

Report of Analysis

Client:	Chemtech Consulting Group		Date Collected:	
Project:	NJ Soil PT		Date Received:	
Client Sample ID:	PB164402BS		SDG No.:	P4495
Lab Sample ID:	PB164402BS		Matrix:	SOIL
Analytical Method:	SW8270SIM		% Solid:	100
Sample Wt/Vol:	30.03	Units: g	Final Vol:	1000 uL
Soil Aliquot Vol:		uL	Test:	SVOCMS Group3
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
BN035094.D	1	10/25/24 09:50	11/14/24 18:59	PB164402

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
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TARGETS

91-20-3	Naphthalene	12.0		0.69	3.30	ug/Kg
91-57-6	2-Methylnaphthalene	12.0		0.80	3.30	ug/Kg
208-96-8	Acenaphthylene	12.1		0.63	3.30	ug/Kg
83-32-9	Acenaphthene	11.6		0.55	3.30	ug/Kg
86-73-7	Fluorene	11.2		0.60	3.30	ug/Kg
85-01-8	Phenanthrene	12.3		0.63	3.30	ug/Kg
120-12-7	Anthracene	12.5		0.79	3.30	ug/Kg
206-44-0	Fluoranthene	10.7		0.67	3.30	ug/Kg
129-00-0	Pyrene	13.6		0.95	3.30	ug/Kg
56-55-3	Benzo(a)anthracene	12.0		0.76	3.30	ug/Kg
218-01-9	Chrysene	12.6		0.86	3.30	ug/Kg
205-99-2	Benzo(b)fluoranthene	13.0		1.20	3.30	ug/Kg
207-08-9	Benzo(k)fluoranthene	12.8		0.98	3.30	ug/Kg
50-32-8	Benzo(a)pyrene	13.0		1.70	3.30	ug/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	11.2		1.30	3.30	ug/Kg
53-70-3	Dibenzo(a,h)anthracene	11.2		1.20	3.30	ug/Kg
191-24-2	Benzo(g,h,i)perylene	10.0		1.10	3.30	ug/Kg

SURROGATES

7297-45-2	2-Methylnaphthalene-d10	0.46		17 - 161	114%	SPK: 0.4
93951-69-0	Fluoranthene-d10	0.32		23 - 138	81%	SPK: 0.4
4165-60-0	Nitrobenzene-d5	0.35		33 - 121	86%	SPK: 0.4
321-60-8	2-Fluorobiphenyl	0.36		32 - 121	89%	SPK: 0.4
1718-51-0	Terphenyl-d14	0.42		21 - 130	105%	SPK: 0.4

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	2570	7.553			
1146-65-2	Naphthalene-d8	6460	10.319			
15067-26-2	Acenaphthene-d10	4850	14.19			
1517-22-2	Phenanthrene-d10	11500	16.932			
1719-03-5	Chrysene-d12	9990	21.133			

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1520-96-3	Perylene-d12	9580	23.303			

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