

Report of Analysis

Client:	CDM Smith	Date Collected:	04/28/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	04/29/25
Client Sample ID:	SS-5	SDG No.:	Q1914
Lab Sample ID:	Q1914-06	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	89.7
Sample Wt/Vol:	30.04	Units:	g
Soil Aliquot Vol:			uL
Extraction Type :		Decanted :	N
Injection Volume :		GPC Factor :	1.0
Prep Method :	SW3541	Final Vol:	1000
		Test:	SVOCMS Group3
		Level :	LOW
		GPC Cleanup :	N
		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BM050106.D	1	04/30/25 11:00	05/05/25 19:37	PB167803

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
91-20-3	Naphthalene	190	U	25.3	190	ug/Kg
208-96-8	Acenaphthylene	190	U	32.2	190	ug/Kg
83-32-9	Acenaphthene	120	J	23.7	190	ug/Kg
86-73-7	Fluorene	250		28.2	190	ug/Kg
85-01-8	Phenanthrene	200		23.3	190	ug/Kg
120-12-7	Anthracene	100	J	37.1	190	ug/Kg
206-44-0	Fluoranthene	360		33.4	190	ug/Kg
129-00-0	Pyrene	360		40.1	190	ug/Kg
56-55-3	Benzo(a)anthracene	110	J	25.6	190	ug/Kg
218-01-9	Chrysene	120	J	22.2	190	ug/Kg
205-99-2	Benzo(b)fluoranthene	150	J	21.2	190	ug/Kg
207-08-9	Benzo(k)fluoranthene	190	U	24.9	190	ug/Kg
50-32-8	Benzo(a)pyrene	110	J	32.8	190	ug/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	190	U	32.4	190	ug/Kg
53-70-3	Dibenzo(a,h)anthracene	190	U	30.5	190	ug/Kg
191-24-2	Benzo(g,h,i)perylene	78.8	J	28.6	190	ug/Kg
SURROGATES						
4165-60-0	Nitrobenzene-d5	41.6		18 - 107	42%	SPK: 100
321-60-8	2-Fluorobiphenyl	40.5		20 - 109	41%	SPK: 100
1718-51-0	Terphenyl-d14	38.1		10 - 105	38%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	358000		7.745		
1146-65-2	Naphthalene-d8	1220000		10.539		
15067-26-2	Acenaphthene-d10	725000		14.398		
1517-22-2	Phenanthrene-d10	1420000		17.145		
1719-03-5	Chrysene-d12	1470000		21.386		
1520-96-3	Perylene-d12	1570000		24.374		

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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