

Report of Analysis

Client:	PSEG	Date Collected:	04/04/25
Project:	PSEG Clay Street	Date Received:	04/04/25
Client Sample ID:	TT-8	SDG No.:	Q1732
Lab Sample ID:	Q1732-01	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	89.7
Sample Wt/Vol:	50.03	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		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BM049846.D	10	04/07/25 08:55	04/07/25 19:18	PB167474

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	2.20	U	1.00	2.20	mg/Kg
108-95-2	Phenol	1.10	U	0.15	1.10	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	1.10	U	0.16	1.10	mg/Kg
95-57-8	2-Chlorophenol	1.10	U	0.16	1.10	mg/Kg
95-48-7	2-Methylphenol	1.10	U	0.20	1.10	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	1.10	U	0.25	1.10	mg/Kg
98-86-2	Acetophenone	1.10	U	0.20	1.10	mg/Kg
65794-96-9	3+4-Methylphenols	2.20	U	0.27	2.20	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.53	U	0.32	0.53	mg/Kg
67-72-1	Hexachloroethane	1.10	U	0.12	1.10	mg/Kg
98-95-3	Nitrobenzene	1.10	U	0.12	1.10	mg/Kg
78-59-1	Isophorone	1.10	U	0.22	1.10	mg/Kg
88-75-5	2-Nitrophenol	1.10	U	0.39	1.10	mg/Kg
105-67-9	2,4-Dimethylphenol	1.10	U	0.43	1.10	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	1.10	U	0.21	1.10	mg/Kg
120-83-2	2,4-Dichlorophenol	1.10	U	0.19	1.10	mg/Kg
91-20-3	Naphthalene	1.10	U	0.15	1.10	mg/Kg
106-47-8	4-Chloroaniline	1.10	U	0.24	1.10	mg/Kg
87-68-3	Hexachlorobutadiene	1.10	U	0.17	1.10	mg/Kg
105-60-2	Caprolactam	2.20	U	0.35	2.20	mg/Kg
59-50-7	4-Chloro-3-methylphenol	1.10	U	0.19	1.10	mg/Kg
91-57-6	2-Methylnaphthalene	1.10	U	0.17	1.10	mg/Kg
77-47-4	Hexachlorocyclopentadiene	2.20	U	0.78	2.20	mg/Kg
88-06-2	2,4,6-Trichlorophenol	1.10	U	0.13	1.10	mg/Kg
95-95-4	2,4,5-Trichlorophenol	1.10	U	0.19	1.10	mg/Kg
92-52-4	1,1-Biphenyl	1.10	U	0.15	1.10	mg/Kg
91-58-7	2-Chloronaphthalene	1.10	U	0.15	1.10	mg/Kg
88-74-4	2-Nitroaniline	1.10	U	0.32	1.10	mg/Kg
131-11-3	Dimethylphthalate	1.10	U	0.18	1.10	mg/Kg

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208-96-8	Acenaphthylene	1.10	U	0.19	1.10	mg/Kg
606-20-2	2,6-Dinitrotoluene	1.10	U	0.22	1.10	mg/Kg
99-09-2	3-Nitroaniline	1.10	U	0.31	1.10	mg/Kg
83-32-9	Acenaphthene	1.10	U	0.14	1.10	mg/Kg
51-28-5	2,4-Dinitrophenol	2.20	U	1.50	2.20	mg/Kg
100-02-7	4-Nitrophenol	2.20	U	0.72	2.20	mg/Kg
132-64-9	Dibenzofuran	1.10	U	0.15	1.10	mg/Kg
121-14-2	2,4-Dinitrotoluene	1.10	U	0.33	1.10	mg/Kg
84-66-2	Diethylphthalate	1.10	U	0.19	1.10	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	1.10	U	0.18	1.10	mg/Kg
86-73-7	Fluorene	1.10	U	0.17	1.10	mg/Kg
100-01-6	4-Nitroaniline	1.10	U	0.43	1.10	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	2.20	U	0.69	2.20	mg/Kg
86-30-6	n-Nitrosodiphenylamine	1.10	U	0.22	1.10	mg/Kg
101-55-3	4-Bromophenyl-phenylether	1.10	U	0.19	1.10	mg/Kg
118-74-1	Hexachlorobenzene	1.10	U	0.17	1.10	mg/Kg
1912-24-9	Atrazine	1.10	U	0.23	1.10	mg/Kg
87-86-5	Pentachlorophenol	2.20	U	0.34	2.20	mg/Kg
85-01-8	Phenanthrene	0.56	J	0.14	1.10	mg/Kg
120-12-7	Anthracene	1.10	U	0.22	1.10	mg/Kg
86-74-8	Carbazole	1.10	U	0.21	1.10	mg/Kg
84-74-2	Di-n-butylphthalate	1.10	U	0.32	1.10	mg/Kg
206-44-0	Fluoranthene	0.88	J	0.20	1.10	mg/Kg
129-00-0	Pyrene	0.72	J	0.24	1.10	mg/Kg
85-68-7	Butylbenzylphthalate	1.10	U	0.48	1.10	mg/Kg
91-94-1	3,3-Dichlorobenzidine	2.20	U	0.25	2.20	mg/Kg
56-55-3	Benzo(a)anthracene	1.10	U	0.15	1.10	mg/Kg
218-01-9	Chrysene	1.10	U	0.13	1.10	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	1.10	U	0.40	1.10	mg/Kg
117-84-0	Di-n-octyl phthalate	2.20	U	0.58	2.20	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.48	J	0.13	1.10	mg/Kg

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207-08-9	Benzo(k)fluoranthene	1.10	U	0.15	1.10	mg/Kg
50-32-8	Benzo(a)pyrene	1.10	U	0.20	1.10	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	1.10	U	0.19	1.10	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	1.10	U	0.18	1.10	mg/Kg
191-24-2	Benzo(g,h,i)perylene	1.10	U	0.17	1.10	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	1.10	U	0.17	1.10	mg/Kg
123-91-1	1,4-Dioxane	1.10	U	0.30	1.10	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	1.10	U	0.18	1.10	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	71.7		18 - 112	48%	SPK: 150
13127-88-3	Phenol-d6	69.5		15 - 107	46%	SPK: 150
4165-60-0	Nitrobenzene-d5	42.5		18 - 107	43%	SPK: 100
321-60-8	2-Fluorobiphenyl	46.7		20 - 109	47%	SPK: 100
118-79-6	2,4,6-Tribromophenol	96.6		10 - 116	64%	SPK: 150
1718-51-0	Terphenyl-d14	42.4		10 - 105	42%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	407000	7.787	
1146-65-2	Naphthalene-d8	1420000	10.581	
15067-26-2	Acenaphthene-d10	930000	14.427	
1517-22-2	Phenanthrene-d10	1820000	17.168	
1719-03-5	Chrysene-d12	1850000	21.409	
1520-96-3	Perylene-d12	2140000	24.421	

TENTATIVE IDENTIFIED COMPOUNDS

	unknown3.504	2500	J		3.50	ug/Kg
000057-10-3	n-Hexadecanoic acid	5000	J		18.1	ug/Kg
000057-11-4	Octadecanoic acid	12600	J		19.4	ug/Kg

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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