

Report of Analysis

Client:	PSEG	Date Collected:	04/25/25
Project:	GW South OP Pole	Date Received:	04/25/25
Client Sample ID:	WC-1	SDG No.:	Q1900
Lab Sample ID:	Q1900-01	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	84.4
Sample Wt/Vol:	50.05	Units:	g
Soil Aliquot Vol:			uL
Extraction Type :		Decanted :	N
Injection Volume :		Level :	LOW
		GPC Factor :	1.0
		GPC Cleanup :	N
		PH :	
Prep Method :	SW3541		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BM050069.D	1	04/29/25 09:30	05/01/25 18:02	PB167779

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.23	U	0.11	0.23	mg/Kg
108-95-2	Phenol	0.12	U	0.016	0.12	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.12	U	0.017	0.12	mg/Kg
95-57-8	2-Chlorophenol	0.12	U	0.017	0.12	mg/Kg
95-48-7	2-Methylphenol	0.12	U	0.021	0.12	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.12	U	0.027	0.12	mg/Kg
98-86-2	Acetophenone	0.12	U	0.021	0.12	mg/Kg
65794-96-9	3+4-Methylphenols	0.23	U	0.029	0.23	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.057	U	0.034	0.057	mg/Kg
67-72-1	Hexachloroethane	0.12	U	0.013	0.12	mg/Kg
98-95-3	Nitrobenzene	0.12	U	0.013	0.12	mg/Kg
78-59-1	Isophorone	0.12	U	0.023	0.12	mg/Kg
88-75-5	2-Nitrophenol	0.12	U	0.041	0.12	mg/Kg
105-67-9	2,4-Dimethylphenol	0.12	U	0.046	0.12	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.12	U	0.022	0.12	mg/Kg
120-83-2	2,4-Dichlorophenol	0.12	U	0.020	0.12	mg/Kg
91-20-3	Naphthalene	0.12	U	0.016	0.12	mg/Kg
106-47-8	4-Chloroaniline	0.12	U	0.025	0.12	mg/Kg
87-68-3	Hexachlorobutadiene	0.12	U	0.018	0.12	mg/Kg
105-60-2	Caprolactam	0.23	U	0.037	0.23	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.12	U	0.020	0.12	mg/Kg
91-57-6	2-Methylnaphthalene	0.12	U	0.018	0.12	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.23	U	0.082	0.23	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.12	U	0.014	0.12	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.12	U	0.021	0.12	mg/Kg
92-52-4	1,1-Biphenyl	0.12	U	0.016	0.12	mg/Kg
91-58-7	2-Chloronaphthalene	0.12	U	0.016	0.12	mg/Kg
88-74-4	2-Nitroaniline	0.12	U	0.034	0.12	mg/Kg
131-11-3	Dimethylphthalate	0.12	U	0.019	0.12	mg/Kg

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Injection Volume :		Level :	LOW
		GPC Factor :	1.0
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208-96-8	Acenaphthylene	0.12	U	0.021	0.12	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.12	U	0.024	0.12	mg/Kg
99-09-2	3-Nitroaniline	0.12	U	0.033	0.12	mg/Kg
83-32-9	Acenaphthene	0.12	U	0.015	0.12	mg/Kg
51-28-5	2,4-Dinitrophenol	0.23	U	0.16	0.23	mg/Kg
100-02-7	4-Nitrophenol	0.23	U	0.076	0.23	mg/Kg
132-64-9	Dibenzofuran	0.12	U	0.016	0.12	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.12	U	0.036	0.12	mg/Kg
84-66-2	Diethylphthalate	0.12	U	0.020	0.12	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.12	U	0.019	0.12	mg/Kg
86-73-7	Fluorene	0.12	U	0.018	0.12	mg/Kg
100-01-6	4-Nitroaniline	0.12	U	0.046	0.12	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.23	U	0.073	0.23	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.12	U	0.023	0.12	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.12	U	0.020	0.12	mg/Kg
118-74-1	Hexachlorobenzene	0.12	U	0.018	0.12	mg/Kg
1912-24-9	Atrazine	0.12	U	0.024	0.12	mg/Kg
87-86-5	Pentachlorophenol	0.23	U	0.036	0.23	mg/Kg
85-01-8	Phenanthrene	0.12	U	0.015	0.12	mg/Kg
120-12-7	Anthracene	0.12	U	0.024	0.12	mg/Kg
86-74-8	Carbazole	0.12	U	0.022	0.12	mg/Kg
84-74-2	Di-n-butylphthalate	0.12	U	0.034	0.12	mg/Kg
206-44-0	Fluoranthene	0.063	J	0.021	0.12	mg/Kg
129-00-0	Pyrene	0.11	J	0.026	0.12	mg/Kg
85-68-7	Butylbenzylphthalate	0.12	U	0.051	0.12	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.23	U	0.026	0.23	mg/Kg
56-55-3	Benzo(a)anthracene	0.14		0.016	0.12	mg/Kg
218-01-9	Chrysene	0.15		0.014	0.12	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.12	U	0.042	0.12	mg/Kg
117-84-0	Di-n-octyl phthalate	0.23	U	0.062	0.23	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.23		0.014	0.12	mg/Kg

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Analytical Method:	SW8270	% Solid:	84.4
Sample Wt/Vol:	50.05	Units:	g
Soil Aliquot Vol:		Final Vol:	1000 uL
Extraction Type :		Test:	SVOC-TCL BNA -20
	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	1.0	GPC Cleanup : N PH :
Prep Method :	SW3541		

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207-08-9	Benzo(k)fluoranthene	0.082	J	0.016	0.12	mg/Kg
50-32-8	Benzo(a)pyrene	0.14		0.021	0.12	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.080	J	0.021	0.12	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.12	U	0.020	0.12	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.095	J	0.018	0.12	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.12	U	0.018	0.12	mg/Kg
123-91-1	1,4-Dioxane	0.12	U	0.032	0.12	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.12	U	0.020	0.12	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	62.3		18 - 112	42%	SPK: 150
13127-88-3	Phenol-d6	67.0		15 - 107	45%	SPK: 150
4165-60-0	Nitrobenzene-d5	37.4		18 - 107	37%	SPK: 100
321-60-8	2-Fluorobiphenyl	35.7		20 - 109	36%	SPK: 100
118-79-6	2,4,6-Tribromophenol	66.6		10 - 116	44%	SPK: 150
1718-51-0	Terphenyl-d14	34.7		10 - 105	35%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	272000	7.745
1146-65-2	Naphthalene-d8	960000	10.539
15067-26-2	Acenaphthene-d10	637000	14.392
1517-22-2	Phenanthrene-d10	1320000	17.139
1719-03-5	Chrysene-d12	1390000	21.38
1520-96-3	Perylene-d12	1450000	24.368

TENTATIVE IDENTIFIED COMPOUNDS

000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	100	A	4.87	ug/Kg
000119-61-9	Benzophenone	130	J	15.8	ug/Kg
074630-43-6	2-Undecene, 6-methyl-, (Z)-	250	J	16.3	ug/Kg
000502-69-2	2-Pentadecanone, 6,10,14-trimethyl	130	J	17.2	ug/Kg
000057-10-3	n-Hexadecanoic acid	320	J	18.0	ug/Kg
	unknown19.045	180	J	19.0	ug/Kg
1000144-64-7	4,4,6-Trimethyl-cyclohex-2-en-1-ol	320	J	19.3	ug/Kg

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007206-21-5	5-Octadecene, (E)-	160	J		20.1	ug/Kg
959085-66-6	Heptadecyl heptafluorobutyrate	670	J		21.1	ug/Kg
074685-30-6	5-Eicosene, (E)-	1600	J		22.2	ug/Kg
000557-59-5	Tetracosanoic acid	77.9	J		22.6	ug/Kg
000638-66-4	Octadecanal	140	J		23.1	ug/Kg
000297-03-0	Cyclotetracosane	720	J		23.6	ug/Kg
000192-97-2	Benzo[e]pyrene	130	J		24.1	ug/Kg
007390-81-0	Oxirane, hexadecyl-	140	J		24.9	ug/Kg
007616-22-0	.gamma.-Tocopherol	200	J		25.2	ug/Kg
000593-49-7	Heptacosane	490	J		25.4	ug/Kg
010192-32-2	1-Tetracosene	350	J		25.5	ug/Kg
007796-19-2	2-Heptacosanone	200	J		25.7	ug/Kg
053584-61-5	17.alpha.(H),21.beta.(H)-Homohopan	1400	J		28.6	ug/Kg

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