

Report of Analysis

Client:	PARSONS Engineering of New York, Inc.		Date Collected:	06/05/25
Project:	Con Ed Non MGP – Atlantic Ave 453957.600024.05		Date Received:	06/06/25
Client Sample ID:	MW-6-20250605DL		SDG No.:	Q2268
Lab Sample ID:	Q2268-07DL		Matrix:	Water
Analytical Method:	8270E		% Solid:	0
Sample Wt/Vol:	960	Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:		uL	Test:	SVOCMS Group1
Extraction Type :		Decanted : N	Level :	LOW
Injection Volume :		GPC Factor : 1.0	GPC Cleanup :	N PH :
Prep Method :	SW3510C			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF142754.D	2	06/10/25 08:10	06/12/25 13:29	PB168376

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
100-52-7	Benzaldehyde	8.10	UD	8.10	20.8	ug/L
108-95-2	Phenol	6.50	JD	1.90	10.4	ug/L
111-44-4	bis(2-Chloroethyl)ether	1.70	UD	1.70	10.4	ug/L
95-57-8	2-Chlorophenol	1.20	UD	1.20	10.4	ug/L
95-48-7	2-Methylphenol	2.30	UD	2.30	10.4	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	2.70	UD	2.70	10.4	ug/L
98-86-2	Acetophenone	1.50	UD	1.50	10.4	ug/L
65794-96-9	3+4-Methylphenols	2.30	UD	2.30	20.8	ug/L
621-64-7	n-Nitroso-di-n-propylamine	2.90	UD	2.90	5.20	ug/L
67-72-1	Hexachloroethane	1.40	UD	1.40	10.4	ug/L
98-95-3	Nitrobenzene	1.60	UD	1.60	10.4	ug/L
78-59-1	Isophorone	1.60	UD	1.60	10.4	ug/L
88-75-5	2-Nitrophenol	3.70	UD	3.70	10.4	ug/L
105-67-9	2,4-Dimethylphenol	130	D	3.90	10.4	ug/L
111-91-1	bis(2-Chloroethoxy)methane	1.40	UD	1.40	10.4	ug/L
120-83-2	2,4-Dichlorophenol	1.10	UD	1.10	10.4	ug/L
91-20-3	Naphthalene	1.00	UD	1.00	10.4	ug/L
106-47-8	4-Chloroaniline	1.80	UD	1.80	10.4	ug/L
87-68-3	Hexachlorobutadiene	1.10	UD	1.10	10.4	ug/L
105-60-2	Caprolactam	2.40	UD	2.40	20.8	ug/L
59-50-7	4-Chloro-3-methylphenol	1.20	UD	1.20	10.4	ug/L
91-57-6	2-Methylnaphthalene	1.20	UD	1.20	10.4	ug/L
77-47-4	Hexachlorocyclopentadiene	7.60	UD	7.60	20.8	ug/L
88-06-2	2,4,6-Trichlorophenol	1.10	UD	1.10	10.4	ug/L
95-95-4	2,4,5-Trichlorophenol	1.30	UD	1.30	10.4	ug/L
92-52-4	1,1-Biphenyl	1.10	UD	1.10	10.4	ug/L
91-58-7	2-Chloronaphthalene	1.30	UD	1.30	10.4	ug/L
88-74-4	2-Nitroaniline	2.60	UD	2.60	10.4	ug/L
131-11-3	Dimethylphthalate	1.30	UD	1.30	10.4	ug/L

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208-96-8	Acenaphthylene	1.60	UD	1.60	10.4	ug/L
606-20-2	2,6-Dinitrotoluene	1.90	UD	1.90	10.4	ug/L
99-09-2	3-Nitroaniline	2.20	UD	2.20	10.4	ug/L
83-32-9	Acenaphthene	1.10	UD	1.10	10.4	ug/L
51-28-5	2,4-Dinitrophenol	12.4	UD	12.4	20.8	ug/L
100-02-7	4-Nitrophenol	5.00	UD	5.00	20.8	ug/L
132-64-9	Dibenzofuran	1.30	UD	1.30	10.4	ug/L
121-14-2	2,4-Dinitrotoluene	2.50	UD	2.50	10.4	ug/L
84-66-2	Diethylphthalate	1.40	UD	1.40	10.4	ug/L
7005-72-3	4-Chlorophenyl-phenylether	1.40	UD	1.40	10.4	ug/L
86-73-7	Fluorene	1.30	UD	1.30	10.4	ug/L
100-01-6	4-Nitroaniline	3.10	UD	3.10	10.4	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	6.00	UD	6.00	20.8	ug/L
86-30-6	n-Nitrosodiphenylamine	1.20	UD	1.20	10.4	ug/L
101-55-3	4-Bromophenyl-phenylether	0.83	UD	0.83	10.4	ug/L
118-74-1	Hexachlorobenzene	1.10	UD	1.10	10.4	ug/L
1912-24-9	Atrazine	2.10	UD	2.10	10.4	ug/L
87-86-5	Pentachlorophenol	3.30	UD	3.30	20.8	ug/L
85-01-8	Phenanthrene	1.00	UD	1.00	10.4	ug/L
120-12-7	Anthracene	1.30	UD	1.30	10.4	ug/L
86-74-8	Carbazole	1.50	UD	1.50	10.4	ug/L
84-74-2	Di-n-butylphthalate	2.50	UD	2.50	10.4	ug/L
206-44-0	Fluoranthene	1.70	UD	1.70	10.4	ug/L
129-00-0	Pyrene	1.00	UD	1.00	10.4	ug/L
85-68-7	Butylbenzylphthalate	4.00	UD	4.00	10.4	ug/L
91-94-1	3,3-Dichlorobenzidine	1.90	UD	1.90	20.8	ug/L
56-55-3	Benzo(a)anthracene	0.94	UD	0.94	10.4	ug/L
218-01-9	Chrysene	0.92	UD	0.92	10.4	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	3.30	UD	3.30	10.4	ug/L
117-84-0	Di-n-octyl phthalate	4.90	UD	4.90	20.8	ug/L
205-99-2	Benzo(b)fluoranthene	1.00	UD	1.00	10.4	ug/L

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207-08-9	Benzo(k)fluoranthene	1.00	UD	1.00	10.4	ug/L
50-32-8	Benzo(a)pyrene	1.10	UD	1.10	10.4	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	1.20	UD	1.20	10.4	ug/L
53-70-3	Dibenzo(a,h)anthracene	1.40	UD	1.40	10.4	ug/L
191-24-2	Benzo(g,h,i)perylene	1.40	UD	1.40	10.4	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	1.10	UD	1.10	10.4	ug/L
123-91-1	1,4-Dioxane	2.10	UD	2.10	10.4	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	1.50	UD	1.50	10.4	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	70.1		23 - 138	47%	SPK: 150
13127-88-3	Phenol-d6	51.0		10 - 134	34%	SPK: 150
4165-60-0	Nitrobenzene-d5	90.7		67 - 132	91%	SPK: 100
321-60-8	2-Fluorobiphenyl	94.7		52 - 132	95%	SPK: 100
118-79-6	2,4,6-Tribromophenol	126		44 - 137	84%	SPK: 150
1718-51-0	Terphenyl-d14	70.1		42 - 152	70%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	64700	6.892			
1146-65-2	Naphthalene-d8	236000	8.175			
15067-26-2	Acenaphthene-d10	121000	9.928			
1517-22-2	Phenanthrene-d10	173000	11.416			
1719-03-5	Chrysene-d12	110000	14.063			
1520-96-3	Perylene-d12	150000	15.557			

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