Remedial Design Work Plan

Farmingdale Plaza Cleaners

Farmingdale Nassau County New York

NYSDEC Site No. 130107



Prepared by New York State Department of Environmental Conservation Division of Environmental Remediation 625 Broadway Albany, NY 12233-7015

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1.0 Introduction

New York State Department of Environmental Conservation (NYSDEC) has developed this work plan to complete the remediation for the Farmingdale Plaza Cleaners (FPC) site (referred to herein as the site). The site is located at 450 Main Street in the Village of Farmingdale, Town of Oyster Bay, Nassau County, New York (see Figure 1).

Site activities are being conducted under the State Superfund Program. The site is divided into two operable units. Operable Unit 1 (OU1) consists of contaminated soil and soil vapor and includes the plaza property and portions of adjacent properties. An Interim Remedial Measure was performed in November 2011 for OU1 to address on-site soil and soil vapor contamination. A Record of Decision, dated March 30, 2012, was issued by NYSDEC for OU1 and required continued operation of a soil vapor extraction system to remediate contamination within onsite soil vapor and soils. OU1 activities are presently being performed by the current site owner. Operable Unit 2 (OU2) addresses on-site and off-site groundwater. A Record of Decision was issued by the NYSDEC on March 25, 2014, which discusses the groundwater contamination migrating from the site and the selected remedial action to install a groundwater extraction well down-gradient of the site to capture the groundwater contamination originating from the site. The potentially responsible parties (PRPs) failed to sign an Order on Consent and the FPC site was referred to the State Superfund for remedial activities. As a result the NYSDEC intends to design and implement the selected remedial action for OU2.

2.0 Background

Site background information regarding physical setting and summary of historical investigation activities at the LIF site and FPC site are presented below.

2.1 Physical Setting & Site Uses

Location: The Farmingdale Plaza Cleaners Site is located at 450 Main Street in Farmingdale, Nassau County. The site is in a suburban area and lies at the intersection of Main Street and Fulton Street. This site is 1,000 feet up-gradient of the Liberty Industrial Finishing (LIF) site, which is a United States Environmental Protection Agency (USEPA) National Priorities List (NPL) site. The sites are shown on Figure 1.

Site Features: The site is a commercial plaza that formerly housed the Farmingdale Plaza Cleaners. The shopping plaza consists of a single 33,000-square foot building and a paved parking lot, surrounded by apartment buildings and other commercial structures.

Current Zoning/Use(s): The former dry cleaner and adjacent former supermarket are no longer present. The site building was recently renovated and a CVS is presently occupying the southern portion of the building. The shopping plaza is zoned for commercial use. The surrounding parcels are zoned residential or commercial.

Past Use(s) of the site: Waldbaum Shopping Plaza was reportedly constructed in 1983, at which time the Farmingdale Plaza Cleaners (FPC) began operation. Dry cleaning operations ceased in August 2003.

2.2 Site Geology

Site investigations encountered the Upper Glacial aquifer (UGA) and Magothy aquifer (MA). The nature of the overburden at the site was characterized and presented in the FPC Remedial Investigation report. The total overburden thickness in the area is unknown as monitoring wells and borings were drilled to a maximum depth of 200 feet below ground surface; however, the deepest overburden water supply well in the area is over 750 feet deep.

Site investigations determined that the UGA was observed to be approximately 85 to 100 feet thick and primarily consisted of gravelly sand, sometimes with a trace of silt. The lower portion of the UGA consisted of fine-grained sand, silt, and clay but sometimes also contained a trace of gravel. The soils changed from brown/tan to dark brown/gray. These finer grained soils of varying thickness were generally recognizable in gamma logs from the profile borings and appear to correlate with a unit identified as the "20-foot-clay" described in regional geologic literature (URS 2000; Perlmutter and Geraghty 1962).

Site investigations determined that the MA soils consisted of interbedded medium- to coarsegrained sands silty sand, clayey silt, and silty clay. Gamma logs showed layers containing a relatively high clay content compared to other zones, which were of variable depth and thickness.

Figure 2 presents a geologic cross-section depicting the subsurface stratigraphy. This cross section presents site information from the north to the south. This section depicts the transition from the UGA to the MA at approximate depths ranging from 100 to 110 feet below ground surface.

2.3 Site Hydrogeology

The Remedial Investigation assessed groundwater flow within the UGA and MA. In March 2012, water level measurements were obtained from 61 monitoring wells. Figures 3 and 4 present the groundwater elevations and interpreted isopleths for the UGA and MA, respectively. Groundwater flow is to the south for both aquifers. In both the UGA and MA, the overall horizontal gradient was measured to be approximately 0.2% to the south. The vertical hydraulic gradient varies within the limits of the investigation area, but in general a downward gradient does appear to be present between the UGA and MA. Vertical migration groundwater will be limited by the fine-grained transitional unit between the UGA and MA.

The LIFS CRI report (URS 2000) indicated that while groundwater movement is predominantly horizontal in the UGA, upward or downward gradients exist within the MA and also between the two aquifers on a seasonal basis.

2.4 Liberty Industrial Finishing Investigation Findings

Investigations were initiated in 1990's by the USEPA at the adjacent LIF site. A 2000 remedial investigation conducted for the USEPA at the LIF site (URS 2000) indicated that the Farmingdale area had been impacted by two Volatile Organic Compound (VOC) contaminant plumes (designated Plume A and Plume B). Plume A consisted of trichloroethene (TCE) and was determined to originate from the LIF site, while the tetrachloroethene (PCE)-dominated Plume B was identified to originate from an up-gradient source that was later identified as the FPC site. Since then, the LIF site has been the subject of a number of environmental investigations over several years by USEPA to define the limits of the plumes. Investigation reports concluded that:

- Based on potentiometric surface elevations, groundwater flow is primarily toward the south with a clear downward vertical hydraulic gradient existing in the UGA near the FPC site; This is further supported by current data presented in Table 1, which was collected at monitoring well clusters.
- Primary VOCs detected during this investigation were PCE, TCE, and cis-1,2-DCE;
- VOC contamination detected within the investigation area originates from multiple potential sources including the LIF site, FPC site, and possibly an unknown source(s) upgradient of the FPC site;

- Plume A consists of cadmium (Cd), chromium (Cr), TCE, and its daughter products extending from the LIF site southward towards the Southern State Parkway. Plume A contamination is located within the UGA and MA. In August 1998, the USEPA issued a unilateral administrative order to the LIF site PRPs to initiate an interim groundwater action. This action ultimately resulted in construction and operation of an on-site groundwater pump and treat system operated as a non-time critical removal action. The off-site groundwater system has numerous components. Groundwater is extracted from the UGA by three recovery wells and from the MA by three additional wells that have an estimated capture width of approximately 940 feet. In addition, at the far end of the cadmium and chromium plume, a single UGA extraction well operates with a design capture width of approximately 500 feet. Figure 5 illustrates the approximate extent of Plume A;
- Plume B (identified as originating from the FPC site and possibly unknown source(s) north of the FPC site) consists of PCE and its daughter products within UGA and MA. Plume B migrates from the FPC site and across the LIF site in a southerly direction with a slightly westward component. Downgradient of the FPC site, higher levels of contamination at depth suggest a general dipping of the contaminant plumes as they move southward along the direction of groundwater flow. The downward dipping trend observed in the plumes is likely the result of the downward hydraulic gradient. Figure 5 illustrates the approximate extent of Plume B;
- Based on the site investigations, Plumes A and B have comingled making delineation difficult (see Figure 5); and
- As a result of data collected during the USEPA investigation, the FPC site was listed as a Class 2 site on the State's Registry of Inactive Hazardous Waste Disposal Sites in December 2002.

2.5 Farmingdale Plaza Cleaners OU 2 Remedial Investigation

In March 2012, the third and final mobilization of the remedial investigation was completed to define the extent of contamination migrating from the FPC site, identified as Plume B. A total of 39 samples were collected from new and existing wells/piezometers, including the six new permanent monitoring wells, 16 existing permanent monitoring wells/piezometer, and 17 existing profile well channels. Each of the existing profile wells sampled (PW-01, PW-03, PW-04, PW-06, PW-08, PW-09, and PW- 11) contains seven "channels" that allow the well to be sampled at varying depths. Only a few channels of each profile well were sampled for this task for comparison with results when the profile wells were installed (YU 2009). Site findings are presented below:

- Groundwater flows in a southerly direction from the site;
- Groundwater migrates vertically downwards through the formation;

- On-site PCE within the UGA was detected at a maximum concentration of 21 parts per billion (ppb) at 40 feet below ground surface (bgs). Most of the PCE concentrations in the on-site groundwater wells are below 5 ppb. The PCE plume shape downgradient of the site was interpreted to be discontinuous based on the locations of the detections exceeding Class GA standards (5 ppb for PCE and TCE) in wells north of Motor Avenue;
- Off-site PCE within the UGA was detected at a maximum concentration of 38 ppb in PW-09 at 94 feet bgs. Most off-site wells did not detect any PCE contamination, with a few wells showing PCE concentrations between 10 ppb to 20 ppb. The portion of the PCE plume with highest concentrations within the UGA appears to be between Motor Avenue and Yoakum Avenue;
- Off-site PCE within the MA was detected at a maximum concentration of 130 ppb in MW-37C at 119 feet bgs. This detection is within the comingled portion of Plumes A and B. The highest concentration of PCE off-site and not located within Plume A was detected at a maximum concentration of 110 ppb in PW-19 at 145 feet bgs, which is located towards the leading edge of the plume. Most of the other groundwater monitoring wells showed PCE concentrations from non-detect to less than 100 ppb;
- The PCE plume widens as it extends southward until the plume ends abruptly at Tomes Avenue (based on the lack of PCE in MW-31C, -31D, and -47C);
- The extent of PCE and TCE contamination in the UGA and MA has decreased compared to data from the sampling completed in 2008 and remains similar to data from LIF site monitoring program in 2011. Reduction in the extent of TCE contamination in the UGA is attributed to the construction and operation of the LIF site groundwater treatment system and the FPC OU-1 vapor extraction system, in addition to natural degradation processes including, but not limited to, dispersion, dilution, and reductive dechlorination. The overall reduction and limitation to the horizontal extent of PCE contamination may be attributed to elimination of the source at the FPC site and natural degradation processes;
- The uncaptured portion of Plume B extends approximately 500 feet east of the capture zone for the LIF site/Plume A at MW- 48C. (see figure 5)
- The base of both the PCE and TCE plumes is interpreted to end at an approximate depth of 200 feet at a clay layer detected in numerous groundwater profile borings;
- PCE is shown to be present up-gradient of the FPC site in monitoring well identified as MW-3. The concentration was 21 ppb in the UGA;
- One monitoring well (EPA-MW-1B) is installed in the MA near the site and sample results from 2007 show the PCE levels to be below Class GA standards in this area;
- The investigation showed that there is no on-going source at the site, since most of the higher PCE concentrations are found at the lead edge of the plume;

- Other breakdown products were also observed within the UGA and MA, but the data are inconclusive as to whether their source is the FPC site or the nearby LIF site; and
- Figure 2 represents a generally north-south cross section. This figure shows the extent of PCE in the UGA and MA from the site to the leading edge of the Plume B.

3.0 Pre-Design Investigation

On August 25, 2015, a sampling event was completed to provide current groundwater flow and VOC concentrations at the lead edge of the plume. This information will be used to design the FPC site groundwater extraction system.

The depth to groundwater was recorded at 19 wells during the sampling event (see Table 1). The groundwater flow direction was determined to be towards the south (see Figure 6).

Groundwater samples were collected at five monitoring wells, identified as MW-28C, MW-28D, MW-37C, MW-46C, and MW-47C. Samples were obtained using passive diffusion bags. Samples were analyzed for VOCs via method 8260 (see Appendix B for laboratory results). Table 2 provides a summary of the analytical results. Highlighted Values indicate exceedances of NYSDEC Groundwater Standards. Samples show levels above Class GA standards for both PCE and TCE (see Figure 7). The analytical results are similar to 2012 findings for these wells, which identified the leading edge of the plume to be near MW-47C.

4.0 Remedial Action Plan

According to the OU 2 Record of Decision, the selected remedial action was groundwater extraction and treatment. This decision was based on FPC site Remedial Investigation Report and the criteria identified for evaluation of remedial alternatives. The FPC groundwater extraction system will be designed and installed so that the capture zone is sufficient to intercept the areal and vertical extent of the elevated PCE contamination that is not currently remediated by the LIF site groundwater extraction system. The LIF site information was used to develop the conceptual design for the FPC site.

A comparison of soil log from LIF extraction wells RW-8 and RW-9 and gamma logs along Tomes Ave (MW-25C, MW-31D, and MW-47C) indicate that the soil conditions are similar. See Table below for log interpretations. The gamma logs show similar material within the upper 140 feet of the logs. MW-25 did not extend beyond that depth, but a soil log was available for RW-9 which extended to 195 feet. Based on the soil log the increased clay zone at approximately 145 feet was not encountered at the LIF recovery well location.

CPS Gamma Readings	well 25C	Well 31D	Well 47C
~ 20 CPS	0 feet	0 feet	0 feet
~ 35 CPS	70 feet	70 feet	75 feet
~ 70 CPS	*	140 feet	150 feet
~ 25 CPS	*	163 feet	194feet

Soil Logs

*Gamma logs ended at ~140 feet.

CPS (counts per second).

Two different gamma loggers were used.

The soil boring logs for these wells are provided in Appendix A. (MW-25 is also identified as OS-15)

In August 1998, the USEPA issued a unilateral administrative order to the LIF site PRPs to initiate an interim groundwater action. The LIF extraction well system was designed based on modelling by Ecology and Environment Engineering, P.C. This design required three extraction wells located within the MA with screen depths ranging from approximately 90 to 180 feet below ground surface (ft bgs) and a total purge rate of 85 gallons per minute, see table below for individual extraction well specifications. The designed capture width was approximately 940 feet. The influence of the LIF groundwater extraction system is shown on Figure 8.

Extraction Well	Design Flow	Screen Top	Screen Bottom	Screen Length
	(gpm)	(ft bgs)	(ft bgs)	(ft)
RW-8	30	94	139	45
RW-9	20	154	184	30
RW-10	35	99	139	40

The FPC groundwater extraction system will be designed to capture the PCE and associated breakdown products at the leading edge of the FPC site plume (Plume B) to minimize any further down-gradient migration of the contaminated groundwater. Based on the LIF pump test results, a similar system shall be considered for the FPC site. The proposed extraction well shall be located near MW-47C and is anticipated to provide a capture width of 400 feet, see Figure 8. The depth of the proposed extraction well is approximately 150 feet bgs and the length of the extraction well screen will be approximately 20 feet (see Figure 9 Vertical Placement of Groundwater Extraction Well). The base of the well screen will be located just above the fine grained soils.

The FPC groundwater extraction system will be integrated with the existing LIF site groundwater extraction and treatment system, or discharged to the Nassau County sewer system. The proposed FPC groundwater extraction well is located approximately 1,500 feet from the existing LIF site extraction well system. The underground piping conveyance is anticipated to run from the FPC extraction well adjacent to MW-47C, along Tomes Avenue, to either the LIF site extraction well system (see Figure 10) or a Nassau County sewer main. This will permit groundwater discharged from the proposed FPC extraction well to be conveyed to an acceptable treatment facility.

5.0 Remedial Design Tasks

Based on the remedial investigation and evaluation of remedial alternatives, the NYSDEC selected groundwater extraction and treatment as the selected remedial action. The FPC groundwater extraction system will be incorporated into the existing LIF site groundwater extraction and treatment system or discharged to the Nassau County sewer system. Discussions with Nassau County regarding discharge requirements shall be performed during the design process. This will capture PCE and associated breakdown products at the leading edge of Plume B that is not presently being captured by the LIF groundwater extraction system. The installation of the FPC extraction well will minimize any further down-gradient migration of the PCE contaminated groundwater. A NYSDEC Engineering Contractor shall be retained to develop design documents. The design elements will include: groundwater recovery (extraction) well and vault construction, pump specifications, initial pump test of extraction well, utilities, and pipe conveyances. Extracted groundwater will not be treated but simply conveyed to a treatment system.

6.0 References

Ecology and Environment Engineering, P.C. February 2008. *Final Engineering Design Report for the Liberty Industrial Finishing Superfund Site Farmingdale, Nassau County, New York.*

Ecology and Environment Engineering, P.C. February 2012. Final Remedial Investigation Report for the Farmingdale Plaza Cleaners Site, Operable Unit 2 Farmingdale, Nassau County, New York.

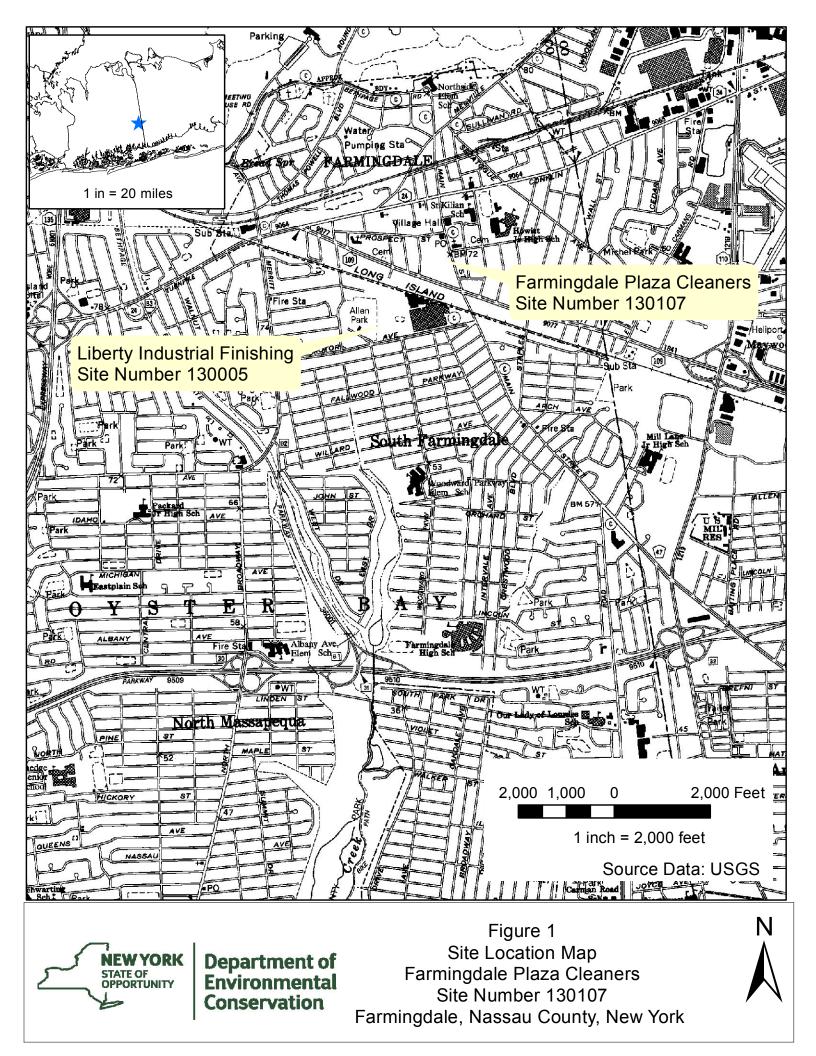
New York State Department of Environmental Conservation March 2014. *Record of Decision Farmingdale Plaza Cleaners Operable Unit Number 02: On-site and Offs-site Groundwater State Superfund Project Farmingdale, Nassau County.*

New York State Department of Environmental Conservation March 2014. *Record of Decision Farmingdale Plaza Cleaners Operable Unit Number 01: Remedial Program – Onsite Area for Soil and Soil Vapor State Superfund Project Farmingdale, Nassau County.*

URS. July 2000. Final Continued Remedial Investigation Report for the Liberty Industrial Finishing Site, Farmingdale, New York. Index No. II, CERCLA 97-0203.

YU & Associates, Inc. 2009. *Final Report, Immediate Investigation, Farmingdale Plaza Cleaners Site, Farmingdale, New York, Site Number: 130107*, prepared for NYSDEC on behalf of AECOM Technical Services Northeast, Inc.

Appendix A







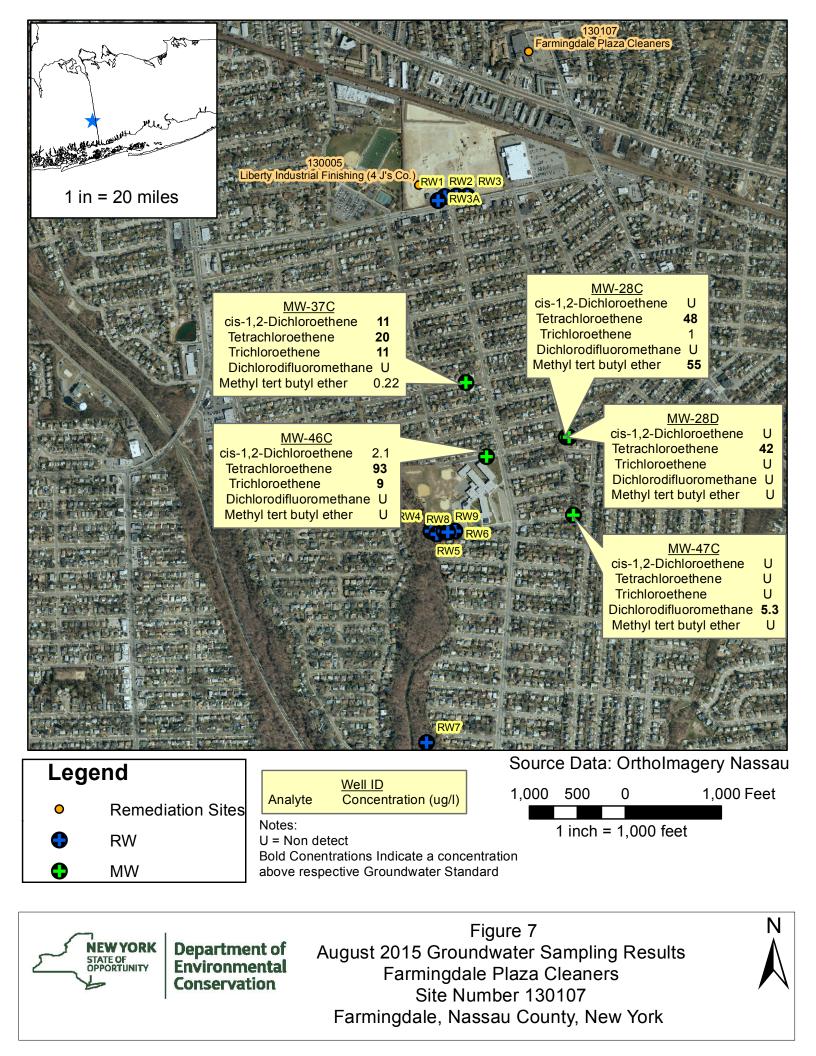
Magothy Aquifer

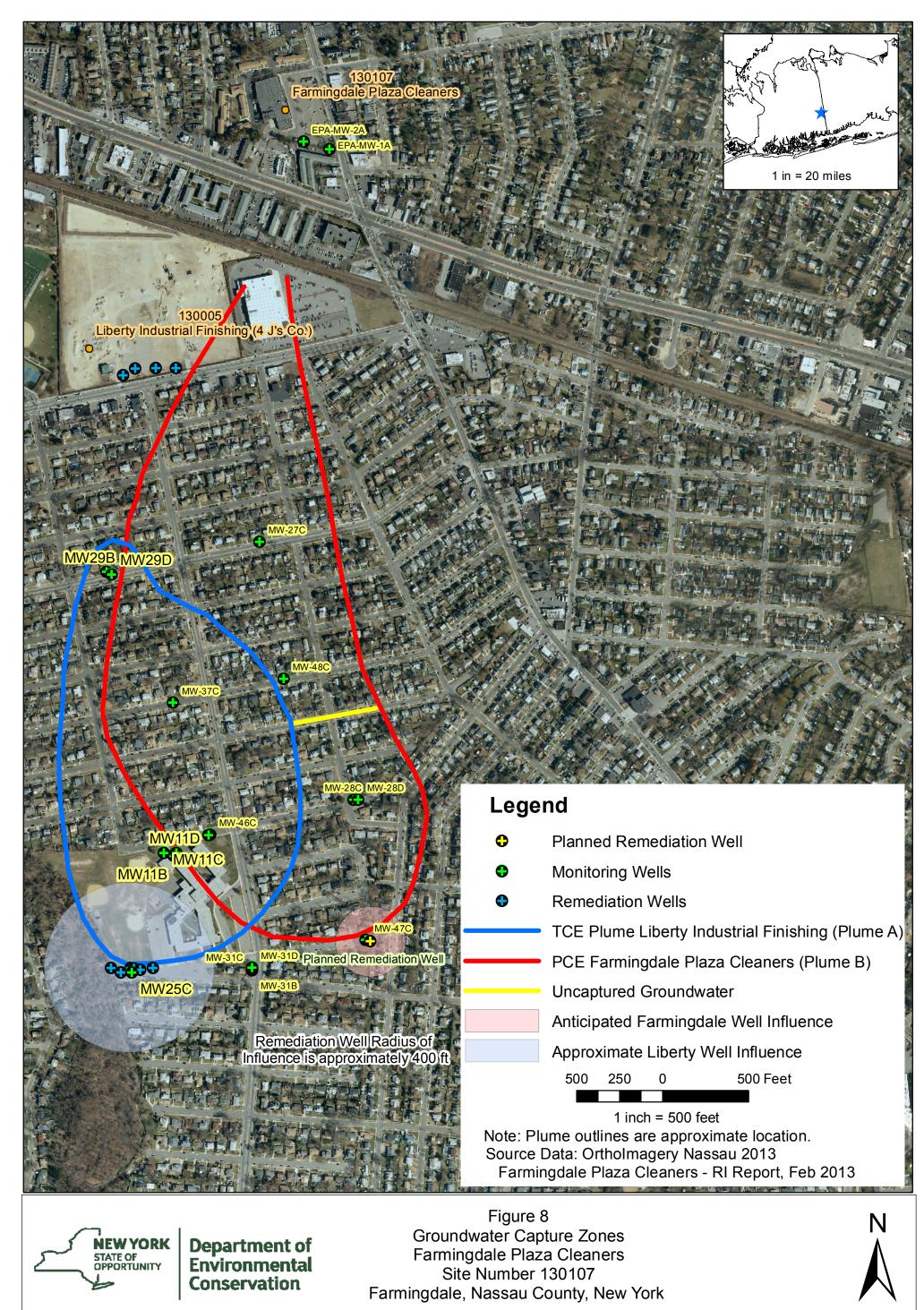


Department of Environmental Conservation

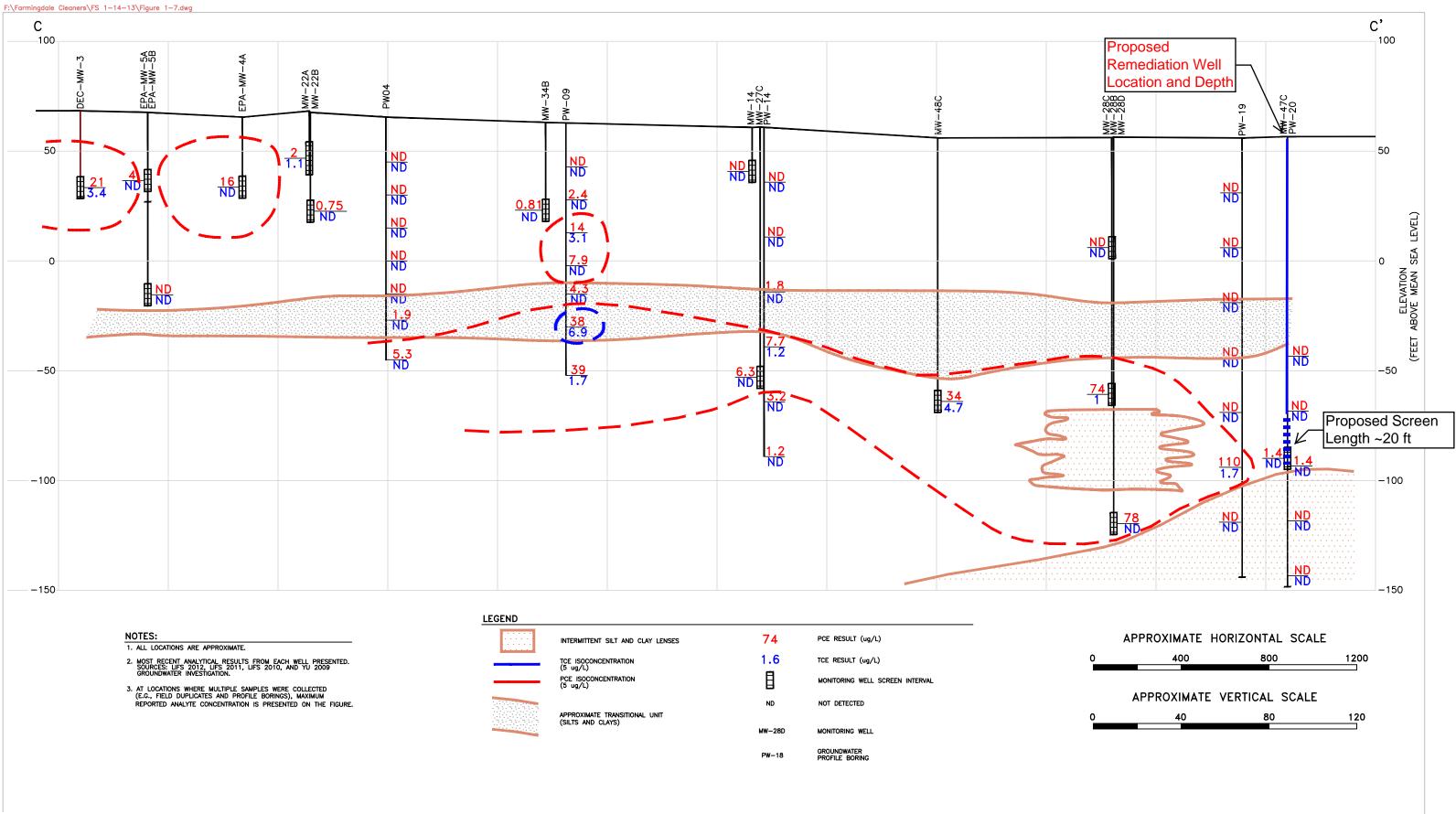
Farmingdale Plaza Cleaners Site Number 130107 Farmingdale, Nassau County, New York

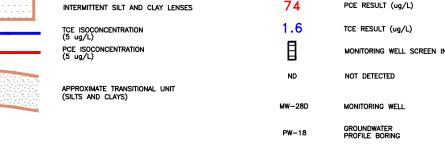




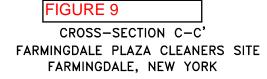








Gecology and environment P.C. -



Original Figure from 2014 Farmingdale Plaza ROD



Department of Environmental Conservation

Piping Conveyance Plan Farmingdale Plaza Cleaners Site Number 130107 Farmingdale, Nassau County, New York

Table 1. Farminguale Haza cleaners Monitoring Weir Elevation Data										
	Latitude		Reference	Top of bag	Depth to	Total	Elevation (ft			
Well ID		Longitude	Elevation	Elevation (ft) 8/25/15	Water (ft)	Depth (ft)	amsl)			
			(ft amsl)	(ft amsl) (11) 8/23/13		8/25/15	8/25/15			
EPA-MW-1A	-73.444948	40.728524	65.50	*	18.50	39.70	47.00			
EPA-MW-2A	-73.445489	40.728650	67.00	*	18.09	37.50	48.91			
MW-11B	-73.448449	40.717324	50.12	*	9.58	*	40.54			
MW-11C	-73.448561	40.717303	50.06	*	10.87	*	39.19			
MW-11D	-73.448396	40.717330	50.19	*	11.10	*	39.09			
MW-25C	-73.449406	40.715479	44.41	*	6.02	46.60	38.39			
MW-27C	-73.446560	40.722273	61.00	*	17.06	*	43.94			
MW-28C	-73.444688	40.718127	56.06	116	15.12	*	40.94			
MW-28D	-73.444592	40.718125	56.41	175	15.35	*	41.06			
MW-29B	-73.449758	40.721821	60.47	*	17.24	50.90	43.23			
MW-29D	-73.449727	40.721829	60.61	*	18.04	195.00	42.57			
MW-31B	-73.446880	40.715433	53.48	*	14.10	*	39.38			
MW-31C	-73.446877	40.715451	53.44	*	15.46	*	37.98			
MW-31D	-73.446870	40.715478	53.30	*	15.66	*	37.64			
MW-37C	-73.448439	40.719727	55.44	113	13.72	*	41.72			
MW-46C	-73.447706	40.717608	54.45	144	14.52	*	39.93			
MW-46D	-73.447739	40.717603	54.39	*	14.71	*	39.68			
MW-47C	-73.444483	40.715883	56.64	145	17.91	*	38.73			
MW-48C	-73.446100	40.720081	56.04	*	13.40	*	42.64			

Table 1: Farmingdale Plaza Cleaners Monitoring Well Elevation Data

Notes:

1.) ft amsl = feet above mean sea level

2.) Length of diffusion bag used was 2 ft

3.) * = no available data

Table 2: Groundwater Analytical Results for Farmingdale Plaza Cleaners for August 25th 2015								
		Monito						
VOC Analyte Concentration	MW-28C	MW-28D	MW-37C	MW-46C	MW-			
	10100-200				47C	Groundwater Standard		
1,1-Dichloroethene	U	0.61	U	1.7	U	5.0		
2-Butanone	1.4	U	U	1.7	U	50.0		
cis-1,2-Dichloroethene	U	U	11	2.1	U	5.0		
Tetrachloroethene	48	42	20	93	U	5.0		
Trichloroethene	1.0	U	11	9.0	U	5.0		
Dichlorodifluoromethane	U	U	U	U	5.3	5.0		
1,2-Dichlorobenzene	U	U	1.8	U	U	3.0		
Chlorobenzene	U	U	2.0	U	U	5.0		
Methyl tert butyl ether	55	U	0.22	U	U	10.0		
Toluene	U	U	1.6	U	U	5.0		
1,1,2-Trichloroethane	U	0.61	U	U	U	1.0		
1,1,-Dichloroethane	U	1.6	U	U	U	5.0		
Chloroform	U	0.45	U	U	U	7.0		
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Note: 1.) VOC = Volatile Organic Compound

2.) Units are in micrograms per liter (ug/l)

3.) Acetone was assumed to be a lab contaminant and thus was not included in the table

4.) U = Analyte was not detected above the laboratory method detection limit

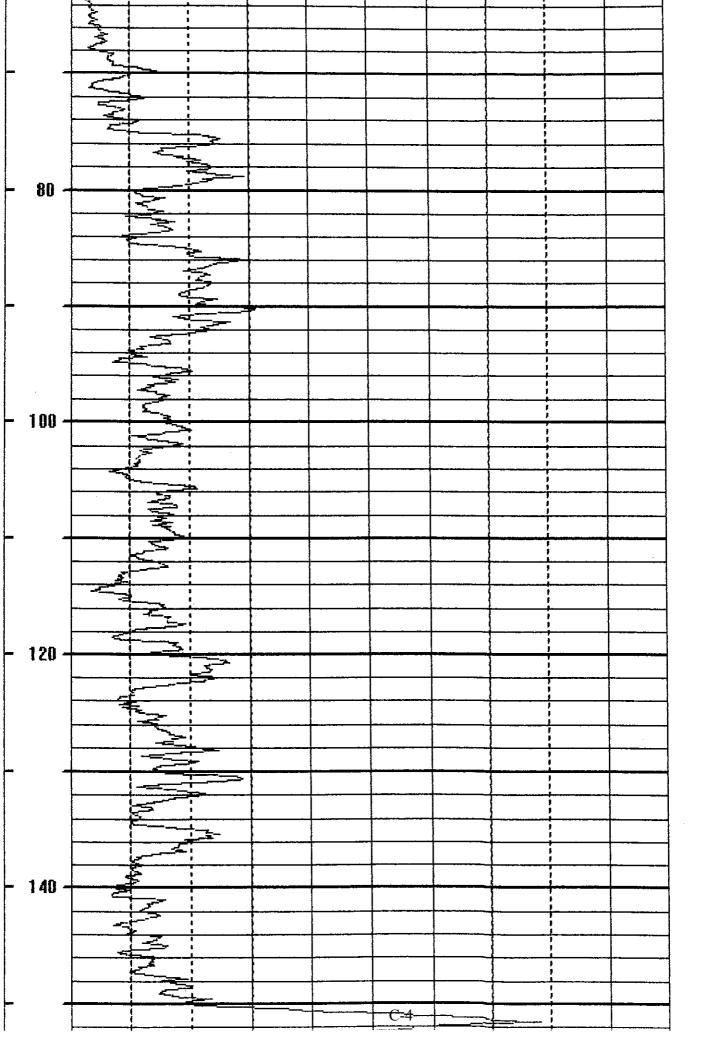
5.) Highlighted Values indicate excedences of NYSDEC Groundwater Standards

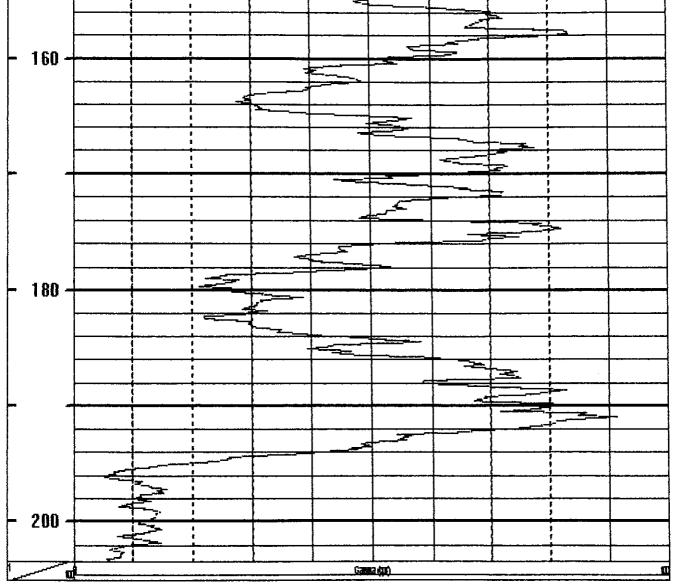
Appendix A

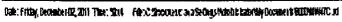
Gamma Logs

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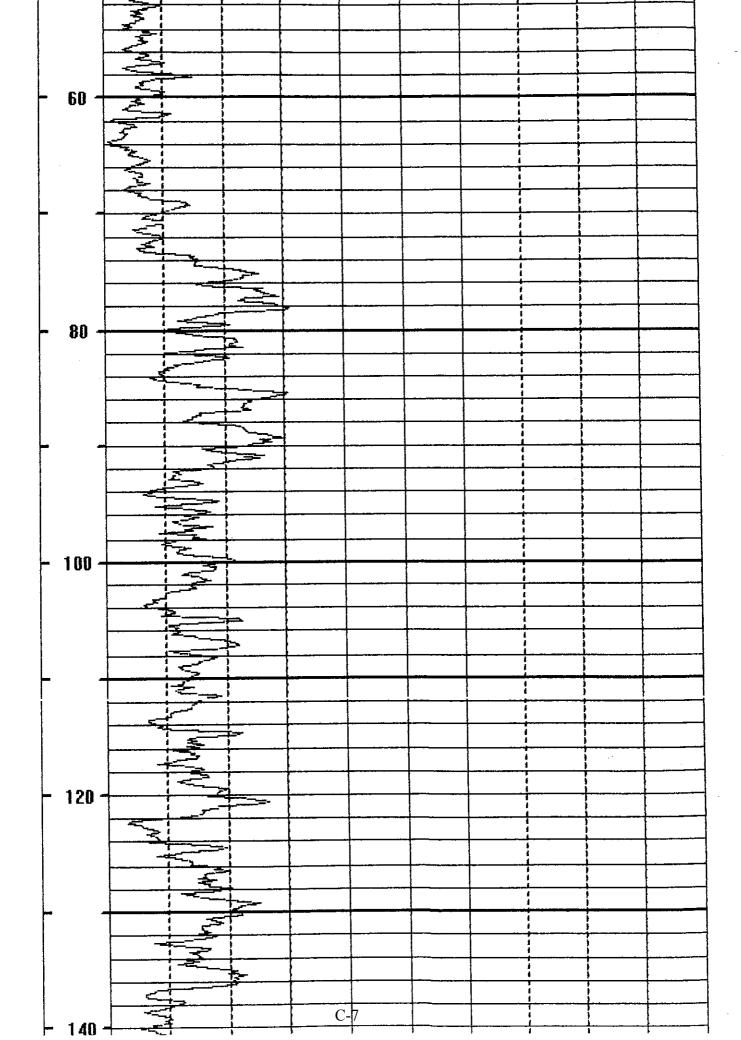


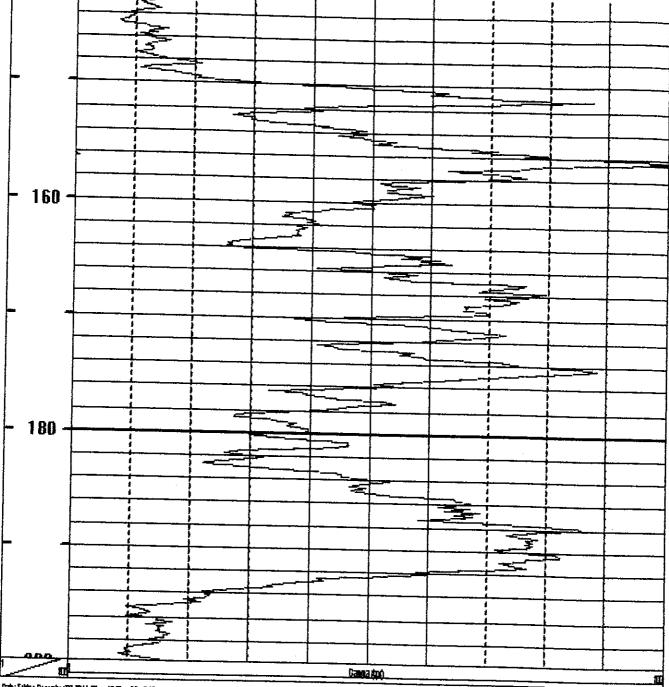


MW47C UP

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		Location	TOMES	STREET					
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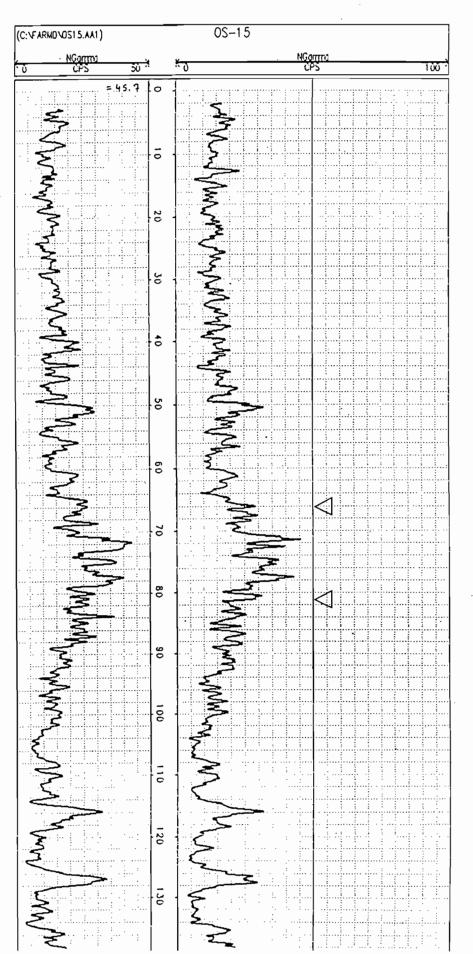
Date: Fritter, December 19, 2011, They 10.28 F.D.; C. LANSARD 206 STEVEN Shirts of the Browner's HILD MINISTOR OF



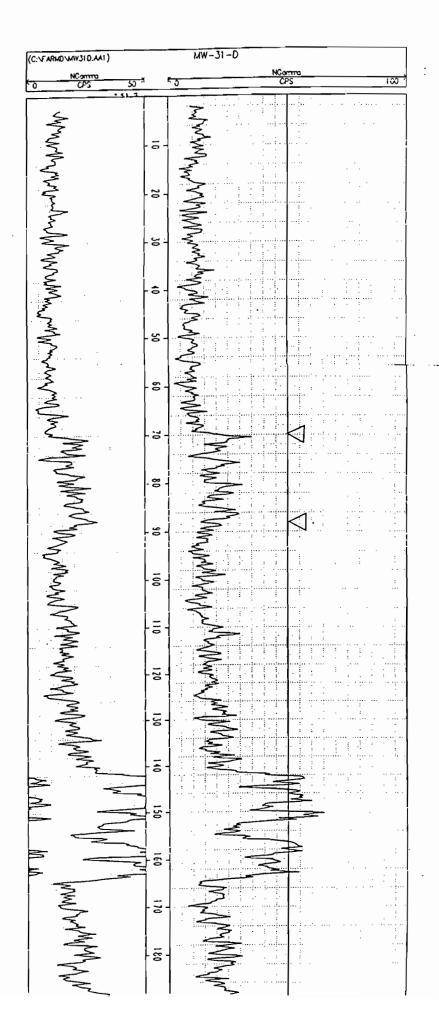


Date: Friday, December 192, 2014 Tipe: 1023 First Calory or the 2010 Strate Mark Strate 194 Boomer's Hill Hill Will Capud

.



OS-15 is also shown as GW-25



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1

Well Logs

DRILLING LOG FOR ______ Farmingdale Plage Cleaners Project Name Water Level (TOIC) Date Time Level(Feet) Sits Location Farmylile, NY Date Started/Finished 12 13 14 - 12/15 14 Dilling Company Masor Drilling Driller's Name_Adam Well Location Sketch 3 B X Bu Geologist's Name ____ 46 A CERT Geologist's Signature_ ZRig Type (s) ____ Rato ANT ROAD - Sonic Bro Drilling Method (s) Spielman \$← MW 28D Bit Size (s) Ave NA. _____ Auger Size (s) NA Augen/Split Spoon Refusal Cilliada 181 Total Depth of Borehole Is _____ NA ıR, Total Depth of Corehols Is____ 5 Look Number Stieleum. æ

				aucivupit	l.
i	SCREENED WELL	Inner Casing Material	OPEN-HOLE WELL	Inner Gesing Material	
	Stick-upft	Inner Casing Inside Diameter 2 inches		Inner Casing Inside DiameterInches	
	Top of Grout	GROUND SURFACE Quantity of Material Used: Bentonite Pellets		Outer Casing DiameterInches	
	Top of Seal at 163-5 ft	Coment Borehole inches Diameter		Borshols Dismeterft	
	Top of Sand Pack 1075 1	Cementi/ Bentonite		Bedrockft	
Y	Top of Screen at 17-1 R	Grout		Bettom of Rock Socket/ Outer Casing1	
	Bottom of 181	Screen Slot Size <u>+020</u> Screen Type		Bottom of Inner Casingtt	
	streen at1	K PVC Stainless Steel		Corshols Diameter	
	Bottom of Hole at <u>191</u> fi Bottom of Sandpack at <u>(21</u>	Pack Type/Size: CKSand 22 Gravel Natural		Bottom of Coreholeft	
					ε.

NOTE: See pages 136 and 137 for well construction diagrams

DRILLING LOG FOR PW23/MW450 Plaza Chanes Project Name Fancine Water Level (TOIC) Time Level(Feet) Date Site Location _ Farming date ĸ Date Started/Finished Drilling Company _ Well Location Sketch Driller's Nama Geologist's Name Geologist's Signature Ц Rig Type (s) ø Sovia Dritting Method (s) Size (s) Bit Size (s) NW 45D NA Auger/Split Spoon Refusal Œ Total Depth of Borehole is Total Depth of Corehole Is, MW-45 D Lock Number____ Stick-up____ tt SCREENED WELL OPEN-HOLE WELL Inner Casing PUC Inner Casing Material_ Material Inner Casino Inside Diameter 2_ inches Inner Casing Inside Stick-up, Diameter inches' GROUND SURFACE Quantity of Material Used: Top of Grout Outer Casing Bentonite Pellets Diameter inches ik-Ge Coment Borehole Topof Seal at 152,5 tt Diameter ŧ Borehole 6 inches Diameter Bedrock_ Top of Sand Pack 156.5 ft Cement/ Bentonite Bottom of Rock Socket/ Grout Top of Outer Casing _ft Screen at 100_1 Ection of Inner Casing_ ė Bottom of Screen Type _ 170 Screen at S PVC_ Corehcie Stainless Steel Diameter Pack Type/Size: Silien #2 Bottom of 171 Hole at ÷ Bottom of Gravel. Corehole, Bottom of Sandpack at ______/7/ Natural

NOTE: See pages 138 and 137 for well construction diagrams

Time	Loval (Feat)
skétch 199	i HII I J
1 1 1 1	
De	50 L
3 8.0	
durantive .	
TILLI	5, 117-
	Wagener
	Skelich Pr Skelich Pr Dog Dog Dog Dog Dog Dog Dog Dog Dog Dog

	SCREENED WELL	Lock Number	OPEN-HOLE WELL	Stick-upft
	Stick-upft	Inner Casing Inside Diameter inches GROUND SURFACE	T	Inner Casing Inside Diamotorinches
2	Top of Grout	Quantly of Material Used: Bentonite Polets		Outer Casing Diameterinches
	Seal at 12-9 ft	Cement		Borehole Diameterti
	Top of Sand Pack 134 tt	Cement/ Benionite		Bedrookft
	Screen at 140 ft	Gou(Bottom of Rock Socket/ Outer Casingft
		Screen Slot Size Screen Type DVFVC		Bottom of Inner Casingft Corehole
	Hole at 19619 It	Stainless Steel Pack Type/Size: S. Sand Gravel		Diameter Botiom of Coreholeft
	Bottom of Sandpack at 15から	n diagrams		

DRILLING LOG FOR _______ Þ Project Name Farming Le Plaza Changes Water Level (TOIC) Time Date Level(Feet) Farmingdele. ٨V Site Location - MMH Data Started/Finished _42 Drifting Company Driller's Name A\$ Well oction Sketch 65 Ben Geologist's Name _ 5n Calo SPielman Ave Geologist's Signature Ъ, MW-45D Rig Type (s) OWITA Ð Drilling Method (s) " rods w/ 20' core brined Bit Size (5) 6." وختت 1 cessig p. ger Size (s) NA Auger/Split Spoon Refusal 195 H Total Depth of Borehole Is School NA-Total Depth of Corehole is_ Lock Number, Stick-up_ SCREENED WELL OPEN-HOLE WELL Inner Casing Inner Casing 000 Material flush ma. Material. inner Casing inside Diameter <u>2.</u> inches Inner Casing Inside Stick-up, Diameter: inches GROUND SURFACE Quantity of Material Used: Top of Grout Outer Casing Bentonile 4 Diameter inches Pellets, Cement Borshole Top of Seelat 176 Diameter ÷ -11 6 Inches Borehole Diameter Bedrock_ ŤI. Top of Sand Pack Cement/ Bentonite, Bottom of Rock Socket/ Gaut Outer Casing_ -tt Top of Screen at 185 at

Screen Slot Size 0,020

PVC 2"× 10'

Steinless Steel

Screen Type _

Pack Type/Size: & Sand ≠2

Gravel

Natural.

风 Sand_ Bottom of Inner Casing_

Corehole Diameter

Bottom of Corehole

ft

NOTE: See pages 136 and 137 for well construction diagrams

ŧ

Bottom of

Screen at

Bettom of

Hole at_

195

195

Bottom of Sandpack et _____195

Water Level (TO)	C)
Time	Lovel(Feet)
on Sketch	*1
Tomes	
-	
A	
MW-476	- e T
O I I]
-19 11-	*
Tomes 2	Tomes
	In Skeich Time Tomes Tomes MW-470

	Lock Number		Stick-upt
SCREENE	Material	OPEN-HOLE WELL	Inner Casing Material
81ick-up -0.5 tt	inner Casing Inside Diameter 🛃 , inches		Inner Casing Inside Diameter Inches
	GROUND SURFACE		
Top of Grout	Quantity of Material Used: Bentonite Poliots_Quilk - Ccl		Outer Casing "Dismeterinches
Top of 130 ti	Carrent		Borehole
Sealar 1.22 II	Borehole inches Diameter		Diemeterft
Top of Sand Pack 136 ti	Cement/ Bertonke		Bedrockft
Top of Screen at 140 t 11/1.5	Grout		Bottom of Rock Socket/ Outer Casingft
Screen at 140 tt 141.5 Screen at 150 151.5	Screen Slot Size < 0.20		Bottom of loner
Sottom of HC 151.5	Screen Type	·	Casingft
	Steinless Stoel		-Corehole Diameter
tolast 153	Pack Type/Size Sand US Silice #2		
Bottom of Sandpeck at 153	Gravel	•	Bottom of Coreholeft

Project Name Farmingdale Plage Clean	reis-	Water Level (TO	(C)	7
Sile Location Farminghete, NY	Date	Time	Loval(Feet)	-
Date Started/Finished 12/15/11 - 12/16/0	-			
Dilling Company Major Drilling				
Driller's Name Alan Cuso	Well Location S	liketoh	103	
Geologist's Name Ben Cile		1854	F Put &	F
Geologist's Signature	A line		1 Finite	8
Rig Type (s) Brut A source			1 20	æ
Drilling Method (s) Noto Somic		Kmw 48	pc .	
Bit Size (s)		SIH AVE.	·	-
Auger/Split Spoon Refusal NA		pierrave		
			1	
Total Depth of Corehole IsIA		6 PLitt	1	1

·				
		Lock Number		Stick-upft
	SCREENED WELL	Inner Casing Materialpv.c	OPEN-HOLE WELL	Inner Casing Material
Stick-upft	T	Inner Casing Inside Diameter 2 inches	T	Inner Casing inside DiameterInches
Top of Grout	1 F	GROUND SURFACE Quartity of Material Used: Benionite Pelleta 45 bass		Outer Casing Diameterinches
Top of Sealet /01 r.		Cement Borehole inches Diametor		Borehole Diameterft
Top of Band Pack ///, 8 ft		Cement/ . Bentonite		Bedrockft
Top of . Screen at <u>/1.5</u> ft		Grout	· 📓 📓	Bottom of Rock Socket/ Outer Casingft
Bottom of Screen atfi		Screen Slot Size _ , 0 2 0 Screen Type		Bottom of Inner Casingfl Corehole Diameter
Bottom of 12-5 n Hole at 12-5 n Bottom of Sandpack at 7		Pack Type/Siza: C Sand _##-2 Gravel Natural		Bottom of Coreholeft
NOTE: See pages 136 and 1	197 for well construct	on diagrams		

Appendix B



ANALYTICAL REPORT

Job Number: 480-86254-1

Job Description: DEC Farmingdale Plaza Cleaners #130107

For: New York State D.E.C. 625 Broadway 9th Floor Albany, NY 12233-7258

Attention: Mr. Brian Jankauskas

Approved for release. Sarah E Brown Project Management Assistant II 9/8/2015 5:55 PM

Designee for Melissa Haas, Project Manager I 777 New Durham Road, Edison, NJ, 08817 (203)944-1310 melissa.haas@testamericainc.com 09/08/2015

The test results in this report meet all NELAP requirements for analytes for which accreditation is required or available. Any exceptions to the NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory. All questions regarding this test report should be directed to the TestAmerica Project Manager who has signed this report. TestAmerica Buffalo NELAC Certifications: CADPH 01169CA, FLDOH E87672, ILEPA 200003, KSDOH E-10187, LADEQ 30708, MDH 036-999-337, NHELAP 2973, NJDEP NY455, NHDOH 10026, ORELAP NY200003, PADEP 68-00281, TXCEQ T-104704412-10-1

TestAmerica Laboratories, Inc.TestAmerica Buffalo10 Hazelwood Drive, Amherst, NY14228-2298Tel (716) 691-2600Fax (716) 691-7991www.testamericainc.com



Job Description: DEC Farmingdale Plaza Cleaners #130107

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed within the body of this report. Release of the data contained in this sample data package and in the electronic data deliverable has been authorized by the Laboratory Manager or his/her designee, as verified by the following signature.

grahe. Brow

Approved for release. Sarah E Brown Project Management Assistant II 9/8/2015 5:55 PM

Designee for Melissa Haas

CASE NARRATIVE

Client: New York State D.E.C.

Project: DEC Farmingdale Plaza Cleaners #130107

Report Number: 480-86254-1

This case narrative is in the form of an exception report, where only the anomalies related to this report, method specific performance and/or QA/QC issues are discussed. If there are no issues to report, this narrative will include a statement that documents that there are no relevant data issues.

It should be noted that samples with elevated Reporting Limits (RLs) as a result of a dilution may not be able to satisfy customer reporting limits in some cases. Such increases in the RLs are unavoidable but acceptable consequence of sample dilution that enables quantification of target analytes or interferences which exceed the calibration range of the instrument.

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

RECEIPT

The samples were received on 8/27/2015 1:45 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 1.3° C.

Note: All samples which require thermal preservation are considered acceptable if the arrival temperature is within 2C of the required temperature or method specified range. For samples with a specified temperature of 4C, samples with a temperature ranging from just above freezing temperature of water to 6C shall be acceptable. Samples that are hand delivered immediately following collection may not meet these criteria, however they will be deemed acceptable according to NELAC standards, if there is evidence that the chilling process has begun, such as arrival on ice, etc.

VOLATILE ORGANICS

Samples MW-46C-82015 (480-86254-1), MW-47C-82015 (480-86254-2), MW-37C-82015 (480-86254-3), MW-28C-82015 (480-86254-4), MW-28D-82015 (480-86254-5) and TRIP BLANK (480-86254-6) were analyzed for Volatile organics in accordance with EPA SW-846 Methods 8260C. The samples were analyzed on 09/03/2015 and 09/04/2015.

The initial calibration curve analyzed in batch 480-261126 was outside method criteria for the low point linear regression fit of 30% read back criteria for the following analytes: Bromoform, 1,2-Dibromo-3-Chloropropane. As indicated in the reference method, sample analysis may proceed; however, any detection or non-detection for the affected analytes is considered an estimated concentration.

The continuing calibration verification (CCV) associated with batch 480-262028 recovered above the upper control limit for the following analyte: Cyclohexane. The sample associated with this CCV were non-detect for the affected analyte; therefore, the data have been reported.

Refer to the QC report for details.

No other difficulties were encountered during the Volatile organics analysis.

All other quality control parameters were within the acceptance limits.

EXECUTIVE SUMMARY - Detections

Client: New York State D.E.C.

Lab Sample ID	Client Sample ID		a 110	Reporting		
Analyte		Result	Qualifier	Limit	Units	Method
480-86254-1	MW-46C-82015	4 7		1.0		00000
1,1-Dichloroethene		1.7		1.0	ug/L	8260C
2-Butanone (MEK)		1.7	J	10	ug/L	8260C
Acetone		35		10	ug/L	8260C
cis-1,2-Dichloroethe	ne	2.1		1.0	ug/L	8260C
Tetrachloroethene		93		1.0	ug/L	8260C
Trichloroethene		9.0		1.0	ug/L	8260C
480-86254-2	MW-47C-82015					
Acetone		14		10	ug/L	8260C
Dichlorodifluorometh	ane	5.3		1.0	ug/L	8260C
Diemoroamaoromea		0.0		1.0	ug/L	02000
480-86254-3	MW-37C-82015					
1,2-Dichlorobenzene	9	1.8		1.0	ug/L	8260C
Acetone		28		10	ug/L	8260C
Chlorobenzene		2.0		1.0	ug/L	8260C
cis-1,2-Dichloroethe	ne	11		1.0	ug/L	8260C
Methyl tert-butyl ethe	er	0.22	J	1.0	ug/L	8260C
Tetrachloroethene		20		1.0	ug/L	8260C
Toluene		1.6		1.0	ug/L	8260C
Trichloroethene		11		1.0	ug/L	8260C
400 00054 4	MM/ 200 02045					
480-86254-4	MW-28C-82015	1.4	J	10	ug/L	8260C
2-Butanone (MEK)		28	J	10	-	8260C 8260C
Acetone	or	28 55		10	ug/L	8260C 8260C
Methyl tert-butyl ethe Tetrachloroethene					ug/L	
Trichloroethene		48 1.0		1.0	ug/L	8260C
Inchioroethene		1.0		1.0	ug/L	8260C
480-86254-5	MW-28D-82015					
1,1,2-Trichloroethan		0.61	J	1.0	ug/L	8260C
1,1-Dichloroethane		1.6		1.0	ug/L	8260C
1,1-Dichloroethene		0.61	J	1.0	ug/L	8260C
Acetone		24		10	ug/L	8260C
Chloroform		0.45	J	1.0	ug/L	8260C
Tetrachloroethene		42	-	1.0	ug/L	8260C
					~ 3 . –	

METHOD SUMMARY

Client: New York State D.E.C.

Job Number: 480-86254-1

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS	TAL BUF	SW846 8260C	
Purge and Trap	TAL BUF		SW846 5030C

Lab References:

TAL BUF = TestAmerica Buffalo

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Method	Analyst	Analyst ID
SW846 8260C	Feldman, Lance J	LJF
SW846 8260C	Goliszek, Gregory T	GTG

SAMPLE SUMMARY

Client: New York State D.E.C.

			Date/Time	Date/Time
Lab Sample ID	Client Sample ID	Client Matrix	Sampled	Received
480-86254-1	MW-46C-82015	Water	08/25/2015 1310	08/27/2015 0145
480-86254-2	MW-47C-82015	Water	08/25/2015 1330	08/27/2015 0145
480-86254-3	MW-37C-82015	Water	08/25/2015 1350	08/27/2015 0145
480-86254-4	MW-28C-82015	Water	08/25/2015 1420	08/27/2015 0145
480-86254-5	MW-28D-82015	Water	08/25/2015 1430	08/27/2015 0145
480-86254-6	TRIP BLANK	Water	08/25/2015 0000	08/27/2015 0145

SAMPLE RESULTS

Analytical Data

Client Sample ID:	MW-46C-82015				
Lab Sample ID: Client Matrix:	480-86254-1 Water				mpled: 08/25/2015 1310 ceived: 08/27/2015 0145
	8	260C Volatile Organi	c Compounds by	GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2101 09/03/2015 2101	Analysis Batch: Prep Batch:	480-261891 N/A	Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	
Analyte		Result (u	g/L) Qualit	fier MDL	RL
1,1,1-Trichloroetha	ane	1.0	U	0.82	1.0
1,1,2,2-Tetrachloro	bethane	1.0	U	0.21	1.0
1,1,2-Trichloro-1,2	,2-trifluoroethane	1.0	U	0.31	1.0
1,1,2-Trichloroetha	ane	1.0	U	0.23	1.0
1,1-Dichloroethane	e	1.0	U	0.38	1.0
1,1-Dichloroethene	e	1.7		0.29	1.0
1,2,3-Trichloroben	zene	1.0	U	0.41	1.0
1,2,4-Trichloroben	zene	1.0	U	0.41	1.0
1,2-Dibromo-3-Chl	oropropane	1.0	U	0.39	1.0
1,2-Dichlorobenze	ne	1.0	U	0.79	1.0
1,2-Dichloroethane	e	1.0	U	0.21	1.0
1,2-Dichloropropa		1.0	U	0.72	1.0
1,3-Dichlorobenze		1.0	U	0.78	1.0
1,4-Dichlorobenze	ne	1.0	U	0.84	1.0
1,4-Dioxane		40	U	9.3	40
2-Butanone (MEK)		1.7	J	1.3	10
2-Hexanone		5.0	U	1.2	5.0
4-Methyl-2-pentan	one (MIBK)	5.0	U	2.1	5.0
Acetone		35		3.0	10
Benzene		1.0	U	0.41	1.0
Bromoform		1.0	U	0.26	1.0
Bromomethane		1.0	U	0.69	1.0
Carbon disulfide	4-	1.0	U	0.19	1.0
Carbon tetrachlorid	de	1.0	U	0.27	1.0
Chlorobenzene		1.0	U	0.75	1.0
Chlorobromometha		1.0	U	0.87	1.0
Chlorodibromomet Chloroethane	nane	1.0 1.0	U	0.32 0.32	1.0 1.0
Chloroform		1.0	U U	0.32	1.0
Chloromethane		1.0	U	0.34	1.0
cis-1,2-Dichloroeth	ana	2.1	0	0.81	1.0
cis-1,3-Dichloropro		1.0	U	0.36	1.0
Cyclohexane	pene	1.0	U	0.18	1.0
Dichlorobromomet	hane	1.0	U	0.39	1.0
Dichlorodifluorome		1.0	Ŭ	0.68	1.0
Ethylbenzene		1.0	U	0.74	1.0
Ethylene Dibromid	e	1.0	Ŭ	0.73	1.0
Isopropylbenzene	•	1.0	Ŭ	0.79	1.0
Methyl acetate		2.5	Ŭ	1.3	2.5
Methyl tert-butyl et	her	1.0	Ŭ	0.16	1.0
Methylcyclohexane		1.0	Ŭ	0.16	1.0
Methylene Chlorid		1.0	Ŭ	0.44	1.0
m-Xylene & p-Xyle		2.0	Ŭ	0.66	2.0
o-Xylene	-	1.0	U	0.76	1.0
Styrene		1.0	Ŭ	0.73	1.0
Tetrachloroethene		93	-	0.36	1.0

Analytical Data

Client Sample ID:	MW-46C-82015							
Lab Sample ID: Client Matrix:	480-86254-1 Water						pled: 08/25/20 eived: 08/27/20	
	82	260C Volatile Organi	ic Compoun	ds by	GC/MS			
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2101 09/03/2015 2101	Analysis Batch: Prep Batch:	480-26189 N/A	1	Instrument ID: Lab File ID: Initial Weight/\ Final Weight/\	/olume:	HP5973S S2162.D 5 mL 5 mL	
Analyte		Result (u	ıg/L)	Quali	fier MDL		RL	
Toluene		1.0		U	0.51		1.0	
trans-1,2-Dichloro	ethene	1.0		U	0.90		1.0	
trans-1,3-Dichloro	propene	1.0		U	0.37		1.0	
Trichloroethene		9.0			0.46		1.0	
Trichlorofluoromet	hane	1.0		U	0.88		1.0	
Vinyl chloride		1.0		U	0.90		1.0	
Surrogate		%Rec		Quali	fier A	Acceptant	ce Limits	
1,2-Dichloroethane	e-d4 (Surr)	105			6	6 - 137		
4-Bromofluoroben	zene (Surr)	101			7	' 3 - 120		
Dibromofluoromet	hane (Surr)	105			6	60 - 140		
Toluene-d8 (Surr)		99			7	' 1 - 126		

Analytical Data

Job Number: 480-86254-1

Client Sample ID:	MW-47C-82015				
Lab Sample ID: Client Matrix:	480-86254-2 Water				mpled: 08/25/2015 1330 ceived: 08/27/2015 0145
		8260C Volatile Organi	c Compounds by	GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2124 09/03/2015 2124	Analysis Batch: Prep Batch:	480-261891 N/A	Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	
Analyte		Result (u	g/L) Quali	fier MDL	RL
1,1,1-Trichloroetha	ane	1.0	U	0.82	1.0
1,1,2,2-Tetrachlor	pethane	1.0	U	0.21	1.0
1,1,2-Trichloro-1,2	2,2-trifluoroethane	1.0	U	0.31	1.0
1,1,2-Trichloroetha	ane	1.0	U	0.23	1.0
1,1-Dichloroethan	9	1.0	U	0.38	1.0
1,1-Dichloroethene	e	1.0	U	0.29	1.0
1,2,3-Trichloroben	zene	1.0	U	0.41	1.0
1,2,4-Trichloroben	zene	1.0	U	0.41	1.0
1,2-Dibromo-3-Ch	loropropane	1.0	U	0.39	1.0
1,2-Dichlorobenze	ne	1.0	U	0.79	1.0
1,2-Dichloroethane	e	1.0	U	0.21	1.0
1,2-Dichloropropa	ne	1.0	U	0.72	1.0
1,3-Dichlorobenze	ne	1.0	U	0.78	1.0
1,4-Dichlorobenze	ne	1.0	U	0.84	1.0
1,4-Dioxane		40	U	9.3	40
2-Butanone (MEK)	10	U	1.3	10
2-Hexanone		5.0	U	1.2	5.0
4-Methyl-2-pentan	one (MIBK)	5.0	U	2.1	5.0
Acetone		14		3.0	10
Benzene		1.0	U	0.41	1.0
Bromoform		1.0	U	0.26	1.0
Bromomethane		1.0	U	0.69	1.0
Carbon disulfide		1.0	U	0.19	1.0
Carbon tetrachlori	de	1.0	U	0.27	1.0
Chlorobenzene		1.0	U	0.75	1.0
Chlorobromometh	ane	1.0	U	0.87	1.0
Chlorodibromome	thane	1.0	U	0.32	1.0
Chloroethane		1.0	U	0.32	1.0
Chloroform		1.0	U	0.34	1.0
Chloromethane		1.0	U	0.35	1.0
cis-1,2-Dichloroeth	nene	1.0	U	0.81	1.0
cis-1,3-Dichloropro	opene	1.0	U	0.36	1.0
Cyclohexane		1.0	U	0.18	1.0
Dichlorobromomet	thane	1.0	U	0.39	1.0
Dichlorodifluorome	ethane	5.3		0.68	1.0
Ethylbenzene		1.0	U	0.74	1.0
Ethylene Dibromid	le	1.0	U	0.73	1.0
Isopropylbenzene		1.0	U	0.79	1.0
Methyl acetate		2.5	U	1.3	2.5
Methyl tert-butyl et		1.0	U	0.16	1.0
Methylcyclohexan	е	1.0	U	0.16	1.0
Methylene Chlorid		1.0	U	0.44	1.0
m-Xylene & p-Xyle	ene	2.0	U	0.66	2.0
o-Xylene		1.0	U	0.76	1.0
Styrene		1.0	U	0.73	1.0
Tetrachloroethene		1.0	U	0.36	1.0

TestAmerica Buffalo

Client Sample ID: MW-47C-82015

Analytical Data

Lab Sample ID: Client Matrix:	480-86254-2 Water					npled: 08/25/2015 1330 eived: 08/27/2015 0145
		8260C Volatile Organi	c Compounds	by GC/I	MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2124 09/03/2015 2124	Analysis Batch: Prep Batch:	480-261891 N/A	La Ini	strument ID: ab File ID: itial Weight/Volume: nal Weight/Volume:	HP5973S S2163.D 5 mL 5 mL
Analyte		Result (u	g/L) G	ualifier	MDL	RL
Toluene		1.0	L		0.51	1.0
trans-1,2-Dichloro	ethene	1.0	U	l	0.90	1.0
trans-1,3-Dichloro	propene	1.0	U	l	0.37	1.0
Trichloroethene		1.0	U	l	0.46	1.0
Trichlorofluoromet	hane	1.0	U	l	0.88	1.0
Vinyl chloride		1.0	L		0.90	1.0
Surrogate		%Rec	G	ualifier	Acceptan	ce Limits
1,2-Dichloroethan	e-d4 (Surr)	106			66 - 137	
4-Bromofluoroben	zene (Surr)	103			73 - 120	
Dibromofluoromet	hane (Surr)	104			60 - 140	
Toluene-d8 (Surr)		98			71 - 126	

Analytical Data

Job Number: 480-86254-1

Client Sample ID:	MW-37C-82015				
Lab Sample ID: Client Matrix:	480-86254-3 Water				mpled: 08/25/2015 1350 ceived: 08/27/2015 0145
		8260C Volatile Organi	c Compounds by	GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2147 09/03/2015 2147	Analysis Batch: Prep Batch:	480-261891 N/A	Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	
Analyte		Result (u	ıg/L) Quali	ifier MDL	RL
1,1,1-Trichloroetha	ane	1.0	U	0.82	1.0
1,1,2,2-Tetrachlor	pethane	1.0	U	0.21	1.0
1,1,2-Trichloro-1,2	2,2-trifluoroethane	1.0	U	0.31	1.0
1,1,2-Trichloroetha	ane	1.0	U	0.23	1.0
1,1-Dichloroethane	9	1.0	U	0.38	1.0
1,1-Dichloroethene	e	1.0	U	0.29	1.0
1,2,3-Trichloroben	zene	1.0	U	0.41	1.0
1,2,4-Trichloroben	zene	1.0	U	0.41	1.0
1,2-Dibromo-3-Ch	loropropane	1.0	U	0.39	1.0
1,2-Dichlorobenze	ne	1.8		0.79	1.0
1,2-Dichloroethane	e	1.0	U	0.21	1.0
1,2-Dichloropropa	ne	1.0	U	0.72	1.0
1,3-Dichlorobenze	ne	1.0	U	0.78	1.0
1,4-Dichlorobenze	ne	1.0	U	0.84	1.0
1,4-Dioxane		40	U	9.3	40
2-Butanone (MEK))	10	U	1.3	10
2-Hexanone		5.0	U	1.2	5.0
4-Methyl-2-pentan	one (MIBK)	5.0	U	2.1	5.0
Acetone		28		3.0	10
Benzene		1.0	U	0.41	1.0
Bromoform		1.0	U	0.26	1.0
Bromomethane		1.0	U	0.69	1.0
Carbon disulfide		1.0	U	0.19	1.0
Carbon tetrachlorie	de	1.0	U	0.27	1.0
Chlorobenzene		2.0		0.75	1.0
Chlorobromometh	ane	1.0	U	0.87	1.0
Chlorodibromomet	thane	1.0	U	0.32	1.0
Chloroethane		1.0	U	0.32	1.0
Chloroform		1.0	U	0.34	1.0
Chloromethane		1.0	U	0.35	1.0
cis-1,2-Dichloroeth		11		0.81	1.0
cis-1,3-Dichloropro	opene	1.0	U	0.36	1.0
Cyclohexane		1.0	U	0.18	1.0
Dichlorobromomet		1.0	U	0.39	1.0
Dichlorodifluorome	ethane	1.0	U	0.68	1.0
Ethylbenzene		1.0	U	0.74	1.0
Ethylene Dibromid	e	1.0	U	0.73	1.0
Isopropylbenzene		1.0	U	0.79	1.0
Methyl acetate		2.5	U	1.3	2.5
Methyl tert-butyl et		0.22	J	0.16	1.0
Methylcyclohexan		1.0	U	0.16	1.0
Methylene Chlorid		1.0	U	0.44	1.0
m-Xylene & p-Xyle	ene	2.0	U	0.66	2.0
o-Xylene		1.0	U	0.76	1.0
Styrene		1.0	U	0.73	1.0
Tetrachloroethene		20		0.36	1.0

TestAmerica Buffalo

Analytical Data

Lab Sample ID: Client Matrix:	480-86254-3 Water					npled: 08/25/2015 1350 eived: 08/27/2015 0145
		8260C Volatile Organi	c Compounds	by GC/N	MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2147 09/03/2015 2147	Analysis Batch: Prep Batch:	480-261891 N/A	Lal Init	strument ID: b File ID: tial Weight/Volume: nal Weight/Volume:	HP5973S S2164.D 5 mL 5 mL
Analyte		Result (u	g/L) G	Qualifier	MDL	RL
Toluene		1.6			0.51	1.0
trans-1,2-Dichloro	ethene	1.0	L		0.90	1.0
trans-1,3-Dichloro	propene	1.0	L	J	0.37	1.0
Trichloroethene		11			0.46	1.0
Trichlorofluoromet	hane	1.0	L		0.88	1.0
Vinyl chloride		1.0	L	J	0.90	1.0
Surrogate		%Rec	G	Qualifier	Acceptan	ce Limits
1,2-Dichloroethan	e-d4 (Surr)	97			66 - 137	
4-Bromofluoroben	zene (Surr)	105			73 - 120	
Dibromofluoromet	hane (Surr)	98			60 - 140	
Toluene-d8 (Surr)		100			71 - 126	

Analytical Data

Job Number: 480-86254-1

Client Sample ID:	MW-28C-82015				
Lab Sample ID: Client Matrix:	480-86254-4 Water				mpled: 08/25/2015 1420 ceived: 08/27/2015 0145
		8260C Volatile Organi	c Compounds by	GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2210 09/03/2015 2210	Analysis Batch: Prep Batch:	480-261891 N/A	Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	
Analyte		Result (u	g/L) Quali	fier MDL	RL
1,1,1-Trichloroetha	ane	1.0	U	0.82	1.0
1,1,2,2-Tetrachlor	pethane	1.0	U	0.21	1.0
1,1,2-Trichloro-1,2	2,2-trifluoroethane	1.0	U	0.31	1.0
1,1,2-Trichloroetha	ane	1.0	U	0.23	1.0
1,1-Dichloroethan	e	1.0	U	0.38	1.0
1,1-Dichloroethene	e	1.0	U	0.29	1.0
1,2,3-Trichloroben		1.0	U	0.41	1.0
1,2,4-Trichloroben		1.0	U	0.41	1.0
1,2-Dibromo-3-Ch	loropropane	1.0	U	0.39	1.0
1,2-Dichlorobenze		1.0	U	0.79	1.0
1,2-Dichloroethan	Э	1.0	U	0.21	1.0
1,2-Dichloropropa		1.0	U	0.72	1.0
1,3-Dichlorobenze		1.0	U	0.78	1.0
1,4-Dichlorobenze	ne	1.0	U	0.84	1.0
1,4-Dioxane		40	U	9.3	40
2-Butanone (MEK)	1.4	J	1.3	10
2-Hexanone		5.0	U	1.2	5.0
4-Methyl-2-pentan	one (MIBK)	5.0	U	2.1	5.0
Acetone		28		3.0	10
Benzene		1.0	U	0.41	1.0
Bromoform		1.0	U	0.26	1.0
Bromomethane		1.0	U	0.69	1.0
Carbon disulfide		1.0	U	0.19	1.0
Carbon tetrachlori	de	1.0	U	0.27	1.0
Chlorobenzene		1.0	U	0.75	1.0
Chlorobromometh	ane	1.0	U	0.87	1.0
Chlorodibromome	thane	1.0	U	0.32	1.0
Chloroethane		1.0	U	0.32	1.0
Chloroform		1.0	U	0.34	1.0
Chloromethane		1.0	U	0.35	1.0
cis-1,2-Dichloroeth	nene	1.0	U	0.81	1.0
cis-1,3-Dichloropro	opene	1.0	U	0.36	1.0
Cyclohexane		1.0	U	0.18	1.0
Dichlorobromomet	thane	1.0	U	0.39	1.0
Dichlorodifluorome	ethane	1.0	U	0.68	1.0
Ethylbenzene		1.0	U	0.74	1.0
Ethylene Dibromid	le	1.0	U	0.73	1.0
Isopropylbenzene		1.0	U	0.79	1.0
Methyl acetate		2.5	U	1.3	2.5
Methyl tert-butyl e	ther	55		0.16	1.0
Methylcyclohexan	е	1.0	U	0.16	1.0
Methylene Chlorid	e	1.0	U	0.44	1.0
m-Xylene & p-Xyle		2.0	U	0.66	2.0
o-Xylene		1.0	U	0.76	1.0
Styrene		1.0	U	0.73	1.0
Tetrachloroethene		48		0.36	1.0

TestAmerica Buffalo

Analytical Data

Client Sample ID:	: MW-28C-82015					
Lab Sample ID: Client Matrix:	480-86254-4 Water					npled: 08/25/2015 1420 eived: 08/27/2015 0145
	8	260C Volatile Organi	c Compounds	by GC/MS	;	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2210 09/03/2015 2210	Analysis Batch: Prep Batch:	480-261891 N/A	Lab F Initial	ument ID: File ID: I Weight/Volume: Weight/Volume:	HP5973S S2165.D 5 mL 5 mL
Analyte		Result (u	g/L) G	Jualifier	MDL	RL
Toluene		1.0	L	l	0.51	1.0
trans-1,2-Dichloro	ethene	1.0	U	I	0.90	1.0
trans-1,3-Dichloro	propene	1.0	U	I	0.37	1.0
Trichloroethene		1.0			0.46	1.0
Trichlorofluoromet	hane	1.0	U	I	0.88	1.0
Vinyl chloride		1.0	L	l	0.90	1.0
Surrogate		%Rec	G	Jualifier	Acceptar	ce Limits
1,2-Dichloroethan	e-d4 (Surr)	101			66 - 137	
4-Bromofluoroben	zene (Surr)	100			73 - 120	
Dibromofluoromet	hane (Surr)	100			60 - 140	
Toluene-d8 (Surr)		96			71 - 126	

Analytical Data

Client Sample ID:	MW-28D-82015				
Lab Sample ID: Client Matrix:	480-86254-5 Water				npled: 08/25/2015 1430 ceived: 08/27/2015 0145
		8260C Volatile Organi	c Compounds by	GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/03/2015 2233 09/03/2015 2233	Analysis Batch: Prep Batch:	480-261891 N/A	Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	HP5973S S2166.D 5 mL 5 mL
Analyte		Result (u	ıg/L) Qual	ifier MDL	RL
1,1,1-Trichloroetha	ane	1.0	U	0.82	1.0
1,1,2,2-Tetrachlor	pethane	1.0	U	0.21	1.0
1,1,2-Trichloro-1,2	2,2-trifluoroethane	1.0	U	0.31	1.0
1,1,2-Trichloroetha	ane	0.61	J	0.23	1.0
1,1-Dichloroethan	е	1.6		0.38	1.0
1,1-Dichloroethene	e	0.61	J	0.29	1.0
1,2,3-Trichloroben		1.0	U	0.41	1.0
1,2,4-Trichloroben	zene	1.0	U	0.41	1.0
1,2-Dibromo-3-Ch		1.0	U	0.39	1.0
1,2-Dichlorobenze		1.0	U	0.79	1.0
1,2-Dichloroethan		1.0	U	0.21	1.0
1,2-Dichloropropa		1.0	U	0.72	1.0
1,3-Dichlorobenze		1.0	U	0.78	1.0
1,4-Dichlorobenze		1.0	U	0.84	1.0
1,4-Dioxane	-	40	U	9.3	40
2-Butanone (MEK)	10	U	1.3	10
2-Hexanone		5.0	U	1.2	5.0
4-Methyl-2-pentan	one (MIBK)	5.0	U	2.1	5.0
Acetone		24		3.0	10
Benzene		1.0	U	0.41	1.0
Bromoform		1.0	Ŭ	0.26	1.0
Bromomethane		1.0	Ŭ	0.69	1.0
Carbon disulfide		1.0	Ŭ	0.19	1.0
Carbon tetrachlori	de	1.0	Ŭ	0.27	1.0
Chlorobenzene		1.0	U	0.75	1.0
Chlorobromometh	ane	1.0	Ŭ	0.87	1.0
Chlorodibromome		1.0	Ŭ	0.32	1.0
Chloroethane		1.0	Ŭ	0.32	1.0
Chloroform		0.45	J	0.34	1.0
Chloromethane		1.0	Ŭ	0.35	1.0
cis-1,2-Dichloroeth	nene	1.0	U	0.81	1.0
cis-1,3-Dichloropro		1.0	Ŭ	0.36	1.0
Cyclohexane		1.0	Ŭ	0.18	1.0
Dichlorobromomet	hane	1.0	Ŭ	0.39	1.0
Dichlorodifluorome		1.0	Ŭ	0.68	1.0
Ethylbenzene		1.0	Ŭ	0.74	1.0
Ethylene Dibromid	le	1.0	Ŭ	0.73	1.0
Isopropylbenzene		1.0	U	0.79	1.0
Methyl acetate		2.5	Ŭ	1.3	2.5
Methyl tert-butyl e	ther	1.0	Ŭ	0.16	1.0
Methylcyclohexan		1.0	Ŭ	0.16	1.0
Methylene Chlorid		1.0	U	0.44	1.0
m-Xylene & p-Xyle		2.0	Ŭ	0.66	2.0
o-Xylene		1.0	Ű	0.76	1.0
Styrene		1.0	Ŭ	0.73	1.0
Tetrachloroethene		42	0	0.36	1.0
		74		0.00	1.0

Analytical Data

Lab Sample ID: 480-86254-5 Client Matrix: Water			npled: 08/25/2015 1430
		Date Red	ceived: 08/27/2015 0145
8260C Volatile Organic Compound	s by GO	C/MS	
Analysis Method:8260CAnalysis Batch:480-261891Prep Method:5030CPrep Batch:N/ADilution:1.0Analysis Date:09/03/2015 2233Prep Date:09/03/2015 2233		Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	HP5973S S2166.D 5 mL 5 mL
Analyte Result (ug/L)	Qualifie	er MDL	RL
Toluene 1.0	U	0.51	1.0
trans-1,2-Dichloroethene 1.0	U	0.90	1.0
trans-1,3-Dichloropropene 1.0	U	0.37	1.0
Trichloroethene 1.0	U	0.46	1.0
Trichlorofluoromethane 1.0	U	0.88	1.0
Vinyl chloride 1.0	U	0.90	1.0
Surrogate %Rec	Qualifie	er Acceptar	nce Limits
1,2-Dichloroethane-d4 (Surr) 103		66 - 137	
4-Bromofluorobenzene (Surr) 97		73 - 120	
Dibromofluoromethane (Surr) 103		60 - 140	
Toluene-d8 (Surr) 95		71 - 126	

Analytical Data

Job Number: 480-86254-1

Client Sample ID:	TRIP BLANK				
Lab Sample ID: Client Matrix:	480-86254-6 Water				ate Sampled: 08/25/2015 0000 ate Received: 08/27/2015 0145
	8	260C Volatile Organi	ic Compounds I	by GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/04/2015 1325 09/04/2015 1325	Analysis Batch: Prep Batch:	480-262028 N/A	Instrument ID: Lab File ID: Initial Weight/Vo Final Weight/Vo	
Analyte		Result (u	ıg/L) Qu	alifier MDL	RL
1,1,1-Trichloroetha	ane	1.0	U	0.82	1.0
1,1,2,2-Tetrachloro		1.0	U	0.21	1.0
1,1,2-Trichloro-1,2		1.0	U	0.31	1.0
1,1,2-Trichloroetha	ane	1.0	U	0.23	1.0
1,1-Dichloroethane	e	1.0	U	0.38	1.0
1,1-Dichloroethene	9	1.0	U	0.29	1.0
1,2,3-Trichloroben		1.0	U	0.41	1.0
1,2,4-Trichloroben	zene	1.0	U	0.41	1.0
1,2-Dibromo-3-Chl	loropropane	1.0	U	0.39	1.0
1,2-Dichlorobenze	ne	1.0	U	0.79	1.0
1,2-Dichloroethane	e	1.0	U	0.21	1.0
1,2-Dichloropropa	ne	1.0	U	0.72	1.0
1,3-Dichlorobenze	ne	1.0	U	0.78	1.0
1,4-Dichlorobenze	ne	1.0	U	0.84	1.0
1,4-Dioxane		40	U	9.3	40
2-Butanone (MEK))	10	U	1.3	10
2-Hexanone		5.0	U	1.2	5.0
4-Methyl-2-pentan	one (MIBK)	5.0	U	2.1	5.0
Acetone		10	U	3.0	10
Benzene		1.0	U	0.41	1.0
Bromoform		1.0	U	0.26	1.0
Bromomethane		1.0	U	0.69	1.0
Carbon disulfide		1.0	U	0.19	1.0
Carbon tetrachlorid	de	1.0	U	0.27	1.0
Chlorobenzene		1.0	U	0.75	1.0
Chlorobromometh		1.0	U	0.87	1.0
Chlorodibromomet	thane	1.0	U	0.32	1.0
Chloroethane		1.0	U	0.32	1.0
Chloroform		1.0	U	0.34	1.0
Chloromethane		1.0	U	0.35 0.81	1.0
cis-1,2-Dichloroeth		1.0	U	0.81	1.0
cis-1,3-Dichloropro	ppene	1.0 1.0	U U	0.36	1.0 1.0
Cyclohexane Dichlorobromomet	hana	1.0	U	0.18	1.0
Dichlorodifluorome		1.0	U	0.68	1.0
Ethylbenzene		1.0	U	0.74	1.0
Ethylene Dibromid	0	1.0	U	0.74	1.0
Isopropylbenzene		1.0	U	0.79	1.0
Methyl acetate		2.5	U	1.3	2.5
Methyl tert-butyl et	ther	1.0	U	0.16	1.0
Methylcyclohexane		1.0	U	0.16	1.0
Methylene Chlorid		1.0	Ű	0.44	1.0
m-Xylene & p-Xyle		2.0	U	0.66	2.0
o-Xylene		1.0	U	0.76	1.0
Styrene		1.0	U	0.73	1.0
Tetrachloroethene		1.0	Ű	0.36	1.0
		1.0	0	0.00	

Analytical Data

Client: New York State D.E.C.

Dibromofluoromethane (Surr)

Toluene-d8 (Surr)

Job Number: 480-86254-1

60 - 140

71 - 126

Client Sample ID	: TRIP BLANK				
Lab Sample ID: Client Matrix:	480-86254-6 Water				mpled: 08/25/2015 0000 eceived: 08/27/2015 0145
	82	60C Volatile Organi	c Compounds by	GC/MS	
Analysis Method: Prep Method: Dilution: Analysis Date: Prep Date:	8260C 5030C 1.0 09/04/2015 1325 09/04/2015 1325	Analysis Batch: Prep Batch:	480-262028 N/A	Instrument ID: Lab File ID: Initial Weight/Volume: Final Weight/Volume:	
Analyte		Result (u	ıg/L) Qual	ifier MDL	RL
Toluene		1.0	U	0.51	1.0
trans-1,2-Dichloro	ethene	1.0	U	0.90	1.0
trans-1,3-Dichloro	propene	1.0	U	0.37	1.0
Trichloroethene		1.0	U	0.46	1.0
Trichlorofluoromet	thane	1.0	U	0.88	1.0
Vinyl chloride		1.0	U	0.90	1.0
Surrogate		%Rec	Qual	ifier Accepta	nce Limits
1,2-Dichloroethan	e-d4 (Surr)	90		66 - 137	,
4-Bromofluoroben	zene (Surr)	97		73 - 120)
D ¹¹ d		• •			

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DATA REPORTING QUALIFIERS

Client: New York State D.E.C.

Lab Section	Qualifier	Description
GC/MS VOA		
	U	Analyzed for but not detected.
	J	Indicates an estimated value.

QUALITY CONTROL RESULTS

Job Number: 480-86254-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:480-26	51891				
_CS 480-261891/6	Lab Control Sample	Т	Water	8260C	
VB 480-261891/8	Method Blank	Т	Water	8260C	
480-86254-1	MW-46C-82015	Т	Water	8260C	
480-86254-2	MW-47C-82015	Т	Water	8260C	
480-86254-3	MW-37C-82015	Т	Water	8260C	
480-86254-4	MW-28C-82015	Т	Water	8260C	
480-86254-5	MW-28D-82015	Т	Water	8260C	
Analysis Batch:480-26	52028				
_CS 480-262028/4	Lab Control Sample	Т	Water	8260C	
VB 480-262028/6	Method Blank	Т	Water	8260C	
180-86236-X-8 MS	Matrix Spike	Т	Water	8260C	
80-86236-X-8 MSD	Matrix Spike Duplicate	Т	Water	8260C	
80-86254-6	TRIP BLANK	Т	Water	8260C	

Report Basis

T = Total

Quality Control Results

Job Number: 480-86254-1

Surrogate Recovery Report

8260C Volatile Organic Compounds by GC/MS

Client Matrix: Water

		DCA	BFB	DBFM	TOL
Lab Sample ID	Client Sample ID	%Rec	%Rec	%Rec	%Rec
480-86254-1	MW-46C-82015	105	101	105	99
480-86254-2	MW-47C-82015	106	103	104	98
480-86254-3	MW-37C-82015	97	105	98	100
480-86254-4	MW-28C-82015	101	100	100	96
480-86254-5	MW-28D-82015	103	97	103	95
480-86254-6	TRIP BLANK	90	97	94	94
MB 480-261891/8		100	101	98	101
MB 480-262028/6		87	95	92	93
LCS 480-261891/6		99	102	102	99
LCS 480-262028/4		84	98	91	92

Surrogate	Acceptance Limits
DCA = 1,2-Dichloroethane-d4 (Surr)	66-137
BFB = 4-Bromofluorobenzene (Surr)	73-120
DBFM = Dibromofluoromethane (Surr)	60-140
TOL = Toluene-d8 (Surr)	71-126

Job Number: 480-86254-1

Surrogate Recovery Report

8260C Volatile Organic Compounds by GC/MS

Client Matrix: Water

		DCA	BFB	TOL	
Lab Sample ID	Client Sample ID	%Rec	%Rec	%Rec	
480-86236-X-8 MS		86	103	94	
480-86236-X-8 MSD		85	99	93	

Surrogate	Acceptance Limits
DCA = 1,2-Dichloroethane-d4 (Surr)	66-137
BFB = 4-Bromofluorobenzene (Surr)	73-120
TOL = Toluene-d8 (Surr)	71-126

Method Blank - Batch: 480-261891

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	MB 480-261891/8 Water 1.0 09/03/2015 1527 09/03/2015 1527 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-261891 N/A N/A ug/L			HP5973S S2148.D 5 mL 5 mL	
Analyte		Res	ult	Qual	MDL	RL	
1,1,1-Trichloroeth	ane	1.0		U	0.82	1.0	
1,1,2,2-Tetrachlor	roethane	1.0		U	0.21	1.0	
1,1,2-Trichloro-1,	2,2-trifluoroethane	1.0		U	0.31	1.0	
1,1,2-Trichloroeth	ane	1.0		U	0.23	1.0	
1 1 Dichlereether	-	1.0			0.20	1.0	

1,1,2-Trichloro-1,2,2-trifluoroethane 1.0 U 0.31 1,1,2-Trichloroethane 1.0 U 0.23 1,1-Dichloroethane 1.0 U 0.38 1,1-Dichloroethane 1.0 U 0.38 1,1-Dichloroethane 1.0 U 0.29 1,2-Trichlorobenzene 1.0 U 0.41 1,2,3-Trichlorobenzene 1.0 U 0.41 1,2,4-Trichlorobenzene 1.0 U 0.41 1,2-Dibromo-3-Chloropropane 1.0 U 0.39 1,2-Dichlorobenzene 1.0 U 0.79 1,2-Dichloropenzene 1.0 U 0.72 1,2-Dichloropropane 1.0 U 0.72 1,3-Dichlorobenzene 1.0 U 0.72 1,3-Dichlorobenzene 1.0 U 0.84 1,4-Dichlorobenzene 1.0 U 9.3 1,4-Dichlorobenzene 1.0 U 1.3 2-Butanone (MEK) 10 U 1.2 4-Methyl-2-pentanone (MIBK) 5.0 U 2.1 Ace	1.0 1.0 1.0 1.0
1,1,2-Trichloroethane 1.0 U 0.23 1,1-Dichloroethane 1.0 U 0.38 1,1-Dichloroethane 1.0 U 0.29 1,2,3-Trichlorobenzene 1.0 U 0.41 1,2,4-Trichlorobenzene 1.0 U 0.41 1,2-Dibromo-3-Chloropropane 1.0 U 0.39 1,2-Dichloroethane 1.0 U 0.39 1,2-Dichlorobenzene 1.0 U 0.79 1,2-Dichloroptopane 1.0 U 0.79 1,2-Dichloroptopane 1.0 U 0.72 1,3-Dichloroptopane 1.0 U 0.72 1,3-Dichlorobenzene 1.0 U 0.78 1,4-Dichlorobenzene 1.0 U 0.84 1,4-Dichlorobenzene 1.0 U 9.3 2-Butanone (MEK) 10 U 1.2 2-Hexanone 5.0 U 1.2 4-Methyl-2-pentanone (MIBK) 5.0 U 2.1 Acetone 10 U 3.0	1.0
1,1-Dichloroethane1.0U0.381,1-Dichloroethene1.0U0.291,2,3-Trichlorobenzene1.0U0.411,2,4-Trichlorobenzene1.0U0.411,2-Dibromo-3-Chloropropane1.0U0.391,2-Dichlorobenzene1.0U0.791,2-Dichloroethane1.0U0.791,2-Dichloropropane1.0U0.721,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dichlorobenzene1.0U1.32-Butanone (MEK)10U1.24-Methyl-2-pentanone (MIBK)5.0U2.14-Methyl-2-pentanone (MIBK)5.0U3.0	
1,1-Dichloroethene1.0U0.291,2,3-Trichlorobenzene1.0U0.411,2,4-Trichlorobenzene1.0U0.411,2-Dibromo-3-Chloropropane1.0U0.391,2-Dichlorobenzene1.0U0.791,2-Dichloroptopane1.0U0.791,2-Dichloroptopane1.0U0.791,2-Dichloroptopane1.0U0.721,3-Dichloroptopane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dichlorobenzene1.0U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	10
1,2,3-Trichlorobenzene1.0U0.411,2,4-Trichlorobenzene1.0U0.411,2-Dibromo-3-Chloropropane1.0U0.391,2-Dichlorobenzene1.0U0.791,2-Dichlorobenzene1.0U0.791,2-Dichloropropane1.0U0.791,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dichlorobenzene1.0U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	
1,2,3-Trichlorobenzene 1.0 U 0.41 1,2,4-Trichlorobenzene 1.0 U 0.41 1,2-Dibromo-3-Chloropropane 1.0 U 0.39 1,2-Dichlorobenzene 1.0 U 0.79 1,2-Dichloroethane 1.0 U 0.79 1,2-Dichloropropane 1.0 U 0.72 1,2-Dichloroptopane 1.0 U 0.72 1,2-Dichloroptopane 1.0 U 0.72 1,3-Dichlorobenzene 1.0 U 0.78 1,4-Dichlorobenzene 1.0 U 0.84 1,4-Dichlorobenzene 1.0 U 9.3 2-Butanone (MEK) 10 U 1.3 2-Hexanone 5.0 U 1.2 4-Methyl-2-pentanone (MIBK) 5.0 U 2.1 Acetone 10 U 3.0	1.0
1,2-Dibromo-3-Chloropropane1.0U0.391,2-Dichlorobenzene1.0U0.791,2-Dichloroethane1.0U0.211,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dichlorobenzene40U9.32-Butanone (MEK)10U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
1,2-Dichlorobenzene1.0U0.791,2-Dichloroethane1.0U0.211,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dichlorobenzene40U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
1,2-Dichlorobenzene1.0U0.791,2-Dichloroethane1.0U0.211,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dichlorobenzene40U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
1,2-Dichloroethane1.0U0.211,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dioxane40U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
1,2-Dichloropropane1.0U0.721,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dioxane40U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
1,3-Dichlorobenzene1.0U0.781,4-Dichlorobenzene1.0U0.841,4-Dioxane40U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
1,4-Dichlorobenzene1.0U0.841,4-Dioxane40U9.32-Butanone (MEK)10U1.32-Hexanone5.0U1.24-Methyl-2-pentanone (MIBK)5.0U2.1Acetone10U3.0	1.0
2-Butanone (MEK) 10 U 1.3 2-Hexanone 5.0 U 1.2 4-Methyl-2-pentanone (MIBK) 5.0 U 2.1 Acetone 10 U 3.0	1.0
2-Butanone (MEK) 10 U 1.3 2-Hexanone 5.0 U 1.2 4-Methyl-2-pentanone (MIBK) 5.0 U 2.1 Acetone 10 U 3.0	40
2-Hexanone 5.0 U 1.2 4 4-Methyl-2-pentanone (MIBK) 5.0 U 2.1 4 Acetone 10 U 3.0 4	10
Acetone 10 U 3.0	5.0
Acetone 10 U 3.0	5.0
	10
Benzene 1.0 U 0.41	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
	1.0
,	1.0
	1.0
	2.5
	1.0
	1.0
	1.0
	2.0
	1.0
	1.0

Client: New York State D.E.C.

Method Blank - Batch: 480-261891

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	MB 480-261891/8 Water 1.0 09/03/2015 1527 09/03/2015 1527 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-261891 N/A N/A ug/L			HP5973S S2148.D 5 mL 5 mL	
Analyte		Res	ult	Qual	MDL	RL	
Tetrachloroethen	e	1.0		U	0.36	1.0	
Toluene		1.0		U	0.51	1.0	
trans-1,2-Dichloro	bethene	1.0		U	0.90	1.0	
trans-1,3-Dichloro	propene	1.0		U	0.37	1.0	
Trichloroethene		1.0		U	0.46	1.0	
Trichlorofluorome	thane	1.0		U	0.88	1.0	
Vinyl chloride		1.0		U 0.90		1.0	
Surrogate		%	Rec		Acceptance Lir	nits	
1,2-Dichloroethar	ne-d4 (Surr)	1	00		66 - 137		
4-Bromofluorober	nzene (Surr)	101		73 - 120			
Dibromofluorome	thane (Surr)	98		60 - 140			
Toluene-d8 (Surr))	1	01				

Lab Control Sample - Batch: 480-261891

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	LCS 480-261891/6 Water 1.0 09/03/2015 1437 09/03/2015 1437 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-261891 N/A N/A ug/L	Instrument IE Lab File ID: Initial Weight Final Weight	/Volume:	HP5973S S2146.D 5 mL 5 mL	
Analyte		Spike Amount	Result	% Rec.	Limit		Qual
1,1,1-Trichloroeth	nane	25.0	24.2	97	73 -	126	
1,1,2,2-Tetrachlo		25.0	23.7	95	70 -	126	
1,1,2-Trichloro-1,	2,2-trifluoroethane	25.0	23.2	93	52 -	148	
1,1,2-Trichloroeth		25.0	24.6	98	76 -	122	
1,1-Dichloroethar	ne	25.0	22.7	91	71 -	129	
1,1-Dichloroether	ne	25.0	21.2	85	58 -	121	
1,2,3-Trichlorobe	nzene	25.0	23.3	93	63 -	138	
1,2,4-Trichlorobe		25.0	23.4	93	70 -	122	
1,2-Dibromo-3-Cl	hloropropane	25.0	23.9	96	56 -	134	
1,2-Dichlorobenz		25.0	23.4	94	80 -	124	
1,2-Dichloroethar	ne	25.0	21.0	84	75 -	127	
1,2-Dichloropropa	ane	25.0	23.8	95	76 -	120	
1,3-Dichlorobenz	ene	25.0	23.6	95	77 -	120	
1,4-Dichlorobenz	ene	25.0	23.0	92	75 -	120	
1,4-Dioxane		500	607	121	50 -	174	
2-Butanone (ME	<)	125	132	105	57 -	140	
2-Hexanone		125	131	105	65 -	127	
4-Methyl-2-penta	none (MIBK)	125	125	100	71 -	125	
Acetone		125	138	110	56 -	142	
Benzene		25.0	23.4	94	71 -	124	
Bromoform		25.0	20.8	83	52 -	132	
Bromomethane		25.0	19.8	79	55 -	144	
Carbon disulfide		25.0	24.1	96	59 -	134	
Carbon tetrachlor	ride	25.0	23.3	93	72 -	134	
Chlorobenzene		25.0	22.2	89	72 -		
Chlorobromometh	hane	25.0	23.5	94	72 -		
Chlorodibromome	ethane	25.0	21.6	87	75 -	125	
Chloroethane		25.0	20.9	83	69 -	136	
Chloroform		25.0	22.7	91	73 -		
Chloromethane		25.0	20.2	81	68 -		
cis-1,2-Dichloroet		25.0	22.7	91	74 -		
cis-1,3-Dichlorop	ropene	25.0	25.3	101	74 -		
Cyclohexane		25.0	22.9	92	59 -		
Dichlorobromome		25.0	24.9	99	80 -		
Dichlorodifluorom	nethane	25.0	20.2	81	59 -		
Ethylbenzene		25.0	23.1	93	77 - 1		
Ethylene Dibromi		25.0	23.7	95	77 - 1		
Isopropylbenzene	9	25.0	24.6	99	77 - 1		
Methyl acetate		125	119	95	74 -		
Methyl tert-butyl		25.0	24.8	99	64 -		
Methylcyclohexar	ne	25.0	24.0	96	61 -	138	

Lab Control Sample - Batch: 480-261891

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	LCS 480-261891/6 Water 1.0 09/03/2015 1437 09/03/2015 1437 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-261891 N/A N/A ug/L	Instrument II Lab File ID: Initial Weigh Final Weight	S2146.[t/Volume: 5 mL	-
Analyte		Spike Amount	Result	% Rec.	Limit	Qual
Methylene Chlori	de	25.0	24.5	98	57 - 132	
m-Xylene & p-Xyl	ene	25.0	23.4	94	76 - 122	
o-Xylene		25.0	24.3	97	76 - 122	
Styrene		25.0	25.3	101	70 - 130	
Tetrachloroethen	e	25.0	22.5	90	74 - 122	
Toluene		25.0	23.0	92	80 - 122	
trans-1,2-Dichloro	bethene	25.0	23.3	93	73 - 127	
trans-1,3-Dichloro	opropene	25.0	24.2	97	72 - 123	
Trichloroethene		25.0	22.9	92	74 - 123	
Trichlorofluorome	ethane	25.0	23.6	95	62 - 152	
Vinyl chloride		25.0	20.3	81	65 - 133	
Surrogate		%	Rec	Acc	ceptance Limits	
1,2-Dichloroethar	ne-d4 (Surr)	9	9		66 - 137	
4-Bromofluorober	nzene (Surr)	1	02		73 - 120	
Dibromofluorome	thane (Surr)	1	02		60 - 140	
Toluene-d8 (Surr))	9	9		71 - 126	

Method Blank - Batch: 480-262028

Method: 8260C Preparation: 5030C

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	MB 480-262028/6 Water 1.0 09/04/2015 1214 09/04/2015 1214 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-262028 N/A N/A ug/L	Lab Fi Initial	nent ID: le ID: Weight/Volume: Veight/Volume:	HP5975T T8290.D 5 mL 5 mL	
Analyte		Res	ult	Qual	MDL	RL	
1,1,1-Trichloroeth	ane	1.0		U	0.82	1.0	
1,1,2,2-Tetrachlor		1.0		U	0.21	1.0	
	2,2-trifluoroethane	1.0		U	0.31	1.0	
1,1,2-Trichloroeth	-	1.0		U	0.23	1.0	
1,1-Dichloroethan		1.0		U	0.38	1.0	
1,1-Dichloroethen		1.0		Ŭ	0.29	1.0	
1,2,3-Trichlorober		1.0		U	0.41	1.0	
1,2,4-Trichlorober		1.0		Ŭ	0.41	1.0	
1,2-Dibromo-3-Ch		1.0		U	0.39	1.0	
1,2-Dichlorobenze		1.0		U	0.79	1.0	
1,2-Dichloroethan		1.0		U	0.21	1.0	
1,2-Dichloropropa		1.0		U	0.72	1.0	
1,3-Dichlorobenze		1.0		Ŭ	0.78	1.0	
1,4-Dichlorobenze		1.0		U	0.84	1.0	
1,4-Dioxane		40		U	9.3	40	
2-Butanone (MEK	()	10		U	1.3	10	
2-Hexanone	-)	5.0		Ŭ	1.2	5.0	
4-Methyl-2-pentar	none (MIBK)	5.0		U	2.1	5.0	
Acetone		10		Ŭ	3.0	10	
Benzene		1.0		U	0.41	1.0	
Bromoform		1.0		U	0.26	1.0	
Bromomethane		1.0		U	0.69	1.0	
Carbon disulfide		1.0		U	0.19	1.0	
Carbon tetrachlor	ide	1.0		U	0.27	1.0	
Chlorobenzene		1.0		Ŭ	0.75	1.0	
Chlorobromometh	nane	1.0		U	0.87	1.0	
Chlorodibromome		1.0		Ŭ	0.32	1.0	
Chloroethane		1.0		U	0.32	1.0	
Chloroform		1.0		U	0.34	1.0	
Chloromethane		1.0		U	0.35	1.0	
cis-1,2-Dichloroet	hene	1.0		U	0.81	1.0	
cis-1,3-Dichloropr		1.0		U	0.36	1.0	
Cyclohexane	•	1.0		U	0.18	1.0	
Dichlorobromome	ethane	1.0		U	0.39	1.0	
Dichlorodifluorom		1.0		U	0.68	1.0	
Ethylbenzene		1.0		U	0.74	1.0	
Ethylene Dibromi	de	1.0		U	0.73	1.0	
Isopropylbenzene		1.0		U	0.79	1.0	
Methyl acetate		2.5		U	1.3	2.5	
Methyl tert-butyl e	ether	1.0		U	0.16	1.0	
Methylcyclohexar		1.0		U	0.16	1.0	
Methylene Chlorid		1.0		U	0.44	1.0	
m-Xylene & p-Xyl		2.0		U	0.66	2.0	
o-Xylene		1.0		Ŭ	0.76	1.0	
Styropo		1.0			0.73	1.0	

Styrene

1.0

U

0.73

1.0

Client: New York State D.E.C.

Method Blank - Batch: 480-262028

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	MB 480-262028/6 Water 1.0 09/04/2015 1214 09/04/2015 1214 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-262028 N/A N/A ug/L			HP5975T T8290.D 5 mL 5 mL	
Analyte		Res	ult	Qual	MDL	RL	
Tetrachloroethen	е	1.0		U	0.36	1.0	
Toluene		1.0		U	0.51	1.0	
trans-1,2-Dichloro	bethene	1.0		U	0.90	1.0	
trans-1,3-Dichloro	opropene	1.0		U	0.37	1.0	
Trichloroethene		1.0		U	0.46	1.0	
Trichlorofluorome	ethane	1.0		U	0.88	1.0	
Vinyl chloride		1.0		U	0.90	1.0	
Surrogate		%	Rec		Acceptance Lir	nits	
1,2-Dichloroethar	ne-d4 (Surr)	8	7		66 - 137		
4-Bromofluorober	nzene (Surr)	95		73 - 120			
Dibromofluorome	thane (Surr)	9	2	60 - 140			
Toluene-d8 (Surr)	93					

Client: New York State D.E.C.

Lab Control Sample - Batch: 480-262028

Analyte Spike Amount Result % Rec. Limit Qual 1,1,1-Trichloroethane 25.0 24.0 96 73.126 1,1,2-Trichloroethane 25.0 26.2 105 52.148 1,1,2-Trichloroethane 25.0 26.2 105 52.148 1,12-Trichloroethane 25.0 25.2 101 58.121 1,1-Dichloroethane 25.0 25.2 101 58.121 1,2-Trichlorobenzene 25.0 24.2 97 70.122 1,2-Dichlorobenzene 25.0 23.4 94 75.127 1,2-Dichlorobenzene 25.0 25.4 106 76.120 1,2-Dichlorobenzene 25.0 25.4 102 77.120 1,2-Dichlorobenzene 25.0 25.6 102 77.120 1,4-Diokane 25.0 25.6 102 77.120 1,4-Diokane 25.0 25.6 102 77.120 1,4-Diokane 25.0 25.8 102 71.125	Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	LCS 480-262028/4 Water 1.0 09/04/2015 1014 09/04/2015 1014 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-262028 N/A N/A ug/L	Instrument II Lab File ID: Initial Weigh Final Weight	٦ t/Volume: t	HP5975T F8285.D 5 mL 5 mL	
1,1,2,2-Tetrachloroethane 25.0 24.9 100 70 - 126 1,1,2-Trichloroethane 25.0 25.1 100 76 - 122 1,1-Dichloroethane 25.0 25.8 103 71 - 129 1,1-Dichloroethane 25.0 25.2 101 58 - 121 1,2-Dichloroethane 25.0 25.2 101 58 - 121 1,2-Dichloroethane 25.0 24.2 97 70 - 122 1,2-Dichloroethane 25.0 25.3 101 80 - 124 1,2-Dichloroptapane 25.0 25.4 94 63 - 134 1,2-Dichloroptapane 25.0 25.4 106 76 - 120 1,2-Dichloroptapane 25.0 25.6 102 77 - 120 1,2-Dichloroptapane 25.0 25.4 106 76 - 120 1,4-Dichlorobenzene 25.0 25.4 102 77 - 120 1,4-Dichlorobenzene 125 121 97 71 - 125 1,4-Dichlorobenzene 25.0 25.6 102 71 - 125 1,4-Dichlorobenzene 25.0 25.6 102 71 +	Analyte		Spike Amount	Result	% Rec.	Limit		Qual
1,1,2-Trichloro-1,2,2-trifluoroethane 25.0 25.1 105 52 + 148 1,1-Dichloroethane 25.0 25.1 100 76 - 122 1,1-Dichloroethane 25.0 25.2 101 58 + 121 1,2,3-Trichlorobenzene 25.0 25.2 101 58 + 121 1,2,4-Trichlorobenzene 25.0 24.2 97 70 - 122 1,2-Dichloroporpane 25.0 25.3 101 80 - 124 1,2-Dichlorobenzene 25.0 25.4 102 77 - 122 1,2-Dichlorobenzene 25.0 25.4 102 77 - 120 1,2-Dichlorobenzene 25.0 25.4 102 77 - 120 1,2-Dichlorobenzene 25.0 25.4 102 77 - 120 1,4-Dioxane 250 25.4 102 77 - 120 1,4-Dioxane 125 134 108 57 - 140 2-Hutanone (MIBK) 125 121 97 71 - 125 Acetone 125 148 119 56 - 142 Benzene 25.0 25.6 102 72 - 120	1,1,1-Trichloroeth	ane	25.0	24.0	96	73 - 12	26	
1,1,2-Trichloroethane 25.0 25.1 100 76 - 122 1,1-Dichloroethane 25.0 25.8 103 71 - 129 1,1-Dichloroethane 25.0 25.2 101 58 - 121 1,2-A-Trichlorobenzene 25.0 23.6 94 63 - 138 1,2-Ditromo-3-Chloropropane 25.0 22.2 89 56 - 134 1,2-Ditromo-3-Chloropropane 25.0 23.4 94 75 - 127 1,2-Ditromo-3-Chloropropane 25.0 25.6 102 77 - 120 1,4-Dichlorobenzene 25.0 25.6 102 77 - 120 1,4-Dichlorobenzene 25.0 25.4 102 75 - 120 1,4-Dichlorobenzene 25.0 25.4 102 75 - 120 1,4-Dichlorobenzene 25.0 25.4 102 75 - 120 1,4-Dichlorobenzene 125 127 101 65 - 127 4-Methyl-2-pentanone (MEK) 125 121 97 71 - 125 Actone 25.0 25.6 102 71 - 124 Bromoform 25.0 25.5 102	1,1,2,2-Tetrachlo	roethane	25.0	24.9	100	70 - 12	26	
1.1-Dichloroethane 25.0 25.8 103 71-129 1.1-Dichloroethene 25.0 25.2 101 58-121 1.2.3-Trichlorobenzene 25.0 23.6 94 63-138 1.2.4-Trichlorobenzene 25.0 24.2 97 70-122 1.2-Dichlorobenzene 25.0 22.2 89 56-134 1.2-Dichlorobenzene 25.0 25.4 101 80-124 1.2-Dichlorobenzene 25.0 25.4 102 77-120 1.4-Dichlorobenzene 25.0 25.4 102 75-120 1.4-Dichlorobenzene 25.0 25.4 102 75-120 1.4-Dichlorobenzene 25.0 25.4 102 75-120 1.4-Dichlorobenzene 25.0 25.4 102 77-120 1.4-Dichlorobenzene 25.0 25.6 102 77-140 2-Hexanone 125 121 97 71-125 Acetone 125 121 97 71-124 Bromoform 25.0 22.8 91 52-132 Bromoform 25	1,1,2-Trichloro-1,	2,2-trifluoroethane	25.0	26.2	105	52 - 14	8	
1.1-Dicklorosentene25.025.210158 - 1211.2.4-Trichlorobenzene25.023.69463 - 1381.2.4-Trichlorobenzene25.022.28956 - 1341.2-Dichlorobenzene25.023.310180 - 1241.2-Dichlorobenzene25.023.49475 - 1271.2-Dichlorobenzene25.025.610277 - 1201.3-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.012710165 - 1271.4-Dichlorobenzene1251219771 - 1252-Butanone (MEK)12514811956 - 1422-Hexanone1251219771 - 124Acetone25.025.610271 - 124Bromoform25.025.510272 - 134Carbon disulfide25.025.510272 - 134Chlorobranene25.025.510272 - 130Chloroforomomethane25.025.510272 - 130Chloroforomomethane25.024.69873 - 127Chloroforomomethane25.025.710371 - 124Chloroforomomethane25.024.698	1,1,2-Trichloroeth	ane	25.0	25.1	100	76 - 12	22	
1.2.3-Trichlorobenzene 25.0 23.6 94 63 - 138 1.2.4-Trichlorobenzene 25.0 24.2 97 70 - 122 1.2-Ditoropopane 25.0 25.3 101 80 - 124 1.2-Ditoropopane 25.0 25.3 101 80 - 124 1.2-Dichloropopane 25.0 25.4 102 77 - 120 1.3-Dichloropopane 25.0 25.4 102 77 - 120 1.4-Dickane 500 552 116 50 - 174 2-Butanone (MEK) 125 134 108 57 - 140 2-Hexanone 125 121 97 71 - 125 Acetone 125 148 119 56 - 142 Benzene 25.0 25.6 102 71 - 125 Acetone 125 148 119 56 - 142 Benzene 25.0 25.6 102 71 - 124 Bromoform 25.0 25.5 102 72 - 134 Chlorobenzene 25.0 25.5 102 72 - 134 Chlorodiforomethane 25.0 25.5	1,1-Dichloroethar	ne	25.0	25.8	103	71 - 12	29	
1.2.4-Trichlorobenzene 25.0 24.2 97 70 - 122 1.2-Dichloropo-3-Chloropropane 25.0 22.2 89 56 - 134 1.2-Dichlorobenzene 25.0 23.4 94 75 - 127 1.2-Dichloropropane 25.0 25.6 102 77 - 120 1.3-Dichlorobenzene 25.0 25.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.4 102 75 - 120 1.4-Dichlorobenzene 25.0 25.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.7 134 108 57 - 140 2-Butanone (MEK) 125 121 97 71 - 125 4-Methyl-2-pentanone (MIBK) 125 124 97 71 - 124 Bromoform 25.0 22.8 91 52 - 132 Bromoform 25.0 25.5 102 72 - 134 Carbon disulfde 25.0 25.5 102 <td< td=""><td>1,1-Dichloroether</td><td>ne</td><td>25.0</td><td>25.2</td><td>101</td><td>58 - 12</td><td>21</td><td></td></td<>	1,1-Dichloroether	ne	25.0	25.2	101	58 - 12	21	
1.2-Dibromo-3-Chloropropane 25.0 22.2 89 56 - 134 1.2-Dichlorobetnaene 25.0 25.3 101 80 - 124 1.2-Dichlorobetnaene 25.0 23.4 94 75 - 127 1.2-Dichlorobetnaene 25.0 26.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.4 102 75 - 120 1.4-Dichlorobenzene 25.0 25.4 102 75 - 120 1.4-Dichlorobenzene 25.0 25.4 102 75 - 120 1.4-Dichlorobenzene 125 134 108 57 - 140 2-Hexanone 125 121 97 71 - 125 Acetone 125 148 119 56 - 142 Benzene 25.0 25.6 102 71 - 124 Bromoform 25.0 22.2 89 55 - 144 Carbon disulfide 25.0 25.5 102 72 - 130 Chlorobenzene 25.0 25.5 102 72 - 130 Chlorobromo	1,2,3-Trichlorobe	nzene	25.0	23.6	94	63 - 13	8	
1.2-Dichlorobenzene25.025.310180 - 1241.2-Dichloroethane25.023.49475 - 1271.2-Dichloropopane25.026.410676 - 1201.3-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.025.410275 - 1201.4-Dicoane12513410857 - 1402-Butanone (MEK)1251219771 - 1252-Hexanone1251219771 - 125Acetone12512811966 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromoform25.022.89152 - 132Bromomethane25.023.59472 - 134Carbon disulfide25.023.59472 - 134Chlorobenzene25.023.89575 - 125Chlorobenmethane25.023.89575 - 125Chlorobenzene25.023.89575 - 125Chlorobenzene25.023.89575 - 125Chlorobenzene25.024.910074 - 124Chlorobenzene25.024.910074 - 124Chlorofirom25.024.910074 - 124Chloroform25.026.710799 - 135Chloroform25.026.710377 - 120Chloroformomethane25.026.710377	1,2,4-Trichlorobe	nzene	25.0	24.2	97	70 - 12	22	
1.2-Dichloroptnane 25.0 23.4 94 75 - 127 1.2-Dichloroptnane 25.0 26.4 106 76 - 120 1.3-Dichlorobenzene 25.0 25.6 102 77 - 120 1.4-Dichlorobenzene 25.0 25.4 102 75 - 120 1.4-Dichlorobenzene 25.0 25.4 102 75 - 120 1.4-Dichlorobenzene 25.0 25.4 102 77 - 120 1.4-Dichlorobenzene 25.0 25.1 116 50 - 174 2-Butanone (MEK) 125 121 97 71 - 125 Acetone 125 148 119 56 - 127 Acetone 25.0 25.6 102 71 - 124 Bromoform 25.0 22.2 89 55 - 144 Carbon disulfide 25.0 25.5 102 72 - 134 Chlorobenzene 25.0 25.5 102 72 - 130 Chlorobromoethane 25.0 26.2 105 59 - 134 Carbon disulfide 25.0 26.5 102 72 - 130 Chlorobromomethane </td <td>1,2-Dibromo-3-Cl</td> <td>nloropropane</td> <td>25.0</td> <td>22.2</td> <td>89</td> <td>56 - 13</td> <td>34</td> <td></td>	1,2-Dibromo-3-Cl	nloropropane	25.0	22.2	89	56 - 13	34	
1.2-Dichloropropane25.026.410676 - 1201.3-Dichlorobenzene25.025.610277 - 1201.4-Dicxane50058211650 - 1742-Butanone (MEK)12513410857 - 1402-Hexanone12512710165 - 1274-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromoform25.022.28955 - 144Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.025.510272 - 130Chlorobenzene25.025.510272 - 130Chlorobromethane25.025.510272 - 130Chlorobromethane25.025.910469 - 136Chlorobromethane25.024.79972 - 130Chlorobromethane25.024.79973 - 127Chlorobromethane25.024.89575 - 125Chloroform25.024.910074 - 124Cis-1,2-Dichloropropene25.026.710759 - 135Chloroform25.024.910074 - 124Cis-1,2-Dichloropropene25.026.710759 - 135Dichloroffluoromethane25.026.710759 - 135Dichloroffluorom	1,2-Dichlorobenze	ene	25.0	25.3	101	80 - 12	24	
1.3-Dichlorobenzene25.025.610277 - 1201.4-Dichlorobenzene25.025.410275 - 1201.4-Dichlorobenzene50058211650 - 1742-Butanone (MEK)12512410857 - 1402-Hexanone12512710165 - 1274-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone25.025.610271 - 124Benzene25.025.610271 - 124Bromoform25.025.610272 - 132Bromomethane25.025.59472 - 134Carbon tetrachloride25.025.510272 - 134Chlorobenzene25.025.510272 - 134Chlorobendhane25.023.89575 - 125Chlorobendhane25.024.79972 - 130Chlorobendhane25.024.89168 - 124Chloroform25.024.89873 - 127Chloroform25.024.69873 - 127Chloroform25.024.910074 - 124cis-1,2-Dichloropene25.024.910074 - 124cis-1,2-Dichloropene25.026.710759 - 135Dichlorofflano25.026.710759 - 135Dichlorofflano25.026.710467 - 123Dichlorofflano25.026.710475 - 135Dichlorofflanomethane25.0	1,2-Dichloroethar	ne	25.0	23.4	94	75 - 12	27	
1.4-Dichlorobenzene25.025.410275 - 1201.4-Dioxane50058211650 - 1742-Butanone (MEK)12513410857 - 1402-Hexanone12512710165 - 1274-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.026.210559 - 134Carbon disulfide25.023.59472 - 134Chlorobenzene25.025.510272 - 130Chlorobenzene25.025.910469 - 136Chlorobenmethane25.025.910469 - 136Chlorobromomethane25.024.79972 - 120Chlorobromomethane25.025.910469 - 136Chloroform25.024.79973 - 125Chloroform25.024.69873 - 127Chloroform25.024.69873 - 127Chloroform25.024.910074 - 124cis-1,2-Dichloroptene25.026.710759 - 135Dichloroffluoromethane25.026.710759 - 135Dichloroffluoromethane25.026.710759 - 135Dichloroffluoromethane25.026.710377 - 123Dichloroffluoromethane <t< td=""><td>1,2-Dichloropropa</td><td>ane</td><td>25.0</td><td>26.4</td><td>106</td><td>76 - 12</td><td>20</td><td></td></t<>	1,2-Dichloropropa	ane	25.0	26.4	106	76 - 12	20	
1,4-Dioxane50058211650 - 1742-Butanone (MEK)12513410857 - 1402-Hexanone12512710165 - 1274-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.026.210559 - 134Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.025.510272 - 134Chlorobromethane25.025.510272 - 134Chlorobromethane25.025.910469 - 136Chlorobromothane25.025.910469 - 136Chlorobromothane25.024.79972 - 130Chlorobromothane25.025.910469 - 136Chloroform25.024.79973 - 127Chlorobromothane25.024.910074 - 124Cisloroform25.024.910074 - 124Cyclohexane25.024.910080 - 122Dichlorobromothane25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromothane25.026.110477 - 123Ethylene Dibromide25.026.110477 - 123Ethylene Dibromide25.0 <t< td=""><td>1,3-Dichlorobenzo</td><td>ene</td><td>25.0</td><td>25.6</td><td>102</td><td>77 - 12</td><td>20</td><td></td></t<>	1,3-Dichlorobenzo	ene	25.0	25.6	102	77 - 12	20	
2-Butanone (MEK)12513410857 - 1402-Hexanone12512710165 - 1274-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.026.210559 - 134Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.025.510272 - 130Chlorobenzene25.025.510272 - 130Chlorobenzene25.025.910469 - 136Chlorobenzene25.025.910469 - 136Chlorobromotethane25.024.79972 - 130Chloroform25.024.69873 - 127Chloroform25.024.69873 - 127Chloroform25.024.910074 - 124cis-1,3-Dichloropropene25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.710477 - 123Bitylene25.026.710377 - 122Dichlorobromomethane </td <td>1,4-Dichlorobenzo</td> <td>ene</td> <td>25.0</td> <td>25.4</td> <td>102</td> <td>75 - 12</td> <td>20</td> <td></td>	1,4-Dichlorobenzo	ene	25.0	25.4	102	75 - 12	20	
2-Hexanone12512710165 - 1274-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.022.28955 - 144Carbon disulfide25.023.59472 - 134Chlorobenzene25.023.59472 - 134Chlorobromomethane25.023.597 - 71 - 125Chlorobromomethane25.023.89575 - 125Chlorobromomethane25.025.910469 - 136Chlorodibromomethane25.023.89575 - 125Chlorodibromomethane25.023.89575 - 125Chlorodibromomethane25.024.69873 - 127Chloroform25.024.69873 - 127Chloromethane25.023.99674 - 124cis-1,2-Dichloroethene25.023.99674 - 124cis-1,3-Dichloropropene25.026.710759 - 135Dichlorobromomethane25.026.110477 - 123Dichlorobromomethane25.026.110477 - 123Dichlorodifluoromethane25.026.110477 - 123Dichlorobromomethane25.025.710377 - 122Dichlorodifluoromethane25.025.710377 - 122	1,4-Dioxane		500	582	116	50 - 17	'4	
4-Methyl-2-pentanone (MIBK)1251219771 - 125Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.022.28955 - 144Carbon disulfide25.026.210559 - 134Carbon disulfide25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobenzene25.023.89575 - 125Chlorodibromomethane25.023.89575 - 125Chloroform25.024.69873 - 127Chloroform25.024.69873 - 127Chloromethane25.024.910074 - 124cis-1,2-Dichloroptopene25.024.910074 - 124cis-1,3-Dichloroptopene25.024.910074 - 124Cyclohexane25.024.910074 - 124Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.110477 - 123Bitylbenzene25.025.710377 - 122Nethyl acetate25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	2-Butanone (MEk	()	125	134	108	57 - 14	0	
Acetone12514811956 - 142Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.022.28955 - 144Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobenzene25.023.89575 - 125Chlorobenzene25.023.89575 - 125Chlorobenzene25.024.79972 - 130Chlorobenzene25.024.89873 - 127Chlorobenzene25.024.89873 - 125Chlorobenzene25.024.69873 - 127Chlorobenzene25.024.910074 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.110477 - 123Ethylenzene25.026.110477 - 123Ethylenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	2-Hexanone		125	127	101	65 - 12	27	
Benzene25.025.610271 - 124Bromoform25.022.89152 - 132Bromomethane25.022.28955 - 144Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobromomethane25.023.89575 - 125Chlorobromomethane25.025.910469 - 136Chlorobromomethane25.024.69873 - 127Chloroform25.024.69873 - 127Chlorobrhane25.024.910074 - 124cis-1,2-Dichloroethene25.023.99674 - 124cis-1,3-Dichloropropene25.026.710759 - 135Dichlorobromomethane25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.023.99677 - 120Dichlorobromomethane25.023.99677 - 120Dichlorobromomethane25.023.99677 - 120Dichlorobromide25.023.99677 - 120Borpopylbenzene25.025.710377 - 120Borpopylbenzene25.025.710377 - 120Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127 <td>4-Methyl-2-penta</td> <td>none (MIBK)</td> <td>125</td> <td>121</td> <td>97</td> <td>71 - 12</td> <td>25</td> <td></td>	4-Methyl-2-penta	none (MIBK)	125	121	97	71 - 12	25	
Bromoform25.022.89152 - 132Bromomethane25.022.28955 - 144Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobinomethane25.023.89575 - 125Chlorobinomethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloroform25.024.910074 - 124cis-1,2-Dichloroptene25.023.99674 - 124cis-1,3-Dichloroptene25.024.910080 - 122Dichlorobifluoromethane25.024.910080 - 122Dichlorobifluoromethane25.024.910080 - 122Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylbenzene25.025.710377 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Acetone		125	148	119	56 - 14	2	
Bromomethane25.022.28955 - 144Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobromomethane25.024.79972 - 130Chlorobromomethane25.023.89575 - 125Chlorobromomethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloroform25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorobromomethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylbenzene25.025.710377 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Benzene		25.0	25.6	102	71 - 12	24	
Carbon disulfide25.026.210559 - 134Carbon tetrachloride25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobromomethane25.024.79972 - 130Chlorodibromomethane25.023.89575 - 125Chloroethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124Cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.026.710759 - 135Dichlorodifluoromethane25.024.910080 - 122Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.110477 - 123Dichlorobromomethane25.026.110477 - 123Dichlorodifluoromethane25.025.710377 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Bromoform		25.0	22.8	91	52 - 13	32	
Carbon tetrachloride25.023.59472 - 134Chlorobenzene25.025.510272 - 120Chlorobromomethane25.024.79972 - 130Chlorodibromomethane25.023.89575 - 125Chloroethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.023.99674 - 124cis-1,3-Dichloropropene25.026.710759 - 135Dichlorobromomethane25.026.710759 - 135Dichlorobromomethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylpenzene25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Bromomethane		25.0	22.2	89	55 - 14	4	
Chlorobenzene25.025.510272 - 120Chlorobromomethane25.024.79972 - 130Chlorodibromomethane25.023.89575 - 125Chloroethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorodifluoromethane25.024.910080 - 122Dichlorobromomethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylpenzene25.025.710377 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Carbon disulfide		25.0	26.2	105	59 - 13	34	
Chlorobromomethane25.024.79972 - 130Chlorodibromomethane25.023.89575 - 125Chloroethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorobromomethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Carbon tetrachlor	ide	25.0	23.5	94	72 - 13	34	
Chlorodibromomethane25.023.89575 - 125Chloroethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobtromomethane25.024.910080 - 122Dichlorobtromomethane25.026.710759 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 123Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Chlorobenzene		25.0	25.5	102	72 - 12	20	
Chloroethane25.025.910469 - 136Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.025.710377 - 122Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Chlorobromometh	nane	25.0	24.7	99	72 - 13	0	
Chloroform25.024.69873 - 127Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.026.110477 - 123Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.025.710377 - 122Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Chlorodibromome	ethane	25.0	23.8	95	75 - 12	25	
Chloromethane25.021.38568 - 124cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.015.26159 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Chloroethane		25.0	25.9	104	69 - 13	6	
cis-1,2-Dichloroethene25.024.910074 - 124cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.015.26159 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Chloroform		25.0	24.6	98	73 - 12	27	
cis-1,3-Dichloropropene25.023.99674 - 124Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.015.26159 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Chloromethane		25.0	21.3	85	68 - 12	24	
Cyclohexane25.026.710759 - 135Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.015.26159 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	cis-1,2-Dichloroet	hene	25.0	24.9	100	74 - 12	24	
Dichlorobromomethane25.024.910080 - 122Dichlorodifluoromethane25.015.26159 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	cis-1,3-Dichloropi	ropene	25.0	23.9	96	74 - 12	24	
Dichlorodifluoromethane25.015.26159 - 135Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Cyclohexane		25.0	26.7	107	59 - 13	5	
Ethylbenzene25.026.110477 - 123Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Dichlorobromome	ethane	25.0	24.9	100	80 - 12	2	
Ethylene Dibromide25.023.99677 - 120Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Dichlorodifluorom	ethane	25.0	15.2	61	59 - 13	5	
Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	Ethylbenzene		25.0	26.1	104	77 - 12	23	
Isopropylbenzene25.025.710377 - 122Methyl acetate1251189474 - 133Methyl tert-butyl ether25.023.59464 - 127	-	de	25.0	23.9	96	77 - 12	20	
Methyl acetate 125 118 94 74 - 133 Methyl tert-butyl ether 25.0 23.5 94 64 - 127	-		25.0	25.7	103	77 - 12	22	
Methyl tert-butyl ether 25.0 23.5 94 64 - 127				118		74 - 13	3	
	•	ether	25.0		94			
	• •		25.0	24.5	98	61 - 13	8	

Client: New York State D.E.C.

Lab Control Sample - Batch: 480-262028

Method: 8260C Preparation: 5030C

Lab Sample ID: Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	LCS 480-262028/4 Water 1.0 09/04/2015 1014 09/04/2015 1014 N/A	Analysis Batch: Prep Batch: Leach Batch: Units:	480-262028 N/A N/A ug/L	Instrument II Lab File ID: Initial Weigh Final Weight	T8285. t/Volume: 5 mL	-
Analyte		Spike Amount	Result	% Rec.	Limit	Qual
Methylene Chlorid	de	25.0	24.6	98	57 - 132	
m-Xylene & p-Xyl	ene	25.0	25.1	100	76 - 122	
o-Xylene		25.0	25.0	100	76 - 122	
Styrene		25.0	25.9	104	70 - 130	
Tetrachloroethen	e	25.0	25.5	102	74 - 122	
Toluene		25.0	25.9	104	80 - 122	
trans-1,2-Dichloro	bethene	25.0	25.5	102	73 - 127	
trans-1,3-Dichloro	propene	25.0	24.1	96	72 - 123	
Trichloroethene		25.0	25.2	101	74 - 123	
Trichlorofluorome	thane	25.0	26.3	105	62 - 152	
Vinyl chloride		25.0	21.6	86	65 - 133	
Surrogate		%	Rec	Acc	eptance Limits	
1,2-Dichloroethan	ie-d4 (Surr)	8	4		66 - 137	
4-Bromofluorober	nzene (Surr)	9	8		73 - 120	
Dibromofluorome	thane (Surr)	9	1		60 - 140	
Toluene-d8 (Surr)		9	2		71 - 126	

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HP5975T

T8310.D

Method: 8260C

Instrument ID:

Lab File ID:

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Preparation: 5030C

Initial Weight/Volume: 5 mL

Job Number: 480-86254-1

Client: New York State D.E.C.

MS Lab Sample ID: 480-86236-X-8 MS

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Water

Client Matrix:

Dilution:

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 480-262028

Analysis Date: Prep Date: Leach Date:	20 09/04/2015 2010 09/04/2015 2010 N/A	Lea	ch Batch:	N/A		eight/Volume: eight/Volume:	5 mL 5 mL 5 mL	
MSD Lab Sample Client Matrix: Dilution: Analysis Date: Prep Date: Leach Date:	ID: 480-86236-X-8 MSD Water 20 09/04/2015 2033 09/04/2015 2033 N/A	Prep	llysis Batch: p Batch: ch Batch:	480-262028 N/A N/A			HP5975T T8311.D 5 mL 5 mL 5 mL	
			Rec.					
Analyte		MS	MSD	Limit	RPD	RPD Limit	MS Qual	MSD Qual
1,1,1-Trichloroeth	ane	105	97	73 - 126	8	15		
1,1,2,2-Tetrachlor	roethane	101	106	70 - 126	5	15		
1,1,2-Trichloroeth	ane	104	104	76 - 122	0	15		
1,1-Dichloroethar	ne	112	103	71 - 129	9	20		
1,1-Dichloroether	ne	114	101	58 - 121	12	16		
1,2-Dibromo-3-Cł	nloropropane	88	103	56 - 134	15	15		
1,2-Dichlorobenze	ene	104	102	80 - 124	1	20		
1,2-Dichloroethar	ne	101	96	75 - 127	5	20		
1,2-Dichloropropa	ane	112	105	76 - 120	6	20		
2-Butanone (MEK	()	111	105	57 - 140	5	20		
2-Hexanone		100	102	65 - 127	2	15		
4-Methyl-2-pentar	none (MIBK)	109	103	71 - 125	6	35		
Acetone		110	106	56 - 142	4	15		
Benzene		109	103	71 - 124	6	13		
Bromoform		94	91	52 - 132	3	15		
Bromomethane		100	95	55 - 144	6	15		
Carbon disulfide		114	106	59 - 134	7	15		
Carbon tetrachlor	ide	100	94	72 - 134	7	15		
Chlorobenzene		107	104	72 - 120	3	25		
Chlorobromometh	nane	105	102	72 - 130	3	15		
Chlorodibromome	ethane	94	97	75 - 125	3	15		
Chloroethane		112	106	69 - 136	5	15		

Analysis Batch:

Prep Batch:

Leach Batch:

480-262028

N/A

N/A

99 Chloroform 105 73 - 127 Chloromethane 96 89 68 - 124 102 74 - 124 cis-1,2-Dichloroethene 109 cis-1,3-Dichloropropene 104 94 74 - 124 80 - 122 Dichlorobromomethane 108 99 Dichlorodifluoromethane 69 64 59 - 135 Ethylbenzene 114 107 77 - 123 Ethylene Dibromide 100 98 77 - 120 Methylene Chloride 106 101 57 - 132 76 - 122 m-Xylene & p-Xylene 107 100

TestAmerica Buffalo

Toluene-d8 (Surr)

HP5975T

T8310.D

71 - 126

Method: 8260C

Instrument ID:

Lab File ID:

Preparation: 5030C

Job Number: 480-86254-1

Client: New York State D.E.C.

MS Lab Sample ID: 480-86236-X-8 MS

Water

Client Matrix:

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 480-262028

Dilution: Analysis Date: Prep Date: Leach Date:	20 09/04/2015 2010 09/04/2015 2010 N/A	Lea	ach Batch:	N/A		eight/Volume: eight/Volume:	5 mL 5 mL 5 mL	
MSD Lab Sample ID:480-86236-X-8 MSDClient Matrix:WaterDilution:20Analysis Date:09/04/2015 2033Prep Date:09/04/2015 2033Leach Date:N/A		Pre	alysis Batch: p Batch: ach Batch:	480-262028 N/A N/A			HP5975T T8311.D 5 mL 5 mL 5 mL	
		%	Rec.					
Analyte		MS	MSD	Limit	RPD	RPD Limit	MS Qual	MSD Qual
o-Xylene		110	103	76 - 122	6	16		
Styrene		109	104	70 - 130	5	20		
Tetrachloroether	ne	104	100	74 - 122	3	20		
Toluene		109	107	80 - 122	2	15		
trans-1,2-Dichlor	oethene	112	102	73 - 127	9	20		
trans-1,3-Dichlor	opropene	94	96	72 - 123	2	15		
Trichloroethene		109	101	74 - 123	7	16		
Trichlorofluorom	ethane	112	105	62 - 152	6	20		
Vinyl chloride		100	91	65 - 133	10	15		
Surrogate			MS % Rec	MSD 9	% Rec	Acce	eptance Lim	its
1,2-Dichloroetha	ne-d4 (Surr)		86	85		6	6 - 137	
4-Bromofluorobe			103	99			3 - 120	
T-1			04	00		7	4 400	

94

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Analysis Batch: 480-262028

N/A

Prep Batch:

	8	THE LEADER IN ENVIRONMENTAL TESTING TestAmerica Laboratories, Inc. TAL-8210 (0713)	COC No:	Sampler	For Lab Use Only:	/alk-in Client:		ob / SDG No.:		Sample Specific Notes:					are will bulke								· for Months		Therm ID No.:	Bate/Time: B-26-15 OS'45	Date/Time:	Date/Time	13 #1
	039946		ate: <i>&</i> /z <i>S</i> / <i>1S</i>	Carrier:				tin of Custody							7	2	/					osessea il salliples ale le	Disposal by Lab	d	Corr'd:	Company: ALL	Company:	Company:	
	Chain of Custody Record		June Lewill	Lab Contact:	<u> </u>					1	M3 \	8	2		w3				/		Sounds Discover (A feet week for the second s	סמוווטופ טוצעטטמו (א ופפ ווומן עם מאפטאפט וו אמוווטופא מופ ובומוווכת וטווקפו ווומוו ז וווטוווון).	C Return to Client		Cooler Temp (); pas'd	Received by: All	Regiver the fill	Received in Laboratory by:	
•	Chain o		Su lawba	Analvsis Tumaround Time		[0	10		umple Vpe	Matrix Cont.	6- 6W 3 W	6- 76au 7 W	33	5 660 5 N	6 6W 3 W				/			odes for the sample in the	Unknown	DET EDI)		-C Plate/Time: 545 Received	Date/Time:	Daje/Time:	
			Project Manager:	Tei/Fax:		TAT if different		2 days	Sample Sample	Time	Stastis 110	0E1 5/5=/8	8/22/12/20	8/22/220	8/25/8-1230			/	 		33, 5=NaOH, 6= Other	Please List any EPA Waste Codes for the sample in the	Poison B	Cat A, NYSDI	Custody Seal No.:	Company: ルビルビ	Company: TH - HI IS	Company:	
-	Testhinerica Buffalo 18 hazelueod Brive	Amberst, NY 14228 Phone: 716,691.2600 Fax: 716.691.7991	Client Contact		NZIP: Alberry NY	8-402-4626	Project Name: Forming Sele Pazu Cleavel	olite: PO#		Sample Identification	MW-46C-82015	MW-476- 52017	MW-37C-82015	21028-285-WM	\$MW-28D-82015	36	of 33	7			Preservation Used: 1= (ce. 2= HCL 3= H2SO4, 4=HNO3, 5=NaOH, 6= Other Describbe Dereved Manufferetiese:	A Hazardous Waste? dispose of the sample.	Non-Hazard	Special instructions/QC Requirements & Comments:	Custody Seals Intact: 📉 Yes 🔲 No	Relinquished by:	Relinquishearby	Relipduished by:	

Login Number: 86254 List Number: 1 Creator: Williams, Christopher S

Job Number: 480-86254-1

List Source: TestAmerica Buffalo

luestion	Answer	Comment
adioactivity either was not measured or, if measured, is at or below ackground	True	
he cooler's custody seal, if present, is intact.	True	
he cooler or samples do not appear to have been compromised or ampered with.	True	
amples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
s the Field Sampler's name present on COC?	True	
here are no discrepancies between the sample IDs on the containers and ne COC.	False	No: Received Trip Blank(s) not listed on COC
amples are received within Holding Time.	True	
ample containers have legible labels.	True	
containers are not broken or leaking.	True	
ample collection date/times are provided.	True	
ppropriate sample containers are used.	True	
ample bottles are completely filled.	True	
ample Preservation Verified	True	
here is sufficient vol. for all requested analyses, incl. any requested IS/MSDs	True	
'OA sample vials do not have headspace or bubble is <6mm (1/4") in iameter.	True	
necessary, staff have been informed of any short hold time or quick TAT eeds	True	
Iultiphasic samples are not present.	True	
amples do not require splitting or compositing.	True	
ampling Company provided.	True	NYSDEC
amples received within 48 hours of sampling.	True	
amples requiring field filtration have been filtered in the field.	N/A	
Chlorine Residual checked.	N/A	