

Report of Analysis

Client:	PSEG	Date Collected:	05/15/25
Project:	Oradell Gas and Appliance Service MA00006789	Date Received:	05/15/25
Client Sample ID:	OR-02-051525	SDG No.:	Q2055
Lab Sample ID:	Q2055-01	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	89.1
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
BF142436.D	1	05/16/25 10:00	05/16/25 19:40	PB168041

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.10	U	0.10	0.22	mg/Kg
108-95-2	Phenol	0.015	U	0.015	0.11	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.016	U	0.016	0.11	mg/Kg
95-57-8	2-Chlorophenol	0.016	U	0.016	0.11	mg/Kg
95-48-7	2-Methylphenol	0.020	U	0.020	0.11	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.025	U	0.025	0.11	mg/Kg
98-86-2	Acetophenone	0.020	U	0.020	0.11	mg/Kg
65794-96-9	3+4-Methylphenols	0.028	U	0.028	0.22	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.032	U	0.032	0.054	mg/Kg
67-72-1	Hexachloroethane	0.012	U	0.012	0.11	mg/Kg
98-95-3	Nitrobenzene	0.012	U	0.012	0.11	mg/Kg
78-59-1	Isophorone	0.022	U	0.022	0.11	mg/Kg
88-75-5	2-Nitrophenol	0.039	U	0.039	0.11	mg/Kg
105-67-9	2,4-Dimethylphenol	0.044	U	0.044	0.11	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.021	U	0.021	0.11	mg/Kg
120-83-2	2,4-Dichlorophenol	0.019	U	0.019	0.11	mg/Kg
91-20-3	Naphthalene	0.015	U	0.015	0.11	mg/Kg
106-47-8	4-Chloroaniline	0.024	U	0.024	0.11	mg/Kg
87-68-3	Hexachlorobutadiene	0.017	U	0.017	0.11	mg/Kg
105-60-2	Caprolactam	0.035	U	0.035	0.22	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.019	U	0.019	0.11	mg/Kg
91-57-6	2-Methylnaphthalene	0.017	U	0.017	0.11	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.078	U	0.078	0.22	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.013	U	0.013	0.11	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.020	U	0.020	0.11	mg/Kg
92-52-4	1,1-Biphenyl	0.015	U	0.015	0.11	mg/Kg
91-58-7	2-Chloronaphthalene	0.015	U	0.015	0.11	mg/Kg
88-74-4	2-Nitroaniline	0.032	U	0.032	0.11	mg/Kg
131-11-3	Dimethylphthalate	0.018	U	0.018	0.11	mg/Kg

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208-96-8	Acenaphthylene	0.019	U	0.019	0.11	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.023	U	0.023	0.11	mg/Kg
99-09-2	3-Nitroaniline	0.031	U	0.031	0.11	mg/Kg
83-32-9	Acenaphthene	0.046	J	0.014	0.11	mg/Kg
51-28-5	2,4-Dinitrophenol	0.15	U	0.15	0.22	mg/Kg
100-02-7	4-Nitrophenol	0.072	U	0.072	0.22	mg/Kg
132-64-9	Dibenzofuran	0.015	U	0.015	0.11	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.034	U	0.034	0.11	mg/Kg
84-66-2	Diethylphthalate	0.019	U	0.019	0.11	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.018	U	0.018	0.11	mg/Kg
86-73-7	Fluorene	0.074	J	0.017	0.11	mg/Kg
100-01-6	4-Nitroaniline	0.043	U	0.043	0.11	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.069	U	0.069	0.22	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.022	U	0.022	0.11	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.019	U	0.019	0.11	mg/Kg
118-74-1	Hexachlorobenzene	0.017	U	0.017	0.11	mg/Kg
1912-24-9	Atrazine	0.023	U	0.023	0.11	mg/Kg
87-86-5	Pentachlorophenol	0.035	U	0.035	0.22	mg/Kg
85-01-8	Phenanthrene	0.59		0.014	0.11	mg/Kg
120-12-7	Anthracene	0.14		0.022	0.11	mg/Kg
86-74-8	Carbazole	0.11		0.021	0.11	mg/Kg
84-74-2	Di-n-butylphthalate	0.032	U	0.032	0.11	mg/Kg
206-44-0	Fluoranthene	1.00		0.020	0.11	mg/Kg
129-00-0	Pyrene	0.69		0.024	0.11	mg/Kg
85-68-7	Butylbenzylphthalate	0.048	U	0.048	0.11	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.025	U	0.025	0.22	mg/Kg
56-55-3	Benzo(a)anthracene	0.41		0.016	0.11	mg/Kg
218-01-9	Chrysene	0.38		0.013	0.11	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.040	U	0.040	0.11	mg/Kg
117-84-0	Di-n-octyl phthalate	0.058	U	0.058	0.22	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.52		0.013	0.11	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.18		0.015	0.11	mg/Kg
50-32-8	Benzo(a)pyrene	0.39		0.020	0.11	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.17		0.020	0.11	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.051	J	0.018	0.11	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.21		0.017	0.11	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.017	U	0.017	0.11	mg/Kg
123-91-1	1,4-Dioxane	0.030	U	0.030	0.11	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.018	U	0.018	0.11	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	80.7		18 - 112	54%	SPK: 150
13127-88-3	Phenol-d6	76.4		15 - 107	51%	SPK: 150
4165-60-0	Nitrobenzene-d5	57.0		18 - 107	57%	SPK: 100
321-60-8	2-Fluorobiphenyl	55.6		20 - 109	56%	SPK: 100
118-79-6	2,4,6-Tribromophenol	79.9		10 - 116	53%	SPK: 150
1718-51-0	Terphenyl-d14	42.6		10 - 105	43%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	142000	6.904
1146-65-2	Naphthalene-d8	493000	8.181
15067-26-2	Acenaphthene-d10	207000	9.939
1517-22-2	Phenanthrene-d10	317000	11.422
1719-03-5	Chrysene-d12	238000	14.063
1520-96-3	Perylene-d12	192000	15.557

TENTATIVE IDENTIFIED COMPOUNDS

000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	150	A	5.12	ug/Kg
006180-61-6	1-Propanol, 3-phenoxy-	74.4	J	8.49	ug/Kg
000629-59-4	Tetradecane	64.8	J	9.33	ug/Kg
000544-76-3	Hexadecane	87.7	J	9.86	ug/Kg
	unknown10.028	230	J	10.0	ug/Kg
006165-40-8	Pentadecane, 7-methyl-	110	J	10.4	ug/Kg
000629-78-7	Heptadecane	160	J	10.8	ug/Kg

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013287-21-3	Tridecane, 6-methyl-	100	J		11.3	ug/Kg
000233-02-3	Naphtho[2,1-b]thiophene	73.6	J		11.3	ug/Kg
000629-50-5	Tridecane	92.8	J		11.7	ug/Kg
000112-39-0	Hexadecanoic acid, methyl ester	530	J		11.8	ug/Kg
000057-10-3	n-Hexadecanoic acid	380	J		11.9	ug/Kg
000203-64-5	4H-Cyclopenta[def]phenanthrene	200	J		12.0	ug/Kg
093327-56-1	6-Phenylbenzocyclohepten-7-one	87.0	J		12.2	ug/Kg
056554-46-2	12-Octadecenoic acid, methyl ester	2400	J		12.5	ug/Kg
000057-11-4	Octadecanoic acid	83.9	J		12.7	ug/Kg
002381-21-7	Pyrene, 1-methyl-	87.0	J		13.2	ug/Kg
000243-17-4	11H-Benzo[b]fluorene	82.1	J		13.3	ug/Kg
000192-97-2	Benzo[e]pyrene	79.2	J		15.2	ug/Kg
000198-55-0	Perylene	290	J		15.4	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