

Report of Analysis

Client:	PSEG	Date Collected:	05/02/25
Project:	Moorestown	Date Received:	05/02/25
Client Sample ID:	CLEAN-FILL	SDG No.:	Q1952
Lab Sample ID:	Q1952-01	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	74.4
Sample Wt/Vol:	50.06	Units:	g
Soil Aliquot Vol:			uL
Extraction Type :		Decanted :	N
Injection Volume :		GPC Factor :	1.0
Prep Method :	SW3541	Level :	LOW
		Test:	SVOC-TCL BNA -20
		GPC Cleanup :	N
		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BM050121.D	1	05/05/25 09:35	05/06/25 15:59	PB167857

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.13	U	0.13	0.27	mg/Kg
108-95-2	Phenol	0.018	U	0.018	0.14	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.020	U	0.020	0.14	mg/Kg
95-57-8	2-Chlorophenol	0.020	U	0.020	0.14	mg/Kg
95-48-7	2-Methylphenol	0.024	U	0.024	0.14	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.030	U	0.030	0.14	mg/Kg
98-86-2	Acetophenone	0.024	U	0.024	0.14	mg/Kg
65794-96-9	3+4-Methylphenols	0.033	U	0.033	0.27	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.038	U	0.038	0.064	mg/Kg
67-72-1	Hexachloroethane	0.014	U	0.014	0.14	mg/Kg
98-95-3	Nitrobenzene	0.015	U	0.015	0.14	mg/Kg
78-59-1	Isophorone	0.026	U	0.026	0.14	mg/Kg
88-75-5	2-Nitrophenol	0.047	U	0.047	0.14	mg/Kg
105-67-9	2,4-Dimethylphenol	0.052	U	0.052	0.14	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.025	U	0.025	0.14	mg/Kg
120-83-2	2,4-Dichlorophenol	0.023	U	0.023	0.14	mg/Kg
91-20-3	Naphthalene	0.15		0.018	0.14	mg/Kg
106-47-8	4-Chloroaniline	0.029	U	0.029	0.14	mg/Kg
87-68-3	Hexachlorobutadiene	0.020	U	0.020	0.14	mg/Kg
105-60-2	Caprolactam	0.042	U	0.042	0.27	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.023	U	0.023	0.14	mg/Kg
91-57-6	2-Methylnaphthalene	0.021	U	0.021	0.14	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.093	U	0.093	0.27	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.016	U	0.016	0.14	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.023	U	0.023	0.14	mg/Kg
92-52-4	1,1-Biphenyl	0.018	U	0.018	0.14	mg/Kg
91-58-7	2-Chloronaphthalene	0.018	U	0.018	0.14	mg/Kg
88-74-4	2-Nitroaniline	0.039	U	0.039	0.14	mg/Kg
131-11-3	Dimethylphthalate	0.022	U	0.022	0.14	mg/Kg

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Injection Volume :		GPC Factor :	1.0
Prep Method :	SW3541	Level :	LOW
		Final Vol:	1000
		Test:	SVOC-TCL BNA -20
		GPC Cleanup :	N
		PH :	

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208-96-8	Acenaphthylene	0.023	U	0.023	0.14	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.027	U	0.027	0.14	mg/Kg
99-09-2	3-Nitroaniline	0.037	U	0.037	0.14	mg/Kg
83-32-9	Acenaphthene	0.17		0.017	0.14	mg/Kg
51-28-5	2,4-Dinitrophenol	0.18	U	0.18	0.27	mg/Kg
100-02-7	4-Nitrophenol	0.086	U	0.086	0.27	mg/Kg
132-64-9	Dibenzofuran	0.084	J	0.018	0.14	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.040	U	0.040	0.14	mg/Kg
84-66-2	Diethylphthalate	0.023	U	0.023	0.14	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.022	U	0.022	0.14	mg/Kg
86-73-7	Fluorene	0.13	J	0.020	0.14	mg/Kg
100-01-6	4-Nitroaniline	0.052	U	0.052	0.14	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.083	U	0.083	0.27	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.027	U	0.027	0.14	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.022	U	0.022	0.14	mg/Kg
118-74-1	Hexachlorobenzene	0.020	U	0.020	0.14	mg/Kg
1912-24-9	Atrazine	0.027	U	0.027	0.14	mg/Kg
87-86-5	Pentachlorophenol	0.041	U	0.041	0.27	mg/Kg
85-01-8	Phenanthrene	1.60		0.017	0.14	mg/Kg
120-12-7	Anthracene	0.38		0.027	0.14	mg/Kg
86-74-8	Carbazole	0.26		0.025	0.14	mg/Kg
84-74-2	Di-n-butylphthalate	0.039	U	0.039	0.14	mg/Kg
206-44-0	Fluoranthene	3.40	E	0.024	0.14	mg/Kg
129-00-0	Pyrene	2.70	E	0.029	0.14	mg/Kg
85-68-7	Butylbenzylphthalate	0.058	U	0.058	0.14	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.030	U	0.030	0.27	mg/Kg
56-55-3	Benzo(a)anthracene	1.90		0.019	0.14	mg/Kg
218-01-9	Chrysene	1.70		0.016	0.14	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.048	U	0.048	0.14	mg/Kg
117-84-0	Di-n-octyl phthalate	0.070	U	0.070	0.27	mg/Kg
205-99-2	Benzo(b)fluoranthene	2.30	E	0.015	0.14	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.71		0.018	0.14	mg/Kg
50-32-8	Benzo(a)pyrene	1.60		0.024	0.14	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.81		0.023	0.14	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.30		0.022	0.14	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.92		0.021	0.14	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.021	U	0.021	0.14	mg/Kg
123-91-1	1,4-Dioxane	0.036	U	0.036	0.14	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.022	U	0.022	0.14	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	71.9		18 - 112	48%	SPK: 150
13127-88-3	Phenol-d6	75.2		15 - 107	50%	SPK: 150
4165-60-0	Nitrobenzene-d5	41.7		18 - 107	42%	SPK: 100
321-60-8	2-Fluorobiphenyl	30.1		20 - 109	30%	SPK: 100
118-79-6	2,4,6-Tribromophenol	64.8		10 - 116	43%	SPK: 150
1718-51-0	Terphenyl-d14	24.2		10 - 105	24%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	310000	7.746
1146-65-2	Naphthalene-d8	1120000	10.54
15067-26-2	Acenaphthene-d10	742000	14.392
1517-22-2	Phenanthrene-d10	1460000	17.139
1719-03-5	Chrysene-d12	1420000	21.386
1520-96-3	Perylene-d12	1370000	24.38

TENTATIVE IDENTIFIED COMPOUNDS

000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	130	A	4.87	ug/Kg
000119-61-9	Benzophenone	170	J	15.8	ug/Kg
000486-25-9	9H-Fluoren-9-one	64.7	J	16.8	ug/Kg
000132-65-0	Dibenzothiophene	91.3	J	17.0	ug/Kg
002531-84-2	Phenanthrene, 2-methyl-	190	J	18.0	ug/Kg
000057-10-3	n-Hexadecanoic acid	190	J	18.0	ug/Kg
000832-69-9	Phenanthrene, 1-methyl-	320	J	18.1	ug/Kg

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	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	1.0	GPC Cleanup : N PH :
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000610-48-0	Anthracene, 1-methyl-	120	J		18.1	ug/Kg
000203-64-5	4H-Cyclopenta[def]phenanthrene	410	J		18.2	ug/Kg
000949-41-7	1H-Cyclopropa[1]phenanthrene,1a,9b	130	J		18.3	ug/Kg
000612-94-2	Naphthalene, 2-phenyl-	150	J		18.5	ug/Kg
000084-65-1	9,10-Anthracenedione	130	J		18.6	ug/Kg
003674-66-6	Phenanthrene, 2,5-dimethyl-	130	J		18.9	ug/Kg
005737-13-3	Cyclopenta(def)phenanthrenone	160	J		19.1	ug/Kg
000243-17-4	11H-Benzo[b]fluorene	95.3	J		20.1	ug/Kg
000239-35-0	Benzo[b]naphtho[2,1-d]thiophene	71.4	J		21.0	ug/Kg
000225-51-4	Benz[c]acridine	64.7	J		21.1	ug/Kg
000192-97-2	Benzo[e]pyrene	330	J		23.7	ug/Kg
000198-55-0	Perylene	1300	J		24.1	ug/Kg
000205-82-3	Benzo[j]fluoranthene	580	J		24.5	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