

Report of Analysis

Client:	PSEG	Date Collected:	08/27/25
Project:	Trenton Gas and Appliance Service MA00006789	Date Received:	08/27/25
Client Sample ID:	TR-05-08272025	SDG No.:	Q2969
Lab Sample ID:	Q2969-01	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	95.5
Sample Wt/Vol:	50.03 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
BP025617.D	1	08/28/25 09:30	08/28/25 18:42	PB169431

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.098	U	0.098	0.21	mg/Kg
108-95-2	Phenol	0.014	U	0.014	0.11	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.015	U	0.015	0.11	mg/Kg
95-57-8	2-Chlorophenol	0.015	U	0.015	0.11	mg/Kg
95-48-7	2-Methylphenol	0.019	U	0.019	0.11	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.024	U	0.024	0.11	mg/Kg
98-86-2	Acetophenone	0.019	U	0.019	0.11	mg/Kg
65794-96-9	3+4-Methylphenols	0.026	U	0.026	0.21	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.030	U	0.030	0.050	mg/Kg
67-72-1	Hexachloroethane	0.011	U	0.011	0.11	mg/Kg
98-95-3	Nitrobenzene	0.012	U	0.012	0.11	mg/Kg
78-59-1	Isophorone	0.021	U	0.021	0.11	mg/Kg
88-75-5	2-Nitrophenol	0.037	U	0.037	0.11	mg/Kg
105-67-9	2,4-Dimethylphenol	0.041	U	0.041	0.11	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.019	U	0.019	0.11	mg/Kg
120-83-2	2,4-Dichlorophenol	0.018	U	0.018	0.11	mg/Kg
91-20-3	Naphthalene	0.014	U	0.014	0.11	mg/Kg
106-47-8	4-Chloroaniline	0.022	U	0.022	0.11	mg/Kg
87-68-3	Hexachlorobutadiene	0.016	U	0.016	0.11	mg/Kg
105-60-2	Caprolactam	0.033	U	0.033	0.21	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.018	U	0.018	0.11	mg/Kg
91-57-6	2-Methylnaphthalene	0.016	U	0.016	0.11	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.073	U	0.073	0.21	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.012	U	0.012	0.11	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.018	U	0.018	0.11	mg/Kg
92-52-4	1,1-Biphenyl	0.014	U	0.014	0.11	mg/Kg
91-58-7	2-Chloronaphthalene	0.014	U	0.014	0.11	mg/Kg
88-74-4	2-Nitroaniline	0.030	U	0.030	0.11	mg/Kg
131-11-3	Dimethylphthalate	0.017	U	0.017	0.11	mg/Kg

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208-96-8	Acenaphthylene	0.018	U	0.018	0.11	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.021	U	0.021	0.11	mg/Kg
99-09-2	3-Nitroaniline	0.029	U	0.029	0.11	mg/Kg
83-32-9	Acenaphthene	0.013	U	0.013	0.11	mg/Kg
51-28-5	2,4-Dinitrophenol	0.14	U	0.14	0.21	mg/Kg
100-02-7	4-Nitrophenol	0.067	U	0.067	0.21	mg/Kg
132-64-9	Dibenzofuran	0.014	U	0.014	0.11	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.032	U	0.032	0.11	mg/Kg
84-66-2	Diethylphthalate	0.018	U	0.018	0.11	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.017	U	0.017	0.11	mg/Kg
86-73-7	Fluorene	0.016	U	0.016	0.11	mg/Kg
100-01-6	4-Nitroaniline	0.040	U	0.040	0.11	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.065	U	0.065	0.21	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.021	U	0.021	0.11	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.018	U	0.018	0.11	mg/Kg
118-74-1	Hexachlorobenzene	0.016	U	0.016	0.11	mg/Kg
1912-24-9	Atrazine	0.021	U	0.021	0.11	mg/Kg
87-86-5	Pentachlorophenol	0.032	U	0.032	0.21	mg/Kg
85-01-8	Phenanthrene	0.013	U	0.013	0.11	mg/Kg
120-12-7	Anthracene	0.021	U	0.021	0.11	mg/Kg
86-74-8	Carbazole	0.020	U	0.020	0.11	mg/Kg
84-74-2	Di-n-butylphthalate	0.030	U	0.030	0.11	mg/Kg
206-44-0	Fluoranthene	0.019	U	0.019	0.11	mg/Kg
129-00-0	Pyrene	0.023	U	0.023	0.11	mg/Kg
85-68-7	Butylbenzylphthalate	0.045	U	0.045	0.11	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.023	U	0.023	0.21	mg/Kg
56-55-3	Benzo(a)anthracene	0.014	U	0.014	0.11	mg/Kg
218-01-9	Chrysene	0.013	U	0.013	0.11	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.037	U	0.037	0.11	mg/Kg
117-84-0	Di-n-octyl phthalate	0.055	U	0.055	0.21	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.012	U	0.012	0.11	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.014	U	0.014	0.11	mg/Kg
50-32-8	Benzo(a)pyrene	0.019	U	0.019	0.11	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.018	U	0.018	0.11	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.017	U	0.017	0.11	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.016	U	0.016	0.11	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.016	U	0.016	0.11	mg/Kg
123-91-1	1,4-Dioxane	0.028	U	0.028	0.11	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.017	U	0.017	0.11	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	64.4		18 - 112	43%	SPK: 150
13127-88-3	Phenol-d6	62.9		15 - 107	42%	SPK: 150
4165-60-0	Nitrobenzene-d5	40.0		18 - 107	40%	SPK: 100
321-60-8	2-Fluorobiphenyl	41.5		20 - 109	41%	SPK: 100
118-79-6	2,4,6-Tribromophenol	62.9		10 - 116	42%	SPK: 150
1718-51-0	Terphenyl-d14	40.6		10 - 105	41%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	215000	7.772	
1146-65-2	Naphthalene-d8	859000	10.543	
15067-26-2	Acenaphthene-d10	517000	14.39	
1517-22-2	Phenanthrene-d10	983000	17.183	
1719-03-5	Chrysene-d12	1030000	21.63	
1520-96-3	Perylene-d12	1190000	25.001	

TENTATIVE IDENTIFIED COMPOUNDS

000994-05-8	Butane, 2-methoxy-2-methyl-	300	J		3.03	ug/Kg
000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	110	A		4.94	ug/Kg
000544-76-3	Hexadecane	95.9	J		14.3	ug/Kg
000629-59-4	Tetradecane	130	J		15.2	ug/Kg
003892-00-0	Pentadecane, 2,6,10-trimethyl-	62.6	J		15.7	ug/Kg
000119-61-9	Benzophenone	120	J		15.8	ug/Kg
000629-78-7	Heptadecane	160	J		16.1	ug/Kg

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001921-70-6	Pentadecane, 2,6,10,14-tetramethyl	85.6	J		16.2	ug/Kg
055045-08-4	Dodecane, 2-methyl-6-propyl-	130	J		16.9	ug/Kg
000629-50-5	Tridecane	82.9	J		17.0	ug/Kg
000593-49-7	Heptacosane	120	J		17.7	ug/Kg
000057-10-3	n-Hexadecanoic acid	380	J		18.1	ug/Kg
000593-45-3	Octadecane	120	J		18.4	ug/Kg
000629-94-7	Heneicosane	110	J		19.1	ug/Kg
000057-11-4	Octadecanoic acid	200	J		19.4	ug/Kg
035599-77-0	Tridecane, 1-iodo-	95.6	J		19.7	ug/Kg
000629-97-0	Docosane	70.5	J		20.3	ug/Kg
007225-64-1	Heptadecane, 9-octyl-	60.5	J		20.8	ug/Kg
003452-07-1	1-Eicosene	150	J		21.3	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