

Report of Analysis

Client:	PSEG	Date Collected:	04/14/25
Project:	Oradell Gas and Appliance Service MA00006789	Date Received:	04/14/25
Client Sample ID:	OR-03-041425	SDG No.:	Q1806
Lab Sample ID:	Q1806-01	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	89.8
Sample Wt/Vol:	50.05 Units: g	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	SVOC-TCL BNA -20
Extraction Type :	Decanted : N	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
BM049939.D	2	04/15/25 09:35	04/15/25 17:52	PB167598

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.21	U	0.21	0.44	mg/Kg
108-95-2	Phenol	0.030	U	0.030	0.23	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.032	U	0.032	0.23	mg/Kg
95-57-8	2-Chlorophenol	0.033	U	0.033	0.23	mg/Kg
95-48-7	2-Methylphenol	0.040	U	0.040	0.23	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.050	U	0.050	0.23	mg/Kg
98-86-2	Acetophenone	0.039	U	0.039	0.23	mg/Kg
65794-96-9	3+4-Methylphenols	0.055	U	0.055	0.44	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.063	U	0.063	0.11	mg/Kg
67-72-1	Hexachloroethane	0.024	U	0.024	0.23	mg/Kg
98-95-3	Nitrobenzene	0.024	U	0.024	0.23	mg/Kg
78-59-1	Isophorone	0.044	U	0.044	0.23	mg/Kg
88-75-5	2-Nitrophenol	0.078	U	0.078	0.23	mg/Kg
105-67-9	2,4-Dimethylphenol	0.087	U	0.087	0.23	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.041	U	0.041	0.23	mg/Kg
120-83-2	2,4-Dichlorophenol	0.038	U	0.038	0.23	mg/Kg
91-20-3	Naphthalene	0.030	U	0.030	0.23	mg/Kg
106-47-8	4-Chloroaniline	0.047	U	0.047	0.23	mg/Kg
87-68-3	Hexachlorobutadiene	0.034	U	0.034	0.23	mg/Kg
105-60-2	Caprolactam	0.070	U	0.070	0.44	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.038	U	0.038	0.23	mg/Kg
91-57-6	2-Methylnaphthalene	0.034	U	0.034	0.23	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.15	U	0.15	0.44	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.026	U	0.026	0.23	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.039	U	0.039	0.23	mg/Kg
92-52-4	1,1-Biphenyl	0.029	U	0.029	0.23	mg/Kg
91-58-7	2-Chloronaphthalene	0.030	U	0.030	0.23	mg/Kg
88-74-4	2-Nitroaniline	0.064	U	0.064	0.23	mg/Kg
131-11-3	Dimethylphthalate	0.036	U	0.036	0.23	mg/Kg

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208-96-8	Acenaphthylene	0.13	J	0.039	0.23	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.045	U	0.045	0.23	mg/Kg
99-09-2	3-Nitroaniline	0.061	U	0.061	0.23	mg/Kg
83-32-9	Acenaphthene	0.29		0.028	0.23	mg/Kg
51-28-5	2,4-Dinitrophenol	0.31	U	0.31	0.44	mg/Kg
100-02-7	4-Nitrophenol	0.14	U	0.14	0.44	mg/Kg
132-64-9	Dibenzofuran	0.14	J	0.030	0.23	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.067	U	0.067	0.23	mg/Kg
84-66-2	Diethylphthalate	0.038	U	0.038	0.23	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.036	U	0.036	0.23	mg/Kg
86-73-7	Fluorene	0.32		0.034	0.23	mg/Kg
100-01-6	4-Nitroaniline	0.086	U	0.086	0.23	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.14	U	0.14	0.44	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.044	U	0.044	0.23	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.037	U	0.037	0.23	mg/Kg
118-74-1	Hexachlorobenzene	0.034	U	0.034	0.23	mg/Kg
1912-24-9	Atrazine	0.045	U	0.045	0.23	mg/Kg
87-86-5	Pentachlorophenol	0.069	U	0.069	0.44	mg/Kg
85-01-8	Phenanthrene	2.80		0.028	0.23	mg/Kg
120-12-7	Anthracene	0.81		0.045	0.23	mg/Kg
86-74-8	Carbazole	0.30		0.042	0.23	mg/Kg
84-74-2	Di-n-butylphthalate	0.064	U	0.064	0.23	mg/Kg
206-44-0	Fluoranthene	3.90	E	0.040	0.23	mg/Kg
129-00-0	Pyrene	3.10		0.048	0.23	mg/Kg
85-68-7	Butylbenzylphthalate	0.095	U	0.095	0.23	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.049	U	0.049	0.44	mg/Kg
56-55-3	Benzo(a)anthracene	1.40		0.031	0.23	mg/Kg
218-01-9	Chrysene	1.30		0.027	0.23	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.079	U	0.079	0.23	mg/Kg
117-84-0	Di-n-octyl phthalate	0.12	U	0.12	0.44	mg/Kg
205-99-2	Benzo(b)fluoranthene	1.40		0.025	0.23	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.53		0.030	0.23	mg/Kg
50-32-8	Benzo(a)pyrene	1.20		0.039	0.23	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.60		0.039	0.23	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.18	J	0.037	0.23	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.70		0.034	0.23	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.034	U	0.034	0.23	mg/Kg
123-91-1	1,4-Dioxane	0.060	U	0.060	0.23	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.037	U	0.037	0.23	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	77.7		18 - 112	52%	SPK: 150
13127-88-3	Phenol-d6	75.9		15 - 107	51%	SPK: 150
4165-60-0	Nitrobenzene-d5	56.8		18 - 107	57%	SPK: 100
321-60-8	2-Fluorobiphenyl	51.6		20 - 109	52%	SPK: 100
118-79-6	2,4,6-Tribromophenol	84.5		10 - 116	56%	SPK: 150
1718-51-0	Terphenyl-d14	43.3		10 - 105	43%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	414000	7.769
1146-65-2	Naphthalene-d8	1320000	10.563
15067-26-2	Acenaphthene-d10	779000	14.416
1517-22-2	Phenanthrene-d10	1440000	17.157
1719-03-5	Chrysene-d12	1470000	21.398
1520-96-3	Perylene-d12	1640000	24.403

TENTATIVE IDENTIFIED COMPOUNDS

000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	420	A	4.89	ug/Kg
000107-70-0	2-Pentanone, 4-methoxy-4-methyl-	140	J	5.97	ug/Kg
000119-61-9	Benzophenone	230	J	15.8	ug/Kg
000132-65-0	Dibenzothiophene	180	J	17.0	ug/Kg
067388-11-8	4-Methylnaphtho[1,2-b]thiophene	100	J	17.7	ug/Kg
002531-84-2	Phenanthrene, 2-methyl-	310	J	18.0	ug/Kg
000832-69-9	Phenanthrene, 1-methyl-	970	J	18.1	ug/Kg

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000613-12-7	Anthracene, 2-methyl-	200	J		18.2	ug/Kg
000203-64-5	4H-Cyclopenta[def]phenanthrene	810	J		18.2	ug/Kg
000779-02-2	Anthracene, 9-methyl-	180	J		18.3	ug/Kg
000612-94-2	Naphthalene, 2-phenyl-	220	J		18.5	ug/Kg
000084-65-1	9,10-Anthracenedione	220	J		18.6	ug/Kg
003674-66-6	Phenanthrene, 2,5-dimethyl-	250	J		19.0	ug/Kg
005737-13-3	Cyclopenta(def)phenanthrenone	250	J		19.1	ug/Kg
000057-11-4	Octadecanoic acid	560	J		19.3	ug/Kg
002381-21-7	Pyrene, 1-methyl-	110	J		19.9	ug/Kg
000238-84-6	11H-Benzo[a]fluorene	130	J		20.1	ug/Kg
003442-78-2	Pyrene, 2-methyl-	100	J		20.2	ug/Kg
027208-37-3	Cyclopenta[cd]pyrene	93.4	J		21.1	ug/Kg
000192-97-2	Benzo[e]pyrene	1100	J		24.1	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