
	70% PRINCIPAL BUILDING 15% ACCESSORY BUILDING	<70%
	80%	<80%
	45 FT	<45 FT
	12,000 GFA	<12,000 GFA
	1.5 SPACE PER DWELLING UNIT 1 SPACE PER 4 MACHINES	35 SPACES* (35 REQUIRED)
۲	THE 7 PARALLEL SPACES ON CR	AIG STREET

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GENERAL CONSTRUCTION:

- THE CONTRACTOR SHALL PROTECT EXISTING PROPERTY LINE MONUMENTATION. ANY MONUMENTATION DISTURBED OR DESTROYED, AS JUDGED BY THE ENGINEER OR OWNER, SHALL BE REPLACED AT THE CONTRACTOR'S EXPENSE AND UNDER THE SUPERVISION OF A NEW YORK STATE LICENSED LAND SURVEYOR.
- 2. ALL PAVEMENT RESTORATION SHALL MEET AND MATCH EXISTING GRADES. ALL SAWCUT LINES SHALL BE PARALLEL AND CURVILINEAR TO EXISTING OR PROPOSED CURBING AND SHALL BE A CONSTANT
- DISTANCE OF 18" MIN AWAY. ALL ARCHITECTURE IS SUBJECT TO PLANNING BOARD REVIEW. 5. NOTIFY ENGINEER 48 HOURS PRIOR TO INITIALIZATION OF ANY WORK ON SITE.
- 6. THE ENGINEER SHALL BE NOTIFIED IN WRITING OF ANY CONDITIONS THAT VARY FROM THOSE SHOWN ON THE PLANS. THE CONTRACTOR'S WORK SHALL NOT VARY FROM THE PLANS WITHOUT PRIOR REVIEW FROM THE ENGINEER. CONTRACTOR IS RESPONSIBLE FOR EMPLOYING AND MAINTAINING ALL
- TRAFFIC CONTROL AND SAFETY MEASURES DURING CONSTRUCTION. CONTRACTOR IS RESPONSIBLE FOR PROPERLY & SAFELY MAINTAINING
- AREA BETWEEN ALL ADJOINING PROPERTIES. NO WORK, STORAGE OR TRESPASS SHALL BE PERMITTED BEYOND THE SITE PROPERTY LINES OR PUBLIC RIGHT-OF-WAY. 10. ALL EXISTING LAWN AREA, CURBING, PAVING, SIDEWALKS, CULVERTS
- OR OTHER PUBLIC OR PRIVATE PROPERTY DAMAGED BY TRENCHING OR EXCAVATION OPERATIONS SHALL BE REPLACED OR REPAIRED TO A CONDITION EQUAL TO EXISTING, AS DESCRIBED IN CONTRACT DOCUMENTS OR AS ORDERED BY ENGINEER (AOBE). MAILBOXES, SIGN POSTS, ETC SHALL BE PROTECTED OR REMOVED AND REPLACED EXACTLY AS THEY WERE BEFORE BEING DISTURBED. REMOVE AND REPLACE AFFECTED CURBING AND SIDEWALK TO NEAREST JOINT. REMOVE PAVEMENT AND REPLACE TO SAW CUT LINE, SAW CUT IN STRAIGHT LINE TO POINT NEEDED TO BLEND GRADE, REMOVE LAWN AND REPLACE TO MINIMUM LIMIT OF EXCAVATION.

LAYOUT

- BUILDING DIMENSIONS TO BE TAKEN FROM ARCHITECTURAL BUILDING PLANS. NOTIFY THE ENGINEER OF ANY DEVIATION FROM CONDITIONS SHOWN ON THIS PLAN.
- THE CONTRACTOR SHALL BE RESPONSIBLE FOR COORDINATING ALL FIELD LAYOUT. THE CONTRACTOR SHALL TAKE TIES TO ALL UTILITY CONNECTIONS AND PROVIDE MARKED-UP AS BUILT PLANS FOR ALL UTILITIES SHOWING TIES TO CONNECTIONS, BENDS, VALVES, LENGTHS OF LINES AND INVERTS. AS-BUILT PLANS SHALL BE REVIEWED BY THE OWNER AND THE ENGINEER AND THE CONTRACTOR SHALL PROVIDE ANY CORRECTION OR ADDITIONS TO THE SATISFACTION OF THE OWNER AND THE ENGINEER BEFORE UTILITIES WILL BE ACCEPTED.

PAVING:

- NO VEHICULAR TRAFFIC OF ANY SORT SHALL BE PERMITTED ON THE SURFACE OF SUBBASE COURSE MATERIAL ONCE IT HAS BEEN FINE GRADED, COMPACTED, AND IS READY FOR PAVING. SUBBASE MATERIAL SO PREPARED FOR PAVING SHALL BE PAVED WITHIN THREE DAYS OF PREPARATION.
- 2. SUBBASE MATERIAL AND THE VARIOUS ASPHALT CONCRETE MATERIALS CALLED FOR IN THESE DRAWINGS SHALL CONFORM WITH THE REFERENCED SECTION OF THE NEW YORK STATE DEPARTMENT OF TRANSPORTATION STANDARD SPECIFICATIONS FOR CONSTRUCTION AND MATERIALS, DATED "LATEST EDITION". CONSTRUCTION SHALL BE AS FURTHER SET FORTH IN THOSE SPECIFICATIONS AND AS OTHERWISE PROVIDED FOR IN THESE DRAWINGS. 3. PLACE ASPHALT CONCRETE MIXTURE ON PREPARED SURFACE.
- SPREAD AND STRIKE-OFF USING A SELF-PROPELLED PAVING MACHINE, WITH VIBRATING SCREED. PLACEMENT IN INACCESSIBLE AND SMALL AREAS MAY BE BY HAND 4. PROVIDE JOINTS BETWEEN OLD AND NEW PAVEMENTS OR BETWEEN
- SUCCESSIVE DAY'S WORK. 5. TACK COAT WHEN SPECIFIED OR CALLED OUT ON THE DRAWINGS OR REQUIRED BY THE REFERENCED SPECIFICATION SHALL CONFORM WITH THE FOLLOWING
- A. TACK COAT SHALL MEET THE MATERIAL REQUIREMENTS OF 702-90 ASPHALT EMULSION FOR TACK COAT OF THE NEW YORK STATE DEPARTMENT OF TRANSPORTATION STANDARD SPECIFICATIONS FOR CONSTRUCTION AND MATERIALS, DATED "LATEST EDITION", SHALL BE APPLIED IN ACCORDANCE WITH SECTION 407 - TACK COAT SHALL BE IN ACCORDANCE WITH THOSE SPECIFICATIONS AND AS OTHERWISE PROVIDED FOR IN THESE DRAWINGS.
- B. REMOVE LOOSE AND FOREIGN MATERIAL FROM ASPHALT SURFACE BEFORE PAVING NEXT COURSE. USE POWER BROOMS, BLOWERS OR HAND BROOM. C. APPLY TACK COAT TO ASPHALT PAVEMENT SURFACES & AND
- SURFACES OF CURBS, GUTTERS, MANHOLES, AND OTHER STRUCTURES PROJECTING INTO OR ABUTTING PAVEMENT. DRY TO A "TACKY" CONSISTENCY BEFORE PAVING. D. TACK COAT ENTIRE VERTICAL SURFACE OF ABUTTING EXISTING
- PAVEMENT. 6. AFTER COMPLETION OF PAVING AND SURFACING OPERATIONS, CLEAN SURFACES OF EXCESS OR SPILLED ASPHALT, GRAVEL OR STONE

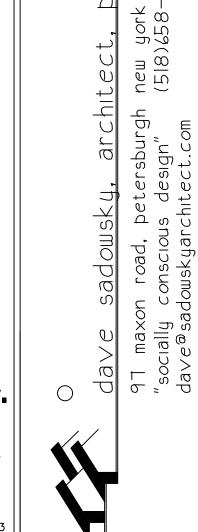
MATERIALS TO THE SATISFACTION OF THE ENGINEER. STRIPING:

- . STRIPE PAVEMENT AS INDICATED ON THE PLANS AND/OR IN ACCORDANCE WITH ALL APPLICABLE LOCAL, STATE AND FEDERAL REQUIREMENTS
- 2. COLOR: DRIVE LANE DIVIDERS WHITE OR AOBE NO PARKING ZONE WARNINGS - WHITE OR AOBE PARKING DIVIDERS - WHITE OR AOBE WALKING LINES - WHITE OR AOBE
 - HANDICAP PARKING LINES & SYMBOL BLUE
 - Dig Safely. New York ☐ Call Before You Dig ☐ Wait The Required Time 🔲 Confirm Utility Respons Respect The Marks Dig With Care 800-962-7962

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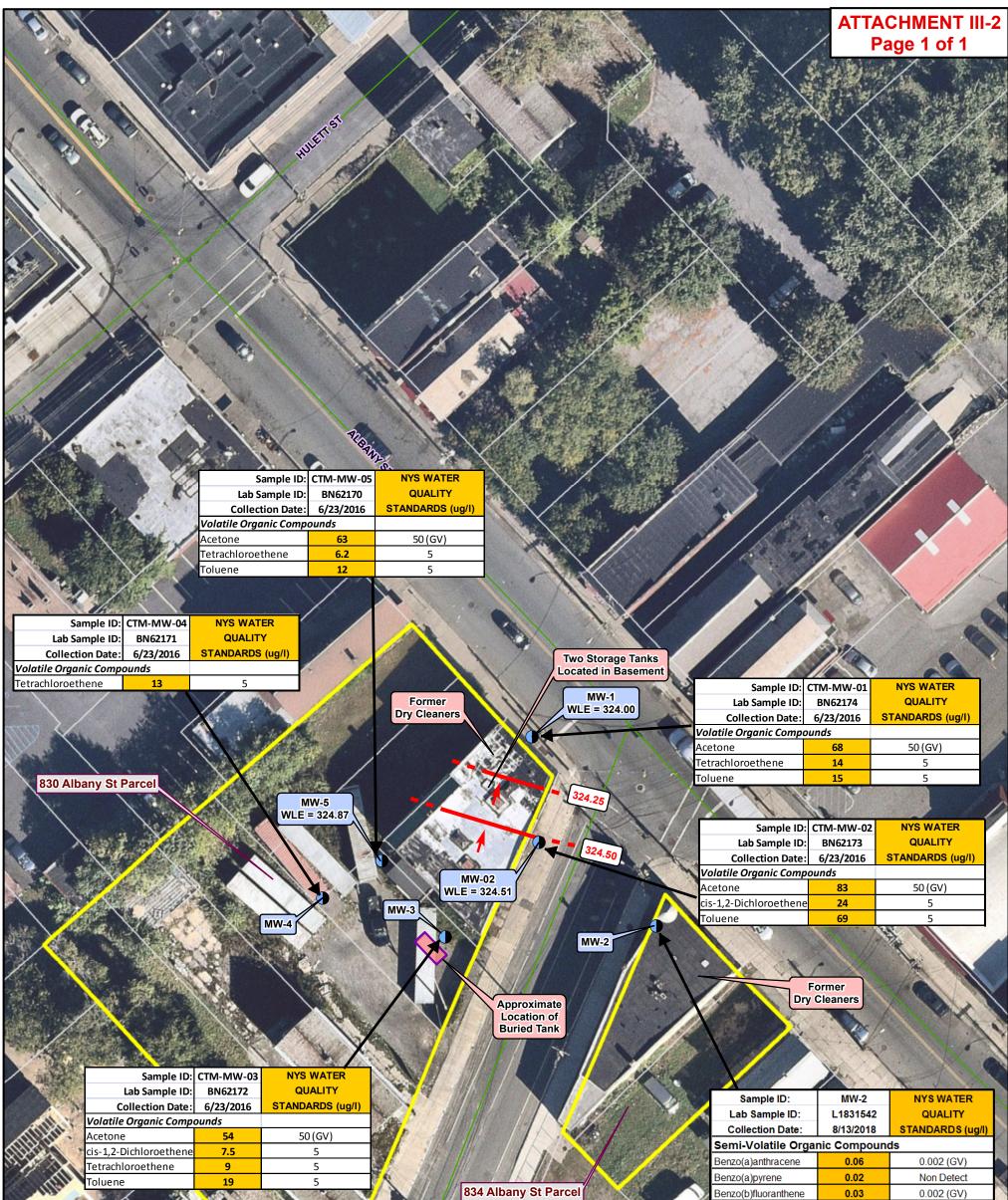
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	Chrysene	0.03	0.002 (GV)
X	Indeno(1,2,3-cd)pyrene	0.02	0.002 (GV)
1	Total Metals		
/	Iron, Total	31,200	300
	Lead, Total	39.48	25
	Manganese, Total	1,361	300
and the second	Sodium, Total	155,000	20000
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Copyright nearmap 2015





Project Number: 16 6334 Data Source: NYSGIS Clearinghouse Projection: State Plane NAD83 NYE (feet) Date: October 22, 2018 File: Fig2HamiltonHill2018_11x17.mxd GIS: C Secor

80 Legend FT Approximate Monitoring Well Locations -Road Centerlines Project Site

CRAIG ST

City of Schenectady Parcels

Map Notes: 1.

- 2. The locations and features depicted on this map are approximate and do not represent a field survey.

FIGURE 4: GROUNDWATER EXCEEDING NYS AWQS PHASE II ESA SAMPLED MONITORING WELLS HAMILTON HILL II - TARGET AREA 1 SITE

City of Schenectady

Schenectady County, New York



C.T. MALE ASSOCIATES ENGINEERING, SURVEYING, ARCHITECTURE & LANDSCAPE ARCHITECTURE, D.AC 50 CENTURY HILL DRIVE, LATHAM, NEW YORK 12110 (518) 786-7400 * FAX (518) 786-7299 * WWW.CTMALE.COM

ATTACHMENT B DIRECT PUSH LOGS

C.T	. M	٩LE	AS	Sociates								
		<u>密</u> ひ][][DIRECT-PUS BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW1 338.62' 6/25/20 1 of 2	DA	ΓUM: ISH D	AMSL ATE: 6/25/20			
PROJI	ECT:		Har	milton Hill II - Target Area 1 BCP Si	te (C447052)	CTM PROJECT NO .:	16.6	6334				
LOCA	TION:		830	& 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	sby			
DEPTH (FT)		NUMBER	SAMPLE CLASSIFICATION						NOTES			
2 4 6 8 10 12 14 16	2 1 4 Fill: Fine brown SAND, little brown silt, trac 2 2 4 gravel 4 3 3 - 6 4 Brown and orange fine to coarse SAND, trace - 10 6 - - - 10 6 - - - 14 7 4 Fine brown SAND and SILT, trace coarse sub - 14 8 Brown SILT and CLAY, trace to coarse sub -					rown silt		t @ ±7	7.5' bgs H bgs			
DRILLIN			OR:	Precision Industrial Services, Inc.			GRC	UNDW/	ATER LEVEL READINGS			
DIRECT-PUSH TYPE: Track-Mounted Geoprobe METHOD OF SAMPLING: 4' Long Macro-Core Sampler with Acetate Liner							DATE	LEVEL	REFERENCE MEASURING POINT			
	J UF 5/	NVIPLIN	I G .	4' Long Macro-Core Sampler with Acetate			6/30/20	13.64	Top of PVC Casing			
IS MAD INFORI AS A S USERS	DE AVA MATIC UBSTI S.	ILABL N AVA TUTE	e to Nlabl For I	RMATION SHOWN HEREON WAS OB AUTHORIZED USERS ONLY THAT TH E TO C.T. MALE. IT IS PRESENTED NVESTIGATIONS, INTERPRETATION	IEY MAY HAVE ACC IN GOOD FAITH, BL OR JUDGMENT OF	ESS TO THE SAME IT IS NOT INTENDED SUCH AUTHORIZED	SAMP	LE CL4	ASSIFICATION BY:			

C.T	. M	4LE	AS	Sociates					
		密 ひ			DIRECT-PUSI BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW1 338.62' 6/25/20 2 of 2	DA	TUM: ISH D	AMSL ATE: 6/25/20
PROJI				milton Hill II - Target Area 1 BCP Sit	, ,	CTM PROJECT NO .:		6334	
LOCA		-) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	isby
DЕРТН (FT)		NUMBER	m RECOVERY (FT)	SAMPLE CLASSIFICATION					NOTES
<u>18</u> <u>20</u> <u>22</u> <u>24</u> <u>26</u> <u>28</u> <u>30</u> <u>32</u>		9	4	Brown CLAY, Some brown Silt, Gray CLAY, little brown silt Boring Termi	inated @ ±20' bg				
DRILLIN	IG CON	TRACT	OR:	Precision Industrial Services, Inc.			GRO		ATER LEVEL READINGS
				Track-Mounted Geoprobe					
METHO	U OF S/	AMPLIN	IG:	4' Long Macro-Core Sampler with Acetate I	Liner		DATE 6/30/20	LEVEL 13.64	REFERENCE MEASURING POINT Top of PVC Casing
IS MAD INFORI AS A S USERS	DE AVA MATIO UBSTI 3.	NILABL N AVA TUTE	e to Nlabi For I	RMATION SHOWN HEREON WAS OB AUTHORIZED USERS ONLY THAT TH LE TO C.T. MALE. IT IS PRESENTED IN NVESTIGATIONS, INTERPRETATION	IEY MAY HAVE ACC IN GOOD FAITH, BU OR JUDGMENT OF	CESS TO THE SAME JT IS NOT INTENDED SUCH AUTHORIZED	SAMP	LE CL/	ASSIFICATION BY:

C.T	. MA	٩LE	AS	Sociates						
		密 び)[][DIRECT-PUSI BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW2 337.73' 6/25/20 1 of 2	DA) TUM: ISH D	AMSL ATE: 6/25/20	
PROJI	ECT:		На	milton Hill II - Target Area 1 B	CP Site (C447052)	CTM PROJECT NO .:	16.6	6334		
LOCA	TION:		830	0 & 834 Albany Street, City/Co	ounty Schenectady	CTM OBSERVER:	Chr	is Orm	ısby	
DЕРТН (FT)	INTERVAL (0	NUMBER	RECOVERY (FT)	SAMPL	E CLASSIFICATIO	DN	NOTES			
_2 _4 _6 _8		1 2 3 4 5	3 4 4	Fill: Fine to coarse brown angular gravel			Moist	t		
<u>10</u> <u>12</u> <u>14</u> 16		6 7 8	3.5	Fine brown SAND and SIL Brown SILT and CLAY, tra			Wet (@ ±1:	3' bgs	
DRILLIN	G CON	RACT	OR:	Precision Industrial Services, Inc.			GRO	איטאוו	ATER LEVEL READINGS	
DIRECT-PUSH TYPE: Track-Mounted Geoprobe								1		
METHO	D OF SA	MPLIN	IG:	4' Long Macro-Core Sampler with A	Acetate Liner		DATE 6/30/20	LEVEL	REFERENCE MEASURING POINT	
IS MAE INFOR AS A S USERS	HE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. I S MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME IFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED S A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED								Top of PVC Casing ASSIFICATION BY: hris Ormsby:	
(Round	#1)\Atta	achme	ent B [Direct-Push Log.xls			Samo		SUXA SAIIDIIDU KESUIS	

C.T	. M <i>i</i>	٩LE	AS	Sociates									
		<u>密</u> ひ][][DIRECT-PUSI BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW2 337.73' 6/25/20 2 of 2	DA	σ ΓUM: ISH D	AMSL ATE: 6/25/20				
PROJE		-		milton Hill II - Target Area 1 BCP Sit		CTM PROJECT NO .:		6334					
LOCA) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	isby				
DЕРТН (FT)	INTERVAL	NUMBER	SAMPLE CLASSIFICATION						NOTES				
		9	2	Brown SILT and CLAY, trace fir	ne brown sand								
18	Gray CLAY, trace brown silt												
	/			Boring Termi	inated @ ±18' bg	IS							
20													
22													
_24													
26													
28													
30													
32													
DRILLIN	G CON	FRACT	OR:	Precision Industrial Services, Inc.			GRO	UNDW	ATER LEVEL READINGS				
				Track-Mounted Geoprobe	Linor		DATE	LEVEL	REFERENCE MEASURING POINT				
METHO	JUF SA	WPLIN	10:	4' Long Macro-Core Sampler with Acetate I	LINEI		6/30/20	12.9	Top of PVC Casing				
IS MAD	E AVA MATIO UBSTI	ILABL N AVA	E TO	RMATION SHOWN HEREON WAS OB AUTHORIZED USERS ONLY THAT TH LE TO C.T. MALE. IT IS PRESENTED I INVESTIGATIONS, INTERPRETATION	IEY MAY HAVE ACC IN GOOD FAITH, BL	CESS TO THE SAME JT IS NOT INTENDED	0.0.45						
		334\E	nv\Ta	raet Area 1 Remedial Investigation 2018	3-2019\Remedial Act	ion\Post-RA Groundwate			ASSIFICATION BY: ^{Thris Ormsby} st RA Sampling Results				

C.T	. M	۹LE	AS	Sociates		h exploration		2				
		<u>密</u> ひ	<u>י</u> ןר כו		BORING NO.: ELEV.: START DATE: SHEET	RAMW3 336.03 6/25/20 1 of 2	DA	TUM: ISH D	AMSL ATE: 6/25/20			
PROJI	ECT:		Ha	milton Hill II - Target Area 1 BCP Sit	te (C447052)	CTM PROJECT NO .:	16.6	6334				
LOCA	_) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	isby			
DЕРТН (FT)	INTERVAL	NUMBER	SAMPLE CLASSIFICATION						NOTES			
2 4 6 8 10 12 14		1 2 3 4 5 6 7 8	2 2 4	Fill: Fine to coarse brown SANE angular gravel Light brown to brown fine to coa Dark brown coarse SAND, trace Brown CLAY, little brown silt, tra sub-angular gravel	arse SAND, little	brown silt		-	1.75' bgs ing (±12.5'-13.5' bgs)			
		TRACT	OR.	Precision Industrial Services Inc								
DRILLING CONTRACTOR: Precision Industrial Services, Inc. DIRECT-PUSH TYPE: Track-Mounted Geoprobe							GRC	UNDWA	ATER LEVEL READINGS			
METHO	D OF S/	MPLIN	IG:		Liner		DATE	LEVEL	REFERENCE MEASURING POINT			
IS MAE INFOR AS A S USERS	IETHOD OF SAMPLING: 4' Long Macro-Core Sampler with Acetate Liner HE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. IT S MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME NFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED S A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED ISERS. \Projects\166334\Env\Target Area 1 Remedial Investigation 2018-2019\Remedial Action\Post-RA Groundwate								Top of PVC Casing			

C.T	. M <i>i</i>	٩LE	AS	Sociates									
		<u>密</u> ひ			DIRECT-PUS BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW3 336.03' 6/25/20 2 of 2	DA	⊐ TUM: ISH D	AMSL ATE: 6/25/20				
PROJE	ECT:		Hai	milton Hill II - Target Area 1 BCP Sit	te (C447052)	CTM PROJECT NO.:	16.6	6334					
LOCA	FION:	•	830) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	nsby				
DЕРТН (FT)	INTERVAL	NUMBER	RECOVERY (FT)	Gray CLAY, little brown silt, trace coarse sub-angular gravel					NOTES				
18		9	2	Gray CLAY, little brown silt, trac	e coarse sub-an	gular gravel							
				Boring Term	inated @ ±18' bo	js							
20													
22													
_24													
26													
_28													
30													
32													
DRILLIN	G CON	RACT	OR:	Precision Industrial Services, Inc.			CPC		ATER LEVEL READINGS				
DIRECT				Track-Mounted Geoprobe									
METHO	D OF SA	MPLIN	IG:	4' Long Macro-Core Sampler with Acetate	Liner		DATE 6/30/20	LEVEL	REFERENCE MEASURING POINT				
IS MAD INFORI AS A S	E AVA MATIO UBSTI	ILABL N AVA	E TO	RMATION SHOWN HEREON WAS OB AUTHORIZED USERS ONLY THAT TH LE TO C.T. MALE. IT IS PRESENTED NVESTIGATIONS, INTERPRETATION	IEY MAY HAVE ACC IN GOOD FAITH, BL	CESS TO THE SAME JT IS NOT INTENDED							
USERS K:\Proie		<u>334\E</u>	<u>nv∖Ta</u>	raet Area 1 Remedial Investigation 2018	3-2019\Remedial Act	ion\Post-RA Groundwate			ASSIFICATION BY: Thris Ormsby St RA Sampling Results				

C.T	. M	ALE	AS	Sociates						
		密 び			DIRECT-PUSI BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW4 334.93' 6/25/20 1 of 1	DA	ish d	AMSL ATE: 6/25/20	
PROJI	ECT:		Ha	milton Hill II - Target Area 1 BCP Si	te (C447052)	CTM PROJECT NO .:	: 16.6334			
LOCA	TION:		830	0 & 834 Albany Street, City/County	Schenectady	CTM OBSERVER:	Chr	is Orm	nsby	
	S	AMPL	E							
DEPTH (FT)	INTERVAL	NUMBER	RECOVERY (FT)	FILL: Fine to coarse brown SAND, little brown silt, trace coarse sub-					NOTES	
		1 2 3	3	FILL: Fine to coarse brown SAI angular gravel	ND, little brown si	lt, trace coarse sub-				
6 8 10		4	3.5	Light brown/orange fine SAND, Dark brown fine to coarse SAN		lt				
<u>12</u> <u>14</u> 16		6 7 8	4	Light brown SILT and CLAY gra brown silt, trace coarse sub-and Boring Term		nd brown CLAY, little	Wet @ ±11' bgs			
		TD 4	05	-		,-				
			UR:	Precision Industrial Services, Inc. Track-Mounted Geoprobe			GRO	UNDW	ATER LEVEL READINGS	
	DIRECT-PUSH TYPE: Track-Mounted Geoprobe METHOD OF SAMPLING: 4' Long Macro-Core Sampler with Acetate Liner							LEVEL	REFERENCE MEASURING POINT	
IS MAD INFOR AS A S	THE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. IT S MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME NFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED AS A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED							10.55	Top of PVC Casing	
USERS									ASSIFICATION BY:	
K:\Proie (Round	cts\16 #1)\At	6334\E tachme	ent B E	raet Area 1 Remedial Investigation 201 Direct-Push Log.xls	8-2019\Remedial Act	ion\Post-RA Groundwate	Samp	lina\Pò	st RA Sampling Results	

C.T	. M	٩LE	AS	Sociates					
		<u>密</u> び			DIRECT-PUS BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW5 335.66' 6/25/20 1 of 2	DA) TUM: ISH D	AMSL ATE: 6/25/20
PROJI	ECT:		Ha	milton Hill II - Target Area 1 B0	CP Site (C447052)	CTM PROJECT NO .:	16.6	6334	
LOCA	TION:		830) & 834 Albany Street, City/Co	unty Schenectady	CTM OBSERVER:	Chr	is Orm	nsby
	S	AMPL	E						
DEPTH (FT)	INTERVAL	NUMBER	RECOVERY (FT)	SAMPLI	E CLASSIFICATIO	N			NOTES
_2		1 2	1.5	FILL: Dark brown fine to co gravel, trace brown silt	oarse SAND, little coa	rse sub-angular			
6		3	2	Brown fine to coarse SANE	D, little brown silt, trac	e fine angular gravel			
<u>10</u> 12		5 6	3	Brown fine SAND, Some b Mixed dark brown/black/ora		e dark brown silt			
<u>14</u> 16		7 8	3.5	Brown SILT and CLAY, tra	ce fine sub-angular g		Wet	@ ±14	4.5' bgs
DRILLIN	G CON	TRACT	OR:	Precision Industrial Services, Inc.					
DIRECT				Track-Mounted Geoprobe			GRC	UNDW	ATER LEVEL READINGS
METHOD OF SAMPLING: 4' Long Macro-Core Sampler with Acetate Liner						DATE	LEVEL	REFERENCE MEASURING POINT	
THE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. IT IS MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME INFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED AS A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED						6/30/20	11.96	Top of PVC Casing	
				raet Area 1 Remedial Investidation Direct-Push Log.xls	n 2018-2019\Remedial Act	ion\Post-RA Groundwate	Samp	lina\Po	Chris Ormsby st RA Sampling Results

C.T	. M	٩LE	AS	Sociates					
		<u>密</u> ひ][][DIRECT-PUSI BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW5 335.66' 6/25/20 2 of 2	DA	ΓUM: ISH D	AMSL ATE: 6/25/20
PROJ	ECT:		Hai	milton Hill II - Target Area 1 BCP Sit	te (C447052)	CTM PROJECT NO.:	16.6	6334	
LOCA	TION:	-	830) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	isby
DЕРТН (FT)	INTERVAL	NUMBER	RECOVERY (FT)	SAMPLE CI	ASSIFICATIO	N	NOTES		
<u>18</u> 20		9 10	4	Brown SILT and CLAY, trace fir Boring Term	inated @ ±20' bç				
<u>22</u> 24									
26	-								
_28	-								
<u>30</u> 32									
DRILLIN	IG CON	TRACT	OR:	Precision Industrial Services, Inc.			GRO		ATER LEVEL READINGS
			10	Track-Mounted Geoprobe					
METHO	ט OF S/	MPLIN	IG:	4' Long Macro-Core Sampler with Acetate	Liner		DATE 6/30/20	LEVEL 11.96	REFERENCE MEASURING POINT
6/30/20 11.96 Top of PVC Casing THE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. IT IS MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME INFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED AS A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED USERS. Chris Ormsby: Chris Ormsby: Chris Ormsby: Chris Ormsby: Chris Ormsby:							ASSIFICATION BY:		

C.T	. M/	٩LE	AS	Sociates					
		<u>密</u> 乙][][DIRECT-PUSI BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW6 336.94' 6/25/20 1 of 2	DA	J TUM: ISH D	AMSL ATE: 6/25/20
PROJI	ECT:		Hai	milton Hill II - Target Area 1 BCP Sit	te (C447052)	CTM PROJECT NO .:	16.0	6334	
LOCA	TION:		830) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	nsby
	S	AMPL	E						
DEPTH (FT)	INTERVAL	NUMBER	RECOVERY (FT)	SAMPLE CL	ASSIFICATIO	N			NOTES
		1	2	FILL: Coarse grey sub-angular (GRAVEL				
2 4		1 2	2	FILL: Dark brown fine to coarse sub-angular gravel	SAND, little brow	wn silt, trace coarse			
<u>6</u> 8		3 4	0.1						
<u>10</u> 12		5	0.5	Mixed dark brown/white/black c	oarse SAND, tra	ce brown silt			
<u>14</u> 16		7 8	4				Wet	@ ±15	5' bgs
DRILLING CONTRACTOR: Precision Industrial Services, Inc.							GRC	DUNDW	ATER LEVEL READINGS
DIRECT-PUSH TYPE: Track-Mounted Geoprobe METHOD OF SAMPLING: 4' Long Macro-Core Sampler with Acetate Liner						DATE	LEVEL	REFERENCE MEASURING POINT	
	0 0 0/						6/30/20	12.83	Top of PVC Casing
6/30/20 12.83 Top of PVC Casing THE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. IT IS MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME INFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED AS A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED USERS. Chris Ormsby, Chris Orms									

C.T	. M	٩LE	AS	Sociates					
K V		<u>密</u> ひ			DIRECT-PUS BORING NO.: ELEV.: START DATE: SHEET	H EXPLORATION RAMW6 336.94 6/25/20 2 of 2	DA	σ ΓUM: ISH DA	AMSL ATE: 6/25/20
PROJECT: Hamilton Hill II - Target Area 1 BCP Site (C447052) CTM PROJECT NO.: 16.6334									
LOCA	TION:	-	830) & 834 Albany Street, City/County S	Schenectady	CTM OBSERVER:	Chr	is Orm	sby
DЕРТН (FT)	INTERVAL 0	NUMBER	RECOVERY (FT)	SAMPLE CI	ASSIFICATIO	NC		I	NOTES
<u>18</u> <u>20</u> <u>22</u> <u>24</u> <u>26</u> <u>28</u> <u>30</u> <u>32</u>		9	3.9	Mixed brown/white/black coarse gravel Brown SILT, Some gray Clay, tr Boring Term		and			
DRILLIN	IG CON	TRACT	OR:	Precision Industrial Services, Inc.			GRO		ATER LEVEL READINGS
DIRECT				Track-Mounted Geoprobe					
METHO	D OF SA	MPLIN	IG:	4' Long Macro-Core Sampler with Acetate	Liner		DATE	LEVEL	REFERENCE MEASURING POINT
IS MAD INFORI AS A S USERS	6/30/20 12.83 Top of PVC Casing THE SUBSURFACE INFORMATION SHOWN HEREON WAS OBTAINED FOR C.T. MALE EVALUATION. IT IS MADE AVAILABLE TO AUTHORIZED USERS ONLY THAT THEY MAY HAVE ACCESS TO THE SAME INFORMATION AVAILABLE TO C.T. MALE. IT IS PRESENTED IN GOOD FAITH, BUT IS NOT INTENDED AS A SUBSTITUTE FOR INVESTIGATIONS, INTERPRETATION OR JUDGMENT OF SUCH AUTHORIZED SAMPLE CLASSIFICATION BY: Chris Ormsby: Chris Ormsby: Chris Ormsby: Chris Ormsby:								ASSIFICATION BY:

ATTACHMENT C

ORGANIC VAPOR HEADSPACE ANALYSIS LOGS



ORGANIC VAPOR HEADSPACE ANALYSIS LOG

PROJECT: Hamilton	n Hill II - Targe	et Area 1 Site	(C447052)	PROJECT #:	16.6334	PAGE 1 OF 3
CLIENT:		DATE				
Location:	y, NY eV	COLLECTED: 6/25/2020				
INSTRUMENT USED:	DATE					
DATE INSTRUMENT			6/25/2020	BY:	Chris Ormsby	ANALYZED: 6/25/2020
TEMPERATURE OF S	SOIL:	am	bient	SAMPLE	BACKGROUND	ANALYST: Chris Ormsby
EXPLORATION	SAMPLE	DEPTH	SAMPLE	READING	READING	
NUMBER	NUMBER	(FT.)***	TYPE	(PPM)**	(PPM)**	REMARKS
RAMW1	1	0-2	Soil	0.2	0.2	No Odors/No Staining
RAMW1	2	2-4	Soil	0.3	0.2	No Odors/No Staining
RAMW1	3	4-6	Soil	0.4	0.2	No Odors/No Staining
RAMW1	4	6-8	Soil	0.4	0.2	No Odors/No Staining
RAMW1	5	8-10	Soil	0.5	0.2	No Odors/No Staining
RAMW1	6	10-12	Soil	0.6	0.2	No Odors/No Staining
RAMW1	7	12-14	Soil	0.5	0.3	No Odors/No Staining
RAMW1	8	14-16	Soil	0.5	0.3	No Odors/No Staining
RAMW1	9	16-18	Soil	0.5	0.3	No Odors/No Staining
RAMW1	10	18-20	Soil	0.5	0.3	No Odors/No Staining
RAMW2	1	0-2	Soil	0.4	0.3	No Odors/No Staining
RAMW2	2	2-4	Soil	0.4	0.3	No Odors/No Staining
RAMW2	3	4-6	Soil	0.4	0.3	No Odors/No Staining
RAMW2	4	6-8	Soil	0.4	0.3	No Odors/No Staining
RAMW2	5	8-10	Soil	0.8	0.3	No Odors/No Staining
RAMW2	6	10-12	Soil	0.8	0.3	No Odors/No Staining
RAMW2	7	12-14	Soil	1.2	0.3	No Odors/No Staining
RAMW2	8	14-16	Soil	1.2	0.3	No Odors/No Staining
RAMW2	9	16-18	Soil	0.5	0.2	No Odors/No Staining

*Instrument was calibrated in accordance with manufacturer's recommended procedure using a calibration gas supplied by the manufacturer. **PPM represents concentration of detectable volatile and gaseous compounds in parts per million of air. *** represents feet below the ground surface



ORGANIC VAPOR HEADSPACE ANALYSIS LOG

PROJECT: Hamilton	n Hill II - Targe	et Area 1 Site	(C447052)	PROJECT #:	16.6334	PAGE 2 OF 3
CLIENT:	Hamilton Hil	l II Limited Pa	rtnership			DATE
Location:	830 & 834 A		City/County of			COLLECTED: 6/25/2020
INSTRUMENT USED:		<i>MiniRae 3000</i>	LAMP	10.6	eV	DATE
DATE INSTRUMENT			6/25/2020	BY:	Chris Ormsby	ANALYZED: 6/25/2020
TEMPERATURE OF S		am	bient	SAMPLE	BACKGROUND	ANALYST: Chris Ormsby
EXPLORATION	SAMPLE	DEPTH	SAMPLE	READING	READING	
NUMBER	NUMBER	(FT.)***	TYPE	(PPM)**	(PPM)**	REMARKS
RAMW3	1	0-2	Soil	0.4	0.3	No Odors/No Staining
RAMW3	2	2-4	Soil	0.4	0.3	No Odors/No Staining
RAMW3	3	4-6	Soil	0.4	0.2	No Odors/No Staining
RAMW3	4	6-8	Soil	0.4	0.2	No Odors/No Staining
RAMW3	5	8-10	Soil	0.5	0.5	No Odors/No Staining
RAMW3	6	10-12	Soil	0.5	0.5	No Odors/No Staining
RAMW3	7	12-14	Soil	71.1	0.4	Black Stain/Petro-Type Odor
RAMW3	8	14-16	Soil	1.1	0.4	No Odors/No Staining
RAMW3	9	16-18	Soil	0.8	0.3	No Odors/No Staining
RAMW4	1	0-2	Soil	0.3	0.2	No Odors/No Staining
RAMW4	2	2-4	Soil	0.3	0.2	No Odors/No Staining
RAMW4	3	4-6	Soil	0.3	0.2	No Odors/No Staining
RAMW4	4	6-8	Soil	0.5	0.2	No Odors/No Staining
RAMW4	5	8-10	Soil	0.5	0.2	No Odors/No Staining
RAMW4	6	10-12	Soil	0.5	0.2	No Odors/No Staining
RAMW4	7	12-14	Soil	0.9	0.2	No Odors/No Staining
RAMW4	8	14-16	Soil	0.9	0.2	No Odors/No Staining

*Instrument was calibrated in accordance with manufacturer's recommended procedure using a calibration gas supplied by the manufacturer. **PPM represents concentration of detectable volatile and gaseous compounds in parts per million of air. *** represents feet below the ground surface



ORGANIC VAPOR HEADSPACE ANALYSIS LOG

PROJECT: Hamiltor				PROJECT #:	16.6334	PAGE 3 OF 3
CLIENT:		I II Limited Pa				DATE
LOCATION:			City/County of			COLLECTED: 6/25/2020
INSTRUMENT USED:		ViniRae 3000	LAMP	10.6	eV	
DATE INSTRUMENT TEMPERATURE OF S			6/25/2020 bient	BA:	Chris Ormsby	ANALYZED: 6/25/2020 ANALYST: Chris Ormsby
ILIVIPERATURE OF 3		am		SAMPLE	BACKGROUND	
EXPLORATION	SAMPLE	DEPTH	SAMPLE	READING	READING	
NUMBER	NUMBER	(FT.)***	TYPE	(PPM)**	(PPM)**	REMARKS
RAMW5	1	0-2	Soil	0.3	0.2	No Odors/No Staining
RAMW5	2	2-4	Soil	0.4	0.2	No Odors/No Staining
RAMW5	3	4-6	Soil	0.4	0.2	No Odors/No Staining
RAMW5	4	6-8	Soil	0.3	0.2	No Odors/No Staining
RAMW5	5	8-10	Soil	0.4	0.1	No Odors/No Staining
RAMW5	6	10-12	Soil	0.8	0.1	No Odors/No Staining
RAMW5	7	12-14	Soil	0.7	0.2	No Odors/No Staining
RAMW5	8	14-16	Soil	0.6	0.2	No Odors/No Staining
RAMW5	9	16-18	Soil	0.7	0.2	No Odors/No Staining
RAMW5	10	18-20	Soil	0.7	0.2	No Odors/No Staining
RAMW6	1	0-2	Soil	0.2	0.2	No Odors/No Staining
RAMW6	2	2-4	Soil	0.2	0.2	No Odors/No Staining
RAMW6	3	4-6	Soil	0.2	0.2	No Odors/No Staining
RAMW6	4	6-8	Soil	0.2	0.2	No Odors/No Staining
RAMW6	5	8-10	Soil	0.2	0.2	No Odors/No Staining
RAMW6	6	10-12	Soil	0.2	0.2	No Odors/No Staining
RAMW6	7	12-14	Soil	0.3	0.2	No Odors/No Staining
RAMW6	8	14-16	Soil	0.3	0.2	No Odors/No Staining
RAMW6	9	16-18	Soil	0.4	0.2	No Odors/No Staining
RAMW6	10	18-20	Soil	0.5	0.2	No Odors/No Staining

*Instrument was calibrated in accordance with manufacturer's recommended procedure using a calibration gas supplied by the manufacturer. **PPM represents concentration of detectable volatile and gaseous compounds in parts per million of air. *** represents feet below the ground surface

ATTACHMENT D

MONITORING WELL CONSTRUCTION LOGS



Protective Enclosure Curb Box Guard Pipe 338.65 ft. elev. 338.44 ft. elev.	Project Name: Hamilton Hill II - Target Area 1 BCP Site 830 & 834 Albany Street Project Number: 16.6334 Well No.: RAMW1
. <u>. 338.62 ft. elev.</u> GROUND SURFACE Concrete Surface Seal ft.* 2- inch diameter drilled hole	Town/City: City of Schenectady County: Schenectady State: NY Installation Date(s): 6/25/2020
Well Casing/Riser 	Drilling Contractor: Precision Environmental Services, Inc. Drilling Method: Direct Push Water Depth From Top of Riser: 13.64 ft 6/30/2020 Date C.T. Mela Observation Observation
	C.T. Male Observer: Chris Ormsby Materials Used: 0.5 Bags of Sand (50lb. bags) Sand Size: 1 Brand: FilPro 0.1 Bags of Bentonite (50lb. bags) Brand: Holeplug 10 ft. of 0.01-Slot, 1-inch diameter 10 ft. of 1-inch diameter 0.5 Bags of Cement/Concrete (80lb. bags) Brand: Quikrete
Gravel Pack Sand Pack Formation Collapse 20.0 ft* 20.0 ft*	Grout Mixture: Bags of Cement [lb. bags] Lbs. of Bentonite Gallons of Water Grout Batches
* Depth below ground surface.	



	Project Name: Hamilton Hill II - Target Area 1 BCP Site
	830 & 834 Albany Street
Protective Enclosure Curb Box Guard Pipe 337.43 ft elev	Project Number: <u>16.6334</u> Well No.: RAMW2 Boring No.: RAMW2
<u> </u>	
ROUND SURFACE	Town/City: City of Schenectady
ft.*	County: <u>Schenectady</u> State: <u>NY</u>
2- inch diameter drilled hole	Installation Date(s): <u>6/25/2020</u>
Well Casing/Riser 1- inch diameter	Drilling Contractor: Precision Environmental Services, Inc.
	Drilling Method: Direct Push
Grout	Water Depth From Top of Riser: <u>12.9</u> ft <u>6/25/2020</u> Date
	C.T. Male Observer: Chris Ormsby
Grade ft* □slurry	<u>Materials Used</u> : 0.5 Bags of Sand (50 lb. bags)
a 3.0 ft* ☐ pellets chips	Sand Size:1Brand:FilPro0.15Bags of Bentonite(50lb. bags)
8.0 ft*	Brand: Holeplug
Well Screen	10 ft. of 0.01-Slot, 1-inch diameter well screen 8 ft. of 1-inch diameter well riser
1 -inch diameter	0.5 Bags of Cement/Concrete(80 lb. bags) Brand: Quikrete
avel Pack	
nd Pack rmation Collapse	Grout Mixture: Bags of Cement (lb. bags)
18.0 ft*	Lbs. of Bentonite Gallons of Water
18.0 ft*	Grout Batches
	Notes:
е.	
	Curb Box Guard Pipe 337.43 ft. elev. ROUND SURFACE a Surface Seal ft.* 2- inch diameter drilled hole Well Casing/Riser 1- inch diameter Backfill Grout Grade ft* Slurry pellets chips 8.0 ft* Well Screen 1 - inch diameter 0.01 slot avel Pack mation Collapse 18.0 ft*



Protective Enclosure Curb Box Guard Pipe 335.97 ft. elev. 336.03 ft. elev.	Project Name: Hamilton Hill II - Target Area 1 BCP Site 830 & 834 Albany Street Project Number: 16.6334 Well No.: RAMW3 Boring No.: RAMW3 Town/City: City of Schenectady County: Schenectady Schenectady State: NY Installation Date(s): 6/25/2020 Drilling Contractor: Precision Environmental Services, Inc. Drilling Method: Direct Push Water Depth From Top of Riser: 11.54 ft 6/25/2020 Date C.T. Male Observer: Chris Ormsby Materials Used: 1 Brand FilPro
Bentonite 2.0 ft* pellets chips 8.0 ft* Well Screen <u>1</u> -inch diameter <u>0.01</u> slot Gravel Pack Sand Pack Formation Collapse <u>18.0</u> ft*	0.5 Bags of Sand (<u>50</u> lb. bags)
* Depth below ground surface.	



Protective Enclosure	Project Name: Hamilton Hill II - Target Area 1 BCP Site 830 & 834 Albany Street Project Number: 16.6334 Well No.: RAMW4 Boring No.: RAMW4 Town/City: City of Schenectady
Concrete Surface Seal <u>ft.*</u> <u>concrete Surface Seal</u> <u>ft.*</u> <u>concrete Surface Seal</u>	County: Schenectady State: NY Installation Date(s): 6/25/2020
Well Casing/Riser 	Drilling Contractor: Precision Environmental Services, Inc. Drilling Method: Direct Push Water Depth From Top of Riser: 10.55 ft 6/25/2020 Date C.T. Male Observer: Chris Ormsby
Grade ft* Bentonite 2.0 ft* Bentonite 2.0 ft* Chips 6.0 ft* Well Screen 1 -inch diameter 0.01 slot Gravel Pack Sand Pack Formation Collapse 16.0 ft*	Materials Used: 0.5 Bags of Sand (50lb. bags) Sand Size: 1 Brand: FilPro 0.1 Bags of Bentonite (50lb. bags) Brand: Holeplug 10 ft. of 0.01-Slot, 1-inch diameter 6 ft. of 1-inch diameter 0.5 Bags of Cement/Concrete (80lb. bags) Brand: Quikrete
* Depth below ground surface.	<u>Notes:</u>



Protective Enclosure	Project Name: Hamilton Hill II - Target Area 1 BCP Site 830 & 834 Albany Street Project Number: 16.6334 Well No.: RAMW5 Boring No.: RAMW5 Town/City: City of Schenectady
Concrete Surface Seal ft.* 2- inch diameter drilled hole Well Casing/Riser	County: Schenectady State: NY Installation Date(s): 6/25/2020 Drilling Contractor: Precision Environmental Services, Inc.
	Drilling Method: Direct Push Water Depth From Top of Riser: <u>11.96 ft</u> <u>6/25/2020</u> Date C.T. Male Observer: <u>Chris Ormsby</u>
Grade ft* Bentonite 2.0 ft* Slurry pellets chips 10.0 ft* Well Screen1 -inch diameter 0.01 slot	Materials Used: 0.5 Bags of Sand (50 lb. bags) Sand Size: 1 Brand: FilPro 0.1 Bags of Bentonite (50 lb. bags) Brand: Holeplug 10 ft. of 0.01-Slot, 1-inch diameter well screen 10 ft. of 1-inch diameter well riser 0.5 Bags of Cement/Concrete (80 lb. bags) Brand: Quikrete
Gravel Pack Sand Pack Formation Collapse 20.0 ft* 20.0 ft*	Grout Mixture: Bags of Cement [] Lbs. of Bentonite Gallons of Water Grout Batches
* Depth below ground surface.	<u>Notes:</u>



	Project Name: Hamilton Hill II - Target Area 1 BCP Site
	830 & 834 Albany Street
Protective Enclosure	
Curb Box	Project Number: 16.6334
$-\underline{337.02} \text{ ft. elev.} \qquad \qquad \square \text{ Guard Pipe}$	
<u>336.82</u> ft. ele	Well No.: RAMW6 Boring No.: RAMW6
GROUND SURFACE	Town/City: City of Schenectady
<u>336.94 ft. elev. </u> GROUND SURFACE	Town/City. City of Scheneclady
ft*	County: Schenectady State: NY
2- inch diameter	
drilled hole	Installation Date(s): 6/25/2020
Well Casing/Riser	Drilling Contractor: <u>Precision Environmental Services, Inc.</u>
<u> </u>	Drilling Method: Direct Push
Backfill	
Grout	Water Depth From Top of Riser: 12.83 ft 6/25/2020
	Date
	C.T. Male Observer: Chris Ormsby
Grade ft*	Materials Used:
slurry	0.5 Bags of Sand (50 lb. bags)
Bentonite 2.0 ft* pellets chips	Sand Size: 1 Brand FilPro
	0.1 Bags of Bentonite (50 lb. bags)
8.0 ft*	Brand: Holeplug
	10 ft. of 0.01-Slot, 1-inch diameter well screen
Well Screen	8 ft. of <u>1-inch diameter</u> well riser 0.5 Bags of Cement/Concrete (80 lb. bags)
	Brand: Quikrete
	Grout Mixture:
Formation Collapse	Bags of Cement (lb. bags)
	Lbs. of Bentonite
18.0 ft*	Gallons of Water
20.0.6*	Grout Batches
<u>20.0</u> ft*	
	Notes:
* Depth below ground surface.	Notes:

C.T. MALE ASSOCIATES

ATTACHMENT E

ANALYTICAL RESULTS SUMMARY TABLE

POST REMEDIAL ACTION GROUNDWATER SAMPLING ANALYTICAL RESULTS SUMMARY HAMILTON HILL II - TARGET AREA 1 BCP SITE (C447052) 830 AND 834 ALBANY STREET CITY AND COUNTY OF SCHENECTADY

		SAMPLE ID:		RAM	WW1-2006	30		RA	MW2-2006	30		RA	MW3-2006	30		RA	MW4-2006	30		RA	MW5-20063	30
		LAB ID:		L2	2027725-04	1		L	2027725-0	5		L	2027725-06	i		L	.2027725-03	3		L	2027725-01	1
		COLLECTION DATE:		6	6/30/2020				6/30/2020				6/30/2020				6/30/2020				6/30/2020	
		SAMPLE MATRIX:			WATER				WATER				WATER				WATER				WATER	
		NY-AWQS ⁽¹⁾																				
ANALYTE	CAS	(ug/l)	Conc	Q	RL	MDL	Conc	Q	RL	MDL	Conc	Q	RL	MDL	Conc	Q	RL	MDL	Conc	Q	RL	MDL
VOLATILE ORGANICS BY O	GC/MS	(3)																				
Acetone	67-64-1	50 (GV)	ND		5	1.5	ND		5	1.5	ND		5	1.5	ND		5	1.5	ND		5	1.5
Bromodichloromethane	75-27-4	50 (GV)	ND		0.5	0.19	0.46	J	0.5	0.19	0.69		0.5	0.19	ND		0.5	0.19	ND		0.5	0.19
Chloroform	67-66-3	7	ND		2.5	0.7	2.8	-	2.5	0.7	1.7	J	2.5	0.7	ND		2.5	0.7	ND		2.5	0.7
Dibromochloromethan	124-48-1	50 (GV)	ND		0.5	0.15	0.26	J	0.5	0.15	ND		0.5	0.15	ND		0.5	0.15	ND		0.5	0.15
Tetrachloroethene	127-18-4	5	7.3		0.5	0.18	4.4		0.5	0.18	6.8		0.5	0.18	8.7		0.5	0.18	1.4		0.5	0.18
Trichloroethene	79-01-6	5	3.2		0.5	0.18	ND		0.5	0.18	ND		0.5	0.18	0.59		0.5	0.18	ND		0.5	0.18
Total VOCs		NS	10.5	-	-	-	7.92	-	-	-	9.19	-	-	-	9.29	-	-	-	1.4	-	-	-
VOLATILE ORGANICS BY C	GC/MS-TIC																					
SEMIVOLATILE ORGANICS	S BY GC/MS																					-
SEMIVOLATILE ORGANICS	BY GC/MS-SIM																					
Anthracene	120-12-7	50 (GV)	ND		0.1	0.01	ND		0.1	0.01	ND		0.1	0.01	ND		0.1	0.01	ND		0.1	0.01
Fluoranthene	206-44-0	50 (GV)	ND		0.1	0.02	ND		0.1	0.02	ND		0.1	0.02	ND		0.1	0.02	ND		0.1	0.02
Fluorene	86-73-7	50 (GV)	ND		0.1	0.01	ND		0.1	0.01	ND		0.1	0.01	ND		0.1	0.01	ND		0.1	0.01
Hexachlorobenzen	118-74-1	0.04	ND		0.8	0.01	ND		0.8	0.01	ND		0.8	0.01	ND		0.8	0.01	ND		0.8	0.01
Phenanthrene	85-01-8	50 (GV)	ND		0.1	0.02	ND		0.1	0.02	ND		0.1	0.02	ND		0.1	0.02	ND		0.1	0.02
Total SVOCs		NS	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
SEMIVOLATILE ORGANICS	S BY GC/MS-TIC																					
Unknown		NS	-		-	-	-		-	-	-		-	-	2.65	J	0	0	2.44	J	0	0
Unknown		NS	2.44	J	0	0	-		-	-	2.4	J	0	0	-	-	-	-	-	-	-	-
Unknown		NS	2.4	J	0	0	-		-	-	1.45	J	0	0	-	-	-	-	-	-	-	-
Unknown Organic Acic		NS	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Total TIC Compounds		NS	4.84	J	0	0	-		-	-	3.85	J	0	0	2.65	J	0	0	2.44	J	0	0

(1) New York Division of Water Technical and Operational Guidance Series (TOGS) 1.1.1 Ambient Water Quality Standards and Guidance Values. June 1998 and Addendu (GV) denotes Guidance Value

NS denotes No Standarc

ND denotes Non-Detec

RL denotes the laboratory's Reporting Lim MDL denotes the laboratory's Method Detection Lirr

Qualifiers in parantheses reflect qualifications made by the data validat

POST REMEDIAL ACTION GROUNDWATER SAMPLING ANALYTICAL RESULTS SUMMARY HAMILTON HILL II - TARGET AREA 1 BCP SITE (C447052) 830 AND 834 ALBANY STREET CITY AND COUNTY OF SCHENECTADY

		SAMPLE ID:		RA	AMW6-20063	30	FD	01-200	630 (RAMW	6-200630)		E	B01-20063	0		TRIP	BLANK-20	0630
		LAB ID:		L	_2027725-02	2		L	_2027725-07	7		L	.2027725-0	8		L	2027725-0	9
		COLLECTION DATE:			6/30/2020				6/30/2020				6/30/2020				6/30/2020	
		SAMPLE MATRIX:			WATER				WATER				WATER				WATER	
		NY-AWQS ⁽¹⁾															MATER	
ANALYTE	CAS	(ug/l)	Conc	Q	RL	MDL	Conc	Q	RL	MDL	Conc	Q	RL	MDL	Conc	Q	RL	MDL
VOLATILE ORGANICS BY GO	C/MS	(3)																
Acetone	67-64-1	50 (GV)	ND		5	1.5	ND		5	1.5	3.8	J	5	1.5	ND		5	1.5
Bromodichloromethane	75-27-4	50 (GV)	ND		0.5	0.19	ND		0.5	0.19	ND		0.5	0.19	ND		0.5	0.19
Chloroform	67-66-3	7	ND		2.5	0.7	ND		2.5	0.7	ND		2.5	0.7	ND		2.5	0.7
Dibromochloromethan	124-48-1	50 (GV)	ND		0.5	0.15	ND		0.5	0.15	ND		0.5	0.15	ND		0.5	0.15
Tetrachloroethene	127-18-4	5	ND		0.5	0.18	ND		0.5	0.18	ND		0.5	0.18	ND		0.5	0.18
Trichloroethene	79-01-6	5	ND		0.5	0.18	ND		0.5	0.18	ND		0.5	0.18	ND		0.5	0.18
Total VOCs		NS	-	-	-	-	-	-	-	-	3.8	-	-	-	-	-	-	-
VOLATILE ORGANICS BY GO	C/MS-TIC																	
SEMIVOLATILE ORGANICS E	BY GC/MS																	
SEMIVOLATILE ORGANICS E	BY GC/MS-SIM																	
Anthracene	120-12-7	50 (GV)	ND		0.1	0.01	ND		0.1	0.01	0.02	J	0.1	0.01	-	-	-	-
Fluoranthene	206-44-0	50 (GV)	ND		0.1	0.02	ND		0.1	0.02	0.01	(U)	0.1	0.02	-	-	-	-
Fluorene	86-73-7	50 (GV)	ND		0.1	0.01	ND		0.1	0.01	0.02	Ĵ	0.1	0.01	-	-	-	-
Hexachlorobenzen	118-74-1	0.04	ND		0.8	0.01	ND		0.8	0.01	0.03	J	0.8	0.01	-	-	-	-
Phenanthrene	85-01-8	50 (GV)	ND		0.1	0.02	ND		0.1	0.02	0.01	(U)	0.1	0.02	-	-	-	-
Total SVOCs		NS	-	-	-	-	-	-	-	-	0.12	-	-	-	-	-	-	
SEMIVOLATILE ORGANICS E	BY GC/MS-TIC																	
Unknown		NS	1.78	J	0	0	1.85	J	0	0	-		-	-	-	-	-	-
Unknown		NS	-	-	-	-	-		-	-	-		-	-	-	-	-	-
Unknown		NS	-	-	-	-	-		-	-	-		-	-	-	-	-	-
Unknown Organic Acid		NS	-	-	-	-	-	-	-	-	1.71	J	0	0	-	-	-	-
Total TIC Compounds		NS	1.78	J	0	0	1.85	J	0	0	1.71	J	0	0	-	-	-	-

(1) New York Division of Water Technical and Operational Guidance Series (TOGS) 1.1.1 Ambient Water Quality Standards and Guidance Values. June 1998 and Addendu (GV) denotes Guidance Value

NS denotes No Standarc

ND denotes Non-Detec

RL denotes the laboratory's Reporting Lim MDL denotes the laboratory's Method Detection Lirr

ATTACHMENT F ANALYTICAL REPORT



ANALYTICAL REPORT

Lab Number:	L2027725
Client:	C.T. Male Associates 50 Century Hill Drive Latham, NY 12110
ATTN: Phone:	Aimee Smith (518) 786-7400
Project Name:	HHII TA1
Project Number:	16.6334
Report Date:	07/08/20

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

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

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



Serial_No:07082013:26

Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2027725-01	RAMW5-200630	WATER	SCHENECTADY NY	06/30/20 08:00	06/30/20
L2027725-02	RAMW6-200630	WATER	SCHENECTADY NY	06/30/20 08:35	06/30/20
L2027725-03	RAMW4-200630	WATER	SCHENECTADY NY	06/30/20 11:35	06/30/20
L2027725-04	RAMW1-200630	WATER	SCHENECTADY NY	06/30/20 11:50	06/30/20
L2027725-05	RAMW2-200630	WATER	SCHENECTADY NY	06/30/20 12:05	06/30/20
L2027725-06	RAMW3-200630	WATER	SCHENECTADY NY	06/30/20 12:55	06/30/20
L2027725-07	FD01-200630	WATER	SCHENECTADY NY	06/30/20 00:00	06/30/20
L2027725-08	EB01-200630	WATER	SCHENECTADY NY	06/30/20 13:05	06/30/20
L2027725-09	TRIP BLANK-200630	WATER	SCHENECTADY NY	06/30/20 00:00	06/30/20



Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Case Narrative

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

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.



Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

Sample Receipt

L2027725-02: The collection date and time on the chain of custody was 30-JUN-20 08:35; however, the collection date and time on the container label was 30-JUN-20 09:20. At the client's request, the collection date and time is reported as 30-JUN-20 08:35.

Semivolatile Organics by SIM

The WG1388230-1 Method Blank, associated with L2027725-01 through -08, has a concentration above the reporting limit for Naphthalene. Since the samples were non-detect to the RL for this target analyte, no further actions were taken. The results of the original analysis are reported.

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

Authorized Signature:

Jufani Morrissey - Tiffani Morrissey

Title: Technical Director/Representative

Date: 07/08/20



ORGANICS



VOLATILES



			Serial_No	0:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-01		Date Collected:	06/30/20 08:00
Client ID:	RAMW5-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water			
Analytical Method:	1,8260C			
Analytical Date:	07/05/20 16:14			
Analyst:	NLK			

Parameter	Result	Qualifier U	nits	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - We	stborough Lab					
Methylene chloride	ND	U	ıg/l	2.5	0.70	1
1,1-Dichloroethane	ND		ıg/l	2.5	0.70	1
Chloroform	ND	L	ıg/l	2.5	0.70	1
Carbon tetrachloride	ND		ıg/l	0.50	0.13	1
1,2-Dichloropropane	ND	U	ıg/l	1.0	0.14	1
Dibromochloromethane	ND	U	ıg/l	0.50	0.15	1
1,1,2-Trichloroethane	ND	U	ıg/l	1.5	0.50	1
Tetrachloroethene	1.4	U	ıg/l	0.50	0.18	1
Chlorobenzene	ND	U	ıg/l	2.5	0.70	1
Trichlorofluoromethane	ND	U	ıg/l	2.5	0.70	1
1,2-Dichloroethane	ND	U	ıg/l	0.50	0.13	1
1,1,1-Trichloroethane	ND	U	ıg/l	2.5	0.70	1
Bromodichloromethane	ND	U	ıg/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND	U	ıg/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND	L	ıg/l	0.50	0.14	1
Bromoform	ND	L	ıg/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND	L	ıg/l	0.50	0.17	1
Benzene	ND	U	ıg/l	0.50	0.16	1
Toluene	ND	L	ıg/l	2.5	0.70	1
Ethylbenzene	ND	u	ıg/l	2.5	0.70	1
Chloromethane	ND	u	ıg/l	2.5	0.70	1
Bromomethane	ND	u	ıg/l	2.5	0.70	1
Vinyl chloride	ND	L	ıg/l	1.0	0.07	1
Chloroethane	ND	L	ıg/l	2.5	0.70	1
1,1-Dichloroethene	ND	L	ıg/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND	L	ıg/l	2.5	0.70	1
Trichloroethene	ND	L	ıg/l	0.50	0.18	1
1,2-Dichlorobenzene	ND	U	ıg/l	2.5	0.70	1



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-01 Date Collected: 06/30/20 08:00 Client ID: RAMW5-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1.0 1 ug/l 5.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 1,2-Dibromoethane ND 0.65 1 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 0.70 1 Freon-113 ug/l ND Methyl cyclohexane ug/l 10 0.40 1

Tentatively Identified Compounds

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		106		70-130	
Toluene-d8		105		70-130	
4-Bromofluorobenzene		111		70-130	
Dibromofluoromethane		101		70-130	



			Serial_No	0:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-02		Date Collected:	06/30/20 08:35
Client ID:	RAMW6-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water			
Analytical Method:	1,8260C			
Analytical Date:	07/05/20 16:39			
Analyst:	NLK			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - We	stborough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-02 Date Collected: 06/30/20 08:35 Client ID: RAMW6-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1 ug/l 5.0 1.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 ND 0.65 1 1,2-Dibromoethane 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1 Freon-113 0.70 ug/l ND Methyl cyclohexane ug/l 10 0.40 1

Tentatively Identified Compounds

Tentatively Identified Compounds	ND	ug/l			1	
Surrogate		% Recovery	Qualifier	Acceptance Criteria		
1,2-Dichloroethane-d4		103		70-130		
Toluene-d8		104		70-130		
4-Bromofluorobenzene		109		70-130		
Dibromofluoromethane		101		70-130		



			Serial_No	p:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID: Client ID:	L2027725-03 RAMW4-200630		Date Collected: Date Received:	06/30/20 11:35 06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix: Analytical Method: Analytical Date: Analyst:	Water 1,8260C 07/05/20 17:05 NLK			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Volatile Organics by GC/MS - Westborough Lab										
Methylene chloride	ND		ug/l	2.5	0.70	1				
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1				
Chloroform	ND		ug/l	2.5	0.70	1				
Carbon tetrachloride	ND		ug/l	0.50	0.13	1				
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1				
Dibromochloromethane	ND		ug/l	0.50	0.15	1				
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1				
Tetrachloroethene	8.7		ug/l	0.50	0.18	1				
Chlorobenzene	ND		ug/l	2.5	0.70	1				
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1				
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1				
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1				
Bromodichloromethane	ND		ug/l	0.50	0.19	1				
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1				
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1				
Bromoform	ND		ug/l	2.0	0.65	1				
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1				
Benzene	ND		ug/l	0.50	0.16	1				
Toluene	ND		ug/l	2.5	0.70	1				
Ethylbenzene	ND		ug/l	2.5	0.70	1				
Chloromethane	ND		ug/l	2.5	0.70	1				
Bromomethane	ND		ug/l	2.5	0.70	1				
Vinyl chloride	ND		ug/l	1.0	0.07	1				
Chloroethane	ND		ug/l	2.5	0.70	1				
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1				
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1				
Trichloroethene	0.59		ug/l	0.50	0.18	1				
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1				



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-03 Date Collected: 06/30/20 11:35 Client ID: RAMW4-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1 ug/l 5.0 1.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 ND 0.65 1 1,2-Dibromoethane 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1

Tentatively Identified Compounds

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		105		70-130	
Toluene-d8		105		70-130	
4-Bromofluorobenzene		109		70-130	
Dibromofluoromethane		101		70-130	

ND

ug/l

ug/l



1

0.70

0.40

10

Freon-113

Methyl cyclohexane

			Serial_No	p:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID: Client ID: Sample Location: Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	L2027725-04 RAMW1-200630 SCHENECTADY NY Water 1,8260C 07/05/20 17:30 NLK		Date Collected: Date Received: Field Prep:	06/30/20 11:50 06/30/20 Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
/olatile Organics by GC/MS - Westborough Lab								
Methylene chloride	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1		
Chloroform	ND		ug/l	2.5	0.70	1		
Carbon tetrachloride	ND		ug/l	0.50	0.13	1		
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1		
Dibromochloromethane	ND		ug/l	0.50	0.15	1		
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1		
Tetrachloroethene	7.3		ug/l	0.50	0.18	1		
Chlorobenzene	ND		ug/l	2.5	0.70	1		
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1		
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1		
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1		
Bromodichloromethane	ND		ug/l	0.50	0.19	1		
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1		
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1		
Bromoform	ND		ug/l	2.0	0.65	1		
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1		
Benzene	ND		ug/l	0.50	0.16	1		
Toluene	ND		ug/l	2.5	0.70	1		
Ethylbenzene	ND		ug/l	2.5	0.70	1		
Chloromethane	ND		ug/l	2.5	0.70	1		
Bromomethane	ND		ug/l	2.5	0.70	1		
Vinyl chloride	ND		ug/l	1.0	0.07	1		
Chloroethane	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1		
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1		
Trichloroethene	3.2		ug/l	0.50	0.18	1		
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1		



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-04 Date Collected: 06/30/20 11:50 Client ID: RAMW1-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1.0 1 ug/l 5.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 1,2-Dibromoethane ND 0.65 1 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1 Freon-113 0.70 ug/l

Tentatively Identified Compounds

Methyl cyclohexane

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		106		70-130	
Toluene-d8		104		70-130	
4-Bromofluorobenzene		109		70-130	
Dibromofluoromethane		102		70-130	

ug/l

10

0.40

ND



			Serial_No	0:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID: Client ID: Sample Location: Sample Depth: Matrix: Analytical Method:	L2027725-05 RAMW2-200630 SCHENECTADY NY Water 1,8260C		Date Collected: Date Received: Field Prep:	06/30/20 12:05 06/30/20 Not Specified
Analytical Date: Analyst:	07/05/20 17:56 NLK			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Volatile Organics by GC/MS - Westl	/olatile Organics by GC/MS - Westborough Lab								
Methylene chloride	ND		ug/l	2.5	0.70	1			
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1			
Chloroform	2.8		ug/l	2.5	0.70	1			
Carbon tetrachloride	ND		ug/l	0.50	0.13	1			
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1			
Dibromochloromethane	0.26	J	ug/l	0.50	0.15	1			
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1			
Tetrachloroethene	4.4		ug/l	0.50	0.18	1			
Chlorobenzene	ND		ug/l	2.5	0.70	1			
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1			
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1			
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1			
Bromodichloromethane	0.46	J	ug/l	0.50	0.19	1			
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1			
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1			
Bromoform	ND		ug/l	2.0	0.65	1			
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1			
Benzene	ND		ug/l	0.50	0.16	1			
Toluene	ND		ug/l	2.5	0.70	1			
Ethylbenzene	ND		ug/l	2.5	0.70	1			
Chloromethane	ND		ug/l	2.5	0.70	1			
Bromomethane	ND		ug/l	2.5	0.70	1			
Vinyl chloride	ND		ug/l	1.0	0.07	1			
Chloroethane	ND		ug/l	2.5	0.70	1			
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1			
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1			
Trichloroethene	ND		ug/l	0.50	0.18	1			
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1			



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-05 Date Collected: 06/30/20 12:05 Client ID: RAMW2-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1 ug/l 5.0 1.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 ND 0.65 1,2-Dibromoethane 2.0 1 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1 Freon-113 0.70 ug/l ND

Tentatively Identified Compounds

Methyl cyclohexane

Tentatively Identified Compounds	ND	ug/l		1	
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		107		70-130	
Toluene-d8		105		70-130	
4-Bromofluorobenzene		107		70-130	
Dibromofluoromethane		103		70-130	

ug/l

10

0.40



			Serial_No	p:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID: Client ID: Sample Location:	L2027725-06 RAMW3-200630 SCHENECTADY NY		Date Collected: Date Received: Field Prep:	06/30/20 12:55 06/30/20 Not Specified
Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	Water 1,8260C 07/05/20 18:21 NLK			·

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
/olatile Organics by GC/MS - Westborough Lab								
Methylene chloride	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1		
Chloroform	1.7	J	ug/l	2.5	0.70	1		
Carbon tetrachloride	ND		ug/l	0.50	0.13	1		
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1		
Dibromochloromethane	ND		ug/l	0.50	0.15	1		
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1		
Tetrachloroethene	6.8		ug/l	0.50	0.18	1		
Chlorobenzene	ND		ug/l	2.5	0.70	1		
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1		
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1		
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1		
Bromodichloromethane	0.69		ug/l	0.50	0.19	1		
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1		
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1		
Bromoform	ND		ug/l	2.0	0.65	1		
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1		
Benzene	ND		ug/l	0.50	0.16	1		
Toluene	ND		ug/l	2.5	0.70	1		
Ethylbenzene	ND		ug/l	2.5	0.70	1		
Chloromethane	ND		ug/l	2.5	0.70	1		
Bromomethane	ND		ug/l	2.5	0.70	1		
Vinyl chloride	ND		ug/l	1.0	0.07	1		
Chloroethane	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1		
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1		
Trichloroethene	ND		ug/l	0.50	0.18	1		
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1		



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-06 Date Collected: 06/30/20 12:55 Client ID: RAMW3-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1 ug/l 5.0 1.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 ND 0.65 1 1,2-Dibromoethane 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1 Freon-113 0.70 ug/l

Tentatively Identified Compounds

Methyl cyclohexane

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		103		70-130	
Toluene-d8		104		70-130	
4-Bromofluorobenzene		111		70-130	
Dibromofluoromethane		99		70-130	

ug/l

10

0.40

ND



			Serial_No	p:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-07		Date Collected:	06/30/20 00:00
Client ID:	FD01-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water			
Analytical Method:	1,8260C			
Analytical Date:	07/05/20 18:46			
Analyst:	NLK			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Volatile Organics by GC/MS - Westborough Lab								
Methylene chloride	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1		
Chloroform	ND		ug/l	2.5	0.70	1		
Carbon tetrachloride	ND		ug/l	0.50	0.13	1		
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1		
Dibromochloromethane	ND		ug/l	0.50	0.15	1		
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1		
Tetrachloroethene	ND		ug/l	0.50	0.18	1		
Chlorobenzene	ND		ug/l	2.5	0.70	1		
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1		
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1		
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1		
Bromodichloromethane	ND		ug/l	0.50	0.19	1		
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1		
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1		
Bromoform	ND		ug/l	2.0	0.65	1		
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1		
Benzene	ND		ug/l	0.50	0.16	1		
Toluene	ND		ug/l	2.5	0.70	1		
Ethylbenzene	ND		ug/l	2.5	0.70	1		
Chloromethane	ND		ug/l	2.5	0.70	1		
Bromomethane	ND		ug/l	2.5	0.70	1		
Vinyl chloride	ND		ug/l	1.0	0.07	1		
Chloroethane	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1		
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1		
Trichloroethene	ND		ug/l	0.50	0.18	1		
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1		



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-07 Date Collected: 06/30/20 00:00 Client ID: FD01-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1 ug/l 5.0 1.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 ND 0.65 1 1,2-Dibromoethane 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l

Tentatively Identified Compounds

o Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		105		70-130	
Toluene-d8		106		70-130	
4-Bromofluorobenzene		110		70-130	
Dibromofluoromethane		101		70-130	

2.5

10

ug/l

ug/l

0.70

0.40

ND

ND



1

1

Freon-113

Methyl cyclohexane

			Serial_No	0:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID: Client ID:	L2027725-08 EB01-200630		Date Collected: Date Received:	06/30/20 13:05 06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix: Analytical Method: Analytical Date: Analyst:	Water 1,8260C 07/05/20 15:49 NLK			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Wes	stborough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-08 Date Collected: 06/30/20 13:05 Client ID: EB01-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 J Acetone 3.8 ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1.0 1 ug/l 5.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 1,2-Dibromoethane ND 0.65 1 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1 Freon-113 0.70 ug/l

Tentatively Identified Compounds

Methyl cyclohexane

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		105		70-130	
Toluene-d8		104		70-130	
4-Bromofluorobenzene		110		70-130	
Dibromofluoromethane		102		70-130	

ug/l

10

0.40

ND



			Serial_N	p:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID: Client ID: Sample Location: Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	L2027725-09 TRIP BLANK-200630 SCHENECTADY NY Water 1,8260C 07/05/20 15:23 NLK		Date Collected: Date Received: Field Prep:	06/30/20 00:00 06/30/20 Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - We	stborough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Serial_No:07082013:26 **Project Name:** HHII TA1 Lab Number: L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-09 Date Collected: 06/30/20 00:00 Client ID: TRIP BLANK-200630 Date Received: 06/30/20 Sample Location: SCHENECTADY NY Field Prep: Not Specified Sample Depth: Qualifier MDL Parameter Result Units RL **Dilution Factor** Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND 2.5 0.70 1 ug/l 1,4-Dichlorobenzene ND ug/l 2.5 0.70 1 Methyl tert butyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND 2.5 0.70 1 ug/l o-Xylene ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 ND Acetone ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 ND 1 2-Butanone ug/l 5.0 1.9 4-Methyl-2-pentanone ND 1 ug/l 5.0 1.0 ND 2-Hexanone 5.0 1.0 1 ug/l Bromochloromethane ND 0.70 1 ug/l 2.5 ND 0.65 1 1,2-Dibromoethane 2.0 ug/l 1,2-Dibromo-3-chloropropane ND 2.5 0.70 1 ug/l Isopropylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND 2.5 0.70 1 ug/l 1,2,4-Trichlorobenzene ND 2.5 0.70 1 ug/l Methyl Acetate ND 2.0 0.23 1 ug/l Cyclohexane ND 10 0.27 1 ug/l 1 1,4-Dioxane ND 250 61. ug/l ND 2.5 1 Freon-113 0.70 ug/l

Tentatively Identified Compounds

Methyl cyclohexane

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4		106		70-130	
Toluene-d8		105		70-130	
4-Bromofluorobenzene		111		70-130	
Dibromofluoromethane		102		70-130	

ug/l

10

0.40

ND



HHII TA1 16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Method Blank Analysis Batch Quality Control

Analytical Method:1,8260CAnalytical Date:07/05/20 14:58Analyst:AJK

Project Name:

Project Number:

arameter	Result	Qualifier Units	RL	MDL
olatile Organics by GC/MS - V	Vestborough Lab	for sample(s): (01-09 Batch:	WG1389227-5
Methylene chloride	ND	ug/l	2.5	0.70
1,1-Dichloroethane	ND	ug/l	2.5	0.70
Chloroform	ND	ug/l	2.5	0.70
Carbon tetrachloride	ND	ug/l	0.50	0.13
1,2-Dichloropropane	ND	ug/l	1.0	0.14
Dibromochloromethane	ND	ug/l	0.50	0.15
1,1,2-Trichloroethane	ND	ug/l	1.5	0.50
Tetrachloroethene	ND	ug/l	0.50	0.18
Chlorobenzene	ND	ug/l	2.5	0.70
Trichlorofluoromethane	ND	ug/l	2.5	0.70
1,2-Dichloroethane	ND	ug/l	0.50	0.13
1,1,1-Trichloroethane	ND	ug/l	2.5	0.70
Bromodichloromethane	ND	ug/l	0.50	0.19
trans-1,3-Dichloropropene	ND	ug/l	0.50	0.16
cis-1,3-Dichloropropene	ND	ug/l	0.50	0.14
Bromoform	ND	ug/l	2.0	0.65
1,1,2,2-Tetrachloroethane	ND	ug/l	0.50	0.17
Benzene	ND	ug/l	0.50	0.16
Toluene	ND	ug/l	2.5	0.70
Ethylbenzene	ND	ug/l	2.5	0.70
Chloromethane	ND	ug/l	2.5	0.70
Bromomethane	ND	ug/l	2.5	0.70
Vinyl chloride	ND	ug/l	1.0	0.07
Chloroethane	ND	ug/l	2.5	0.70
1,1-Dichloroethene	ND	ug/l	0.50	0.17
trans-1,2-Dichloroethene	ND	ug/l	2.5	0.70
Trichloroethene	ND	ug/l	0.50	0.18
1,2-Dichlorobenzene	ND	ug/l	2.5	0.70
1,3-Dichlorobenzene	ND	ug/l	2.5	0.70



Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Method Blank Analysis Batch Quality Control

Analytical Method:1,8260CAnalytical Date:07/05/20 14:58Analyst:AJK

arameter	Result 0	Qualifier Units	RL	MDL
olatile Organics by GC/MS - V	/estborough Lab fe	or sample(s): 01-09	Batch:	WG1389227-5
1,4-Dichlorobenzene	ND	ug/l	2.5	0.70
Methyl tert butyl ether	ND	ug/l	2.5	0.70
p/m-Xylene	ND	ug/l	2.5	0.70
o-Xylene	ND	ug/l	2.5	0.70
cis-1,2-Dichloroethene	ND	ug/l	2.5	0.70
Styrene	ND	ug/l	2.5	0.70
Dichlorodifluoromethane	ND	ug/l	5.0	1.0
Acetone	ND	ug/l	5.0	1.5
Carbon disulfide	ND	ug/l	5.0	1.0
2-Butanone	ND	ug/l	5.0	1.9
4-Methyl-2-pentanone	ND	ug/l	5.0	1.0
2-Hexanone	ND	ug/l	5.0	1.0
Bromochloromethane	ND	ug/l	2.5	0.70
1,2-Dibromoethane	ND	ug/l	2.0	0.65
1,2-Dibromo-3-chloropropane	ND	ug/l	2.5	0.70
Isopropylbenzene	ND	ug/l	2.5	0.70
1,2,3-Trichlorobenzene	ND	ug/l	2.5	0.70
1,2,4-Trichlorobenzene	ND	ug/l	2.5	0.70
Methyl Acetate	ND	ug/l	2.0	0.23
Cyclohexane	ND	ug/l	10	0.27
1,4-Dioxane	ND	ug/l	250	61.
Freon-113	ND	ug/l	2.5	0.70
Methyl cyclohexane	ND	ug/l	10	0.40

Tentatively Identified Compounds

No Tentatively Identified Compounds

ug/l



 Lab Number:
 L2027725

 Report Date:
 07/08/20

Project Name:HHII TA1Project Number:16.6334

Method Blank Analysis Batch Quality Control

Analytical Method:1,8260CAnalytical Date:07/05/20 14:58Analyst:AJK

Parameter	Result	Qualifier	Units	RL	MDL	
Volatile Organics by GC/MS	- Westborough La	ab for sample	e(s): 01-09	Batch:	WG1389227-5	

		Acceptance			
Surrogate	%Recovery Qເ	ualifier Criteria			
1.2-Dichloroethane-d4	107	70-130			
Toluene-d8	104	70-130			
4-Bromofluorobenzene	112	70-130			
Dibromofluoromethane	99	70-130			



Lab Control Sample Analysis Batch Quality Control

Project Name: HHII TA1 **Project Number:** 16.6334

Lab Number: L2027725

Report Date: 07/08/20

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Volatile Organics by GC/MS - Westborough	Lab Associated	sample(s): 01-0	09 Batch: V	VG1389227-3	WG1389227-4				
Methylene chloride	92		90		70-130	2		20	
1,1-Dichloroethane	95		94		70-130	1		20	
Chloroform	97		92		70-130	5		20	
Carbon tetrachloride	96		91		63-132	5		20	
1,2-Dichloropropane	92		90		70-130	2		20	
Dibromochloromethane	88		87		63-130	1		20	
1,1,2-Trichloroethane	86		90		70-130	5		20	
Tetrachloroethene	93		94		70-130	1		20	
Chlorobenzene	88		90		75-130	2		20	
Trichlorofluoromethane	98	_	96		62-150	2		20	
1,2-Dichloroethane	94		94		70-130	0		20	
1,1,1-Trichloroethane	99		96		67-130	3		20	
Bromodichloromethane	92		91		67-130	1		20	
trans-1,3-Dichloropropene	93		91		70-130	2		20	
cis-1,3-Dichloropropene	93		91		70-130	2		20	
Bromoform	84		86		54-136	2		20	
1,1,2,2-Tetrachloroethane	82		84		67-130	2		20	
Benzene	90		88		70-130	2		20	
Toluene	92		91		70-130	1		20	
Ethylbenzene	91		91		70-130	0		20	
Chloromethane	81		82		64-130	1		20	
Bromomethane	82		63		39-139	26	Q	20	
Vinyl chloride	82	_	78		55-140	5		20	



Lab Control Sample Analysis Batch Quality Control

Report Date: 07/08/20

Parameter	LCS %Recovery	Qual		CSD covery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough	Lab Associated	sample(s):	01-09	Batch:	WG1389227-3	WG1389227-4			
Chloroethane	86			85		55-138	1		20
1,1-Dichloroethene	94			94		61-145	0		20
trans-1,2-Dichloroethene	94			93		70-130	1		20
Trichloroethene	89			86		70-130	3		20
1,2-Dichlorobenzene	88			90		70-130	2		20
1,3-Dichlorobenzene	88			91		70-130	3		20
1,4-Dichlorobenzene	89			91		70-130	2		20
Methyl tert butyl ether	87			87		63-130	0		20
p/m-Xylene	90			90		70-130	0		20
o-Xylene	90			90		70-130	0		20
cis-1,2-Dichloroethene	92			90		70-130	2		20
Styrene	85			90		70-130	6		20
Dichlorodifluoromethane	73			75		36-147	3		20
Acetone	86			81		58-148	6		20
Carbon disulfide	93			91		51-130	2		20
2-Butanone	82			87		63-138	6		20
4-Methyl-2-pentanone	84			86		59-130	2		20
2-Hexanone	87			88		57-130	1		20
Bromochloromethane	94			94		70-130	0		20
1,2-Dibromoethane	85			85		70-130	0		20
1,2-Dibromo-3-chloropropane	79			81		41-144	3		20
Isopropylbenzene	93			96		70-130	3		20
1,2,3-Trichlorobenzene	77			85		70-130	10		20



Lab Control Sample Analysis

Batch Quality Control

Project Name: HHII TA1 Project Number: 16.6334 Lab Number: L2027725 Report Date: 07/08/20

LCS LCSD %Recovery RPD %Recovery %Recovery Parameter Qual Qual Limits RPD Qual Limits Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-09 Batch: WG1389227-3 WG1389227-4 1,2,4-Trichlorobenzene 81 87 70-130 7 20 84 86 2 Methyl Acetate 70-130 20 Cyclohexane 98 96 70-130 2 20 1,4-Dioxane 82 90 56-162 9 20 Freon-113 100 96 70-130 4 20 20 Methyl cyclohexane 96 95 70-130 1

Surrogate	LCS %Recovery Qua	LCSD I %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	110	110	70-130
Toluene-d8	104	105	70-130
4-Bromofluorobenzene	111	108	70-130
Dibromofluoromethane	103	103	70-130



Matrix Spike Analysis

Project Name:	HHII TA1	Batch Quality Control	ab

Project Number: 16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	MSD Qual Found	MSD %Recovery	Recovery Qual Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS RAMW5-200630	- Westborough	Lab Assoc	iated sample(s): 01-09 QC	Batch ID: WG13892	27-6 WG138	9227-7 QC Sample	e: L202	7725-01	Client ID:
Methylene chloride	ND	10	8.6	86	10	100	70-130	15		20
1,1-Dichloroethane	ND	10	9.0	90	11	110	70-130	20		20
Chloroform	ND	10	9.3	93	11	110	70-130	17		20
Carbon tetrachloride	ND	10	8.6	86	11	110	63-132	24	Q	20
1,2-Dichloropropane	ND	10	8.8	88	10	100	70-130	13		20
Dibromochloromethane	ND	10	8.0	80	10	100	63-130	22	Q	20
1,1,2-Trichloroethane	ND	10	8.3	83	9.9	99	70-130	18		20
Tetrachloroethene	1.4	10	9.6	82	12	106	70-130	22	Q	20
Chlorobenzene	ND	10	8.4	84	10	100	75-130	17		20
Trichlorofluoromethane	ND	10	9.0	90	12	120	62-150	29	Q	20
1,2-Dichloroethane	ND	10	8.9	89	11	110	70-130	21	Q	20
1,1,1-Trichloroethane	ND	10	9.2	92	11	110	67-130	18		20
Bromodichloromethane	ND	10	8.7	87	11	110	67-130	23	Q	20
trans-1,3-Dichloropropene	ND	10	8.1	81	10	100	70-130	21	Q	20
cis-1,3-Dichloropropene	ND	10	8.1	81	9.9	99	70-130	20		20
Bromoform	ND	10	7.8	78	9.5	95	54-136	20		20
1,1,2,2-Tetrachloroethane	ND	10	7.7	77	9.6	96	67-130	22	Q	20
Benzene	ND	10	8.5	85	10	100	70-130	16		20
Toluene	ND	10	8.6	86	11	110	70-130	24	Q	20
Ethylbenzene	ND	10	8.2	82	11	110	70-130	29	Q	20
Chloromethane	ND	10	8.2	82	10	100	64-130	20		20
Bromomethane	ND	10	4.8	48	7.2	72	39-139	40	Q	20
Vinyl chloride	ND	10	7.9	79	10	100	55-140	23	Q	20



Matrix Spike Analysis

Project Name:	HHII TA1	Batch Quality Control	Lab Numbe
Project Number:	16.6334		Report Dat

L2027725 ber: 07/08/20 Report Date:

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	MSD Qual Found	MSD %Recovery	Recovery Qual Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/M RAMW5-200630	S - Westborough	Lab Assoc	ciated sample	(s): 01-09 QC	Batch ID: WG13892	27-6 WG138	9227-7 QC Sample	e: L202	7725-01	Client ID:
Chloroethane	ND	10	8.8	88	10	100	55-138	13		20
1,1-Dichloroethene	ND	10	9.0	90	11	110	61-145	20		20
trans-1,2-Dichloroethene	ND	10	8.9	89	11	110	70-130	21	Q	20
Trichloroethene	ND	10	8.1	81	10	100	70-130	21	Q	20
1,2-Dichlorobenzene	ND	10	8.0	80	10	100	70-130	22	Q	20
1,3-Dichlorobenzene	ND	10	7.8	78	10	100	70-130	25	Q	20
1,4-Dichlorobenzene	ND	10	7.7	77	10	100	70-130	26	Q	20
Methyl tert butyl ether	ND	10	8.2	82	9.9	99	63-130	19		20
p/m-Xylene	ND	20	16	80	21	105	70-130	27	Q	20
o-Xylene	ND	20	16	80	21	105	70-130	27	Q	20
cis-1,2-Dichloroethene	ND	10	8.6	86	10	100	70-130	15		20
Styrene	ND	20	16	80	20	100	70-130	22	Q	20
Dichlorodifluoromethane	ND	10	6.1	61	8.5	85	36-147	33	Q	20
Acetone	ND	10	8.3	83	10	100	58-148	19		20
Carbon disulfide	ND	10	8.5	85	11	110	51-130	26	Q	20
2-Butanone	ND	10	7.3	73	9.1	91	63-138	22	Q	20
4-Methyl-2-pentanone	ND	10	8.2	82	9.6	96	59-130	16		20
2-Hexanone	ND	10	8.1	81	10	100	57-130	21	Q	20
Bromochloromethane	ND	10	9.6	96	10	100	70-130	4		20
1,2-Dibromoethane	ND	10	8.4	84	9.9	99	70-130	16		20
1,2-Dibromo-3-chloropropane	ND	10	7.6	76	9.3	93	41-144	20		20
Isopropylbenzene	ND	10	8.3	83	11	110	70-130	28	Q	20
1,2,3-Trichlorobenzene	ND	10	6.9	69	Q 9.6	96	70-130	33	Q	20



Matrix Spike Analysis

Project Name:	HHII TA1	Batch Quality Control	Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20

Parameter	Native Sample	MS Added	MS Found	MS %Recover	ry Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS RAMW5-200630	- Westborough	Lab Assoc	iated sample(s	s): 01-09 (QC Batch ID:	WG13892	27-6 WG1389	9227-7	QC Sample	e: L2027	7725-01	Client ID:
1,2,4-Trichlorobenzene	ND	10	6.8	68	Q	9.9	99		70-130	37	Q	20
Methyl Acetate	ND	10	7.6	76		8.8	88		70-130	15		20
Cyclohexane	ND	10	8.7J	87		11	110		70-130	23	Q	20
1,4-Dioxane	ND	500	420	84		470	94		56-162	11		20
Freon-113	ND	10	8.1	81		11	110		70-130	30	Q	20
Methyl cyclohexane	ND	10	7.6J	76		11	110		70-130	37	Q	20

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
1,2-Dichloroethane-d4	111	111	70-130
4-Bromofluorobenzene	105	105	70-130
Dibromofluoromethane	103	103	70-130
Toluene-d8	105	105	70-130



SEMIVOLATILES



			Serial_No	07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-01		Date Collected:	06/30/20 08:00
Client ID:	RAMW5-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	I: EPA 3510C
Analytical Method:	1,8270D		Extraction Date:	07/01/20 18:11
Analytical Date:	07/02/20 18:57			
Analyst:	SZ			
•				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS -	Westborough Lab					
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1
3.3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1
Isophorone	ND		ug/l	5.0	1.2	1
Nitrobenzene	ND		ug/l	2.0	0.77	1
NDPA/DPA	ND		ug/l	2.0	0.42	1
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1
Diethyl phthalate	ND		ug/l	5.0	0.38	1
Dimethyl phthalate	ND		ug/l	5.0	1.8	1
Biphenyl	ND		ug/l	2.0	0.46	1
4-Chloroaniline	ND		ug/l	5.0	1.1	1
2-Nitroaniline	ND		ug/l	5.0	0.50	1
3-Nitroaniline	ND		ug/l	5.0	0.81	1
4-Nitroaniline	ND		ug/l	5.0	0.80	1
Dibenzofuran	ND		ug/l	2.0	0.50	1
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1
Acetophenone	ND		ug/l	5.0	0.53	1
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1



Serial_No:07082013:26 **Project Name:** Lab Number: HHII TA1 L2027725 **Project Number:** Report Date: 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-01 Date Collected: 06/30/20 08:00 Client ID: Date Received: 06/30/20 RAMW5-200630 Sample Location: Field Prep: Not Specified SCHENECTADY NY Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - Wes	stborough Lab					
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1
2-Chlorophenol	ND		ug/l	2.0	0.48	1
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1
2-Nitrophenol	ND		ug/l	10	0.85	1
4-Nitrophenol	ND		ug/l	10	0.67	1
2,4-Dinitrophenol	ND		ug/l	20	6.6	1
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1
Phenol	ND		ug/l	5.0	0.57	1
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1
Carbazole	ND		ug/l	2.0	0.49	1
Atrazine	ND		ug/l	10	0.76	1
Benzaldehyde	ND		ug/l	5.0	0.53	1
Caprolactam	ND		ug/l	10	3.3	1
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1

Total TIC Compounds 2.44 J ug/l 1					
	Total TIC Compounds	2.44	J	ug/l	1
Unknown 2.44 J ug/l 1	Unknown	2.44	J	ug/l	1

2-Fluorophenol 66 21-120 Phenol-d6 54 10-120 Nitrobenzene-d5 93 23-120 2-Fluorobiphenyl 70 15-120
Nitrobenzene-d5 93 23-120
2-Fluorobiphenyl 70 15-120
2,4,6-Tribromophenol 57 10-120
4-Terphenyl-d14 73 41-149



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-01		Date Collected:	06/30/20 08:00
Client ID:	RAMW5-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	: EPA 3510C
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 18:15
Analytical Date:	07/02/20 19:37			
Analyst:	DV			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-	-SIM - Westborough La	b				
Acenaphthene	ND		ug/l	0.10	0.01	1
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1
Fluoranthene	ND		ug/l	0.10	0.02	1
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1
Naphthalene	ND		ug/l	0.10	0.05	1
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1
Chrysene	ND		ug/l	0.10	0.01	1
Acenaphthylene	ND		ug/l	0.10	0.01	1
Anthracene	ND		ug/l	0.10	0.01	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1
Fluorene	ND		ug/l	0.10	0.01	1
Phenanthrene	ND		ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1
Pyrene	ND		ug/l	0.10	0.02	1
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1
Pentachlorophenol	ND		ug/l	0.80	0.01	1
Hexachlorobenzene	ND		ug/l	0.80	0.01	1
Hexachloroethane	ND		ug/l	0.80	0.06	1



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor
Sample Depth:						
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified
Client ID:	RAMW5-200630				Date Received:	06/30/20
Lab ID:	L2027725-01				Date Collected:	06/30/20 08:00
		SAMPI	LE RESULTS	3		
Project Number:	16.6334				Report Date:	07/08/20
Project Name:	HHII TA1				Lab Number:	L2027725
					Serial_N	o:07082013:26

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	71	21-120
Phenol-d6	64	10-120
Nitrobenzene-d5	107	23-120
2-Fluorobiphenyl	89	15-120
2,4,6-Tribromophenol	112	10-120
4-Terphenyl-d14	116	41-149



			Serial_No	07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-02		Date Collected:	06/30/20 08:35
Client ID:	RAMW6-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	I: EPA 3510C
Analytical Method:	1,8270D		Extraction Date:	07/01/20 18:11
Analytical Date:	07/02/20 20:31			
Analyst:	SZ			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Westborough Lab									
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1			
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1			
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1			
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1			
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1			
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1			
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1			
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1			
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1			
Isophorone	ND		ug/l	5.0	1.2	1			
Nitrobenzene	ND		ug/l	2.0	0.77	1			
NDPA/DPA	ND		ug/l	2.0	0.42	1			
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1			
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1			
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1			
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1			
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1			
Diethyl phthalate	ND		ug/l	5.0	0.38	1			
Dimethyl phthalate	ND		ug/l	5.0	1.8	1			
Biphenyl	ND		ug/l	2.0	0.46	1			
4-Chloroaniline	ND		ug/l	5.0	1.1	1			
2-Nitroaniline	ND		ug/l	5.0	0.50	1			
3-Nitroaniline	ND		ug/l	5.0	0.81	1			
4-Nitroaniline	ND		ug/l	5.0	0.80	1			
Dibenzofuran	ND		ug/l	2.0	0.50	1			
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1			
Acetophenone	ND		ug/l	5.0	0.53	1			
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1			



Serial_No:07082013:26 **Project Name:** Lab Number: HHII TA1 L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-02 Date Collected: 06/30/20 08:35 Client ID: Date Received: 06/30/20 RAMW6-200630 Sample Location: Field Prep: SCHENECTADY NY Not Specified Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS	- Westborough Lab					
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1
2-Chlorophenol	ND		ug/l	2.0	0.48	1
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1
2-Nitrophenol	ND		ug/l	10	0.85	1
4-Nitrophenol	ND		ug/l	10	0.67	1
2,4-Dinitrophenol	ND		ug/l	20	6.6	1
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1
Phenol	ND		ug/l	5.0	0.57	1
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1
Carbazole	ND		ug/l	2.0	0.49	1
Atrazine	ND		ug/l	10	0.76	1
Benzaldehyde	ND		ug/l	5.0	0.53	1
Caprolactam	ND		ug/l	10	3.3	1
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1

Total TIC Compounds	1.78	J	ug/l	1
Unknown	1.78	J	ug/l	1
Childiown	1.70	0	ug/i	•

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	69	21-120
Phenol-d6	58	10-120
Nitrobenzene-d5	92	23-120
2-Fluorobiphenyl	70	15-120
2,4,6-Tribromophenol	66	10-120
4-Terphenyl-d14	77	41-149



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-02		Date Collected:	06/30/20 08:35
Client ID:	RAMW6-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth: Matrix:	Water		Extraction Method	: EPA 3510C
Analytical Method: Analytical Date: Analyst:	1,8270D-SIM 07/02/20 19:57 DV		Extraction Date:	07/01/20 18:15

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS-SIM - Westborough Lab									
Acenaphthene	ND		ug/l	0.10	0.01	1			
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1			
Fluoranthene	ND		ug/l	0.10	0.02	1			
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1			
Naphthalene	ND		ug/l	0.10	0.05	1			
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1			
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1			
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1			
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1			
Chrysene	ND		ug/l	0.10	0.01	1			
Acenaphthylene	ND		ug/l	0.10	0.01	1			
Anthracene	ND		ug/l	0.10	0.01	1			
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1			
Fluorene	ND		ug/l	0.10	0.01	1			
Phenanthrene	ND		ug/l	0.10	0.02	1			
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1			
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1			
Pyrene	ND		ug/l	0.10	0.02	1			
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1			
Pentachlorophenol	ND		ug/l	0.80	0.01	1			
Hexachlorobenzene	ND		ug/l	0.80	0.01	1			
Hexachloroethane	ND		ug/l	0.80	0.06	1			



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor
Sample Depth:						
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified
Client ID:	RAMW6-200630				Date Received:	06/30/20
Lab ID:	L2027725-02				Date Collected:	06/30/20 08:35
		SAMP		6		
Project Number:	16.6334				Report Date:	07/08/20
Project Name:	HHII TA1				Lab Number:	L2027725
					Serial_N	0:07082013:26

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	73	21-120
Phenol-d6	66	10-120
Nitrobenzene-d5	108	23-120
2-Fluorobiphenyl	90	15-120
2,4,6-Tribromophenol	108	10-120
4-Terphenyl-d14	117	41-149



		Serial_No	:07082013:26
HHII TA1		Lab Number:	L2027725
16.6334		Report Date:	07/08/20
	SAMPLE RESULTS		
L2027725-03		Date Collected:	06/30/20 11:35
RAMW4-200630		Date Received:	06/30/20
SCHENECTADY NY		Field Prep:	Not Specified
Water		Extraction Method	: EPA 3510C
1.8270D		Extraction Date:	07/01/20 18:11
07/02/20 20:54			
SZ			
	16.6334 L2027725-03 RAMW4-200630 SCHENECTADY NY Water 1,8270D 07/02/20 20:54	16.6334 SAMPLE RESULTS L2027725-03 RAMW4-200630 SCHENECTADY NY Water 1,8270D 07/02/20 20:54	HHII TA1 Lab Number: 16.6334 Report Date: L2027725-03 Date Collected: RAMW4-200630 SCHENECTADY NY Field Prep: Water 1,8270D 07/02/20 20:54 Extraction Method

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Westborough Lab									
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1			
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1			
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1			
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1			
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1			
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1			
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1			
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1			
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1			
Isophorone	ND		ug/l	5.0	1.2	1			
Nitrobenzene	ND		ug/l	2.0	0.77	1			
NDPA/DPA	ND		ug/l	2.0	0.42	1			
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1			
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1			
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1			
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1			
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1			
Diethyl phthalate	ND		ug/l	5.0	0.38	1			
Dimethyl phthalate	ND		ug/l	5.0	1.8	1			
Biphenyl	ND		ug/l	2.0	0.46	1			
4-Chloroaniline	ND		ug/l	5.0	1.1	1			
2-Nitroaniline	ND		ug/l	5.0	0.50	1			
3-Nitroaniline	ND		ug/l	5.0	0.81	1			
4-Nitroaniline	ND		ug/l	5.0	0.80	1			
Dibenzofuran	ND		ug/l	2.0	0.50	1			
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1			
Acetophenone	ND		ug/l	5.0	0.53	1			
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1			



Serial_No:07082013:26 **Project Name:** Lab Number: HHII TA1 L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-03 Date Collected: 06/30/20 11:35 Client ID: Date Received: 06/30/20 RAMW4-200630 Sample Location: Field Prep: SCHENECTADY NY Not Specified Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Westborough Lab									
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1			
2-Chlorophenol	ND		ug/l	2.0	0.48	1			
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1			
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1			
2-Nitrophenol	ND		ug/l	10	0.85	1			
4-Nitrophenol	ND		ug/l	10	0.67	1			
2,4-Dinitrophenol	ND		ug/l	20	6.6	1			
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1			
Phenol	ND		ug/l	5.0	0.57	1			
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1			
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1			
Carbazole	ND		ug/l	2.0	0.49	1			
Atrazine	ND		ug/l	10	0.76	1			
Benzaldehyde	ND		ug/l	5.0	0.53	1			
Caprolactam	ND		ug/l	10	3.3	1			
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1			

Total TIC Compounds	2.65	J	ug/l	1
Unknown	2.65	J	ug/l	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	68	21-120
Phenol-d6	58	10-120
Nitrobenzene-d5	88	23-120
2-Fluorobiphenyl	67	15-120
2,4,6-Tribromophenol	67	10-120
4-Terphenyl-d14	74	41-149



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-03		Date Collected:	06/30/20 11:35
Client ID:	RAMW4-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	: EPA 3510C
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 18:15
Analytical Date:	07/02/20 20:16			
Analyst:	DV			
-				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM - Westborough Lab								
Acenaphthene	ND		ug/l	0.10	0.01	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1		
Fluoranthene	ND		ug/l	0.10	0.02	1		
Hexachlorobutadiene	ND		ug/l	0.50	0.02	1		
Naphthalene	ND		ug/l	0.00	0.05	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.03	1		
Benzo(a)pyrene	ND		•	0.10	0.02			
	ND		ug/l	0.10	0.02	1		
Benzo(b)fluoranthene			ug/l					
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1		
Chrysene	ND		ug/l	0.10	0.01	1		
Acenaphthylene	ND		ug/l	0.10	0.01	1		
Anthracene	ND		ug/l	0.10	0.01	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1		
Fluorene	ND		ug/l	0.10	0.01	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1		
Pyrene	ND		ug/l	0.10	0.02	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1		
Pentachlorophenol	ND		ug/l	0.80	0.01	1		
Hexachlorobenzene	ND		ug/l	0.80	0.01	1		
Hexachloroethane	ND		ug/l	0.80	0.06	1		



Parameter		Result	Qualifier	Units	RL	MDL	Dilution Factor	
Sample Depth:								
Sample Location:	SCHENECTADY NY				Field Prep	:	Not Specified	
Client ID:	RAMW4-200630				Date Rece	ived:	06/30/20	
Lab ID:	L2027725-03				Date Colle	cted:	06/30/20 11:35	
		SAMPI	LE RESULTS	5				
Project Number:	16.6334				Report D	ate:	07/08/20	
Project Name:	HHII TA1				Lab Num	nber:	L2027725	
	Serial_No:07082013:26							

Surrogate	% Recovery	Qualifier	Acceptance Criteria
2-Fluorophenol	72		21-120
Phenol-d6	66		10-120
Nitrobenzene-d5	101		23-120
2-Fluorobiphenyl	85		15-120
2,4,6-Tribromophenol	123	Q	10-120
4-Terphenyl-d14	120		41-149



			Serial_No:07082013:26		
Project Name:	HHII TA1		Lab Number:	L2027725	
Project Number:	16.6334		Report Date:	07/08/20	
		SAMPLE RESULTS			
Lab ID:	L2027725-04		Date Collected:	06/30/20 11:50	
Client ID:	RAMW1-200630		Date Received:	06/30/20	
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified	
Sample Depth:					
Matrix:	Water		Extraction Method	: EPA 3510C	
Analytical Method:	1,8270D		Extraction Date:	07/01/20 18:11	
Analytical Date:	07/02/20 21:18				
Analyst:	SZ				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Westborough Lab								
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1		
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1		
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1		
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1		
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1		
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1		
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1		
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1		
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1		
Isophorone	ND		ug/l	5.0	1.2	1		
Nitrobenzene	ND		ug/l	2.0	0.77	1		
NDPA/DPA	ND		ug/l	2.0	0.42	1		
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1		
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1		
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1		
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1		
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1		
Diethyl phthalate	ND		ug/l	5.0	0.38	1		
Dimethyl phthalate	ND		ug/l	5.0	1.8	1		
Biphenyl	ND		ug/l	2.0	0.46	1		
4-Chloroaniline	ND		ug/l	5.0	1.1	1		
2-Nitroaniline	ND		ug/l	5.0	0.50	1		
3-Nitroaniline	ND		ug/l	5.0	0.81	1		
4-Nitroaniline	ND		ug/l	5.0	0.80	1		
Dibenzofuran	ND		ug/l	2.0	0.50	1		
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1		
Acetophenone	ND		ug/l	5.0	0.53	1		
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1		



Serial_No:07082013:26 Project Name: Lab Number: HHII TA1 L2027725 **Project Number:** Report Date: 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-04 Date Collected: 06/30/20 11:50 Client ID: Date Received: 06/30/20 RAMW1-200630 Sample Location: Field Prep: Not Specified SCHENECTADY NY Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Westborough Lab								
	ND			2.0	0.05	1		
p-Chloro-m-cresol			ug/l	2.0	0.35	•		
2-Chlorophenol	ND		ug/l	2.0	0.48	1		
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1		
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1		
2-Nitrophenol	ND		ug/l	10	0.85	1		
4-Nitrophenol	ND		ug/l	10	0.67	1		
2,4-Dinitrophenol	ND		ug/l	20	6.6	1		
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1		
Phenol	ND		ug/l	5.0	0.57	1		
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1		
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1		
Carbazole	ND		ug/l	2.0	0.49	1		
Atrazine	ND		ug/l	10	0.76	1		
Benzaldehyde	ND		ug/l	5.0	0.53	1		
Caprolactam	ND		ug/l	10	3.3	1		
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1		

Tentatively Identified Compounds				
Total TIC Compounds	4.84	J	ug/l	1
Unknown	2.44	J	ug/l	1
Unknown	2.40	J	ug/l	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
2-Fluorophenol	63	21-120	
Phenol-d6	53	10-120	
Nitrobenzene-d5	84	23-120	
2-Fluorobiphenyl	64	15-120	
2,4,6-Tribromophenol	57	10-120	
4-Terphenyl-d14	71	41-149	



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-04		Date Collected:	06/30/20 11:50
Client ID:	RAMW1-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 18:15
Analytical Date:	07/02/20 20:36			
Analyst:	DV			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS	-SIM - Westborough La	b				
Acenaphthene	ND		ug/l	0.10	0.01	1
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1
Fluoranthene	ND		ug/l	0.10	0.02	1
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1
Naphthalene	ND		ug/l	0.10	0.05	1
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1
Chrysene	ND		ug/l	0.10	0.01	1
Acenaphthylene	ND		ug/l	0.10	0.01	1
Anthracene	ND		ug/l	0.10	0.01	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1
Fluorene	ND		ug/l	0.10	0.01	1
Phenanthrene	ND		ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1
Pyrene	ND		ug/l	0.10	0.02	1
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1
Pentachlorophenol	ND		ug/l	0.80	0.01	1
Hexachlorobenzene	ND		ug/l	0.80	0.01	1
Hexachloroethane	ND		ug/l	0.80	0.06	1



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor
Sample Depth:						
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified
Client ID:	RAMW1-200630				Date Received:	06/30/20
Lab ID:	L2027725-04				Date Collected:	06/30/20 11:50
		SAMP	LE RESULTS	5		
Project Number:	16.6334				Report Date:	07/08/20
Project Name:	HHII TA1				Lab Number:	L2027725
					Serial_N	lo:07082013:26

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	71	21-120
Phenol-d6	62	10-120
Nitrobenzene-d5	101	23-120
2-Fluorobiphenyl	85	15-120
2,4,6-Tribromophenol	108	10-120
4-Terphenyl-d14	116	41-149



			Serial_No:07082013:26		
Project Name:	HHII TA1		Lab Number:	L2027725	
Project Number:	16.6334		Report Date:	07/08/20	
		SAMPLE RESULTS			
Lab ID:	L2027725-05		Date Collected:	06/30/20 12:05	
Client ID:	RAMW2-200630		Date Received:	06/30/20	
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified	
Sample Depth:					
Matrix:	Water		Extraction Method	: EPA 3510C	
Analytical Method:	1,8270D		Extraction Date:	07/01/20 18:11	
Analytical Date:	07/02/20 21:41				
Analyst:	SZ				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS -	Westborough Lab					
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1
Isophorone	ND		ug/l	5.0	1.2	1
Nitrobenzene	ND		ug/l	2.0	0.77	1
NDPA/DPA	ND		ug/l	2.0	0.42	1
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1
Diethyl phthalate	ND		ug/l	5.0	0.38	1
Dimethyl phthalate	ND		ug/l	5.0	1.8	1
Biphenyl	ND		ug/l	2.0	0.46	1
4-Chloroaniline	ND		ug/l	5.0	1.1	1
2-Nitroaniline	ND		ug/l	5.0	0.50	1
3-Nitroaniline	ND		ug/l	5.0	0.81	1
4-Nitroaniline	ND		ug/l	5.0	0.80	1
Dibenzofuran	ND		ug/l	2.0	0.50	1
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1
Acetophenone	ND		ug/l	5.0	0.53	1
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1



Serial_No:07082013:26 **Project Name:** Lab Number: HHII TA1 L2027725 **Project Number:** Report Date: 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-05 Date Collected: 06/30/20 12:05 Client ID: Date Received: 06/30/20 RAMW2-200630 Sample Location: Field Prep: SCHENECTADY NY Not Specified Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Westborough Lab									
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1			
2-Chlorophenol	ND		ug/l	2.0	0.48	1			
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1			
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1			
2-Nitrophenol	ND		ug/l	10	0.85	1			
4-Nitrophenol	ND		ug/l	10	0.67	1			
2,4-Dinitrophenol	ND		ug/l	20	6.6	1			
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1			
Phenol	ND		ug/l	5.0	0.57	1			
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1			
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1			
Carbazole	ND		ug/l	2.0	0.49	1			
Atrazine	ND		ug/l	10	0.76	1			
Benzaldehyde	ND		ug/l	5.0	0.53	1			
Caprolactam	ND		ug/l	10	3.3	1			
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1			

Tentatively Identified Compounds	ND	ug/l			1
Surrogate		% Recovery	Qualifier	Acceptance Criteria	
2-Fluorophenol		51		21-120	
Phenol-d6		42		10-120	
Nitrobenzene-d5		70		23-120	
2-Fluorobiphenyl		56		15-120	
2,4,6-Tribromophenol		54		10-120	
4-Terphenyl-d14		69		41-149	



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-05		Date Collected:	06/30/20 12:05
Client ID:	RAMW2-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	Water 1,8270D-SIM 07/02/20 20:56 DV		Extraction Method Extraction Date:	: EPA 3510C 07/01/20 18:15

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS-SIM - Westborough Lab									
Acenaphthene	ND		ug/l	0.10	0.01	1			
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1			
Fluoranthene	ND		ug/l	0.10	0.02	1			
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1			
Naphthalene	ND		ug/l	0.10	0.05	1			
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1			
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1			
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1			
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1			
Chrysene	ND		ug/l	0.10	0.01	1			
Acenaphthylene	ND		ug/l	0.10	0.01	1			
Anthracene	ND		ug/l	0.10	0.01	1			
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1			
Fluorene	ND		ug/l	0.10	0.01	1			
Phenanthrene	ND		ug/l	0.10	0.02	1			
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1			
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1			
Pyrene	ND		ug/l	0.10	0.02	1			
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1			
Pentachlorophenol	ND		ug/l	0.80	0.01	1			
Hexachlorobenzene	ND		ug/l	0.80	0.01	1			
Hexachloroethane	ND		ug/l	0.80	0.06	1			



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor
Danamatan		Beault	Qualifian	Unito	RL MDL	Dilution Foster
Sample Depth:						
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified
Client ID:	RAMW2-200630				Date Received:	06/30/20
Lab ID:	L2027725-05				Date Collected:	06/30/20 12:05
		SAMP		6		
Project Number:	16.6334				Report Date:	07/08/20
Project Name:	HHII TA1				Lab Number:	L2027725
					Serial_N	lo:07082013:26

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	56	21-120
Phenol-d6	49	10-120
Nitrobenzene-d5	82	23-120
2-Fluorobiphenyl	72	15-120
2,4,6-Tribromophenol	91	10-120
4-Terphenyl-d14	114	41-149



			Serial_No	0:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-06		Date Collected:	06/30/20 12:55
Client ID:	RAMW3-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	d: EPA 3510C
Analytical Method:	1,8270D		Extraction Date:	07/01/20 18:11
Analytical Date:	07/02/20 22:05			
Analyst:	SZ			
•				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Westborough Lab								
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1		
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1		
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1		
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1		
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1		
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1		
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1		
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1		
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1		
Isophorone	ND		ug/l	5.0	1.2	1		
Nitrobenzene	ND		ug/l	2.0	0.77	1		
NDPA/DPA	ND		ug/l	2.0	0.42	1		
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1		
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1		
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1		
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1		
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1		
Diethyl phthalate	ND		ug/l	5.0	0.38	1		
Dimethyl phthalate	ND		ug/l	5.0	1.8	1		
Biphenyl	ND		ug/l	2.0	0.46	1		
4-Chloroaniline	ND		ug/l	5.0	1.1	1		
2-Nitroaniline	ND		ug/l	5.0	0.50	1		
3-Nitroaniline	ND		ug/l	5.0	0.81	1		
4-Nitroaniline	ND		ug/l	5.0	0.80	1		
Dibenzofuran	ND		ug/l	2.0	0.50	1		
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1		
Acetophenone	ND		ug/l	5.0	0.53	1		
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1		



Serial_No:07082013:26 Project Name: Lab Number: HHII TA1 L2027725 **Project Number:** Report Date: 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-06 Date Collected: 06/30/20 12:55 Client ID: Date Received: 06/30/20 RAMW3-200630 Sample Location: Field Prep: Not Specified SCHENECTADY NY Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Westborough Lab									
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1			
2-Chlorophenol	ND		ug/l	2.0	0.48	1			
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1			
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1			
2-Nitrophenol	ND		ug/l	10	0.85	1			
4-Nitrophenol	ND		ug/l	10	0.67	1			
2,4-Dinitrophenol	ND		ug/l	20	6.6	1			
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1			
Phenol	ND		ug/l	5.0	0.57	1			
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1			
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1			
Carbazole	ND		ug/l	2.0	0.49	1			
Atrazine	ND		ug/l	10	0.76	1			
Benzaldehyde	ND		ug/l	5.0	0.53	1			
Caprolactam	ND		ug/l	10	3.3	1			
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1			

Tentatively Identified Compounds				
Total TIC Compounds	3.85	J	ug/l	1
Unknown	2.40	J	ug/l	1
Unknown	1.45	J	ug/l	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
2-Fluorophenol	56	21-120	
Phenol-d6	47	10-120	
Nitrobenzene-d5	75	23-120	
2-Fluorobiphenyl	57	15-120	
2,4,6-Tribromophenol	46	10-120	
4-Terphenyl-d14	61	41-149	



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-06		Date Collected:	06/30/20 12:55
Client ID:	RAMW3-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	I: EPA 3510C
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 18:15
Analytical Date:	07/02/20 21:15			
Analyst:	DV			
-				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS-SIM - Westborough Lab									
Acenaphthene	ND		ug/l	0.10	0.01	1			
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1			
Fluoranthene	ND		ug/l	0.10	0.02	1			
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1			
Naphthalene	ND		ug/l	0.10	0.05	1			
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1			
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1			
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1			
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1			
Chrysene	ND		ug/l	0.10	0.01	1			
Acenaphthylene	ND		ug/l	0.10	0.01	1			
Anthracene	ND		ug/l	0.10	0.01	1			
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1			
Fluorene	ND		ug/l	0.10	0.01	1			
Phenanthrene	ND		ug/l	0.10	0.02	1			
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1			
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1			
Pyrene	ND		ug/l	0.10	0.02	1			
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1			
Pentachlorophenol	ND		ug/l	0.80	0.01	1			
Hexachlorobenzene	ND		ug/l	0.80	0.01	1			
Hexachloroethane	ND		ug/l	0.80	0.06	1			



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor
Sample Depth:						
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified
Client ID:	RAMW3-200630				Date Received:	06/30/20
Lab ID:	L2027725-06				Date Collected:	06/30/20 12:55
		SAMP	LE RESULTS	6		
Project Number:	16.6334				Report Date:	07/08/20
Project Name:	HHII TA1				Lab Number:	L2027725
					Serial_N	lo:07082013:26

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	58	21-120
Phenol-d6	53	10-120
Nitrobenzene-d5	89	23-120
2-Fluorobiphenyl	76	15-120
2,4,6-Tribromophenol	90	10-120
4-Terphenyl-d14	95	41-149



		Serial_No:	07082013:26
HHII TA1		Lab Number:	L2027725
16.6334		Report Date:	07/08/20
	SAMPLE RESULTS		
L2027725-07		Date Collected:	06/30/20 00:00
FD01-200630		Date Received:	06/30/20
SCHENECTADY NY		Field Prep:	Not Specified
Water		Extraction Method:	EPA 3510C
		Extraction Date:	07/01/20 18:11
,			
SZ			
	16.6334 L2027725-07 FD01-200630 SCHENECTADY NY Water 1,8270D 07/02/20 22:29	16.6334 SAMPLE RESULTS L2027725-07 FD01-200630 SCHENECTADY NY Water 1,8270D 07/02/20 22:29	HHII TA1 Lab Number: 16.6334 Report Date: L2027725-07 Date Collected: FD01-200630 SCHENECTADY NY Date Received: SCHENECTADY NY Extraction Method: 1,8270D 07/02/20 22:29

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Westborough Lab								
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1		
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1		
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1		
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1		
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1		
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1		
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1		
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1		
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1		
Isophorone	ND		ug/l	5.0	1.2	1		
Nitrobenzene	ND		ug/l	2.0	0.77	1		
NDPA/DPA	ND		ug/l	2.0	0.42	1		
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1		
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1		
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1		
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1		
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1		
Diethyl phthalate	ND		ug/l	5.0	0.38	1		
Dimethyl phthalate	ND		ug/l	5.0	1.8	1		
Biphenyl	ND		ug/l	2.0	0.46	1		
4-Chloroaniline	ND		ug/l	5.0	1.1	1		
2-Nitroaniline	ND		ug/l	5.0	0.50	1		
3-Nitroaniline	ND		ug/l	5.0	0.81	1		
4-Nitroaniline	ND		ug/l	5.0	0.80	1		
Dibenzofuran	ND		ug/l	2.0	0.50	1		
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1		
Acetophenone	ND		ug/l	5.0	0.53	1		
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1		



Serial_No:07082013:26 **Project Name:** Lab Number: HHII TA1 L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-07 Date Collected: 06/30/20 00:00 Client ID: Date Received: 06/30/20 FD01-200630 Sample Location: Field Prep: Not Specified SCHENECTADY NY

Sample	Depth:
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Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Westborough Lab								
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1		
2-Chlorophenol	ND		ug/l	2.0	0.48	1		
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1		
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1		
2-Nitrophenol	ND		ug/l	10	0.85	1		
4-Nitrophenol	ND		ug/l	10	0.67	1		
2,4-Dinitrophenol	ND		ug/l	20	6.6	1		
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1		
Phenol	ND		ug/l	5.0	0.57	1		
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1		
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1		
Carbazole	ND		ug/l	2.0	0.49	1		
Atrazine	ND		ug/l	10	0.76	1		
Benzaldehyde	ND		ug/l	5.0	0.53	1		
Caprolactam	ND		ug/l	10	3.3	1		
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1		

Total TIC Compounds	1.85	J	ug/l	1
Unknown	1.85	J	ug/l	1

2-Fluorophenol 67 21-120 Phenol-d6 55 10-120
Phenol-d6 55 10-120
Nitrobenzene-d5 84 23-120
2-Fluorobiphenyl 62 15-120
2,4,6-Tribromophenol 61 10-120
4-Terphenyl-d14 73 41-149



			Serial_No:	07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-07		Date Collected:	06/30/20 00:00
Client ID:	FD01-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method:	EPA 3510C
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 18:15
Analytical Date:	07/02/20 21:35			
Analyst:	DV			
•				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM - Westborough Lab								
Acenaphthene	ND		ug/l	0.10	0.01	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.01	1		
Fluoranthene	ND		•	0.20	0.02	1		
			ug/l					
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1		
Naphthalene	ND		ug/l	0.10	0.05	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1		
Chrysene	ND		ug/l	0.10	0.01	1		
Acenaphthylene	ND		ug/l	0.10	0.01	1		
Anthracene	ND		ug/l	0.10	0.01	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1		
Fluorene	ND		ug/l	0.10	0.01	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1		
Pyrene	ND		ug/l	0.10	0.02	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1		
Pentachlorophenol	ND		ug/l	0.80	0.01	1		
Hexachlorobenzene	ND		ug/l	0.80	0.01	1		
Hexachloroethane	ND		ug/l	0.80	0.06	1		



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor
Sample Depth:						
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified
Client ID:	FD01-200630				Date Received:	06/30/20
Lab ID:	L2027725-07				Date Collected:	06/30/20 00:00
		SAMP	LE RESULTS	5		
Project Number:	16.6334				Report Date:	07/08/20
Project Name:	HHII TA1				Lab Number:	L2027725
	Serial_No:07082013:26					lo:07082013:26

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	72	21-120
Phenol-d6	64	10-120
Nitrobenzene-d5	99	23-120
2-Fluorobiphenyl	89	15-120
2,4,6-Tribromophenol	111	10-120
4-Terphenyl-d14	117	41-149



		Serial_No:07082013:26		
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-08		Date Collected:	06/30/20 13:05
Client ID:	EB01-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	d: EPA 3510C
Analytical Method:	1,8270D		Extraction Date:	07/01/20 18:11
Analytical Date:	07/02/20 22:52			
Analyst:	SZ			
-				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Westborough Lab								
Bis(2-chloroethyl)ether	ND		ug/l	2.0	0.50	1		
3,3'-Dichlorobenzidine	ND		ug/l	5.0	1.6	1		
2,4-Dinitrotoluene	ND		ug/l	5.0	1.2	1		
2,6-Dinitrotoluene	ND		ug/l	5.0	0.93	1		
4-Chlorophenyl phenyl ether	ND		ug/l	2.0	0.49	1		
4-Bromophenyl phenyl ether	ND		ug/l	2.0	0.38	1		
Bis(2-chloroisopropyl)ether	ND		ug/l	2.0	0.53	1		
Bis(2-chloroethoxy)methane	ND		ug/l	5.0	0.50	1		
Hexachlorocyclopentadiene	ND		ug/l	20	0.69	1		
Isophorone	ND		ug/l	5.0	1.2	1		
Nitrobenzene	ND		ug/l	2.0	0.77	1		
NDPA/DPA	ND		ug/l	2.0	0.42	1		
n-Nitrosodi-n-propylamine	ND		ug/l	5.0	0.64	1		
Bis(2-ethylhexyl)phthalate	ND		ug/l	3.0	1.5	1		
Butyl benzyl phthalate	ND		ug/l	5.0	1.2	1		
Di-n-butylphthalate	ND		ug/l	5.0	0.39	1		
Di-n-octylphthalate	ND		ug/l	5.0	1.3	1		
Diethyl phthalate	ND		ug/l	5.0	0.38	1		
Dimethyl phthalate	ND		ug/l	5.0	1.8	1		
Biphenyl	ND		ug/l	2.0	0.46	1		
4-Chloroaniline	ND		ug/l	5.0	1.1	1		
2-Nitroaniline	ND		ug/l	5.0	0.50	1		
3-Nitroaniline	ND		ug/l	5.0	0.81	1		
4-Nitroaniline	ND		ug/l	5.0	0.80	1		
Dibenzofuran	ND		ug/l	2.0	0.50	1		
1,2,4,5-Tetrachlorobenzene	ND		ug/l	10	0.44	1		
Acetophenone	ND		ug/l	5.0	0.53	1		
2,4,6-Trichlorophenol	ND		ug/l	5.0	0.61	1		



Serial_No:07082013:26 **Project Name:** Lab Number: HHII TA1 L2027725 **Project Number: Report Date:** 16.6334 07/08/20 SAMPLE RESULTS Lab ID: L2027725-08 Date Collected: 06/30/20 13:05 Client ID: Date Received: 06/30/20 EB01-200630 Sample Location: Field Prep: SCHENECTADY NY Not Specified

Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Westborough Lab									
p-Chloro-m-cresol	ND		ug/l	2.0	0.35	1			
2-Chlorophenol	ND		ug/l	2.0	0.48	1			
2,4-Dichlorophenol	ND		ug/l	5.0	0.41	1			
2,4-Dimethylphenol	ND		ug/l	5.0	1.8	1			
2-Nitrophenol	ND		ug/l	10	0.85	1			
4-Nitrophenol	ND		ug/l	10	0.67	1			
2,4-Dinitrophenol	ND		ug/l	20	6.6	1			
4,6-Dinitro-o-cresol	ND		ug/l	10	1.8	1			
Phenol	ND		ug/l	5.0	0.57	1			
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0	0.48	1			
2,4,5-Trichlorophenol	ND		ug/l	5.0	0.77	1			
Carbazole	ND		ug/l	2.0	0.49	1			
Atrazine	ND		ug/l	10	0.76	1			
Benzaldehyde	ND		ug/l	5.0	0.53	1			
Caprolactam	ND		ug/l	10	3.3	1			
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0	0.84	1			

Total TIC Compounds	1 71	1	ug/l	1
	1.71	J	ug/l	1
Unknown Organic Acid	1.71	J	ug/l	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	64	21-120
Phenol-d6	52	10-120
Nitrobenzene-d5	85	23-120
2-Fluorobiphenyl	64	15-120
2,4,6-Tribromophenol	63	10-120
4-Terphenyl-d14	72	41-149



			Serial_No	:07082013:26
Project Name:	HHII TA1		Lab Number:	L2027725
Project Number:	16.6334		Report Date:	07/08/20
		SAMPLE RESULTS		
Lab ID:	L2027725-08		Date Collected:	06/30/20 13:05
Client ID:	EB01-200630		Date Received:	06/30/20
Sample Location:	SCHENECTADY NY		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method	: EPA 3510C
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 18:15
Analytical Date:	07/02/20 21:55			
Analyst:	DV			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		ug/l	0.10	0.01	1	
2-Chloronaphthalene	ND		ug/l	0.20	0.02	1	
Fluoranthene	0.02	J	ug/l	0.10	0.02	1	
Hexachlorobutadiene	ND		ug/l	0.50	0.05	1	
Naphthalene	ND		ug/l	0.10	0.05	1	
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1	
Benzo(a)pyrene	ND		ug/l	0.10	0.02	1	
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	1	
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	1	
Chrysene	ND		ug/l	0.10	0.01	1	
Acenaphthylene	ND		ug/l	0.10	0.01	1	
Anthracene	0.02	J	ug/l	0.10	0.01	1	
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	1	
Fluorene	0.02	J	ug/l	0.10	0.01	1	
Phenanthrene	0.03	J	ug/l	0.10	0.02	1	
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	1	
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	1	
Pyrene	ND		ug/l	0.10	0.02	1	
2-Methylnaphthalene	ND		ug/l	0.10	0.02	1	
Pentachlorophenol	ND		ug/l	0.80	0.01	1	
Hexachlorobenzene	0.03	J	ug/l	0.80	0.01	1	
Hexachloroethane	ND		ug/l	0.80	0.06	1	



Parameter		Result	Qualifier	Units	RL MDL	Dilution Factor		
Sample Depth:								
Sample Location:	SCHENECTADY NY				Field Prep:	Not Specified		
Client ID:	EB01-200630				Date Received:	06/30/20		
Lab ID:	L2027725-08				Date Collected:	06/30/20 13:05		
SAMPLE RESULTS								
Project Number:	16.6334				Report Date:	07/08/20		
Project Name:	HHII TA1				Lab Number:	L2027725		
					Serial_No:07082013:26			

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	67	21-120
Phenol-d6	61	10-120
Nitrobenzene-d5	97	23-120
2-Fluorobiphenyl	86	15-120
2,4,6-Tribromophenol	118	10-120
4-Terphenyl-d14	122	41-149



Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Method Blank Analysis Batch Quality Control

Analytical Method:	1,
Analytical Date:	07
Analyst:	S

1,8270D 07/02/20 15:07 SZ Extraction Method: EPA 3510C Extraction Date: 07/01/20 18:11

arameter	Result	Qualifier Units	RL		MDL
emivolatile Organics by GC/MS	- Westborough	Lab for sample(s):	01-08	Batch:	WG1388229-1
Bis(2-chloroethyl)ether	ND	ug/l	2.0		0.50
3,3'-Dichlorobenzidine	ND	ug/l	5.0		1.6
2,4-Dinitrotoluene	ND	ug/l	5.0		1.2
2,6-Dinitrotoluene	ND	ug/l	5.0		0.93
4-Chlorophenyl phenyl ether	ND	ug/l	2.0		0.49
4-Bromophenyl phenyl ether	ND	ug/l	2.0		0.38
Bis(2-chloroisopropyl)ether	ND	ug/l	2.0		0.53
Bis(2-chloroethoxy)methane	ND	ug/l	5.0		0.50
Hexachlorocyclopentadiene	ND	ug/l	20		0.69
Isophorone	ND	ug/l	5.0		1.2
Nitrobenzene	ND	ug/l	2.0		0.77
NDPA/DPA	ND	ug/l	2.0		0.42
n-Nitrosodi-n-propylamine	ND	ug/l	5.0		0.64
Bis(2-ethylhexyl)phthalate	ND	ug/l	3.0		1.5
Butyl benzyl phthalate	ND	ug/l	5.0		1.2
Di-n-butylphthalate	ND	ug/l	5.0		0.39
Di-n-octylphthalate	ND	ug/l	5.0		1.3
Diethyl phthalate	ND	ug/l	5.0		0.38
Dimethyl phthalate	ND	ug/l	5.0		1.8
Biphenyl	ND	ug/l	2.0		0.46
4-Chloroaniline	ND	ug/l	5.0		1.1
2-Nitroaniline	ND	ug/l	5.0		0.50
3-Nitroaniline	ND	ug/l	5.0		0.81
4-Nitroaniline	ND	ug/l	5.0		0.80
Dibenzofuran	ND	ug/l	2.0		0.50
1,2,4,5-Tetrachlorobenzene	ND	ug/l	10		0.44
Acetophenone	ND	ug/l	5.0		0.53
2,4,6-Trichlorophenol	ND	ug/l	5.0		0.61
p-Chloro-m-cresol	ND	ug/l	2.0		0.35



Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Method Blank Analysis Batch Quality Control

Analytical Method:1,8Analytical Date:07Analyst:SZ

1,8270D 07/02/20 15:07 SZ Extraction Method: EPA 3510C Extraction Date: 07/01/20 18:11

arameter	Result	Qualifier	Units	RL		MDL
emivolatile Organics by GC/MS	6 - Westborough	h Lab for sa	ample(s):	01-08	Batch:	WG1388229-1
2-Chlorophenol	ND		ug/l	2.0		0.48
2,4-Dichlorophenol	ND		ug/l	5.0		0.41
2,4-Dimethylphenol	ND		ug/l	5.0		1.8
2-Nitrophenol	ND		ug/l	10		0.85
4-Nitrophenol	ND		ug/l	10		0.67
2,4-Dinitrophenol	ND		ug/l	20		6.6
4,6-Dinitro-o-cresol	ND		ug/l	10		1.8
Phenol	ND		ug/l	5.0		0.57
3-Methylphenol/4-Methylphenol	ND		ug/l	5.0		0.48
2,4,5-Trichlorophenol	ND		ug/l	5.0		0.77
Carbazole	ND		ug/l	2.0		0.49
Atrazine	ND		ug/l	10		0.76
Benzaldehyde	ND		ug/l	5.0		0.53
Caprolactam	ND		ug/l	10		3.3
2,3,4,6-Tetrachlorophenol	ND		ug/l	5.0		0.84

Tentatively Identified Compounds							
Total TIC Compounds	5.27	J	ug/l				
Unknown	3.09		-				
		J	ug/l				
Unknown	2.18	J	ug/l				



Serial_No:07082013:26

Project Name: Project Number:	HHII TA1 16.6334		Lab Number: Report Date:	L2027725 07/08/20
		Method Blank Analysis Batch Quality Control		
Analytical Method: Analytical Date: Analyst:	1,8270D 07/02/20 15:07 SZ		Extraction Method: Extraction Date:	EPA 3510C 07/01/20 18:11

Parameter	Result	Qualifier	Units	RL		MDL
Semivolatile Organics by GC/MS -	Westboroug	h Lab for s	ample(s):	01-08	Batch:	WG1388229-1

Surrogate	%Recovery Qua	Acceptance lifier Criteria
2-Fluorophenol	64	21-120
Phenol-d6	51	10-120
Nitrobenzene-d5	84	23-120
2-Fluorobiphenyl	62	15-120
2,4,6-Tribromophenol	64	10-120
4-Terphenyl-d14	78	41-149



 Lab Number:
 L2027725

 Report Date:
 07/08/20

Method Blank Analysis Batch Quality Control

Analytical Method:	1,8270D-SIM
Analytical Date:	07/02/20 18:37
Analyst:	DV

Extraction Method:EPA 3510CExtraction Date:07/01/20 18:15

arameter	Result	Qualifier	Units	RL	MDL	
emivolatile Organics by	GC/MS-SIM - Westbo	rough Lab	for sample(s)	01-08	Batch:	WG1388230-1
Acenaphthene	ND		ug/l	0.10	0.01	
2-Chloronaphthalene	ND		ug/l	0.20	0.02	
Fluoranthene	ND		ug/l	0.10	0.02	
Hexachlorobutadiene	ND		ug/l	0.50	0.05	
Naphthalene	0.28		ug/l	0.10	0.05	
Benzo(a)anthracene	ND		ug/l	0.10	0.02	
Benzo(a)pyrene	ND		ug/l	0.10	0.02	
Benzo(b)fluoranthene	ND		ug/l	0.10	0.01	
Benzo(k)fluoranthene	ND		ug/l	0.10	0.01	
Chrysene	ND		ug/l	0.10	0.01	
Acenaphthylene	0.02	J	ug/l	0.10	0.01	
Anthracene	ND		ug/l	0.10	0.01	
Benzo(ghi)perylene	ND		ug/l	0.10	0.01	
Fluorene	0.02	J	ug/l	0.10	0.01	
Phenanthrene	0.05	J	ug/l	0.10	0.02	
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.01	
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.01	
Pyrene	0.02	J	ug/l	0.10	0.02	
2-Methylnaphthalene	0.02	J	ug/l	0.10	0.02	
Pentachlorophenol	ND		ug/l	0.80	0.01	
Hexachlorobenzene	ND		ug/l	0.80	0.01	
Hexachloroethane	ND		ug/l	0.80	0.06	



Serial_No:07082013:26

Project Name: Project Number:	HHII TA1 16.6334		Lab Number: Report Date:	L2027725 07/08/20
		Method Blank Analysis Batch Quality Control		
Analytical Method: Analytical Date: Analyst:	1,8270D-SIM 07/02/20 18:37 DV		Extraction Method: Extraction Date:	EPA 3510C 07/01/20 18:15

Parameter	Result	Qualifier	Units	RL	MDL	
Semivolatile Organics by GC/MS-S	IM - Westbo	orough Lab	for sample	(s): 01-08	Batch: WG138823	0-1

Surrogate	%Recovery	Qualifier	Acceptance Criteria
2-Fluorophenol	72		21-120
Phenol-d6	62		10-120
Nitrobenzene-d5	103		23-120
2-Fluorobiphenyl	86		15-120
2,4,6-Tribromophenol	126	Q	10-120
4-Terphenyl-d14	123		41-149



Project Name: HHII TA1 Project Number: 16.6334

Lab Number: L2027725

Report Date: 07/08/20

	LCS	o <i>i</i>	LCSD		%Recovery		RPD
Parameter	%Recovery	Qual	%Recovery	Qual	Limits	RPD	Qual Limits
Semivolatile Organics by GC/MS - Westb	orough Lab Associ	ated sample(s):	01-08 Batch	: WG138822	9-2 WG138822	29-3	
Bis(2-chloroethyl)ether	67		65		40-140	3	30
3,3'-Dichlorobenzidine	50		47		40-140	6	30
2,4-Dinitrotoluene	69		68		48-143	1	30
2,6-Dinitrotoluene	64		62		40-140	3	30
4-Chlorophenyl phenyl ether	59		58		40-140	2	30
4-Bromophenyl phenyl ether	56		52		40-140	7	30
Bis(2-chloroisopropyl)ether	80		77		40-140	4	30
Bis(2-chloroethoxy)methane	62		61		40-140	2	30
Hexachlorocyclopentadiene	53		54		40-140	2	30
Isophorone	66		64		40-140	3	30
Nitrobenzene	77		75		40-140	3	30
NDPA/DPA	59		57		40-140	3	30
n-Nitrosodi-n-propylamine	73		71		29-132	3	30
Bis(2-ethylhexyl)phthalate	70		68		40-140	3	30
Butyl benzyl phthalate	69		63		40-140	9	30
Di-n-butylphthalate	61		58		40-140	5	30
Di-n-octylphthalate	76		72		40-140	5	30
Diethyl phthalate	64		62		40-140	3	30
Dimethyl phthalate	57		54		40-140	5	30
Biphenyl	62		61		40-140	2	30
4-Chloroaniline	35	Q	36	Q	40-140	3	30
2-Nitroaniline	65		61		52-143	6	30
3-Nitroaniline	51		50		25-145	2	30



Lab Number: L2027725

Project Number: 16.6334

HHII TA1

Project Name:

Report Date: 07/08/20

arameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	RPD Qual Limits
emivolatile Organics by GC/MS - We	estborough Lab Associa	ated sample(s)	: 01-08 Batch	n: WG138	8229-2 WG13882	29-3	
4-Nitroaniline	48	Q	47	Q	51-143	2	30
Dibenzofuran	61		59		40-140	3	30
1,2,4,5-Tetrachlorobenzene	57		57		2-134	0	30
Acetophenone	66		64		39-129	3	30
2,4,6-Trichlorophenol	58		56		30-130	4	30
p-Chloro-m-cresol	63		60		23-97	5	30
2-Chlorophenol	67		64		27-123	5	30
2,4-Dichlorophenol	65		62		30-130	5	30
2,4-Dimethylphenol	48		44		30-130	9	30
2-Nitrophenol	82		79		30-130	4	30
4-Nitrophenol	45		43		10-80	5	30
2,4-Dinitrophenol	75		72		20-130	4	30
4,6-Dinitro-o-cresol	78		77		20-164	1	30
Phenol	50		48		12-110	4	30
3-Methylphenol/4-Methylphenol	66		64		30-130	3	30
2,4,5-Trichlorophenol	60		56		30-130	7	30
Carbazole	59		57		55-144	3	30
Atrazine	63		60		40-140	5	30
Benzaldehyde	61		62		40-140	2	30
Caprolactam	38		37		10-130	3	30
2,3,4,6-Tetrachlorophenol	59		56		40-140	5	30



Project Name:HHII TA1Project Number:16.6334

Lab Number: L2027725

Report Date: 07/08/20

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Semivolatile Organics by GC/MS - Westbo	rough Lab Associa	ated sample(s	s): 01-08 Batch	: WG138	8229-2 WG13882	29-3			

Surrogate	LCS %Recovery Qua	LCSD al %Recovery Qual	Acceptance Criteria
2-Fluorophenol	60	59	21-120
Phenol-d6	52	52	10-120
Nitrobenzene-d5	84	85	23-120
2-Fluorobiphenyl	55	55	15-120
2,4,6-Tribromophenol	59	57	10-120
4-Terphenyl-d14	58	55	41-149



Project Name: HHII TA1 Project Number: 16.6334

Report Date: 07/08/20

Parameter	LCS %Recovery	LCSD Qual %Recovery	%Recovery Qual Limits	RPD	RPD Qual Limits
Semivolatile Organics by GC/MS-SIM - Wes	stborough Lab As	sociated sample(s): 01-08	Batch: WG1388230-2 W	G1388230-3	
Acenaphthene	70	77	40-140	10	40
2-Chloronaphthalene	66	72	40-140	9	40
Fluoranthene	86	92	40-140	7	40
Hexachlorobutadiene	59	64	40-140	8	40
Naphthalene	61	67	40-140	9	40
Benzo(a)anthracene	87	89	40-140	2	40
Benzo(a)pyrene	92	100	40-140	8	40
Benzo(b)fluoranthene	86	92	40-140	7	40
Benzo(k)fluoranthene	82	90	40-140	9	40
Chrysene	76	84	40-140	10	40
Acenaphthylene	81	89	40-140	9	40
Anthracene	83	90	40-140	8	40
Benzo(ghi)perylene	91	98	40-140	7	40
Fluorene	74	80	40-140	8	40
Phenanthrene	71	78	40-140	9	40
Dibenzo(a,h)anthracene	107	112	40-140	5	40
Indeno(1,2,3-cd)pyrene	108	115	40-140	6	40
Pyrene	86	93	40-140	8	40
2-Methylnaphthalene	66	72	40-140	9	40
Pentachlorophenol	97	106	40-140	9	40
Hexachlorobenzene	67	73	40-140	9	40
Hexachloroethane	61	66	40-140	8	40



Project Name:HHII TA1Project Number:16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

 LCS
 LCSD
 %Recovery
 RPD

 Parameter
 %Recovery
 Qual
 Value
 Limits
 RPD
 Qual
 Limits

 Semivolatile Organics by GC/MS-SIM - Westborough Lab
 Associated sample(s):
 01-08
 Batch:
 WG1388230-2
 WG1388230-3

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
2-Fluorophenol	58	64	21-120
Phenol-d6	52	57	10-120
Nitrobenzene-d5	80	88	23-120
2-Fluorobiphenyl	69	75	15-120
2,4,6-Tribromophenol	104	111	10-120
4-Terphenyl-d14	102	109	41-149



Matrix Spike Analysis Batch Quality Control

		Batch	Qu

Project Name:HHII TA1Project Number:16.6334

_

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery		Recovery Limits	RPD	Qual	RPD Limits
Semivolatile Organics by GC/M ID: RAMW5-200630	IS - Westbor	ough Lab	Associated sar	mple(s): 01-08	QC Batch	n ID: WG1	388229-4 WG	31388229	9-5 QC Sa	imple: L	202772	5-01 Client
Bis(2-chloroethyl)ether	ND	18.2	11	61		15	83		40-140	31	Q	30
3,3'-Dichlorobenzidine	ND	18.2	8.3	46		11	61		40-140	28		30
2,4-Dinitrotoluene	ND	18.2	13	72		19	100		48-143	38	Q	30
2,6-Dinitrotoluene	ND	18.2	12	66		18	99		40-140	40	Q	30
4-Chlorophenyl phenyl ether	ND	18.2	10	55		15	83		40-140	40	Q	30
4-Bromophenyl phenyl ether	ND	18.2	9.6	53		14	77		40-140	37	Q	30
Bis(2-chloroisopropyl)ether	ND	18.2	14	77		20	110		40-140	35	Q	30
Bis(2-chloroethoxy)methane	ND	18.2	11	61		16	88		40-140	37	Q	30
Hexachlorocyclopentadiene	ND	18.2	9.6J	53		14.J	77		40-140	37	Q	30
Isophorone	ND	18.2	12	66		17	94		40-140	34	Q	30
Nitrobenzene	ND	18.2	14	77		19	100		40-140	30		30
NDPA/DPA	ND	18.2	10	55		16	88		40-140	46	Q	30
n-Nitrosodi-n-propylamine	ND	18.2	13	72		18	99		29-132	32	Q	30
Bis(2-ethylhexyl)phthalate	ND	18.2	13	72		20	110		40-140	42	Q	30
Butyl benzyl phthalate	ND	18.2	12	66		19	100		40-140	45	Q	30
Di-n-butylphthalate	ND	18.2	11	61		17	94		40-140	43	Q	30
Di-n-octylphthalate	ND	18.2	14	77		20	110		40-140	35	Q	30
Diethyl phthalate	ND	18.2	12	66		18	99		40-140	40	Q	30
Dimethyl phthalate	ND	18.2	10	55		15	83		40-140	40	Q	30
Biphenyl	ND	18.2	11	61		16	88		40-140	37	Q	30
4-Chloroaniline	ND	18.2	6.6	36	Q	8.2	45		40-140	22		30
2-Nitroaniline	ND	18.2	12	66		18	99		52-143	40	Q	30
3-Nitroaniline	ND	18.2	9.6	53		13	72		25-145	30		30



Matrix Spike Analysis Batch Quality Control

HHII TA1		-

Project Number: 16.6334

Project Name:

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery		Recovery Limits	RPD	Qual	RPD Limits
Semivolatile Organics by GC ID: RAMW5-200630	C/MS - Westbor	ough Lab	Associated sa	mple(s): 01-08	QC Batc	h ID: WG1	1388229-4 WG	138822	29-5 QC Sa	ample: L	202772	5-01 Client
4-Nitroaniline	ND	18.2	9.1	50	Q	13	72		51-143	35	Q	30
Dibenzofuran	ND	18.2	10	55		15	83		40-140	40	Q	30
1,2,4,5-Tetrachlorobenzene	ND	18.2	10	55		14	77		2-134	33	Q	30
Acetophenone	ND	18.2	12	66		17	94		39-129	34	Q	30
2,4,6-Trichlorophenol	ND	18.2	10	55		15	83		30-130	40	Q	30
p-Chloro-m-cresol	ND	18.2	11	61		17	94		23-97	43	Q	30
2-Chlorophenol	ND	18.2	12	66		16	88		27-123	29		30
2,4-Dichlorophenol	ND	18.2	12	66		17	94		30-130	34	Q	30
2,4-Dimethylphenol	ND	18.2	9.4	52		13	72		30-130	32	Q	30
2-Nitrophenol	ND	18.2	14	77		21	120		30-130	40	Q	30
4-Nitrophenol	ND	18.2	11	61		19	100	Q	10-80	53	Q	30
2,4-Dinitrophenol	ND	18.2	15.J	83		22	120		20-130	38	Q	30
4,6-Dinitro-o-cresol	ND	18.2	14	77		21	120		20-164	40	Q	30
Phenol	ND	18.2	9.2	51		13	72		12-110	34	Q	30
3-Methylphenol/4-Methylphenol	ND	18.2	12	66		17	94		30-130	34	Q	30
2,4,5-Trichlorophenol	ND	18.2	11	61		16	88		30-130	37	Q	30
Carbazole	ND	18.2	10	55		16	88		55-144	46	Q	30
Atrazine	ND	18.2	11	61		17	94		40-140	43	Q	30
Benzaldehyde	ND	18.2	11	61		16	88		40-140	37	Q	30
Caprolactam	ND	18.2	7.4J	41		10	55		10-130	30		30
2,3,4,6-Tetrachlorophenol	ND	18.2	10	55		16	88		40-140	46	Q	30



Project Name:HHII TA1Matrix Spike Analysis
Batch Quality ControlLab Number:L2027725Project Number:16.63347/08/2007/08/20

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limit	
Semivolatile Organics by GC/M ID: RAMW5-200630	IS - Westbor	ough Lab	Associated sa	imple(s): 01-08	QC Bate	ch ID: WG1	1388229-4 WG	138822	29-5 QC Sa	ample: L	202772	5-01	Client
				MS			MSD		Accept	ance			

		liide	Acceptance	
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria	
2,4,6-Tribromophenol	54	85	10-120	
2-Fluorobiphenyl	54	79	15-120	
2-Fluorophenol	57	81	21-120	
4-Terphenyl-d14	51	85	41-149	
Nitrobenzene-d5	83	119	23-120	
Phenol-d6	52	74	10-120	



Matrix Spike Analysis

Project Name:	HHII TA1	Batch Quality Control	Lab Nu	ımber:	
			_	_	

Project Number: 16.6334

 Lab Number:
 L2027725

 Report Date:
 07/08/20

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	MSD Qual Found	MSD %Recovery	Recover Qual Limits	/ RPD	RPD Qual Limits
Semivolatile Organics by Client ID: RAMW5-20063		stborough Lab	Associate	ed sample(s): 01-	08 QC Batch ID:	WG1388230-4	WG1388230-5	QC Sam	ple: L2027725-01
Acenaphthene	ND	18.2	10	55	13	72	40-140	26	40
2-Chloronaphthalene	ND	18.2	12	66	14	77	40-140	15	40
Fluoranthene	ND	18.2	12	66	15	83	40-140	22	40
Hexachlorobutadiene	ND	18.2	10	55	12	66	40-140	18	40
Naphthalene	ND	18.2	10	55	12	66	40-140	18	40
Benzo(a)anthracene	ND	18.2	12	66	15	83	40-140	22	40
Benzo(a)pyrene	ND	18.2	12	66	16	88	40-140	29	40
Benzo(b)fluoranthene	ND	18.2	11	61	14	77	40-140	24	40
Benzo(k)fluoranthene	ND	18.2	11	61	15	83	40-140	31	40
Chrysene	ND	18.2	9.7	53	13	72	40-140	29	40
Acenaphthylene	ND	18.2	14	77	16	88	40-140	13	40
Anthracene	ND	18.2	12	66	15	83	40-140	22	40
Benzo(ghi)perylene	ND	18.2	11	61	14	77	40-140	24	40
Fluorene	ND	18.2	11	61	14	77	40-140	24	40
Phenanthrene	ND	18.2	9.6	53	12	66	40-140	22	40
Dibenzo(a,h)anthracene	ND	18.2	13	72	19	100	40-140	38	40
Indeno(1,2,3-cd)pyrene	ND	18.2	14	77	17	94	40-140	19	40
Pyrene	ND	18.2	12	66	15	83	40-140	22	40
2-Methylnaphthalene	ND	18.2	11	61	14	77	40-140	24	40
Pentachlorophenol	ND	18.2	12	66	16	88	40-140	29	40
Hexachlorobenzene	ND	18.2	9.7	53	12	66	40-140	21	40
Hexachloroethane	ND	18.2	11	61	13	72	40-140	17	40



		Matrix Spike Analysis Batch Quality Control		
Project Name:	HHII TA1	Lak	Number:	L2027725
Project Number:	16.6334	Rej	oort Date:	07/08/20

	Native	MS	MS	MS		MSD	MSD		Recovery			RPD
Parameter	Sample	Added	Found	%Recovery	Qual	Found	%Recovery	Qual	Limits	RPD	Qual	Limits
Semivolatile Organics by GC Client ID: RAMW5-200630	/MS-SIM - Wes	stborough Lat	Associate	ed sample(s): 01-	08 QC	Batch ID:	WG1388230-4	WG13	88230-5 C	C Samp	le: L202	27725-01

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
2,4,6-Tribromophenol	88	116	10-120
2-Fluorobiphenyl	63	77	15-120
2-Fluorophenol	57	67	21-120
4-Terphenyl-d14	71	94	41-149
Nitrobenzene-d5	79	94	23-120
Phenol-d6	54	64	10-120



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Sample Receipt and Container Information

Were project specific reporting limits specified?

YES

Cooler Information

Cooler	Custody Seal				
A	Absent				

Container Info	rmation		Initial	Final	Temp			Frozen	
Container ID	Container Type	Cooler	рН	рН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2027725-01A	Vial HCl preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01B	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01D	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01E	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01F	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01G	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01H	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01J	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-01K	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-01L	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-01M	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-01N	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-01P	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-01Q	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-02A	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-02B	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-02C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-02D	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-02E	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2027725-03A	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-03B	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2027725-03C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)

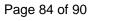


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Container Info	iner Information		Initial	Final	Temp			Frozen		
Container ID	Container Type	Cooler	рН	рН	-	Pres	Seal	Date/Time	Analysis(*)	
L2027725-03D	Amber 250ml unpreserved	A	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-03E	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-04A	Vial HCl preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-04B	Vial HCl preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-04C	Vial HCl preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-04D	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-04E	Amber 250ml unpreserved	A	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-05A	Vial HCI preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-05B	Vial HCI preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-05C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-05D	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-05E	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-06A	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-06B	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-06C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-06D	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-06E	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-07A	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-07B	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-07C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-07D	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-07E	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-08A	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-08B	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-08C	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	
L2027725-08D	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-08E	Amber 250ml unpreserved	А	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)	
L2027725-09A	Vial HCI preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)	



Container Info	ormation		Initial	Final	Temp			Frozen	
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2027725-09B	Vial HCl preserved	А	NA		3.4	Y	Absent		NYTCL-8260-R2(14)





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Project Number: 16.6334

Lab Number: L2027725

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GLOSSARY

Acronyms

Acronyms	
DL	- Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)
EDL	- Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).
EMPC	- Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.
EPA	- Environmental Protection Agency.
LCS	- Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
LCSD	- Laboratory Control Sample Duplicate: Refer to LCS.
LFB	- Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
LOD	- Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)
LOQ	- Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)
	Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)
MDL	- Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
MS	- Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.
MSD	- Matrix Spike Sample Duplicate: Refer to MS.
NA	- Not Applicable.
NC	 Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.
NDPA/DPA NI	- N-Nitrosodiphenylamine/Diphenylamine.
	- Not Ignitable.
NP RL	 Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil. Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL
KL.	includes any adjustments from dilutions, concentrations or moisture content, where applicable.
RPD	- Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.
SRM	- Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.
STLP	- Semi-dynamic Tank Leaching Procedure per EPA Method 1315.
TEF	- Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.
TEQ	- Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.
TIC	- Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.
Footnotes	

Report Format: DU Report with 'J' Qualifiers



Project Name:	HHII TA1	Lab Number:	L2027725
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- The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum. Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Waterpreserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'. Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. If a 'Total' result is requested, the results of its individual components will also be reported.

The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA,this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

- A Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. ND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- **D** Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- J Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- ${f P}$ The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration

Report Format: DU Report with 'J' Qualifiers



Serial_No:07082013:26

Project Name:	HHII TA1	Lab Number:	L2027725
Project Number:	16.6334	Report Date:	07/08/20

Data Qualifiers

Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)

- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- **S** Analytical results are from modified screening analysis.

Report Format: DU Report with 'J' Qualifiers



 Lab Number:
 L2027725

 Report Date:
 07/08/20

REFERENCES

1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

LIMITATION OF LIABILITIES

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

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



Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene
EPA 8260C: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.
EPA 8270D: NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.
SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.
Mansfield Facility
SM 2540D: TSS
EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.
EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene, 1-Methylnaphthalene.
SPA 3C Fixed gases
Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP. Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs **EPA 625.1**: SVOC (Acid/Base/Neutral Extractables), **EPA 600/4-81-045**: PCB-Oil.

Microbiology SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn. **EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn. **EPA 245.1** Hg. **SM2340B**

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Serial_No:07082013:26

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C.T. MALE ASSOCIATES

ATTACHMENT G

DATA USABILITY SUMMARY REPORT

DATA USABILITY SUMMARY REPORT HAMILTON HILL, SCHENECTADY, NEW YORK

Client:	C.T. Male Associates, Latham, New York
SDG:	L2027725
Laboratory:	Alpha Analytical, Westborough, Massachusetts
Site:	Hamilton Hill II TA-1, Schenectady, New York
Date:	June 14, 2020

EDS ID	Client Sample ID	Laboratory Sample ID	Matrix
1	RAMW5-200630	L2027725-01	Water
1MS**	RAMW5-200630MS	L2027725-01MS	Water
1MSD**	RAMW5-200630MSD	L2027725-01MSD	Water
2	RAMW6-200630	L2027725-02	Water
3	RAMW4-200630	L2027725-03	Water
4	RAMW1-200630	L2027725-04	Water
5	RAMW2-200630	L2027725-05	Water
6	RAMW3-200630	L2027725-06	Water
7	FD01-200630	L2027725-07	Water
8	EB01-200630	L2027725-08	Water
9*	TRIP BLANK-200630	L2027725-09	Water

* - VOC only ** - VOC & SVOC only

A Data Usability Summary Review was performed on the analytical data for seven water samples, one aqueous equipment blank sample, and one aqueous trip blank sample collected on June 30, 2020 by CT Male at the Hamilton Hill II-TA1 site in Schenectady, New York. The samples were analyzed under Environmental Protection Agency (USEPA) "Test Methods for the Evaluation of Solid Waste, USEPA SW-846, Third Edition, September 1986, with revisions".

Specific method references are as follows:

<u>Analysis</u>	Method References
VOCs	USEPA SW-846 Method 8260C
SVOCs	USEPA SW-846 Method 8270D
РАН	USEPA SW-846 Method 8270D-SIM

The data have been validated according to the protocols and quality control (QC) requirements of the analytical methods and the USEPA Region II Data Review Standard Operating Procedures (SOPs) as follows:

- SOP Number HW-33A, Revision 1, September 2016: Low/Medium Volatile Data Validation;
- SOP Number HW-35A, Revision 1, September 2016: Semivolatile Data Validation;
- and the reviewer's professional judgment.

The following items/criteria were reviewed for this report:

Organics

- Data Completeness
- Holding times and sample preservation
- Surrogate Spike recoveries
- Matrix Spike/Matrix Spike Duplicate (MS/MSD) recoveries
- Laboratory Control Sample (LCS) recoveries
- Method blank and field blank contamination
- Gas Chromatography (GC)/Mass Spectroscopy (MS) tuning
- Initial and continuing calibration summaries
- Compound Quantitation
- Internal standard area and retention time summary forms
- Tentatively Identified Compounds (TICs)
- Field Duplicate sample precision

Data Usability Assessment

There were no rejections of data.

The data are acceptable for the intended purposes as qualified for the deficiencies detailed in this report.

Please note that any results qualified (U) due to blank contamination may be then qualified (J) due to another action. Therefore, the results may be qualified (UJ) due to the culmination of the blank contaminations and actions from other exceedances of QC criteria.

Data Completeness

• The data is a complete Category B data package as defined under the requirements for the NYS Department of Environmental Conservation Analytical Services Protocol.

Volatile Organic Compounds (VOCs)

Holding Times

• All samples were analyzed within 14 days for preserved water samples.

Surrogate Spike Recoveries

• All samples exhibited acceptable surrogate %R values.

Matrix Spike/Matrix Spike Duplicate (MS/MSD) Recoveries

• The MS/MSD samples that exhibited percent recoveries (%R) outside the QC limits and/or relative percent differences (RPD) above QC limits. A low %R may indicate a potential low bias while a high %R may indicate a potential high bias. For a low %R, positive results are considered estimated and qualified (J) while non-detects are estimated and qualified (UJ). For a high %R, positive results are considered estimated and qualified and qualified (J).

MS/MSD Sample ID	Compound	MS %R/MSD %R/RPD	Qualifier
1	1,2,3-Trichlorobenzene	69%/OK/33	UJ
	1,2,4-Trichlorobenzene	68%/OK/37	UJ
	28 Compounds	OK/OK/High	None for RPD alone

Laboratory Control Samples

• The LCS samples exhibited acceptable %R values.

Method Blank

• The method blanks were free of contamination.

Field Blank

• The following table lists field QC samples with contamination and the samples associated with the blanks that had results qualified as a consequence of the blank contamination. Detected sample concentrations of acetone, 2-butanone and methylene chloride (common laboratory contaminants) less than ten times (10x) the highest associated blank (after taking sample dilution levels, percent moisture and sample volume into account) are negated and qualified with a (U). For all other compounds, an action level of five times (5x) the highest associated blank concentration is used.

Blank ID	Compound	Conc. ug/L	Qualifier	Affected Samples
EB01-200630	Acetone	3.8	None	All Associated ND
TRIP BLANK-200630	None - ND	2.6	2.05	-

GC/MS Tuning

• All criteria were met.

Initial Calibration

• The initial calibrations exhibited acceptable %RSD and/or correlation coefficients and mean RRF values.

Continuing Calibration

• The following table presents compounds that exceeded percent difference (%D) and/or RRF values <0.05 (0.01 for poor performers) in the continuing calibration (CCAL). A low RRF indicates poor instrument sensitivity for these compounds. Positive results for these compounds in the affected samples are considered estimated and qualified (J). Non-detect results for these compounds in the affected samples are rejected (R) and are unusable for project objectives. A high %D may indicate a potential high or low bias. All results for these compounds in affected samples are considered estimated and qualified (J/UJ).

CCAL Date	Compound	%D/RRF	Qualifier	Affected Samples
07/05/20 (1316)	Bromomethane	36.8%	UJ	All Samples

Compound Quantitation

• All criteria were met.

Internal Standard (IS) Area Performance

• All internal standards met response and retention time (RT) criteria.

Tentatively Identified Compounds (TICs)

• TICs were not detected.

Field Duplicate Sample Precision

• Field duplicate results are summarized below. The precision was acceptable.

Compound	RAMW6-200630 ug/L	FD01-200630 ug/L	RPD	Qualifier
None	ND	ND		

Semivolatile Organic Compounds (SVOCs)

Holding Times

• All samples were extracted within 7 days for water samples and analyzed within 40 days for all samples.

Surrogate Spike Recoveries

• All samples exhibited acceptable surrogate %R values.

Matrix Spike/Matrix Spike Duplicate (MS/MSD) Recoveries

• The following table presents MS/MSD samples that exhibited percent recoveries (%R) outside the QC limits and/or relative percent differences (RPD) above QC limits. A low %R may indicate a potential low bias while a high %R may indicate a potential high bias. For a low %R, positive results are considered estimated and qualified (J) while non-detects are estimated and qualified (UJ). For a high %R, positive results are considered estimated and qualified (J).

MS/MSD Sample ID	Compound	MS %R/MSD %R/ RPD	Qualifier
1	Most Compounds	OK/OK/High	None for RPD alone
	4-Chloroaniline	36%/OK/OK	None - See LCS
	4-Nitroaniline	50%/OK/35	None - See LCS
	4-Nitrophenol	OK/100%/53	None - Sample ND

Laboratory Control Samples

• The following table presents LCS percent recoveries (%R) outside the QC limits. A low %R may indicate a potential low bias while a high %R may indicate a potential high bias. For a low %R, positive results are considered estimated and qualified (J) while non-detects are estimated and qualified (UJ). For a high %R, positive results are considered estimated and qualified and qualified (J). Results are valid and usable, however possibly biased.

LCS ID	Compound	%R	Qualifier	Affected Samples
WG1388229-2	4-Chloroaniline	35%	UJ	All Samples
	4-Nitroaniline	48%	UJ	

Method Blank

• The method blanks were free of contamination.

Field Blank

• The field QC results are summarized below.

Blank ID	Compound	Conc. ug/L	Qualifier	Affected Samples
EB01-200630	None - ND	1.21	-	

GC/MS Tuning

• All criteria were met.

Initial Calibration

• The initial calibrations exhibited acceptable %RSD and/or correlation coefficients and mean RRF values.

Continuing Calibration

• The following table presents compounds that exceeded percent difference (%D) and/or RRF values <0.05 in the continuing calibration (CCAL). A low RRF indicates poor instrument sensitivity for these compounds. Positive results for these compounds in the affected samples are considered estimated and qualified (J). Non-detect results for these compounds in the affected samples are rejected (R) and are unusable for project objectives. A high %D may indicate a potential high or low bias. All results for these compounds in affected samples are considered estimated and qualified (J/UJ).

CCAL Date	Compound	%D/RRF	Qualifier	Affected Samples
07/02/20 (1308)	Bis(2-chloroisopropyl)ether	21.8%	UJ	All Samples
	2-Nitrophenol	24.2%	UJ	-

Compound Quantitation

• All criteria were met.

Internal Standard (IS) Area Performance

• All internal standards met response and retention time (RT) criteria.

Tentatively Identified Compounds (TICs)

• All TICs were qualified estimated (J) for unknown compounds.

Field Duplicate Sample Precision

• Field duplicate results are summarized below. The precision was acceptable.

Compound	RAMW6-200630 ug/L	FD01-200630 ug/L	RPD	Qualifier
None	ND	ND	1.51	(.

Polycyclic Aromatic Hydrocarbons (PAHs)

Holding Times

• All samples were extracted within 7 days for water samples and analyzed within 40 days for all samples.

Surrogate Spike Recoveries

• The following table presents surrogate percent recoveries (%R) outside the QC limits. A low %R may indicate a potential low bias while a high %R may indicate a potential high bias. For a low %R, positive results are considered estimated and qualified (J) while non-detects are estimated and qualified (UJ). For a high %R, positive results are considered estimated and qualified (J).

EDS Sample ID	Surrogate	%R	Qualifier
3	2,4,6-Tribromophenol	123%	None for one out per fraction

Matrix Spike/Matrix Spike Duplicate (MS/MSD) Recoveries

• MS/MSD samples were not analyzed.

Laboratory Control Samples

• The LCS samples exhibited acceptable %R values.

Method Blank

• The following table lists method blanks with contamination and the samples associated with the blanks that had results qualified as a consequence of the blank contamination. Detected sample concentrations less than five times (5x) the highest associated blank (after taking sample dilution levels, percent moisture and sample volume into account) are negated and qualified with a (U).

Blank ID	Compound	Conc. ug/L	Qualifier	Affected Samples
WG1388230-1	Naphthalene	0.28	None	All Associated ND
	Acenaphthylene	0.02	None	
	Fluorene	0.02	U	8
	Phenanthrene	0.05	U	
	Pyrene	0.02	None	All Associated ND
	2-Methylnaphthalene	0.02	None	

Field Blank

The following table lists field QC samples with contamination and the samples associated • with the blanks that had results qualified as a consequence of the blank contamination. Detected sample concentrations less than five times (5x) the highest associated blank (after taking sample dilution levels, percent moisture and sample volume into account) are negated and qualified with a (U).

Blank ID	Compound	Conc. ug/L	Qualifier	Affected Samples
EB01-200630	Fluoranthene	0.02	None	All Associated ND
	Anthracene	0.02	None	
	Hexachlorobenzene	0.03	None	

Initial Calibration

All %RSD and/or correlation coefficient criteria were met. •

Continuing Calibration

All %D criteria were met. .

Compound Quantitation

All criteria were met.

Field Duplicate Sample Precision

• Field duplicate results are summarized below. The precision was acceptable.

Compound	RAMW6-200630 ug/L	FD01-200630 ug/L	RPD	Qualifier
None	ND	ND	-	

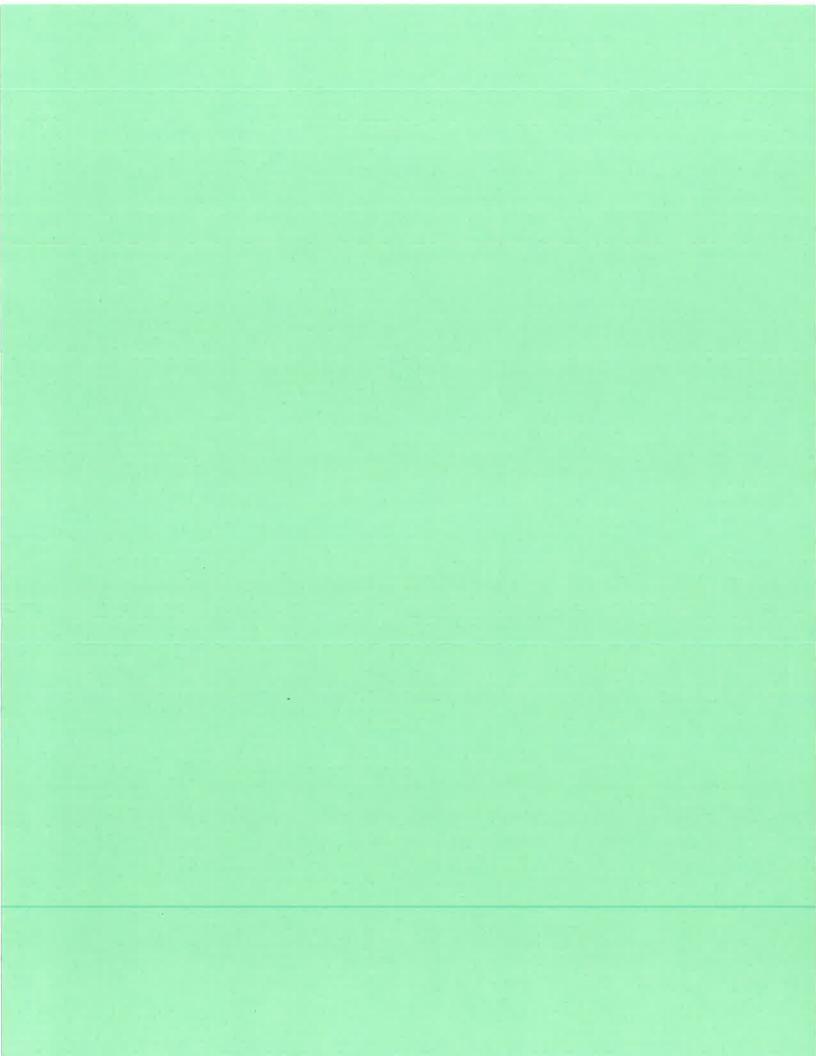
Please contact the undersigned at (757) 564-0090 if you have any questions or need further information.

Signed:

<u>Nancy Weaver</u> Dated: <u>7/17/20</u>

Senior Chemist

Data Qualifier	Definition
U	The analyte was analyzed for, but was not detected above the level of the reported sample quantitation limit.
J	The analyte is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
J+	The result is an estimated quantity, but the result may be biased high.
J-	The result is an estimated quantity, but the result may be biased low.
NJ	The analysis has been "tentatively identified" or "presumptively" as present and the associated numerical value is the estimated concentration in the samples.
UJ	The analyte was analyzed for but was not detected. The reported quantitation limits is approximate and may be inaccurate or imprecise.
R	The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the samples.



Results Summary Form 1 Volatile Organics by GC/MS

Client	: C.T. Male Associates	Lab Number : L2027725
Project Name	: HHII TA1	Project Number 16.6334
Lab ID	: L2027725-01	Date Collected : 06/30/20 08:00
Client ID	: RAMW5-200630	Date Received : 06/30/20
Sample Location	: SCHENECTADY NY	Date Analyzed : 07/05/20 16:14
Sample Matrix	: WATER	Dilution Factor 1
Analytical Method	: 1,8260C	Analyst NLK
Lab File ID	: VG200705A08	Instrument ID : GONZO
Sample Amount	: 10 ml	GC Column : RTX-502.2
Level	: LOW	%Solids : N/A
Extract Volume (MeO	H) : N/A	Injection Volume : N/A

	Parameter		ug/L		
CAS NO.		Results	RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	1.4	0.50	0.18	
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	NUS
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



M1 7/14/20

Results Summary Form 1 Volatile Organics by GC/MS

Client	: C.T. Male Associates	Lab Number : L2027725	
Project Name	: HHII TA1	Project Number : 16.6334	
Lab ID	: L2027725-01	Date Collected : 06/30/20 08:0	00
Client ID	RAMW5-200630	Date Received : 06/30/20	
Sample Location	SCHENECTADY NY	Date Analyzed : 07/05/20 16:	14
Sample Matrix	: WATER	Dilution Factor : 1	
Analytical Method	: 1,8260C	Analyst : NLK	
Lab File ID	; VG200705A08	Instrument ID : GONZO	
Sample Amount	: 10 ml	GC Column : RTX-502.2	
Level	: LOW	%Solids : N/A	
Extract Volume (MeO	H) 🗄 N/A	Injection Volume : N/A	

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U ⁵⁵
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	Jr US
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	N US
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U



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Client	: C.T. Male Associates		Lab Nun	nber	: L2027725
Project Name	: HHII TA1		Project N	lumber	: 16.6334
Lab ID	: L2027725-01		Date Co	llected	: 06/30/20 08:00
Client ID	: RAMW5-200630		Date Re	ceived	: 06/30/20
Sample Location	: SCHENECTADY NY		Date An	alyzed	07/05/20 16:14
Sample Matrix	: WATER		Dilution	Factor	±1
Analytical Method	: 1,8260C		Analyst		: NLK
Lab File ID	: VG200705A08		Instrume	ent ID	GONZO
Sample Amount	: 10 ml		GC Colu	ımn	: RTX-502.2
Level	: LOW		%Solids		: N/A
Extract Volume (M	leOH) : N/A		Injection	Volume	: N/A
			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U



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Client	: C.T. Male Associates	Lab Number : L2027725	
Project Name	: HHII TA1	Project Number : 16.6334	
Lab ID	: L2027725-02	Date Collected : 06/30/20 08:3	35
Client ID	: RAMW6-200630	Date Received : 06/30/20	
Sample Location	SCHENECTADY NY	Date Analyzed : 07/05/20 16:3	39
Sample Matrix	: WATER	Dilution Factor : 1	
Analytical Method	: 1,8260C	Analyst 📰 NLK	
Lab File ID	: VG200705A09	Instrument ID : GONZO	
Sample Amount	: 10 ml	GC Column : RTX-502.2	
Level	LOW	%Solids : N/A	
Extract Volume (MeOI	H) 🗄 N/A	Injection Volume : N/A	

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	201
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



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Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-02	Date Collected	: 06/30/20 08:35
Client ID	: RAMW6-200630	Date Received	: 06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed	: 07/05/20 16:39
Sample Matrix	: WATER	Dilution Factor	ž1
Analytical Method	: 1,8260C	Analyst	: NLK
Lab File ID	: VG200705A09	Instrument ID	: GONZO
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	🗄 N/A
Extract Volume (MeO	H) : N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether 🔗	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichlorœthene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U



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Client	: C.T. Male Associates		Lab Nun	nber	: L2027725
Project Name	: HHII TA1		Project I	Number	: 16.6334
Lab ID	: L2027725-02		Date Co	llected	: 06/30/20 08:35
Client ID	: RAMW6-200630		Date Re	ceived	: 06/30/20
Sample Location	: SCHENECTADY NY		Date An	alyzed	: 07/05/20 16:39
Sample Matrix	: WATER		Dilution	Factor	81
Analytical Method	: 1,8260C		Analyst		: NLK
Lab File ID	: VG200705A09		Instrume	ent ID	: GONZO
Sample Amount	: 10 ml		GC Colu	ımn	: RTX-502.2
Level	: LOW		%Solids		: N/A
Extract Volume (M	leOH):N/A		Injection	Volume	: N/A
			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U



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Client	 C.T. Male Associates HHII TA1 L2027725-03 RAMW4-200630 SCHENECTADY NY WATER 1,8260C VG200705A10 10 ml LOW 	Lab Number	: L2027725
Project Name		Project Number	: 16.6334
Lab ID		Date Collected	: 06/30/20 11:35
Client ID		Date Received	: 06/30/20
Sample Location		Date Analyzed	: 07/05/20 17:05
Sample Matrix		Dilution Factor	: 1
Analytical Method		Analyst	: NLK
Lab File ID		Instrument ID	: GONZO
Sample Amount		GC Column	: RTX-502.2
Level		%Solids	: N/A
Level Extract Volume (MeOH)		%Solids Injection Volume	

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	8.7	0.50	0.18	
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Тошепе	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	NUS
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



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Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-03	Date Collected	: 06/30/20 11:35
Client ID	: RAMW4-200630	Date Received	: 06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed	: 07/05/20 17:05
Sample Matrix	: WATER	Dilution Factor	1
Analytical Method	: 1,8260C	Analyst	: NLK
Lab File ID	: VG200705A10	Instrument ID	: GONZO
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH	l) : N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	0.59	0.50	0.18	
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	ម
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U



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Client	: C.T. Male Associates		Lab Nun	nber	: L2027725
Project Name	: HHII TA1		Project N	lumber	: 16.6334
Lab ID	L2027725-03		Date Co	llected	: 06/30/20 11:35
Client ID	: RAMW4-200630		Date Re	ceived	: 06/30/20
Sample Location	SCHENECTADY NY		Date Ana	alyzed	1 07/05/20 17:05
Sample Matrix	WATER		Dilution	Factor	÷ 1
Analytical Method	: 1,8260C		Analyst		: NLK
Lab File ID	: VG200705A10		Instrume	nt ID	: GONZO
Sample Amount	: 10 ml	(GC Colu	mn	: RTX-502.2
Level	: LOW		%Solids		: N/A
Extract Volume (N	leOH) : N/A	I	njection	Volume	: N/A
			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U



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Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-04	Date Collected	: 06/30/20 11:50
Client ID	: RAMW1-200630	Date Received	: 06/30/20
Sample Location	: SCHENECTADY NY	Date Analyzed	: 07/05/20 17:30
Sample Matrix	: WATER	Dilution Factor	: 1
Analytical Method	: 1,8260C	Analyst	: NLK
Lab File ID	: VG200705A11	Instrument ID	: GONZO
Sample Amount	: 10 mł	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	7.3	0.50	0.18	
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	y US
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



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Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	16.6334
Lab ID	: L2027725-04	Date Collected	: 06/30/20 11:50
Client ID	: RAMW1-200630	Date Received	: 06/30/20
Sample Location	: SCHENECTADY NY	Date Analyzed	: 07/05/20 17:30
Sample Matrix	: WATER	Dilution Factor	š 1
Analytical Method	: 1,8260C	Analyst	: NLK
Lab File ID	: VG200705A11	Instrument ID	: GONZO
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeO	H) : N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	3.2	0.50	0.18	
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
57-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	υ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	υ



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108-87-2	Methyl cyclohexane		ND	10	0.40	U
76-13-1	Freon-113		ND	2.5	0.70	U
CAS NO.	Parameter		Results	RL	MDL	Qualifier
				ug/L		
Extract Vo	blume (MeOH) : N/A			Injection	Volume	: N/A
Level	: LOW			%Solids		; N/A
Sample A				GC Colu	mn	: RTX-502.2
Lab File II	C : VG2007054	A11		Instrume	nt ID	GONZO
Analytical	Method : 1,8260C			Analyst		: NLK
Sample M	atrix 🗄 WATER			Dilution I	Factor	1
Sample L	ocation : SCHENEC	TADY NY		Date Ana	alyzed	: 07/05/20 17:30
Client ID	: RAMW1-20	0630		Date Red	ceived	: 06/30/20
Lab ID	: L2027725-0	04		Date Col	lected	: 06/30/20 11:50
Project Na	ame 💠 🗄 HHII TA1			Project N	lumber	16.6334
Client	: C.T. Male A	Associates		Lab Nurr	ıber	: L2027725



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Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method	 C.T. Male Associates HHII TA1 L2027725-05 RAMW2-200630 SCHENECTADY NY WATER 1,8260C VG200705 A12 	Lab Number Project Number Date Collected Date Received Date Analyzed Dilution Factor Analyst	: L2027725 : 16.6334 : 06/30/20 12:05 : 06/30/20 : 07/05/20 17:56 : 1 : NLK : CONZO
Lab File ID Sample Amount	: VG200705A12 : 10 ml	Instrument ID GC Column	: GONZO : RTX-502.2
Levei Extract Volume (MeOH)	: LOW	%Solids	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	2.8	2.5	0.70	
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	0.26	0.50	0.15	J
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	4.4	0.50	0.18	
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	0.46	0.50	0.19	L
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	N VS
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



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Client	: C.T. Male Associates	Lab Number	E2027725
Project Name	: HHII TA1	Project Number	16.6334
Lab ID	: L2027725-05	Date Collected	: 06/30/20 12:05
Client ID	: RAMW2-200630	Date Received	: 06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed	: 07/05/20 17:56
Sample Matrix	: WATER	Dilution Factor	: 1
Analytical Method	: 1,8260C	Analyst	: NLK
Lab File ID	: VG200705A12	Instrument ID	: GONZO
Sample Amount	🗄 10 mi	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH) : N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
57-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	υ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U



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Client	: C.T. Male Associates	1	.ab Nun	nber	: L2027725
Project Name	: HHII TA1		Project N	lumber	: 16.6334
Lab ID	: L2027725-05	1	Date Co	llected	: 06/30/20 12:05
Client ID	: RAMW2-200630	1	Date Re	ceived	: 06/30/20
Sample Location	: SCHENECTADY NY	1	Date Ana	alyzed	: 07/05/20 17:56
Sample Matrix	WATER	1	Dilution	Factor	ŝ 1
Analytical Method	: 1,8260C		Analyst		: NLK
Lab File ID	: VG200705A12	1	nstrume	ent ID	: GONZO
Sample Amount	: 10 ml	(GC Colu	mn	: RTX-502.2
Level	: LOW	c	%Solids		: N/A
Extract Volume (N	leOH):N/A	I	njection	Volume	: N/A
			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U
100-07-2		ND	10	0.40	0



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Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-06	Date Collected	: 06/30/20 12:55
Client ID	: RAMW3-200630	Date Received	: 06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed	: 07/05/20 18:21
Sample Matrix	: WATER	Dilution Factor	¥ 1
Analytical Method	: 1,8260C	Analyst	: NLK
Lab File ID	: VG200705A13	Instrument ID	: GONZO
Sample Amount	💈 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH): N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	1.7	2.5	0.70	J
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	6.8	0.50	0.18	
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	0.69	0.50	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	JU US
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



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Client	: C.T. Male Associates	Lab Number : L2027725
Project Name	; HHII TA1	Project Number : 16.6334
Lab ID	: L2027725-06	Date Collected : 06/30/20 12:55
Client ID	: RAMW3-200630	Date Received : 06/30/20
Sample Location	: SCHENECTADY NY	Date Analyzed : 07/05/20 18:21
Sample Matrix	: WATER	Dilution Factor : 1
Analytical Method	: 1,8260C	Analyst : NLK
Lab File ID	: VG200705A13	Instrument ID : GONZO
Sample Amount	: 10 ml	GC Column : RTX-502.2
Level	: LOW	%Solids : N/A
Extract Volume (MeO	H) : N/A	Injection Volume : N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	NÐ	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	on U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	υ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U



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Client	: C.T. Male Associates		Lab Num	ber	: L2027725
Project Name	: HHII TA1		Project N	lumber	: 16.6334
Lab ID	: L2027725-06		Date Col	lected	: 06/30/20 12:55
Client ID	: RAMW3-200630		Date Red	eived	: 06/30/20
Sample Location	: SCHENECTADY NY		Date Ana	lyzed	: 07/05/20 18:21
Sample Matrix	: WATER		Dilution	actor	; 1
Analytical Method	: 1,8260C		Analyst		: NLK
Lab File ID	: VG200705A13		Instrume	nt ID	GONZO
Sample Amount	: 10 ml		GC Colu	mn	: RTX-502.2
Level	: LOW		%Solids		: N/A
Extract Volume (Me	DH) : N/A		Injection	Volume	: N/A
			ug/L		
CAS NO. P	arameter	Results	RL	MDL	Qualifier
76-13-1 Fr	eon-113	ND	2.5	0.70	U
108-87-2 M	ethyl cyclohexane	ND	10	0.40	U



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Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Level Extract Volume (I	: WATER : 1,8260C : VG200705A14 : 10 ml : LOW		Lab Num Project N Date Coll Date Rec Date Ana Dilution F Analyst Instrume GC Colu %Solids Injection	lumber lected eived lyzed factor nt ID mn	: L2027725 : 16.6334 : 06/30/20 00:00 : 06/30/20 : 07/05/20 18:46 : 1 : NLK : GONZO : RTX-502.2 : N/A : N/A
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	N US
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U
		110	0.30	0.17	



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Client Project Name Lab ID Client ID Sample Locatio Sample Matrix Analytical Meth Lab File ID Sample Amoun Level Extract Volume	: WATER od : 1,8260C : VG200705A14		Lab Num Project N Date Col Date Rec Date Ana Dilution F Analyst Instrume GC Colu %Solids Injection	lumber lected ceived lyzed Factor nt ID mn	: L2027725 : 16.6334 : 06/30/20 00:00 : 06/30/20 : 07/05/20 18:46 : 1 : NLK : GONZO : RTX-502.2 : N/A : N/A
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U
				- 325	-



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Client	: C.T. Male Associates		Lab Nun	nber	: L2027725
Project Name	: HHII TA1		Project N	lumber	: 16.6334
Lab ID	: L2027725-07		Date Col	lected	: 06/30/20 00:00
Client ID	: FD01-200630		Date Re	ceived	: 06/30/20
Sample Location	SCHENECTADY NY		Date Ana	alyzed	07/05/20 18:46
Sample Matrix	: WATER		Dilution	Factor	11
Analytical Method	: 1,8260C		Analyst		: NLK
Lab File ID	: VG200705A14		Instrume	nt ID	: GONZO
Sample Amount	: 10 ml		GC Colu	mn	: RTX-502.2
Level	: LOW		%Solids		: N/A
Extract Volume (M	eOH) : N/A		Injection	Volume	: N/A
			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U
108-87-2	Metnyi cyclonexane	ND	10	0.40	U



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Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Level Extract Volume (N	: VG200705A07 : 10 ml : LOW		Lab Num Project N Date Col Date Red Date Ana Dilution F Analyst Instrume GC Colu %Solids Injection	lumber lected ceived lyzed Factor nt ID mn	: L2027725 : 16.6334 : 06/30/20 13:05 : 06/30/20 : 07/05/20 15:49 : 1 : NLK : GONZO : RTX-502.2 : N/A : N/A
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	1 mail
75-01-4	Vinyl chloride	ND	_		JV US
75-00-3			1.0	0.07	U
	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



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Client	: C.T. Male Associates	Lab Number : L2027725
Project Name	: HHII TA1	Project Number : 16.6334
Lab ID	: L2027725-08	Date Collected : 06/30/20 13:05
Client ID	EB01-200630	Date Received : 06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed : 07/05/20 15:49
Sample Matrix	: WATER	Dilution Factor : 1
Analytical Method	: 1,8260C	Analyst : NLK
Lab File ID	: VG200705A07	Instrument ID : GONZO
Sample Amount	: 10 ml	GC Column : RTX-502.2
Level	: LOW	%Solids : N/A
Extract Volume (MeO	0H) : N/A	Injection Volume : N/A

1.23			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	บ	
79-01-6	Trichloroethene	ND	0.50	0.18	U	
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U	
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U	
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U	
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U	
179601-23-1	p/m-Xylene	ND	2.5	0.70	U	
95-47-6	o-Xylene	ND	2.5	0.70	U	
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U	
100-42-5	Styrene	ND	2.5	0.70	U	
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U	
67-64-1	Acetone	3.8	5.0	1.5	J	
75-15-0	Carbon disulfide	ND	5.0	1.0	U	
78-93-3	2-Butanone	ND	5.0	1.9	U	
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U	
591-78-6	2-Hexanone	ND	5.0	1.0	U	
74-97-5	Bromochloromethane	ND	2.5	0.70	U	
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U	
98-82-8	Isopropylbenzene	ND	2.5	0.70	U	
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U	
79-20-9	Methyl Acetate	ND	2.0	0.23	U	
110-82-7	Cyclohexane	ND	10	0.27	U	
123-91-1	1,4-Dioxane	ND	250	61.	U	



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108-87-2	Methyl cyclohexane	ND	10	0.40	U
76-13-1	Freon-113	ND	2.5	0.70	U
CAS NO.	Parameter	Results	RL	MDL	Qualifier
			ug/L		
Extract Volum	ne (MeOH) : N/A		Injection	Volume	: N/A
Level	: LOW		%Solids		: N/A
Sample Amou	ınt : 10 ml		GC Colu		: RTX-502.2
Lab File ID	: VG200705A07		Instrume	ent ID	: GONZO
Analytical Me	thod : 1,8260C		Analyst		: NLK
Sample Matri	K WATER		Dilution	Factor	÷1
Sample Locat	ion : SCHENECTADY NY		Date An	alyzed	: 07/05/20 15:49
Client ID	: EB01-200630		Date Re	ceived	: 06/30/20
Lab ID	: L2027725-08		Date Co	llected	: 06/30/20 13:05
Project Name	: HHII TA1		Project N	lumber	: 16.6334
Client	C.T. Male Associates		Lab Nun	nber	: L2027725



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Client	: C.T. Male Associates	Lab Number	L2027725
Project Name	: HHII TA1	Project Number	16.6334
Lab ID	: L2027725-09	Date Collected	06/30/20 00:00
Client ID	: TRIP BLANK-200630	Date Received :	06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed	07/05/20 15:23
Sample Matrix	: WATER	Dilution Factor	1
Analytical Method	: 1,8260C	Analyst :	NLK
Lab File ID	: VG200705A06	Instrument ID	GONZO
Sample Amount	: 10 ml	GC Column	RTX-502.2
Level	: LOW	%Solids	N/A
Extract Volume (MeO	H) : N/A	Injection Volume	N/A

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
75-09-2	Methylene chloride	ND	2.5	0.70	U	
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U	
67-66-3	Chloroform	ND	2.5	0.70	U	
56-23-5	Carbon tetrachloride	ND	0.50	0.13	υ	
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U	
124-48-1	Dibromochloromethane	ND	0.50	0.15	U	
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U	
127-18-4	Tetrachloroethene	ND	0.50	0.18	U	
108-90-7	Chlorobenzene	ND	2.5	0.70	U	
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U	
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U	
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U	
75-27-4	Bromodichloromethane	ND	0.50	0.19	U	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U	
75-25-2	Bromoform	ND	2.0	0.65	U	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U	
71-43-2	Benzene	ND	0.50	0.16	U	
108-88-3	Toluene	ND	2.5	0.70	U	
100-41-4	Ethylbenzene	NĎ	2.5	0.70	U	
74-87-3	Chloromethane	ND	2.5	0.70	U	
74-83-9	Bromomethane	ND	2.5	0.70	NUS	
75-01-4	Vinyl chloride	ND	1.0	0.07	U	
75-00-3	Chloroethane	ND	2.5	0.70	U	
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U	



mr 7/14/20

Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-09	Date Collected	: 06/30/20 00:00
Client ID	: TRIP BLANK-200630	Date Received	: 06/30/20
Sample Location	: SCHENECTADY NY	Date Analyzed	: 07/05/20 15:23
Sample Matrix	: WATER	Dilution Factor	š 1
Analytical Method	: 1,8260C	Analyst	🗄 NLK
Lab File ID	: VG200705A06	Instrument ID	: GONZO
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeO	H) : N/A	Injection Volume	: N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xyłene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
74-97-5	Bromochloromethane	ND	2.5	0.70	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
87-61-6	1,2,3-Trichlorobenzene	ND	2.5	0.70	U
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
123-91-1	1,4-Dioxane	ND	250	61.	U

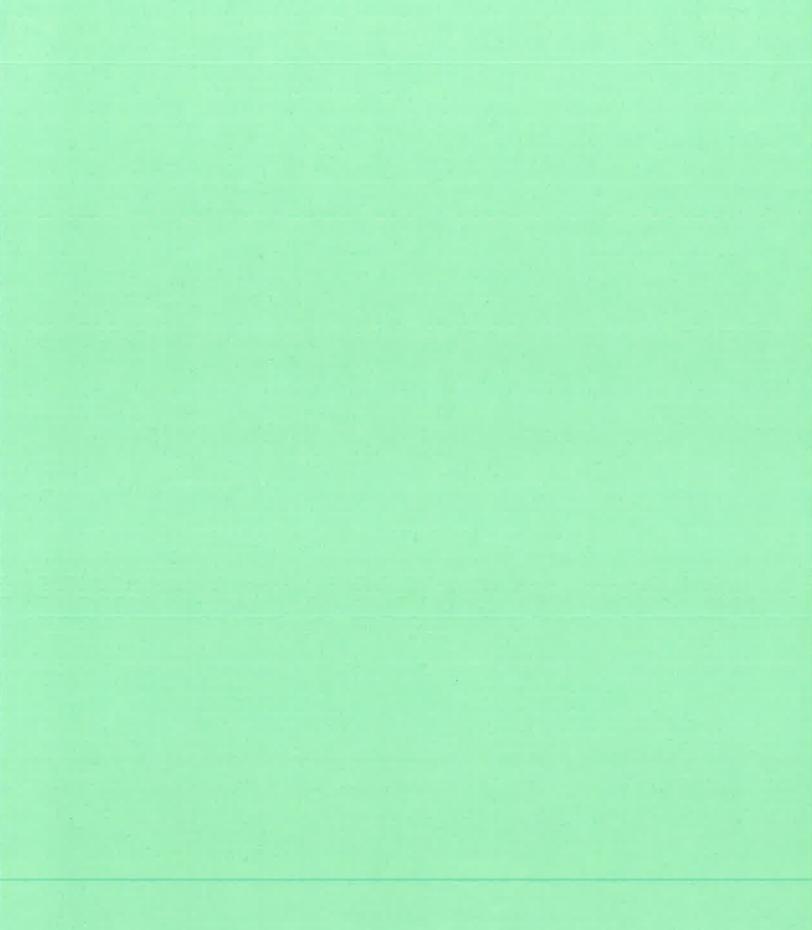


MT 7/14/20

: C.T. Male Associates		Lab Nun	nber	: L2027725
: HHII TA1		Project I	Number	: 16.6334
: L2027725-09		Date Co	llected	: 06/30/20 00:00
: TRIP BLANK-200630		Date Re	ceived	: 06/30/20
: SCHENECTADY NY		Date An	alvzed	: 07/05/20 15:23
: WATER		Dilution	Factor	8 1
: 1,8260C		Analyst		: NLK
: VG200705A06			nt ID	: GONZO
: 10 ml		GC Colu	mn	: RTX-502.2
: LOW		%Solids		: N/A
DH) : N/A		: N/A		
		ug/L		
Irameter	Results	RL	MDL	Qualifier
eon-113	ND	2.5	0.70	U
ethyl cyclohexane	ND	10	0.40	U
	: HHII TA1 : L2027725-09 : TRIP BLANK-200630 : SCHENECTADY NY : WATER : 1,8260C : VG200705A06 : 10 ml : LOW DH) : N/A arameter eon-113	: HHII TA1 : L2027725-09 : TRIP BLANK-200630 : SCHENECTADY NY : WATER : 1,8260C : VG200705A06 : 10 ml : LOW DH) : N/A arameter Results	: HHII TA1Project N: L2027725-09Date Co: TRIP BLANK-200630Date Re: SCHENECTADY NYDate And: WATERDilution: 1,8260CAnalyst: VG200705A06Instrume: 10 mlGC Colu: LOW%SolidsDH) : N/AInjectionarameterug/Leon-113ND2.5	: HHII TA1 Project Number : L2027725-09 Date Collected : TRIP BLANK-200630 Date Received : SCHENECTADY NY Date Analyzed : WATER Dilution Factor : 1,8260C Analyst : VG200705A06 Instrument ID : 10 ml GC Column : LOW %Solids DH) : N/A Injection Volume arameter ug/L eon-113 ND 2.5 0.70



MT 2/14/20



Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Methor Lab File ID Sample Amount Extraction Metho Extract Volume GPC Cleanup	: WATER d : 1,8270D : 27725-01 : 275 ml	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume		: 06/30/20 08:00 : 06/30/20 : 07/02/20 18:57 : 07/01/20 : 1 : SZ : SV124 : RTX5-MS : N/A		
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier	
	1 412/116161	nesuls	ΠL		Quainter	
111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U	
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U	
121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U	
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U	
7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U	
101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U	
108-60-1	Bis(2-chloroisopropyl)ether	ND	2.0	0.53	JU UJ	
111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U	
77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U	
78-59-1	Isophorone	ND	5.0	1.2	U	
98-95-3	Nitrobenzene	ND	2.0	0.77	U	
86-30-6	NDPA/DPA	ND	2.0	0.42	U	
621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U	
117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U	
85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U	
84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U	
117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U	
84-66-2	Diethyl phthalate	ND	5.0	0.38	U	
131-11-3	Dimethyl phthalate	ND	5.0	1.8	U	
92-52-4	Biphenyl	ND	2.0	0.46	U	
106-47-8	4-Chloroaniline	ND	5.0	1.1	NOS 2	



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-01 : 275 ml	Projec Date C Date F Date F Date F Dilutio Analys Instrum GC Cc %Solid	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume		27725 334 0/20 08:00 0/20 2/20 18:57 1/20 24 5-MS
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
-				1	
99-09-2	3-Nitroaniline	ND	5.0	0.81	U
100-01-6	4-Nitroaniline	ND	5.0	0.80	y us
132-64-9	Dibenzofuran	ND	2.0	0.50	U
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
98-86-2	Acetophenone	ND	5.0	0.53	U
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
95-57-8	2-Chlorophenol	ND	2.0	0.48	U
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
88-75-5	2-Nitrophenol	ND	10	0.85	xUS
100-02-7	4-Nitrophenol	ND	10	0.67	υ
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
108-95-2	Phenol	ND	5.0	0.57	U
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
86-74-8	Carbazole	ND	2.0	0.49	U
1912-24-9	Atrazine	ND	10	0.76	U
100-52-7	Benzaldehyde	ND	5.0	0.53	U
105-60-2	Caprolactam	ND	10	3.3	U
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



MT 1/14/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

Client	: C.T. Male Associates	Lab Number	L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-01	Date Collected	: 06/30/20 08:00
Client ID	: RAMW5-200630	Date Received	: 06/30/20
Sample Location	: SCHENECTADY NY	Date Analyzed	: 07/02/20 18:57
Sample Matrix	: WATER	Date Extracted	: 07/01/20
Analytical Method	: 1,8270D	Dilution Factor	å 1
Lab File ID	: 27725-01	Analyst	: SZ
Sample Amount	: 275 ml	Instrument ID	: SV124
Extraction Method	: EPA 3510C	GC Column	:
Extract Volume	: 1000 uL	%Solids	: N/A
GPC Cleanup	: N	Injection Volume	: 1 uL

Number TICS found: 2

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualifier	
	Unknown	10.95	2.44	55	
	Total TIC Compounds		2.44J	13	



MT 1/11/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-02 : 275 ml	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume		: 06/30 : 06/30 : 07/02 : 07/02 : 1 : SZ : SV12 : RTX : N/A	334 0/20 08:35 0/20 2/20 20:31 1/20 24
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U
121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U
7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U
101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U
108-60-1	Bis(2-chlorolsopropyl)ether	ND	2.0	0.53	VUS
111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U
77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U
78-59-1	Isophorone	ND	5.0	1.2	U
98-95-3	Nitrobenzene	ND	2.0	0.77	U
86-30-6	NDPA/DPA	ND	2.0	0.42	U
621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U
117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U
85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U
84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U
117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U
84-66-2	Diethyl phthalate	ND	5.0	0.38	U
131-11-3	Dimethyl phthalate	ND	5.0	1.8	U
92-52-4	Biphenyl	ND	2.0	0.46	U
106-47-8	4-Chloroaniline	ND	5.0	1.1	VUS
88-74-4	2-Nitroaniline	ND	5.0	0.50	U



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-02 : 275 ml	Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solid	t Number collected neceived nalyzed xtracted n Factor t nent ID lumn ls on Volume	: 06/30/20 08:35 : 06/30/20 : 07/02/20 20:31 : 07/01/20 : 1 : SZ : SV124 : RTX5-MS : N/A		
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier	
99-09-2	3-Nitroaniline	ND	5.0	0.81	U	
100-01-6	4-Nitroaniline	ND	5.0	0.80	JU US	
132-64-9	Dibenzofuran	ND	2.0	0.50	U	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U	
98-86-2	Acetophenone	ND	5.0	0.53	U	
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U	
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U	
95-57-8	2-Chlorophenol	ND	2.0	0.48	U	
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U	
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U	
88-75-5	2-Nitrophenol	ND	10	0.85	y US	
100-02-7	4-Nitrophenol	ND	10	0.67	U	
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U	
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U	
108-95-2	Phenol	ND	5.0	0.57	U	
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U	
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U	
86-74-8	Carbazole	ND	2.0	0.49	U	
1912-24-9	Atrazine	ND	10	0.76	U	
100-52-7	Benzaldehyde	ND	5.0	0.53	U	
105-60-2	Caprolactam	ND	10	3.3	U	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U	



MT 1/14/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

Number TICS found: 2

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualifier	
	Unknown	10.95	1.78	13	· · · · · · · · · · · · · · · · · · ·
	Total TIC Compounds		1.78J	13	



MT 7/14/20

CAS NO. Parameter Interval in the second se	Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-03 : 275 ml	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume		: 06/30 : 06/30 : 07/0 : 07/0 : 1 : SZ : SV12 : RTX : N/A	334 0/20 11:35 0/20 2/20 20:54 1/20
91-94-1 3,3'-Dichlorobenzidine ND 5.0 1.6 U 121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U 606-20-2 2,6-Dinitrotoluene ND 5.0 0.33 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chlorolsopropyl)ether ND 2.0 0.53	CAS NO.	Parameter	Results		MDL	Qualifier
121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U 606-20-2 2,6-Dinitrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chlorosiopropyl)ether ND 2.0 0.53 U/U) 111-91-1 Bis(2-chlorositoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 1.2 U 98-85-3 Nitrobenzene ND 2.0 0.69 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 812-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.42 U 817-84-7 Bis(2-ethylhexyl)phthalate ND 5.0 1.2 U 817-84-7 Di-hotylyphthalate ND 5.0 1.2 U 817-84-7 Di-hotylyphthalate ND 5.0 1.3 U 817-84-7	111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U
606-20-2 2,6-Dintrototuene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chlorosboropyl)ether ND 2.0 0.53	91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U
Yous-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 U/U) 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 1.2 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 821-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.44 U 8117-81-7 Bis(2-ethylhexyl)phthalate ND 5.0 0.44 U 84-66-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-66-2 Diehyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4	121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U
101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 U/U) 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.42 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 821-64-7 n-Nitrosodi-n-propylamine ND 5.0 1.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 92-52-4 Biphenyl ND 5.0 1.8 U 92-52-4 Choroaniline	606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U
108-60-1 Bis(2-chlorosisopropy)ether ND 2.0 0.53 V V 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.77 U 821-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 5.0 1.2 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate	7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U
111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 1.5 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.1 J// // // // // // 106-47-8 4-Chloroaniline ND 5.0 1.1 J// // <td>101-55-3</td> <td>4-Bromophenyl phenyl ether</td> <td>ND</td> <td>2.0</td> <td>0.38</td> <td>U</td>	101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U
T7-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 1.5 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 117-84-0 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 J// // // // // // // // // // // // //	108-60-1	Bis(2-chloroisopropyl)ether	ND	2.0	0.53	y UJ
78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.4 J/ J/	111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U
98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 J/ / / / / / /	77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U
86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 J// // // // // //	78-59-1	Isophorone	ND	5.0	1.2	U
621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 J/ / / / /	98-95-3	Nitrobenzene	ND	2.0	0.77	U
117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V V.S	86-30-6	NDPA/DPA	ND	2.0	0.42	U
85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U
84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 JUUS	117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U
117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U
84-66-2 Diethyl phthafate ND 5.0 0.38 U 131-11-3 Dimethyl phthafate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V V	84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U
131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U
92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V V	84-66-2	Diethyl phthalate	ND	5.0	0.38	U
106-47-8 4-Chloroaniline ND 5.0 1.1 V	131-11-3	Dimethyl phthalate	ND	5.0	1.8	U
	92-52-4	Biphenyl	ND	2.0	0.46	U
88-74-4 2-Nitroaniline ND 5.0 0.50 U	106-47-8	4-Chloroaniline	ND	5.0	1.1	W VS
	88-74-4	2-Nitroaniline	ND	5.0	0.50	U



Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-03 : 275 ml	Project Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solid	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume		27725 334 0/20 11:35 0/20 2/20 20:54 1/20 24 5-MS
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
	i unanistei	nesuits	116	MDL	
99-09-2	3-Nitroaniline	ND	5.0	0.81	U
100-01-6	4-Nitroaniline	ND	5.0	0.80	WV3
132-64-9	Dibenzofuran	ND	2.0	0.50	U
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
98-86-2	Acetophenone	ND	5.0	0.53	U
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
95-57-8	2-Chlorophenol	ND	2.0	0.48	U
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
88-75-5	2-Nitrophenol	ND	10	0.85	W US
100-02-7	4-Nitrophenol	ND	10	0.67	U
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
108-95-2	Phenol	ND	5.0	0.57	U
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
86-74-8	Carbazole	ND	2.0	0.49	U
1912-24-9	Atrazine	ND	10	0.76	U
100-52-7	Benzaldehyde	ND	5.0	0.53	U
105-60-2	Caprolactam	ND	10	3.3	U
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



WAT 7/14/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-03 : 275 ml	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume	: L2027725 : 16.6334 : 06/30/20 11:35 : 06/30/20 : 07/02/20 20:54 : 07/01/20 : 1 : SZ : SV124 : : N/A : 1 uL
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Number TICS found: 2

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualifier	
	Unknown	2.38	2.65	63	
	Total TIC Compounds		2.65J	13	



M 7/14/20

CAS NO. Parameter ug/L NDL NDL Outliner 111144-4 Bis(2-chloroethyl)ether ND 2.0 0.50 U 91-94-1 3,3-Dichlorobenzidine ND 5.0 1.6 U 121-14-2 2,4-Dinitrotoluene ND 5.0 0.33 U 606-20-2 2,6-Dinitrotoluene ND 2.0 0.49 U 101-55-3 4-Chlorophenyl phenyl ether ND 2.0 0.33 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.33 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.50 U 111-91-1 Bis(2-chloroethoxy)methane ND 2.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 2.0 0.50 U 78-59-1 Isophorone ND 5.0 0.51 U 88-30-4 NDPA/DPA ND 5.0 0.64 U 8164-81/Discolthophynhynhibalae	Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-04 : 275 ml	Date C Date F Date A Date E Dilutio Analys Instrur GC Cc %Solid	t Number Collected Received Analyzed Extracted In Factor St Inent ID Solumn	: 06/3 : 06/3 : 07/0 : 07/0 : 1 : SZ : SV1 : RTX : N/A	334 0/20 11:50 0/20 2/20 21:18 1/20 24
111-44-4 Bis(2-chloroethyl)ether ND 2.0 0.50 U 91-94-1 3,3'-Dichlorobenzidine ND 5.0 1.6 U 121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U 606-20-2 2,6-Dinitrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chlorostopropyl)ether ND 2.0 0.50 U 77-47-4 Hexachlorocyclopentadisne ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 5.0 1.2 U 86-30-6 NDPA/DPA ND 2.0 0.77 U 86-31-7 Bis(2-ethylhexyl)phthalate ND 5.0 1.2 U 87-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-71-7 Bis(2-ethylhexyl)phthalate ND 5.0 1.2 U 84-66-7	CARNO	Devenuelev				• • • •
ND 5.0 1.6 U 121-14-2 2,4-D initrotoluene ND 5.0 1.2 U 666-20-2 2,6-D initrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.53 JUUS 108-60-1 Bis(2-chlorosiopropyl)ether ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 0.50 U 78-59-1 Isophorone ND 2.0 0.69 U 86-30-6 NDPA/DPA ND 2.0 0.69 U 98-95-3 Nitrobenzene ND 5.0 1.2 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 86-30-6 NDPA/DPA ND 3.0 1.5 U 86-30-6 NDPA/DPA ND 5.0 0.44 U 117-61-7 Bis(2-ethylhexyl)phthalate ND 5.0 1.2 U <td>CAS NO.</td> <td>Parameter</td> <td>Results</td> <td>RL</td> <td>MDL</td> <td>Qualifier</td>	CAS NO.	Parameter	Results	RL	MDL	Qualifier
121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U 606-20-2 2,6-Dinitrotoluene ND 5.0 0.83 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 106-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 \checkmark U 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 1.2 U 86-30-6 NDPA/DPA ND 2.0 0.69 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 812-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 817-74-2 Bis(2-ethylhexyliphthalate ND 5.0 0.64 U 82-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 84-66-2 <td>111-44-4</td> <td>Bis(2-chloroethyl)ether</td> <td>ND</td> <td>2.0</td> <td>0.50</td> <td>U</td>	111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U
B66-20-2 2,6-Dinitrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chlorosiopropyl)ether ND 2.0 0.53 V V 111-91-1 Bis(2-chlorositocytopyl)ether ND 5.0 0.50 U V 77-47-4 Hexachlorocyclopentadiene ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 86-30-6 NDPA/DPA ND 5.0 0.42 U 86-30-6 NDPA/DPA ND 5.0 0.42 U 81-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.44 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 84-66-2	91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U
Total Addition NR State C 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 J// // // // // // // // // // // // //	121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U
101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 J/ U J 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 2.0 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 88-30-6 NDPA/DPA ND 2.0 0.42 U 86-30-6 NDPA/DPA ND 2.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl	606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U
Instruction Instruction <thinstruction< th=""> <thinstruction< th=""></thinstruction<></thinstruction<>	7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U
111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 117-84-0 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.1 J// / / / / /	101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U
77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 0.38 U 117-84-0 Di-n-ottylphthalate ND 5.0 0.38 U 117-84-0 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.1 JV X 106-47-8 4-Chloroaniline <td>108-60-1</td> <td>Bis(2-chloroisopropyl)ether</td> <td>ND</td> <td>2.0</td> <td>0.53</td> <td>NUS</td>	108-60-1	Bis(2-chloroisopropyl)ether	ND	2.0	0.53	NUS
78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 1.8 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V V	111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U
98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.4 V 92-52-4 A-Chloroaniline ND 5.0 1.1 V	77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U
B6-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.2 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	78-59-1	Isophorone	ND	5.0	1.2	U
621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 J// / / / /	98-95-3	Nitrobenzene	ND	2.0	0.77	U
117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.1 V 106-47-8 4-Chloroaniline ND 5.0 1.1 V	86-30-6	NDPA/DPA	ND	2.0	0.42	บ
85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U
84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U
117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U
84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U
131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 V	117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U
92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 1000000000000000000000000000000000000	84-66-2	Diethyl phthalate	ND	5.0	0.38	U
106-47-8 4-Chloroaniline ND 5.0 1.1 4	131-11-3	Dimethyl phthalate	ND	5.0	1.8	U
	92-52-4	Biphenyl	ND	2.0	0.46	U
88-74-4 2-Nitroaniline ND 5.0 0.50 U	106-47-8	4-Chloroaniline	ND	5.0	1.1	y US
	88-74-4	2-Nitroaniline	ND	5.0	0.50	U



MT 7/14/10

CAS NO. Parameter ug/L Results nL MDL Oualifier 99-09-2 3-Nitroaniline ND 5.0 0.81 U 100-01-6 4-Nitroaniline ND 5.0 0.80 V (X) 132-64-9 Dibenzofuran ND 2.0 0.50 U 95-94-3 1,2,4,5-Teitrachlorobenzane ND 10 0.44 U 98-86-2 Acatophenone ND 5.0 0.53 U 98-96-2 2,4,6-Trichlorophenol ND 5.0 0.61 U 98-96-2 2,4,6-Trichlorophenol ND 5.0 0.61 U 98-96-2 2,4,6-Trichlorophenol ND 2.0 0.35 U 98-96-2 2,4,6-Trichlorophenol ND 2.0 0.41 U 98-96-7 p-Chloro-m-resol ND 2.0 0.48 U 120-85-2 2,4-Dichlorophenol ND 5.0 0.41 U 120-85-2 2,4-Dichlorophenol ND <th>Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup</th> <th>: 27725-04 : 275 ml</th> <th>Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solid</th> <th>t Number collected teceived analyzed extracted n Factor st nent ID olumn</th> <th>: 06/30 : 07/03 : 07/0 : 1 : SZ : SV12 : RTX : N/A</th> <th>334 D/20 11:50 D/20 2/20 21:18 1/20</th>	Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-04 : 275 ml	Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solid	t Number collected teceived analyzed extracted n Factor st nent ID olumn	: 06/30 : 07/03 : 07/0 : 1 : SZ : SV12 : RTX : N/A	334 D/20 11:50 D/20 2/20 21:18 1/20
99-09-2 3-Nitroaniline ND 5.0 0.81 U 100-01-6 4-Nitroaniline ND 5.0 0.80 V V S 132-64-9 Dibenzoturan ND 2.0 0.50 U 95-94-3 1,2,4,5-Tetrachlorobenzene ND 10 0.44 U 98-86-2 Acetophenone ND 5.0 0.53 U 95-94-3 1,2,4,5-Trichlorophenol ND 5.0 0.53 U 98-86-2 Acetophenone ND 5.0 0.61 U 95-50-7 p-Chloro-m-cresol ND 2.0 0.35 U 95-57-8 2-Chlorophenol ND 2.0 0.48 U 120-83-2 2,4-Dirichlorophenol ND 5.0 1.8 U 105-67-9 2,4-Dimethylphenol ND 10 0.85 V 100-02-7 4-Nitrophenol ND 10 0.66 U 51-28-5 2,4-Dinitro-o-cresol <t< td=""><td>CAS NO</td><td>Parameter</td><td>Results</td><td></td><td>MDI</td><td>Qualifier</td></t<>	CAS NO	Parameter	Results		MDI	Qualifier
100-01-6 4-Nitroaniline ND 5.0 0.80 V V 132-64-9 Dibenzofuran ND 2.0 0.50 U 95-94-3 1,2,4,5-Tetrachlorobenzene ND 10 0.44 U 98-66-2 Acetophenone ND 5.0 0.53 U 88-06-2 2,4,6-Trichlorophenol ND 5.0 0.61 U 95-50-7 p-Chloro-m-cresol ND 2.0 0.35 U 95-57-8 2-Chlorophenol ND 5.0 0.41 U 100-02-7 2,4-Dichlorophenol ND 5.0 0.41 U 100-02-7 4-Nitrophenol ND 10 0.85 V 100-02-7 4-Nitrophenol ND 10 0.57 U 108-95-2 2-Abinitro-cresol ND 10 0.57 U 108-95-2 Q-Phenol ND 10 0.57 U 108-95-2 Phenol ND 5.0 0.57			neaulta	116	WDL	waannei
Dibenzofuran ND 2.0 0.50 U 95-94-3 1,2,4,5-Tetrachlorobenzene ND 10 0.44 U 98-86-2 Acetophenone ND 5.0 0.53 U 88-06-2 2,4,6-Trichlorophenol ND 5.0 0.61 U 59-50-7 p-Chloro-m-cresol ND 2.0 0.35 U 95-57-8 2-Chlorophenol ND 2.0 0.48 U 120-83-2 2,4-Dichlorophenol ND 5.0 0.41 U 105-57-9 2,4-Dichlorophenol ND 5.0 1.8 U 100-02-7 2,4-Dimethylphenol ND 10 0.67 U 51-28-5 2,4-Dintrophenol ND 10 0.67 U 51-28-5 2,4-Dintrophenol ND 10 0.57 U 51-28-5 2,4-Dintrophenol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U	99-09-2	3-Nitroaniline	ND	5.0	0.81	U
132-64-9DibenzofuranND2.00.50U95-94-31,2,4,5-TetrachlorobenzeneND100.44U98-66-2AcetophenoneND5.00.53U88-06-22,4,6-TrichlorophenolND5.00.61U95-50-7p-Chloro-m-cresolND2.00.35U95-57-82-ChlorophenolND2.00.48U120-83-22,4-DichlorophenolND5.00.41U105-67-92,4-DichlorophenolND5.01.8U105-67-92,4-DichlorophenolND100.67U100-02-74-NitrophenolND100.67U101-02-74-NitrophenolND100.67U102-02-74-NitrophenolND100.67U104-02-74-NitrophenolND101.8U104-02-74-NitrophenolND101.8U104-02-74,6-Dinitro-o-cresolND101.8U104-02-74,6-Dinitro-o-cresolND5.00.57U104-03-4/106-44-53-Methylphenol/4-MethylphenolND5.00.77U104-03-4/106-44-53-MethylphenolND5.00.77U104-03-4/106-44-53-MethylphenolND1.00.49U104-03-4/106-44-53-MethylphenolND1.00.49U104-03-4CarbazoleND1.00	100-01-6	4-Nitroaniline	ND	5.0	0.80	YUS
B8-86-2 Acetophenone ND 5.0 0.53 U B8-06-2 2,4,6-Trichlorophenol ND 5.0 0.61 U 59-50-7 p-Chloro-m-cresol ND 2.0 0.35 U 95-57-8 2-Chlorophenol ND 2.0 0.48 U 120-83-2 2,4-Dichlorophenol ND 5.0 0.41 U 105-67-9 2,4-Dinethylphenol ND 5.0 1.8 U 105-67-9 2,4-Dinethylphenol ND 5.0 1.8 U 105-67-9 2,4-Dinethylphenol ND 10 0.85 ✓ 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 20 6.6 U 534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.77 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.	132-64-9	Dibenzofuran	ND	2.0	0.50	
B8-06-2 2,4,6-Trichlorophenol ND 5.0 0.61 U 59-50-7 p-Chloro-m-cresol ND 2.0 0.35 U 95-57-8 2-Chlorophenol ND 2.0 0.48 U 102-83-2 2,4-Dichlorophenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 1.8 U 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitro-c-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-95-2 Phenol ND 5.0 0.77 U 108-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 108-95-9 Phenol ND 0.49 U	95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
S9-50-7 p-Chloro-m-cresol ND 2.0 0.35 U 95-57-8 2-Chlorophenol ND 2.0 0.48 U 120-83-2 2,4-Dichlorophenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 1.8 U 88-75-5 2-Nitrophenol ND 10 0.85 J/J/J 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 10 0.67 U 534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.77 U 108-39-4/106-44-5 2-4,5-Trichlorophenol ND 5.0 0.77 U 108-57-4 Carbazole ND	98-86-2	Acetophenone	ND	5.0	0.53	U
95-57-8 2-Chlorophenol ND 2.0 0.48 U 120-83-2 2,4-Dichlorophenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 0.41 U 88-75-5 2,4-Dimethylphenol ND 5.0 1.8 U 88-75-5 2-Nitrophenol ND 10 0.85 ✓ 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dimitrophenol ND 10 0.67 U 51-28-5 2,4-Dimitro-o-cresol ND 10 1.8 U 534-52-1 4,6-Dinitro-o-cresol ND 5.0 0.57 U 108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 <t< td=""><td>88-06-2</td><td>2,4,6-Trichlorophenol</td><td>ND</td><td>5.0</td><td>0.61</td><td>U</td></t<>	88-06 -2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
120-83-2 2,4-Dichlorophenol ND 5.0 0.41 U 105-67-9 2,4-Dimethylphenol ND 5.0 1.8 U 88-75-5 2-Nitrophenol ND 10 0.85 JUC 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 10 0.67 U 534-52-1 4,6-Dinitro-ocresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-95-2 Phenol ND 5.0 0.57 U 108-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U	59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
105-67-9 2,4-Dimethylphenol ND 5.0 1.8 U 88-75-5 2-Nitrophenol ND 10 0.85 U/US 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 20 6.6 U 534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 100-52-7 Benzaldehyde ND 5.0 0.53 <td>95-57-8</td> <td>2-Chlorophenol</td> <td>ND</td> <td>2.0</td> <td>0.48</td> <td>U</td>	95-57-8	2-Chlorophenol	ND	2.0	0.48	U
B8-75-5 2-Nitrophenol ND 10 0.85 V 100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 20 6.6 U 534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
100-02-7 4-Nitrophenol ND 10 0.67 U 51-28-5 2,4-Dinitrophenol ND 20 6.6 U 534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
51-28-5 2,4-Dinitrophenol ND 20 6.6 U 534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 100-52-7 Caprolactam ND 5.0 0.53 U	88-75-5	2-Nitrophenol	ND	10	0.85	105
534-52-1 4,6-Dinitro-o-cresol ND 10 1.8 U 108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Capolactam ND 10 3.3 U	100-02-7	4-Nitrophenol	ND	10	0.67	U
108-95-2 Phenol ND 5.0 0.57 U 108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
108-39-4/106-44-5 3-Methylphenol/4-Methylphenol ND 5.0 0.48 U 95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
95-95-4 2,4,5-Trichlorophenol ND 5.0 0.77 U 86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	108-95-2	Phenol	ND	5.0	0.57	U
86-74-8 Carbazole ND 2.0 0.49 U 1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
1912-24-9 Atrazine ND 10 0.76 U 100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
100-52-7 Benzaldehyde ND 5.0 0.53 U 105-60-2 Caprolactam ND 10 3.3 U	86-74-8	Carbazole	ND	2.0	0.49	U
105-60-2 Caprolactam ND 10 3.3 U	1912-24-9	Atrazine	ND	10	0.76	U
	100-52-7	Benzaldehyde	ND	5.0	0.53	U
	105-60-2	Caprolactam	ND	10	3.3	U
	58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



M-T -1/1/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

Sample Amount: 2/5 miInstrument ID: SV124Extraction Method: EPA 3510CGC Column:Extract Volume: 1000 uL%Solids: N/AGPC Cleanup: NInjection Volume: 1 uL	Extract Volume	: 1000 uL	%Solids	
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Number TICS found: 3

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualif	ier
	Unknown	2.38	2.44	3	1
	Unknown	10.95	2.4	5	1
	Total TIC Compounds		4.84J	8	4

MT 7/14/20



CAS NO. Parameter ug/L Ug/L Qualifier	r
111-44-4 Bis(2-chloroethyl)ether ND 2.0 0.50 U	
91-94-1 3,3'-Dichlorobenzidine ND 5.0 1.6 U	
121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U	
606-20-2 2,6-Dinitrotoluene ND 5.0 0.93 U	
7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U	
101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U	
108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53	J
111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U	
77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U	
78-59-1 Isophorone ND 5.0 1.2 U	
98-95-3 Nitrobenzene ND 2.0 0.77 U	
86-30-6 NDPA/DPA ND 2.0 0.42 U	
621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U	
117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U	
85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U	
84-74-2 Di-n-butylphthalate ND 5.0 0.39 U	
117-84-0 Di-n-octylphthalate ND 5.0 1.3 U	
84-66-2 Diethyl phthalate ND 5.0 0.38 U	
131-11-3 Dimethyl phthalate ND 5.0 1.8 U	
92-52-4 Biphenyl ND 2.0 0.46 U	
106-47-8 4-Chloroaniline ND 5.0 1.1 🌿 (ろ
88-74-4 2-Nitroaniline ND 5.0 0.50 U	



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-05 : 275 ml	Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solid	Number ollected eceived nalyzed xtracted Factor t nent ID lumn	: 06/3 : 07/0 : 07/0 : 1 : SZ : SV1 : RTX : N/A	334 0/20 12:05 0/20 2/20 21:41 1/20 24 5-MS
CAS NO.	Parameter	Beaulte	ug/L RL	MDL	Qualifiar
CAS NO.	Farameler	Results	RL	MDL	Qualifier
99-09-2	3-Nitroaniline	ND	5.0	0.81	U
100-01-6	4-Nitroaniline	ND	5.0	0.80	YUS
132-64-9	Dibenzofuran	ND	2.0	0.50	U
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
98-86-2	Acetophenone	ND	5.0	0.53	U
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
95-57-8	2-Chlorophenol	ND	2.0	0.48	U
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
88-75-5	2-Nitrophenol	ND	10	0.85	VUS
100-02-7	4-Nitrophenol	ND	10	0.67	U
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
108-95-2	Phenol	ND	5.0	0.57	U
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
86-74-8	Carbazole	ND	2.0	0.49	U
1912-24-9	Atrazine	ND	10	0.76	U
100-52-7	Benzaldehyde	ND	5.0	0.53	U
105-60-2	Caprolactam	ND	10	3.3	U
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



MT 2/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: WATER d : 1,8270D : 27725-06 : 275 ml	Date C Date R Date A Date E Dilution Analys Instrun GC Co %Solid	t Number collected received malyzed xtracted n Factor t nent ID dumn	: 06/30 : 06/30 : 07/02 : 07/0 : 1 : SZ : SV12 : RTX : N/A	334 0/20 12:55 0/20 2/20 22:05 1/20 24
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U
121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U
7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U
101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U
108-60-1	Bis(2-chloroisopropyl)ether	ND	2.0	0.53	w us
111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U
77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U
78-59-1	Isophorone	ND	5.0	1.2	U
98-95-3	Nitrobenzene	ND	2.0	0.77	U
86-30-6	NDPA/DPA	ND	2.0	0.42	U
621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U
117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U
85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U
84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U
117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U
84-66-2	Diethyl phthalate	ND	5.0	0.38	U
131-11-3	Dimethyl phthalate	ND	5.0	1.8	U
92-52-4	Biphenyl	ND	2.0	0.46	U
106-47-8	4-Chloroaniline	ND	5.0	1.1	W VS
88-74-4	2-Nitroaniline				



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-06 : 275 ml	Projec Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solic	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column %Solids Injection Volume		7725 334 0/20 12:55 0/20 2/20 22:05 1/20 24 5-MS
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
99-09-2	3-Nitroaniline	ND	5.0	0.81	U
100-01-6	4-Nitroaniline	ND	5.0	0.80	× 15
132-64-9	Dibenzofuran	ND	2.0	0.50	U
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
98-86-2	Acetophenone	ND	5.0	0.53	U
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
95-57-8	2-Chlorophenol	ND	2.0	0.48	U
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
88-75-5	2-Nitrophenol	ND	10	0.85	NO VS
100-02-7	4-Nitrophenol	ND	10	0.67	U
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
108-95-2	Phenol	ND	5.0	0.57	U
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
86-74-8	Carbazole	ND	2.0	0.49	U
1912-24-9	Atrazine	ND	10	0.76	U
100-52-7	Benzaldehyde	ND	5.0	0.53	U
105-60-2	Caprolactam	ND	10	3.3	U
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



MA 7/14/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

Client	 C.T. Male Associates HHII TA1 L2027725-06 RAMW3-200630 SCHENECTADY NY WATER 1,8270D 27725-06 275 ml EPA 3510C 1000 uL 	Lab Number	: L2027725
Project Name		Project Number	: 16.6334
Lab ID		Date Collected	: 06/30/20 12:55
Client ID		Date Received	: 06/30/20
Sample Location		Date Analyzed	: 07/02/20 22:05
Sample Matrix		Date Extracted	: 07/01/20
Analytical Method		Dilution Factor	: 1
Lab File ID		Analyst	: SZ
Sample Amount		Instrument ID	: SV124
Extraction Method		GC Column	:
Extract Volume		%Solids	: N/A
Extract Volume	: 1000 uL	%Solids	: N/A
GPC Cleanup	: N	Injection Volume	: 1 uL

Number TICS found: 3

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualifier	
	Unknown	2.39	2.4	17	
	Unknown	10.95	1.45	4	
	Total TIC Compounds		3.85J	1 1	





Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-07 : 275 ml	Date C Date F Date A Date E Dilutio Analys Instrur GC Cc %Solid	t Number Collected Received Analyzed Extracted In Factor St In Factor	: 06/3 : 06/3 : 07/0 : 07/0 : 1 : SZ : SV1 : RTX : N/A	334 0/20 00:00 0/20 2/20 22:29 1/20 24
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U
121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U
7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U
101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U
108-60-1	Bis(2-chloroisopropyl)ether	ND	2.0	0.53	WV3
111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U
77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U
78-59-1	Isophorone	ND	5.0	1.2	U
98-95-3	Nitrobenzene	ND	2.0	0.77	U
86-30-6	NDPA/DPA	ND	2.0	0.42	U
621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U
117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U
85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U
84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U
117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U
84-66-2	Diethyl phthalate	ND	5.0	0.38	U
131-11-3	Dimethyl phthalate	ND	5.0	1.8	U
92-52-4	Biphenyl	ND	2.0	0.46	U
106-47-8	4-Chloroaniline	ND	5.0	1.1	JE US
88-74-4	2-Nitroaniline	ND	5.0	0.50	U



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-07 : 275 ml	Date C Date F Date E Dilutio Analys Instrum GC Cc %Solic	t Number collected received analyzed xtracted n Factor t nent ID Jumn	: 06/30 : 07/02 : 07/0 : 1 : SZ : SV12 : RTX : N/A	334 0/20 00:00 0/20 2/20 22:29 1/20
CAS NO	Baramatar	Populto	ug/L	MDL	Qualifier
CAS NO.	Parameter	Results	RL	MUL	wudimei
99-09-2	3-Nitroaniline	ND	5.0	0.81	U
100-01-6	4-Nitroaniline	ND	5.0	0.80	VUS
132-64-9	Dibenzofuran	ND	2.0	0.50	U
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
98-86-2	Acetophenone	ND	5.0	0.53	U
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
95-57-8	2-Chlorophenol	ND	2.0	0.48	U
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
88-75-5	2-Nitrophenol	ND	10	0.85	W UJ
100-02-7	4-Nitrophenol	ND	10	0.67	U
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
108-95-2	Phenol	ND	5.0	0.57	บ
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
86-74-8	Carbazole	ND	2.0	0.49	U
1912-24-9	Atrazine	ND	10	0.76	U
100-52-7	Benzaldehyde	ND	5.0	0.53	U
105-60-2	Caprolactam	ND	10	3.3	U
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



MT 2/14/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method	: 27725-07 : 275 ml : EPA 3510C	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column % Solide	: L2027725 : 16.6334 : 06/30/20 00:00 : 06/30/20 : 07/02/20 22:29 : 07/01/20 : 1 : SZ : SV124 : N/A
Extraction Method Extract Volume GPC Cleanup	: 1000 uL : N	%Solids Injection Volume	: : N/A : 1 uL

Number TICS found: 2

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualifier	
	Unknown	10.95	1.85	5 3	
	Total TIC Compounds		1.85J	13	

MT 7/11/20



CASNO.ParameterIngleRulRLMDLOutlitier11144-4Bis(2-chloroethyl)etherND2.00.50U81-94-13,3-DichlorobenzidineND5.01.6U121-14-22,4-DintrotolueneND5.01.2U606-20-22,6-DinikrotolueneND5.00.33U105-72-34-Chlorophenyl phenyl etherND2.00.49U105-72-34-Chlorophenyl phenyl etherND2.00.50U108-80-1Bis(2-chloroisopropyl)etherND2.00.50U119-11Bis(2-chlorosbynymthaneND5.00.50U77-47-4HexachlorocyclopentadieneND5.00.50U78-59-1IsophoroneND5.00.64U86-30-6NDPA/DPAND5.00.64U81-64-7n-Mitrosoft-propylamineND5.00.38U81-64-7n-Mitrosoft-propylamineND5.00.38U81-64-7Buyl benzyl phthalateND5.00.38U81-74-1Bis(2-chlorophthalateND5.00.38U81-74-1Buyl benzyl phthalateND5.00.38U81-74-1Buyl benzyl phthalateND5.00.38U81-74-1Bis(2-chloroethoxyl phthalateND5.00.38U81-74-1Buyl benzyl phthalateND5.00.38U<	Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-08 : 275 ml	Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solic	t Number collected teceived malyzed extracted n Factor st nent ID olumn	: 06/3 : 07/0 : 07/0 : 1 : SZ : SV1 : RTX : N/A	334 0/20 13:05 0/20 2/20 22:52 1/20 24 5-MS
111-44-4 Bis(2-chloroethyl)ether ND 2.0 0.50 U 81-94-1 3,3'-Dichlorobenzidine ND 5.0 1.6 U 121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U 606-20-2 2,6-Dinitrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 U/US 111-91-1 Bis(2-chloroisopropyl)ether ND 2.0 0.59 U 77-47-4 Hexachlorocyclopentadiene ND 2.0 0.69 U 78-59-1 Isophorone ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 117-81-7 Bis(2-ethylhexylphthalate ND 3.0 1.5 U 85-68-7 Butyl berzyl phthalate ND 5.0 0.38 U 84-74-2 <th>CASNO</th> <th>Paramotor</th> <th>Populto</th> <th></th> <th>MD</th> <th>Qualifier</th>	CASNO	Paramotor	Populto		MD	Qualifier
ND S.0 1.6 U 121-14-2 2,4-Dinitrotoluene ND S.0 1.2 U 666-20-2 2,6-Dinitrotoluene ND S.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chlorolsopropyl)ether ND 2.0 0.59 U 111-91-1 Bis(2-chlorolsopropyl)ether ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 0.59 U 88-95-3 Nitrobenzene ND 5.0 0.64 U 86-30-6 NDPA/DPA ND 5.0 0.64 U 117-81-7 Bis(2-chtylhexyl)phthalate ND 5.0 0.64 U 84-74-2 D-n-butylphthalate ND 5.0 0.39 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0		רמומוושנטו	nesuits	nL	WDL	Quaimer
121-14-2 2,4-Dinitrotoluene ND 5.0 1.2 U 606-20-2 2,6-Dinitrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroshoropyl)ether ND 2.0 0.50 U 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 1.2 U 86-80-6 NDPA/DPA ND 2.0 0.69 U 86-80-6 NDPA/DPA ND 2.0 0.77 U 86-80-6 NDPA/DPA ND 5.0 1.2 U 86-80-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 815(2-ethylphthalate ND 5.0 1.2 U 82-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-68-2 Di-n-butylphthalate ND </td <td>111-44-4</td> <td>Bis(2-chloroethyl)ether</td> <td>ND</td> <td>2.0</td> <td>0.50</td> <td>U</td>	111-44-4	Bis(2-chloroethyl)ether	ND	2.0	0.50	U
606-20-2 2,6-Dinktrotoluene ND 5.0 0.93 U 7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53	91-94-1	3,3'-Dichlorobenzidine	ND	5.0	1.6	U
7005-72-3 4-Chlorophenyl phenyl ether ND 2.0 0.49 U 101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 U 111-91-1 Bis(2-chloroisopropyl)ether ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 5.0 0.59 U 78-59-1 Isophorone ND 5.0 0.77 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 5.0 0.64 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-66-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 92-52-4 Biphenyl ND 5.0 1.8 U 92-52-4 Biphenyl	121-14-2	2,4-Dinitrotoluene	ND	5.0	1.2	U
101-55-3 4-Bromophenyl phenyl ether ND 2.0 0.38 U 108-60-1 Bis(2-chloroisopropyl)ether ND 2.0 0.53 U/US 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 2.0 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 5.0 1.5 U 8117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 1.2 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND	606-20-2	2,6-Dinitrotoluene	ND	5.0	0.93	U
108-60-1 Bis(2-chlorosisopropy)ether ND 2.0 0.53 V/V 111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 66-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 1.2 U 84-66-2 Diethyl phthalate ND 5.0 1.3 U 82-66-2 Diethyl phthalate ND 5.0 1.3 U 82-66-2 Diethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.	7005-72-3	4-Chlorophenyl phenyl ether	ND	2.0	0.49	U
111-91-1 Bis(2-chloroethoxy)methane ND 5.0 0.50 U 77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 1.5 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 5.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 1.2 U 84-66-2 Diemtyl phthalate ND 5.0 1.3 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.4 ✔ ✔ 106-47-8 4- Choroaniline <	101-55-3	4-Bromophenyl phenyl ether	ND	2.0	0.38	U
77-47-4 Hexachlorocyclopentadiene ND 20 0.69 U 78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 5.0 1.1 Af US 106-47-8 4-Chloroaniline ND 5.0 1.1 Af US	108-60-1	Bis(2-chloroisopropyl)ether	ND	2.0	0.53	V15
78-59-1 Isophorone ND 5.0 1.2 U 98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylarnine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 ✓ ✓	111-91-1	Bis(2-chloroethoxy)methane	ND	5.0	0.50	U
98-95-3 Nitrobenzene ND 2.0 0.77 U 86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 ✓ ✓	77-47-4	Hexachlorocyclopentadiene	ND	20	0.69	U
86-30-6 NDPA/DPA ND 2.0 0.42 U 621-64-7 n-Nitrosodi-n-propylarnine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 0.39 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 0.39 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 🗸 🗸	78-59-1	Isophorone	ND	5.0	1.2	U
621-64-7 n-Nitrosodi-n-propylamine ND 5.0 0.64 U 117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 1.3 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 1.8 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 🗸 🗸 🏹	98-95-3	Nitrobenzene	ND	2.0	0.77	U
117-81-7 Bis(2-ethylhexyl)phthalate ND 3.0 1.5 U 85-68-7 Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 I	86-30-6	NDPA/DPA	ND	2.0	0.42	U
Butyl benzyl phthalate ND 5.0 1.2 U 84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 Image: Control of the second s	621-64-7	n-Nitrosodi-n-propylamine	ND	5.0	0.64	U
84-74-2 Di-n-butylphthalate ND 5.0 0.39 U 117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 Image: Contract of the second	117-81-7	Bis(2-ethylhexyl)phthalate	ND	3.0	1.5	U
117-84-0 Di-n-octylphthalate ND 5.0 1.3 U 84-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 Image: Construction of the second	85-68-7	Butyl benzyl phthalate	ND	5.0	1.2	U
B4-66-2 Diethyl phthalate ND 5.0 0.38 U 131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 Image: Construction of the second	84-74-2	Di-n-butylphthalate	ND	5.0	0.39	U
131-11-3 Dimethyl phthalate ND 5.0 1.8 U 92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 Image: Comparison of the company of the com	117-84-0	Di-n-octylphthalate	ND	5.0	1.3	U
92-52-4 Biphenyl ND 2.0 0.46 U 106-47-8 4-Chloroaniline ND 5.0 1.1 Image: Comparison of the second sec	84-66-2	Diethyl phthalate	ND	5.0	0.38	U
106-47-8 4-Chloroaniline ND 5.0 1.1 1	131-11-3	Dimethyl phthalate	ND	5.0	1.8	U
	92-52-4	Biphenyl	ND	2.0	0.46	U
88-74-4 2-Nitroaniline ND 5.0 0.50 U	106-47-8	4-Chloroaniline	ND	5.0	1.1	15
	88-74-4	2-Nitroaniline	ND	5.0	0.50	U

MT 2/14/20



Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: WATER d : 1,8270D : 27725-08 : 275 ml	Projec Date C Date F Date F Date E Dilutio Analys Instrur GC Cc %Solid	ment ID blumn	: 06/30 : 06/30 : 07/02 : 07/01 : 1 : SZ : SV12 : RTX5 : N/A	334 0/20 13:05 0/20 2/20 22:52 1/20
			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
99-09-2	3-Nitroaniline	ND	5.0	0.81	U
100-01-6	4-Nitroaniline	ND	5.0	0.80	NUS
132-64-9	Dibenzofuran	ND	2.0	0.50	U
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	10	0.44	U
98-86-2	Acetophenone	ND	5.0	0.53	U
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.61	U
59-50-7	p-Chloro-m-cresol	ND	2.0	0.35	U
95-57-8	2-Chlorophenol	ND	2.0	0.48	U
120-83-2	2,4-Dichlorophenol	ND	5.0	0.41	U
105-67-9	2,4-Dimethylphenol	ND	5.0	1.8	U
88-75-5	2-Nitrophenol	ND	10	0.85	JU US
100-02-7	4-Nitrophenol	ND	10	0.67	U
51-28-5	2,4-Dinitrophenol	ND	20	6.6	U
534-52-1	4,6-Dinitro-o-cresol	ND	10	1.8	U
108-95-2	Phenol	ND	5.0	0.57	U
108-39-4/106-44-5	3-Methylphenol/4-Methylphenol	ND	5.0	0.48	U
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.77	U
86-74-8	Carbazole	ND	2.0	0.49	U
1912-24-9	Atrazine	ND	10	0.76	U
100-52-7	Benzaldehyde	ND	5.0	0.53	U
105-60-2	Caprolactam	ND	10	3.3	U
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.84	U



MT 7/14/20

Tentatively Identified Compounds Form 1 Semivolatile Organics by GC/MS

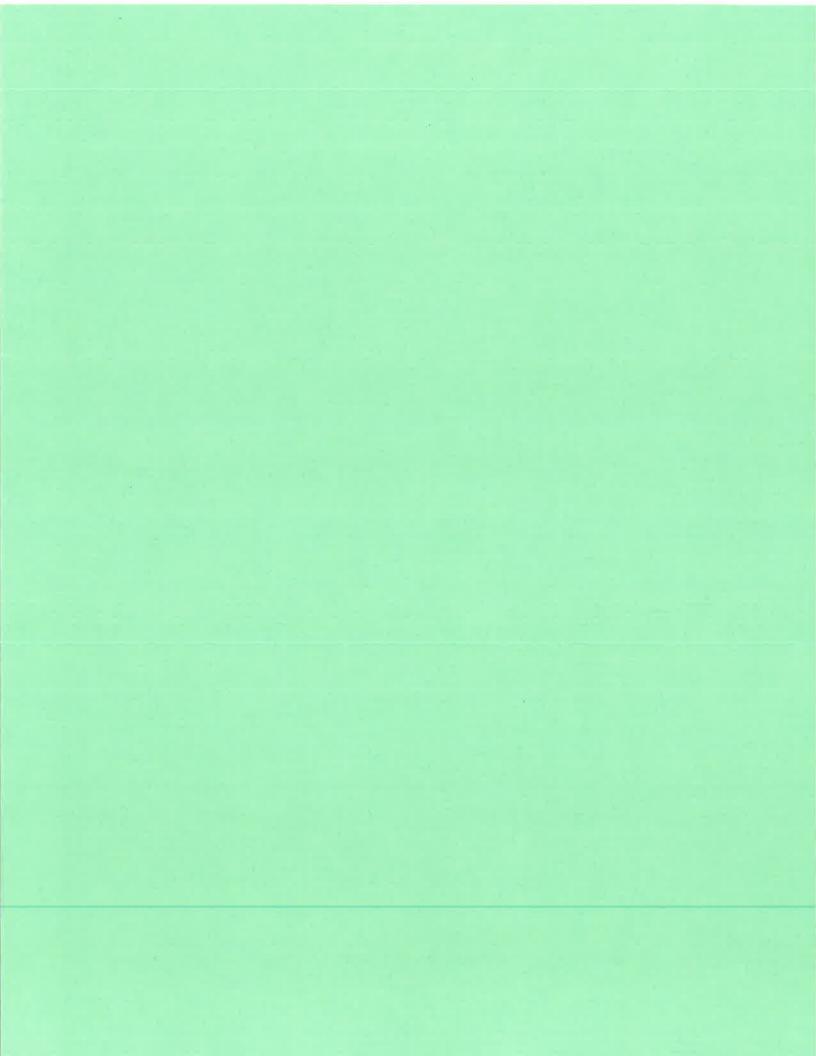
Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method	: 27725-08 : 275 ml	Lab Number Project Number Date Collected Date Received Date Analyzed Date Extracted Dilution Factor Analyst Instrument ID GC Column	: L2027725 : 16.6334 : 06/30/20 13:05 : 06/30/20 : 07/02/20 22:52 : 07/01/20 : 1 : SZ : SV124 :
•			: : N/A

Number TICS found: 2

Concentration Units: ug/L

CAS Number	Compound Name	RT	EST. CONC.	Qualifier	
	Unknown Organic Acid	10.95	1.71	r 5	
	Total TIC Compounds		1.71J	13	

MT 7/14/20



Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-01 : 275 ml	Date C Date R Date A Date E Dilution Analys Instrun GC Co %Solid	t Number collected acceived nalyzed xtracted n Factor t nent ID lumn Is on Volume	: 06/30 : 06/30	334 0/20 08:00 0/20 2/20 19:37 1/20 25
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
83-32-9	Acenaphthene	ND	0.10	0.01	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.02	U
206-44-0	Fluoranthene	ND	0.10	0.02	U
87-68-3	Hexachlorobutadiene	ND	0.50	0.05	U
91-20-3	Naphthalene	ND	0.10	0.05	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.02	U
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.01	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U
218-01-9	Chrysene	ND	0.10	0.01	U
208-96-8	Acenaphthylene	ND	0.10	0.01	U
120-12-7	Anthracene	ND	0.10	0.01	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U
86-73-7	Fluorene	ND	0.10	0.01	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U
129-00-0	Pyrene	ND	0.10	0.02	U
91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U
87-86-5	Pentachlorophenol	ND	0.80	0.01	U
118-74-1	Hexachlorobenzene	ND	0.80	0.01	U
67-72-1	Hexachloroethane	ND	0.80	0.06	U



MT 1/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-02 : 275 ml	Date Co Date Ro Date Au Date Ex Dilution Analyst Instrum GC Col %Solids	Number ollected eceived nalyzed ktracted Factor ent ID umn s	: 06/30	334 0/20 08:35 0/20 2/20 19:57 1/20
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
	F & 61116161	inesuits	nL	MUL	wuqiillei
83-32-9	Acenaphthene	ND	0.10	0.01	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.02	U
206-44-0	Fluoranthene	ND	0.10	0.02	U
87-68-3	Hexachlorobutadiene	ND	0.50	0.05	U
91-20-3	Naphthalene	ND	0.10	0.05	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.02	U
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.01	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U
218-01-9	Chrysene	ND	0.10	0.01	U
208-96-8	Acenaphthylene	ND	0.10	0.01	U
120-12-7	Anthracene	ND	0.10	0.01	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U
86-73-7	Fluorene	ND	0.10	0.01	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U
129-00-0	Pyrene	ND	0.10	0.02	U
91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U
87-86-5	Pentachlorophenol	ND	0.80	0.01	U
118-74-1	Hexachlorobenzene	ND	0.80	0.01	U
67-72-1	Hexachloroethane	ND	0.80	0.06	U



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-03 : 275 ml	Date C Date F Date E Dilutio Analys Instrur GC Cc %Solid	t Number Collected Received Analyzed Extracted In Factor St In St In St I	: 06/30 : 07/02 : 07/01 : 1 : DV : SV12 : RXI-5 : N/A	334 0/20 11:35 0/20 2/20 20:16 1/20 25
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
83-32-9	Acenaphthene	ND	0.10	0.01	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.02	U
206-44-0	Fluoranthene	ND	0.10	0.02	U
87-68-3	Hexachlorobutadiene	ND	0.50	0.05	U
91-20-3	Naphthalene	ND	0.10	0.05	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.02	U
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.01	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U
218-01-9	Chrysene	ND	0.10	0.01	U
208-96-8	Acenaphthylene	ND	0.10	0.01	U
120-12-7	Anthracene	ND	0.10	0.01	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U
86-73-7	Fluorene	ND	0.10	0.01	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U
129-00-0	Pyrene	ND	0.10	0.02	U
91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U
87-86-5	Pentachlorophenol	ND	0.80	0.01	U
118-74-1	Hexachlorobenzene	ND	0.80	0.01	U
67-72-1	Hexachloroethane	ND	0.80	0.06	U



M-1 -1/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-04 : 275 ml	Date C Date F Date A Date E Dilutio Analys Instrur GC Cc %Solid	t Number Collected Received Analyzed Extracted n Factor st ment ID blumn ds on Volume	: 06/3 : 06/3 : 07/0 : 07/0 : 1 : DV : SV1 : RXI- : N/A	334 0/20 11:50 0/20 2/20 20:36 1/20 25
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
83.30.0	Assumption		0.40		
83-32-9	Acenaphthene	ND	0.10	0.01	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.02	U
206-44-0	Fluoranthene	ND	0.10	0.02	U
87-68-3 91-20-3	Hexachlorobutadiene	ND	0.50	0.05	U
	Naphthalene	ND	0.10	0.05	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.02	U
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.01	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U
218-01-9	Chrysene	ND	0.10	0.01	U
208-96-8	Acenaphthylene	ND	0.10	0.01	U
120-12-7	Anthracene	ND	0.10	0.01	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U
86-73-7	Fluorene	ND	0.10	0.01	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U
129-00-0	Pyrene	ND	0.10	0.02	U
91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U
87-86-5	Pentachlorophenol	ND	0.80	0.01	U
118-74-1	Hexachlorobenzene	ND	0.80	0.01	U
67-72-1	Hexachloroethane	ND	0.80	0.06	U



MT 7/14/20

AS NO. Parameter ug/L ug/L 83-32-9 Acenaphthene ND 0.01 0.01 U 81-58-7 2-Chioronaphthalene ND 0.20 0.02 U 206-44-0 Fluoranthene ND 0.10 0.02 U 87-68-3 Hexachlorobutadione ND 0.50 0.05 U 91-20-3 Naphthalene ND 0.10 0.02 U 56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 205-99-2 Benzo(bifuoranthene ND 0.10 0.01 U 207-08-9 Benzo(bifuoranthene ND 0.10 0.01 U 208-98-2 Benzo(bifuoranthene ND 0.10 0.01 U 207-08-9 Benzo(bifuoranthene ND 0.10 0.01 U 208-98-2 Benzo(bifuoranthene ND 0.10 0.01 U 208-99-2 Benzo(bifuoranthene ND 0.10 0.01	Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-05 : 275 ml	Date C Date R Date A Date E Dilutior Analys Instrum GC Co %Solid	t Number ollected ecceived nalyzed xtracted n Factor t ment ID lumn ts on Volume	: 06/30 : 07/02 : 07/0 ⁻ : 1 : DV : SV12 : RXI-5 : N/A	334 0/20 12:05 0/20 2/20 20:56 1/20 25
91-58-7 2-Chioronaphthalene ND 0.20 0.02 U 206-44-0 Fluoranthene ND 0.10 0.02 U 87-68-3 Hexachlorobutadiene ND 0.50 0.05 U 91-20-3 Naphthalene ND 0.10 0.05 U 56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 50-32-8 Benzo(a)pyrene ND 0.10 0.02 U 205-99-2 Benzo(h)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 208-98-2 Benzo(k)fluoranthene ND 0.10 0.01 U 208-98-8 Accenaphthylene ND 0.10 0.01 U 208-98-8 Accenaphthylene ND 0.10 0.01 U 208-98-8 Accenaphthylene ND 0.10 0.01 U 208-19-7 Anthracene ND 0.10 0.01 U 91-24-2 Benzo(ghi)perylene ND	CAS NO.	Parameter	Results		MDL	Qualifier
91-58-7 2-Chioronaphthalene ND 0.20 0.02 U 206-44-0 Fluoranthene ND 0.10 0.02 U 87-68-3 Hexachlorobutadiene ND 0.50 0.05 U 91-20-3 Naphthalene ND 0.10 0.05 U 56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 50-32-8 Benzo(a)pyrene ND 0.10 0.02 U 205-99-2 Benzo(h)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 208-98-2 Benzo(k)fluoranthene ND 0.10 0.01 U 208-98-8 Accenaphthylene ND 0.10 0.01 U 208-98-8 Accenaphthylene ND 0.10 0.01 U 208-98-8 Accenaphthylene ND 0.10 0.01 U 208-19-7 Anthracene ND 0.10 0.01 U 91-24-2 Benzo(ghi)perylene ND						
206-44-0 Fluoranthene ND 0.10 0.02 U 87-68-3 Hexachlorobutadiene ND 0.50 0.05 U 91-20-3 Naphthalene ND 0.10 0.02 U 56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 50-32-8 Benzo(a)pyrene ND 0.10 0.01 U 207-08-9 Benzo(b)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 208-73-7 Fluorene ND 0.10 <						
87-68-3 Hexachlorobutadiene ND 0.50 0.05 U 91-20-3 Naphthalene ND 0.10 0.05 U 56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 50-32-8 Benzo(a)pyrene ND 0.10 0.02 U 205-99-2 Benzo(t)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(t)fluoranthene ND 0.10 0.01 U 208-96-2 Benzo(t)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(t)fluoranthene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 208-97-7 Anthracene ND 0.10 0.01 U 91-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 91-27-7 Fluorene ND 0.10 <td></td> <td>•</td> <td></td> <td></td> <td></td> <td></td>		•				
91-20-3 Naphthalene ND 0.10 0.05 U 56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 50-32-8 Benzo(a)pyrene ND 0.10 0.02 U 205-99-2 Benzo(b)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 218-01-9 Chrysene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 20-12-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(gh)perylene ND 0.10 0						
56-55-3 Benzo(a)anthracene ND 0.10 0.02 U 50-32-8 Benzo(a)pyrene ND 0.10 0.02 U 205-99-2 Benzo(b)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 218-01-9 Chrysene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 201-2-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>						
50-32-8 Benzo(a)pyrene ND 0.10 0.02 U 205-99-2 Benzo(b)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 218-01-9 Chrysene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 201-27 Anthracene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.02 U 35-01-8 Phenanthrene ND 0.10 0.01						
205-99-2 Benzo(b)fluoranthene ND 0.10 0.01 U 207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 218-01-9 Chrysene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 120-12-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 </td <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>						
207-08-9 Benzo(k)fluoranthene ND 0.10 0.01 U 218-01-9 Chrysene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 120-12-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 85-01-8 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.02 U 129-00-0 Pyrene ND 0.10 0.02 U 129-00-0 Pyrene ND 0.10 0.02 U 14-57-6 2-Methylnaphthalene ND 0.10 0.01						
218-01-9 Chrysene ND 0.10 0.01 U 208-96-8 Acenaphthylene ND 0.10 0.01 U 120-12-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.02 U 129-00-0 Pyrene ND 0.10 0.02 U 118-74-1 Hexachlorophenol ND 0.10 0.02 U	205-99-2	Benzo(b)fluoranthene		0.10	0.01	
208-96-8 Acenaphthylene ND 0.10 0.01 U 120-12-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.02 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.10 0.02 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U
120-12-7 Anthracene ND 0.10 0.01 U 191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.02 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorobenzene ND 0.80 0.01 U	218-01-9	Chrysene	ND	0.10	0.01	U
191-24-2 Benzo(ghi)perylene ND 0.10 0.01 U 86-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 129-00-0 Pyrene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.10 0.02 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	208-96-8	Acenaphthylene	ND	0.10	0.01	U
B6-73-7 Fluorene ND 0.10 0.01 U 85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	120-12-7	Anthracene	ND	0.10	0.01	U
85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U
53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.01 U 193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	86-73-7	Fluorene	ND	0.10	0.01	U
193-39-5 Indeno(1,2,3-cd)pyrene ND 0.10 0.01 U 129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	85-01-8	Phenanthrene	ND	0.10	0.02	U
129-00-0 Pyrene ND 0.10 0.02 U 91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U
91-57-6 2-Methylnaphthalene ND 0.10 0.02 U 87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U
87-86-5 Pentachlorophenol ND 0.80 0.01 U 118-74-1 Hexachlorobenzene ND 0.80 0.01 U	129-00-0	Pyrene	ND	0.10	0.02	U
118-74-1 Hexachlorobenzene ND 0.80 0.01 U	91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U
	87-86-5	Pentachlorophenol	ND	0.80	0.01	U
67-72-1 Hexachloroethane ND 0.80 0.06 U	118-74-1	Hexachlorobenzene	ND	0.80	0.01	U
	67-72-1	Hexachloroethane	ND	0.80	0.06	U



MT 7/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-06 : 275 ml	Date C Date R Date A Date E Dilutior Analys Instrun GC Co %Solid	t Number collected received nalyzed xtracted n Factor t nent ID lumn Is on Volume	: 06/30 : 07/02 : 07/07 : 1 : DV : SV12 : RXI-1 : N/A	334 D/20 12:55 D/20 2/20 21:15 1/20 25
CAS NO.	Parameter	Results	ug/L RL	MDL	Qualifier
				_	
83-32-9	Acenaphthene	ND	0.10	0.01	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.02	U
206-44-0	Fluoranthene	ND	0.10	0.02	U
87-68-3	Hexachlorobutadiene	ND	0.50	0.05	U
91-20-3	Naphthalene	ND	0.10	0.05	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.02	U
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.01	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U
218-01-9	Chrysene	ND	0.10	0.01	U
208-96-8	Acenaphthylene	ND	0.10	0.01	U
120-12-7	Anthracene	ND	0.10	0.01	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U
86-73-7	Fluorene	ND	0.10	0.01	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U
129-00-0	Pyrene	ND	0.10	0.02	U
91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U
87-86-5	Pentachlorophenol	ND	0.80	0.01	U
118-74-1	Hexachlorobenzene	ND	0.80	0.01	U
67-72-1	Hexachloroethane	ND	0.80	0.06	U



MT 2/14/20

Client	: C.T. Male Associates	Lab Number	: L2027725
Project Name	: HHII TA1	Project Number	: 16.6334
Lab ID	: L2027725-07	Date Collected	: 06/30/20 00:00
Client ID	: FD01-200630	Date Received	: 06/30/20
Sample Location	SCHENECTADY NY	Date Analyzed	: 07/02/20 21:35
Sample Matrix	: WATER	Date Extracted	: 07/01/20
Analytical Method	: 1,8270D-SIM	Dilution Factor	1
Lab File ID	: 27725-07	Analyst	: DV
Sample Amount	: 275 mi	Instrument ID	: SV125
Extraction Method	: EPA 3510C	GC Column	: RXI-5SilM
Extract Volume	: 1000 uL	%Solids	: N/A
GPC Cleanup	: N	Injection Volume	: 1 uL
		ug/L	
5 NO.	Parameter	Results RL	MDL Qualifier

			uy/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.01	U	
91-58-7	2-Chloronaphthalene	ND	0.20	0.02	U	
206-44-0	Fluoranthene	ND	0.10	0.02	U	
87-68-3	Hexachlorobutadiene	ND	0.50	0.05	U	
91-20-3	Naphthalene	ND	0.10	0.05	U	
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.02	U	
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.01	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.01	U	
218-01-9	Chrysene	ND	0.10	0.01	U	
208-96-8	Acenaphthylene	ND	0.10	0.01	U	
120-12-7	Anthracene	ND	0.10	0.01	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.01	U	
86-73-7	Fluorene	ND	0.10	0.01	U	
85-01-8	Phenanthrene	ND	0.10	0.02	U	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.01	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.01	U	
129-00-0	Pyrene	ND	0.10	0.02	U	
91-57-6	2-Methylnaphthalene	ND	0.10	0.02	U	
87-86-5	Pentachlorophenol	ND	0.80	0.01	U	
118-74-1	Hexachlorobenzene	ND	0.80	0.01	U	
67-72-1	Hexachloroethane	ND	0.80	0.06	U	



MT 1/14/20

Client Project Name Lab ID Client ID Sample Location Sample Matrix Analytical Method Lab File ID Sample Amount Extraction Method Extract Volume GPC Cleanup	: 27725-08 : 275 ml		Date C Date R Date A Date E Dilution Analys Instrum GC Co %Solid	t Number collected teceived analyzed extracted n Factor at nent ID olumn	: 06/30 : 06/30 : 07/0 : 07/0 : 1 : DV : SV12 : RXI- : N/A	334 0/20 13:05 0/20 2/20 21:55 1/20 25
CAS NO.	Parameter		Results	ug/L RL	MDL	Qualifier
			nesults			
83-32-9	Acenaphthene		ND	0.10	0.01	U
91-58-7	2-Chloronaphthalene		ND	0.20	0.02	U
206-44-0	Fluoranthene		0.02	0.10	0.02	J
87-68-3	Hexachlorobutadiene		ND	0.50	0.05	U
91-20-3	Naphthalene		ND	0.10	0.05	U
56-55-3	Benzo(a)anthracene		ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene		ND	0.10	0.02	U
205-99-2	Benzo(b)fluoranthene		ND	0.10	0.01	U
207-08-9	Benzo(k)fluoranthene		ND	0.10	0.01	U
218-01-9	Chrysene		ND	0.10	0.01	U
208-96-8	Acenaphthylene		ND	0.10	0.01	U
120-12-7	Anthracene		0.02	0.10	0.01	J
191-24-2	Benzo(ghi)perylene		ND	0.10	0.01	U
86-73-7	Fluorene	0.10	0.02	0.10	0.01	~ U
85-01-8	Phenanthrene	0.10	0.03	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene		ND	0.10	0.01	U
193-39-5	Indeno(1,2,3-cd)pyrene		ND	0.10	0.01	U
129-00-0	Pyrene	_	ND	0.10	0.02	U
91-57-6	2-Methylnaphthalene		ND	0.10	0.02	U
87-86-5	Pentachiorophenol		ND	0.80	0.01	U
118-74-1	Hexachlorobenzene		0.03	0.80	0.01	J
67-72-1	Hexachloroethane		ND	0.80	0.06	U



MT 7/14/2

EXHIBIT 1

DEPARTMENT APPROVED POST REMEDIAL ACTION GROUNDWATER SAMPLING PLAN

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

Division of Environmental Remediation, Region 4 1130 North Westcott Road, Schenectady, NY 12306-2014 P: (518) 357-2045 | F: (518) 357-2460 www.dec.ny.gov

Ms. Susan McCann Hamilton Hill II Limited Partnership 90 State Street Suite 602 Albany, NY 12207 <u>smccann@tcbinc.org</u> *(Sent via email only)*

RE: Post Remedial Action Groundwater Sampling Plan Hamilton Hill II – Target Area 1 Site (Site No. C447052) 830 and 834 Albany Street, City/County of Schenectady

Dear Ms. McCann:

The New York State Department of Environmental Conservation has reviewed the subject work plan submitted electronically on April 27, 2020 by C.T. Male Associates on behalf of Hamilton Hill II Limited Partnership. The revised work plan is approved with the following modification: all drill cuttings originating below the depth of the remedial excavation should be drummed and disposed of following the Remedial Investigation Work Plan.

Please contact me at 518-357-2008 or joshua.haugh@dec.ny.gov if you have any questions about this letter.

Sincerely,

Josh Haugh Professional Geologist 1

ec: A. Fleck, DEC G. Burke, DEC K. Moline, C.T. Male S. Bieber, C.T. Male



Department of Environmental Conservation

Engineering, Surveying, Architecture, Landscape Architecture & Geology, D.P.C.

50 Century Hill Drive, Latham, NY 12110 518.786.7400 FAX 518.786.7299 www.ctmale.com

April 27, 2020

A700_ _____

Via Email

Joshua Haugh New York State Department of Environmental Conservation Division of Environmental Remediation 1130 North Westcott Road Schenectady, NY 12306 joshua.haugh@dec.ny.gov

RE: Post Remedial Action Groundwater Sampling Plan Hamilton Hill II – Target Area 1 Site (BCP Site No. C447052) 830 and 834 Albany Street City and County of Schenectady

Dear Mr. Haugh:

On behalf of Hamilton Hill II Limited Partnership, C.T. Male Associates Engineering, Surveying, Architecture, Landscape Architecture & Geology, D.P.C. (C.T. Male) has prepared this work plan for the Department's review for the collection of post-remedial action groundwater samples at the above referenced BCP Site No. C447052. The groundwater sampling is being conducted as a component of the November 2019 Decision Document for the Site.

The major elements of the remedial action were completed in April 2020. These included the remediation (by removal) of identified source areas of fill/soil mixtures containing contaminants at concentrations exceeding Unrestricted Use Soil Cleanup Objectives (SCOs), permanent closure (by removal) of underground bulk storage tanks, and backfilling of the fill/soil and tank excavations with imported fill material meeting Unrestricted Use SCOs.

The following presents the methods that will be employed for the installation of monitoring wells and collection of groundwater samples for laboratory analyses.

- Six (6) monitoring wells identified as RAMW1 to RAMW6 are proposed to be installed at the approximate locations shown on the attached two (2) figures. The first figure depicts the locations of the wells in relation to groundwater and soil vapor contaminants identified during the Remedial Investigation (RI). The second figure depicts the locations of the monitoring wells in relation to the proposed Site redevelopment, which is currently in full scale construction.

April 27, 2020 Joshua Haugh Page - 2

- Six (6) soil borings will be advanced at the approximate locations shown on the figures employing direct-push (Geoprobe) methods. Based on average groundwater depths obtained during the RI, the borings will each likely extend to approximately 15 to 20 feet bgs.

- One (1)-inch diameter monitoring wells consisting of Schedule 40 PVC screen (0.01 slot) and solid riser pipe will be installed in each boring such that the screened portion of each well straddles the water table. A sand pack will be placed around the well screens and the wells will be finished to grade with hydrated bentonite. The wells will be finished at the ground surface and will be protected with flush mounted curb boxes set in concrete pads.

- Each well will be developed and purged/sampled in accordance with the Field Sampling Plan presented as Appendix A in the Department approved December 2018 (Revised February 2019) Remedial Investigation Work Plan for the Site.

The following presents an explanation on the well locations compared to analytical detections identified within the remedial investigation phase to be above SCGs.

830 Albany Street Portion of the Site

- Proposed monitoring well RAMW1 will be installed in the vicinity of RI-installed monitoring well RIMW2 located in the west-central portion of the 830 Albany Street parcel. Both chloroform (8 ppb vs. SCG of 7 ppb) and tetrachloroethene (6.1 ppb vs. SCG of 5 ppb) were detected in groundwater at concentrations exceeding SCGs at RIMW2. RAMW1 will also serve as an upgradient monitoring well for the 830 Albany Street parcel.

- Proposed monitoring well RAMW2 will be installed in the vicinity of both RI-installed monitoring well RIGP1 and soil vapor probe RISV1 located in the northern portion of the 830 Albany Street parcel. Because of the new building footprint being over RIGP1, the replacement well location was positioned south of RIGP1. Chloroform (9.4 ppb vs. SCG of 7 ppb) was detected in groundwater above its SCG at RIGP1. Tetrachloroethene was detected in soil vapor at a concentration of 426 ug/m³ at RISV1, but was detected below its SCG of 5 ppb in groundwater at monitoring well RIGP1.

- Proposed monitoring well RAMW3 will be installed in the vicinity of RI-installed monitoring wells RIGP2 and RIGP3 in the eastern portion of the 830 Albany Street parcel. Both chloroform (9.4 ppb vs. SCG of 7 ppb) and tetrachloroethene (6.9 ppb vs. SCG of 5 ppb) were detected in groundwater at concentrations exceeding SCGs at RIGP2. Tetrachloroethene (8.6 ppb vs. SCG of 5 ppb) was detected in groundwater at a concentration exceeding its SCG at RIGP3.

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- Proposed monitoring well RAMW4 will be installed in the eastern portion of the 830 Albany Street parcel in the vicinity of MW 2, which was installed during previous Phase II ESAs conducted on the Site. Chloroform (9.8 ppb vs. SCG of 7 ppb) and seven (7) SVOCs (0.03 ppb to 0.11 ppb concentration range) were detected in groundwater above SCGs when this well was sampled during the RI. RAMW4 will also serve as a downgradient monitoring well for the 830 Albany Street parcel.

834 Albany Street Portion of the Site

- Proposed monitoring well RAMW5 will be installed in the vicinity of RI-installed RIGP4 and RIGP5 in the northern portion of the 834 Albany Street parcel. Five (5) SVOCs (0.02 ppb to 0.04 ppb concentration range) were detected in groundwater at concentrations exceeding SCGs at RIGP5. There were no SVOCs detected above SCGs at RIGP4. RAMW5 will serve as an upgradient well for the 834 Albany Street parcel.

- Proposed monitoring well RAMW6 will be installed in the vicinity RI-installed RIMW5 along the eastern boundary of the 834 Albany Street parcel. Three (3) SVOCs (0.02 ppb to 0.06 ppb concentration range) were detected at concentrations exceeding SCGs at this well location. RAMW6 will serve as a downgradient well for the 834 Albany Street parcel.

Analytical Requirements

- The groundwater samples will be analyzed for the Target Compound List (TCL) for volatile organic compounds (VOCs) and semi-volatile organic compounds (SVOCs).**

** C.T. Male requests the Department to waive the requirement for testing for metals during this post remedial action sampling event. The RI clearly demonstrated that metals were not contaminants of concern that drove the remedial action, and therefore should not require further testing especially with the removal of soils with contaminants above unrestricted use (and groundwater protection) SCOs. As shown on the first figure attached herein, sodium (which is attributed to the application of snow/ice melt on surrounding roadways) was the only metal predominantly detected above SCGs during the RI. Antimony (4.24 ppb) was detected slightly above its SCG of 3 ppb in one (1) groundwater sample only collected from RI sampled monitoring well RIMW4. The isolated manganese and iron detections above SCGs at RI-sampled monitoring wells RIMW4D, RIMW6D, MW1 and MW2 were viewed as naturally occurring and not contaminants of concern.

- One (1) set of quality control (QC) samples will be collected. These samples will include a trip blank, duplicate sample, matrix spike, matrix spike duplicate and equipment blank.

April 27, 2020 Joshua Haugh Page - 4

- The analytical results will be reported as a Category B deliverable for DUSR validation by an independent third party data validator.

Well Decommissioning and Reporting

Upon receipt of the analytical results from this sampling of the additional wells described herein, C.T. Male will prepare a summary report presenting the validated analytical data in tabular format. The results will be compared to NYSDEC Ambient Water Quality Standards for an evaluation on the overall groundwater quality. This comparison will be used to determine when the monitoring wells can be decommissioned.

The monitoring wells will be decommissioned in accordance with DEC Policy CP-43: Groundwater Monitoring Well Decommissioning Policy, dated November 3, 2009. The preferred method for decommissioning will be removal of the protective enclosures followed by casing pulling/grouting. If the well casing cannot be pulled, then the well will be grouted to the surface. Department approval will be obtained prior to decommissioning of the monitoring wells.

Waste Management

Wastes that are anticipated to be generated from installation of monitoring wells and groundwater sampling include drill cuttings, groundwater development and purge water, solid wastes consisting of disposable macro-core acetate liners, and personal protective (i.e., nitrile gloves) and groundwater sampling (i.e., tubing, wipes, etc.) equipment. The Site has been remediated to Unrestricted Use SCOs and soil imported onto the Site as backfill has been tested and meets Unrestricted Use SCOs.

Drill cuttings that are not saturated with groundwater will be disposed of at the ground surface in the vicinity of the soil boring that they were generated from. Drill cuttings that are saturated with groundwater will be placed in a 55-gallon drum for future profiling and ultimate off-site disposal at a permitted treatment, storage and disposal facility (TSDF). Groundwater development and purge water will be placed in a 55-gallon drum for future profiling and ultimate off-site disposal at a permitted TSDF. The drums containing the drill cuttings and water will be staged at a location within the Site that is not anticipated to be disturbed by redevelopment activities.

The disposable sampling and personal protective equipment will be disposed of as solid waste.

April 27, 2020 Joshua Haugh Page - 5

Please contact the undersigned should you have any questions regarding this work plan. At this point, we are anticipating the installation of the monitoring wells to take place after the building foundations have been constructed for the proposed building within the 830 Albany Street parcel. Excavation for the installation of the foundations would likely destroy proposed monitoring wells RAMW2, RAMW3 and RAMW4 if these were installed beforehand.

Respectfully, C.T. MALE ASSOCIATES

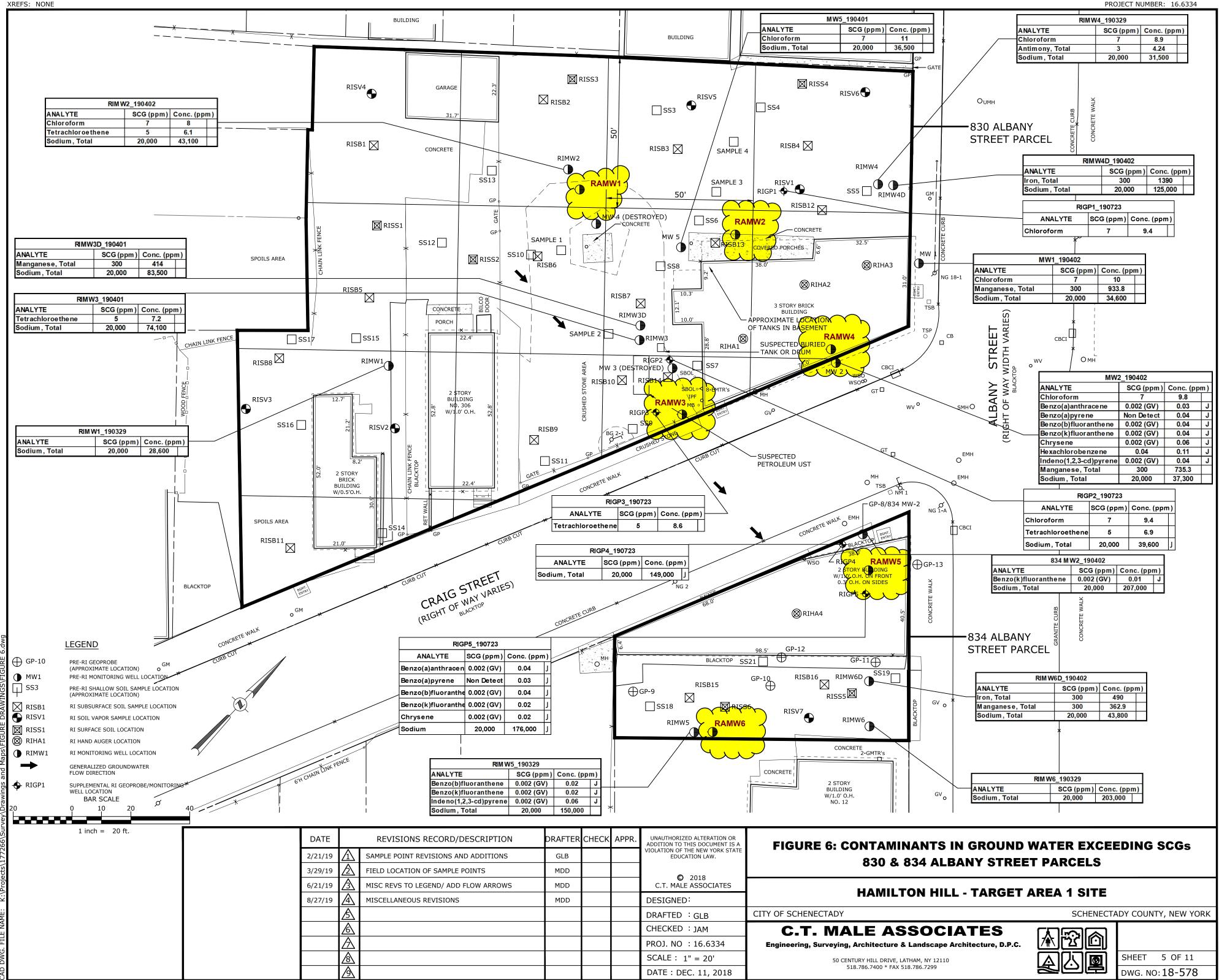
Step

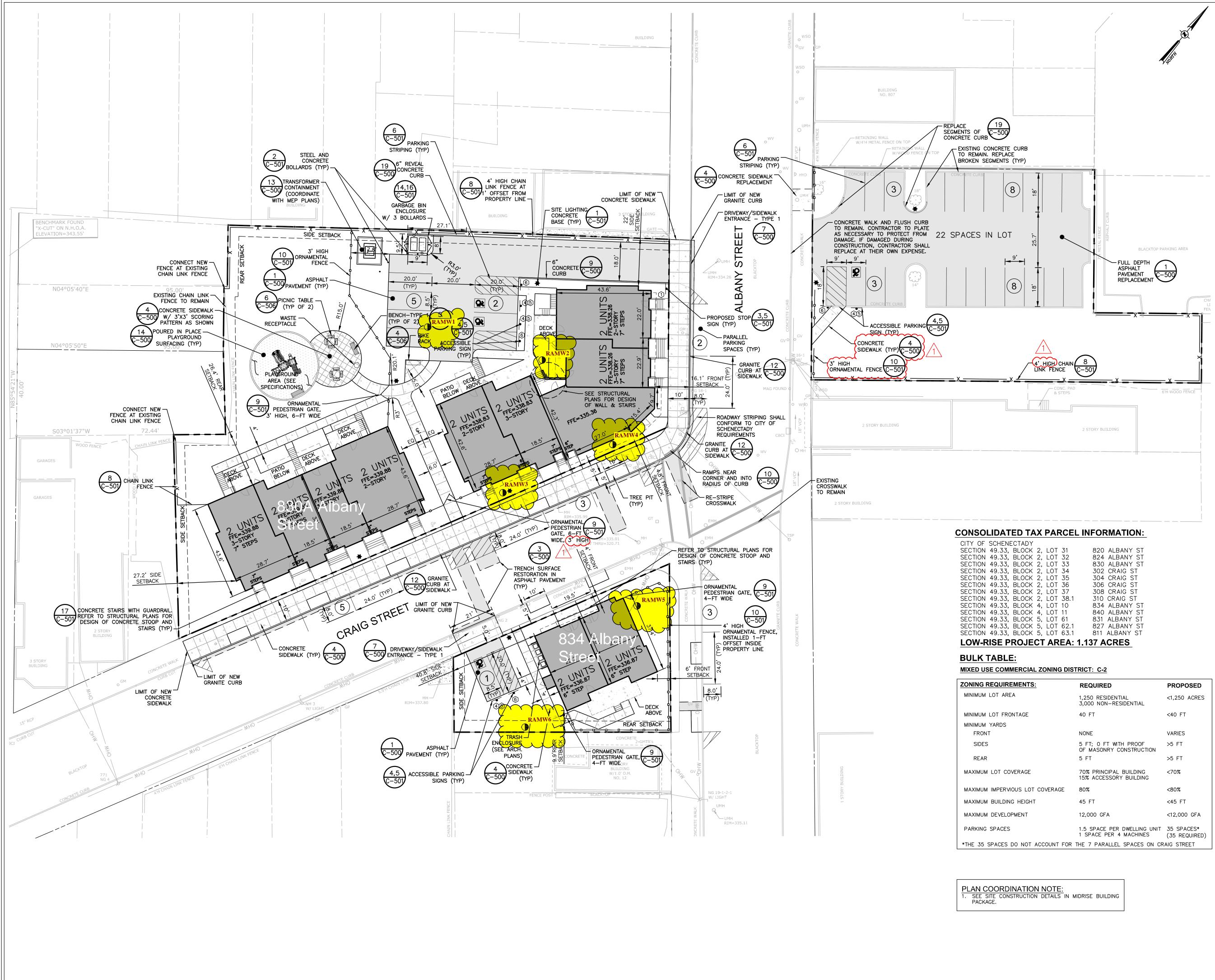
Stephen Bieber, CHMM Project Scientist

Attachments

ec: Kelly Melarango, Hamilton Hill II Limited Partnership Janis Stewart, Hamilton Hill II Limited Partnership Jeff Marx, P.E., C.T. Male Associates Kirk Moline, P.G., C.T. Male Associates







	820 /	ALBANY ST
	824 /	ALBANY ST
	830 /	ALBANY ST
	302 (CRAIG ST
	304 (CRAIG ST
	306 (CRAIG ST
	308 (CRAIG ST
1	310 (CRAIG ST
	834 /	ALBANY ST
	840 /	ALBANY ST
	831 A	ALBANY ST
1	827 /	ALBANY ST
1	811 A	LBANY ST
•	1.137 AC	RES

	REQUIRED	PROPOSED			
	1,250 RESIDENTIAL 3,000 NON-RESIDENTIAL	<1,250 ACRES			
	40 FT	<40 FT			
	NONE	VARIES			
	5 FT; 0 FT WITH PROOF OF MASONRY CONSTRUCTION	>5 FT			
	5 FT	>5 FT			
	70% PRINCIPAL BUILDING 15% ACCESSORY BUILDING	<70%			
	80%	<80%			
	45 FT	<45 FT			
	12,000 GFA	<12,000 GFA			
	1.5 SPACE PER DWELLING UNIT 1 SPACE PER 4 MACHINES	35 SPACES* (35 REQUIRED)			
۲	THE 7 PARALLEL SPACES ON CRAIG STREET				

SITE LEGEND: $O \cup O$ ()>00 $O \to O$ PROPERTY LINE -----_ _ _ <u>SETBACK LINE</u> __ _ SETBACK LINE ர*்ன்* வல ⊔щ́≦́щ́⊇ BUILDING DRAU BURAU BURAU BURAU BURAU PAVEMENT SIDEWALK CURB CHAIN LINK FENCE CONCRETE SURFACE ROAD LINING & STRIPING -----(2) SIGN PARKING COUNT, SHOWN FOR DESIGN PURPOSES ONLY PROPOSED LIGHT POLE SITE PLAN NOTES:

GENERAL CONSTRUCTION:

- THE CONTRACTOR SHALL PROTECT EXISTING PROPERTY LINE MONUMENTATION. ANY MONUMENTATION DISTURBED OR DESTROYED, AS JUDGED BY THE ENGINEER OR OWNER, SHALL BE REPLACED AT THE CONTRACTOR'S EXPENSE AND UNDER THE SUPERVISION OF A NEW YORK STATE LICENSED LAND SURVEYOR.
- 2. ALL PAVEMENT RESTORATION SHALL MEET AND MATCH EXISTING GRADES. ALL SAWCUT LINES SHALL BE PARALLEL AND CURVILINEAR TO EXISTING OR PROPOSED CURBING AND SHALL BE A CONSTANT
- DISTANCE OF 18" MIN AWAY. ALL ARCHITECTURE IS SUBJECT TO PLANNING BOARD REVIEW. 5. NOTIFY ENGINEER 48 HOURS PRIOR TO INITIALIZATION OF ANY WORK ON SITE.
- 6. THE ENGINEER SHALL BE NOTIFIED IN WRITING OF ANY CONDITIONS THAT VARY FROM THOSE SHOWN ON THE PLANS. THE CONTRACTOR'S WORK SHALL NOT VARY FROM THE PLANS WITHOUT PRIOR REVIEW FROM THE ENGINEER. CONTRACTOR IS RESPONSIBLE FOR EMPLOYING AND MAINTAINING ALL
- TRAFFIC CONTROL AND SAFETY MEASURES DURING CONSTRUCTION. CONTRACTOR IS RESPONSIBLE FOR PROPERLY & SAFELY MAINTAINING
- AREA BETWEEN ALL ADJOINING PROPERTIES. NO WORK, STORAGE OR TRESPASS SHALL BE PERMITTED BEYOND THE SITE PROPERTY LINES OR PUBLIC RIGHT-OF-WAY. 10. ALL EXISTING LAWN AREA, CURBING, PAVING, SIDEWALKS, CULVERTS
- OR OTHER PUBLIC OR PRIVATE PROPERTY DAMAGED BY TRENCHING OR EXCAVATION OPERATIONS SHALL BE REPLACED OR REPAIRED TO A CONDITION EQUAL TO EXISTING, AS DESCRIBED IN CONTRACT DOCUMENTS OR AS ORDERED BY ENGINEER (AOBE). MAILBOXES, SIGN POSTS, ETC SHALL BE PROTECTED OR REMOVED AND REPLACED EXACTLY AS THEY WERE BEFORE BEING DISTURBED. REMOVE AND REPLACE AFFECTED CURBING AND SIDEWALK TO NEAREST JOINT. REMOVE PAVEMENT AND REPLACE TO SAW CUT LINE, SAW CUT IN STRAIGHT LINE TO POINT NEEDED TO BLEND GRADE, REMOVE LAWN AND REPLACE TO MINIMUM LIMIT OF EXCAVATION.

LAYOUT

- BUILDING DIMENSIONS TO BE TAKEN FROM ARCHITECTURAL BUILDING PLANS. NOTIFY THE ENGINEER OF ANY DEVIATION FROM CONDITIONS SHOWN ON THIS PLAN.
- THE CONTRACTOR SHALL BE RESPONSIBLE FOR COORDINATING ALL FIELD LAYOUT. THE CONTRACTOR SHALL TAKE TIES TO ALL UTILITY CONNECTIONS AND PROVIDE MARKED-UP AS BUILT PLANS FOR ALL UTILITIES SHOWING TIES TO CONNECTIONS, BENDS, VALVES, LENGTHS OF LINES AND INVERTS. AS-BUILT PLANS SHALL BE REVIEWED BY THE OWNER AND THE ENGINEER AND THE CONTRACTOR SHALL PROVIDE ANY CORRECTION OR ADDITIONS TO THE SATISFACTION OF THE OWNER AND THE ENGINEER BEFORE UTILITIES WILL BE ACCEPTED.

PAVING:

- NO VEHICULAR TRAFFIC OF ANY SORT SHALL BE PERMITTED ON THE SURFACE OF SUBBASE COURSE MATERIAL ONCE IT HAS BEEN FINE GRADED, COMPACTED, AND IS READY FOR PAVING. SUBBASE MATERIAL SO PREPARED FOR PAVING SHALL BE PAVED WITHIN THREE DAYS OF PREPARATION.
- 2. SUBBASE MATERIAL AND THE VARIOUS ASPHALT CONCRETE MATERIALS CALLED FOR IN THESE DRAWINGS SHALL CONFORM WITH THE REFERENCED SECTION OF THE NEW YORK STATE DEPARTMENT OF TRANSPORTATION STANDARD SPECIFICATIONS FOR CONSTRUCTION AND MATERIALS, DATED "LATEST EDITION". CONSTRUCTION SHALL BE AS FURTHER SET FORTH IN THOSE SPECIFICATIONS AND AS OTHERWISE PROVIDED FOR IN THESE DRAWINGS. 3. PLACE ASPHALT CONCRETE MIXTURE ON PREPARED SURFACE.
- SPREAD AND STRIKE-OFF USING A SELF-PROPELLED PAVING MACHINE, WITH VIBRATING SCREED. PLACEMENT IN INACCESSIBLE AND SMALL AREAS MAY BE BY HAND 4. PROVIDE JOINTS BETWEEN OLD AND NEW PAVEMENTS OR BETWEEN
- SUCCESSIVE DAY'S WORK. 5. TACK COAT WHEN SPECIFIED OR CALLED OUT ON THE DRAWINGS OR REQUIRED BY THE REFERENCED SPECIFICATION SHALL CONFORM WITH THE FOLLOWING
- A. TACK COAT SHALL MEET THE MATERIAL REQUIREMENTS OF 702-90 ASPHALT EMULSION FOR TACK COAT OF THE NEW YORK STATE DEPARTMENT OF TRANSPORTATION STANDARD SPECIFICATIONS FOR CONSTRUCTION AND MATERIALS, DATED "LATEST EDITION", SHALL BE APPLIED IN ACCORDANCE WITH SECTION 407 - TACK COAT SHALL BE IN ACCORDANCE WITH THOSE SPECIFICATIONS AND AS OTHERWISE PROVIDED FOR IN THESE DRAWINGS.
- B. REMOVE LOOSE AND FOREIGN MATERIAL FROM ASPHALT SURFACE BEFORE PAVING NEXT COURSE. USE POWER BROOMS, BLOWERS OR HAND BROOM. C. APPLY TACK COAT TO ASPHALT PAVEMENT SURFACES & AND
- SURFACES OF CURBS, GUTTERS, MANHOLES, AND OTHER STRUCTURES PROJECTING INTO OR ABUTTING PAVEMENT. DRY TO A "TACKY" CONSISTENCY BEFORE PAVING. D. TACK COAT ENTIRE VERTICAL SURFACE OF ABUTTING EXISTING
- PAVEMENT. 6. AFTER COMPLETION OF PAVING AND SURFACING OPERATIONS, CLEAN SURFACES OF EXCESS OR SPILLED ASPHALT, GRAVEL OR STONE

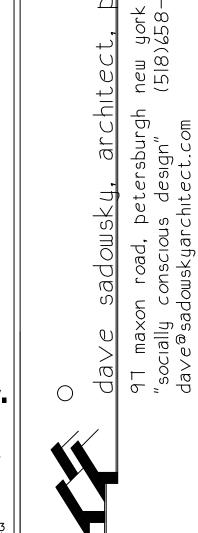
MATERIALS TO THE SATISFACTION OF THE ENGINEER. STRIPING:

- STRIPE PAVEMENT AS INDICATED ON THE PLANS AND/OR IN ACCORDANCE WITH ALL APPLICABLE LOCAL, STATE AND FEDERAL REQUIREMENTS
- 2. COLOR: DRIVE LANE DIVIDERS WHITE OR AOBE NO PARKING ZONE WARNINGS - WHITE OR AOBE PARKING DIVIDERS - WHITE OR AOBE WALKING LINES - WHITE OR AOBE
 - HANDICAP PARKING LINES & SYMBOL BLUE
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ORIGINAL SCALE IN INCHES



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