

Report of Analysis

Client:	PARSONS Engineering of New York, Inc.		Date Collected:	05/08/25	
Project:	Con Ed Non-MGP - East River M0SF 454453		Date Received:	05/09/25	
Client Sample ID:	MW-3-20250508MSD		SDG No.:	Q1993	
Lab Sample ID:	Q1993-03MSD		Matrix:	Water	
Analytical Method:	8260D		% Solid:	0	
Sample Wt/Vol:	5	Units: mL	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group1	
GC Column:	DB-624UI	ID : 0.18	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046120.D	1		05/09/25 17:10	VX050925

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
74-87-3	Chloromethane	53.7		0.32	1.00	ug/L
75-01-4	Vinyl Chloride	52.2		0.26	1.00	ug/L
74-83-9	Bromomethane	49.0		1.40	5.00	ug/L
75-00-3	Chloroethane	59.6		0.47	1.00	ug/L
75-35-4	1,1-Dichloroethene	52.0		0.23	1.00	ug/L
67-64-1	Acetone	250		1.50	5.00	ug/L
75-15-0	Carbon Disulfide	56.7		0.21	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	53.9		0.16	1.00	ug/L
75-09-2	Methylene Chloride	49.5		0.28	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	52.8		0.23	1.00	ug/L
75-34-3	1,1-Dichloroethane	53.7		0.23	1.00	ug/L
78-93-3	2-Butanone	270		0.98	5.00	ug/L
56-23-5	Carbon Tetrachloride	52.6		0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	53.3		0.19	1.00	ug/L
67-66-3	Chloroform	54.0		0.25	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	54.2		0.20	1.00	ug/L
71-43-2	Benzene	52.9		0.15	1.00	ug/L
107-06-2	1,2-Dichloroethane	53.0		0.22	1.00	ug/L
79-01-6	Trichloroethene	51.6		0.090	1.00	ug/L
78-87-5	1,2-Dichloropropane	55.0		0.20	1.00	ug/L
75-27-4	Bromodichloromethane	53.9		0.22	1.00	ug/L
108-10-1	4-Methyl-2-Pentanone	270		0.68	5.00	ug/L
108-88-3	Toluene	53.5		0.14	1.00	ug/L
10061-02-6	t-1,3-Dichloropropene	54.4		0.17	1.00	ug/L
10061-01-5	cis-1,3-Dichloropropene	54.1		0.16	1.00	ug/L
79-00-5	1,1,2-Trichloroethane	53.6		0.21	1.00	ug/L
591-78-6	2-Hexanone	280		0.89	5.00	ug/L
124-48-1	Dibromochloromethane	55.6		0.18	1.00	ug/L
127-18-4	Tetrachloroethene	47.6		0.23	1.00	ug/L
108-90-7	Chlorobenzene	50.7		0.12	1.00	ug/L

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100-41-4	Ethyl Benzene	52.1		0.13	1.00	ug/L
179601-23-1	m/p-Xylenes	100		0.24	2.00	ug/L
95-47-6	o-Xylene	52.7		0.12	1.00	ug/L
100-42-5	Styrene	53.8		0.15	1.00	ug/L
75-25-2	Bromoform	52.8		0.19	1.00	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	51.7		0.26	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	51.3		74 - 125	103%	SPK: 50
1868-53-7	Dibromofluoromethane	51.1		75 - 124	102%	SPK: 50
2037-26-5	Toluene-d8	50.7		86 - 113	101%	SPK: 50
460-00-4	4-Bromofluorobenzene	51.2		77 - 121	102%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	86600	5.549			
540-36-3	1,4-Difluorobenzene	154000	6.757			
3114-55-4	Chlorobenzene-d5	139000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	65100	12.018			

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products