

Report of Analysis

Client:	Nobis Group	Date Collected:	07/17/25
Project:	Raymark Superfund Site	Date Received:	07/17/25
Client Sample ID:	RT2286MS	SDG No.:	Q2639
Lab Sample ID:	Q2635-01MS	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	84.8
Sample Wt/Vol:	50.07	Units:	g
Soil Aliquot Vol:			uL
Extraction Type :		Decanted :	N
Injection Volume :		GPC Factor :	1.0
Prep Method :	SW3541	Level :	LOW
		Test:	SVOCMS Group3
		Final Vol:	1000
		GPC Cleanup :	N
		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF143180.D	1	07/21/25 09:30	07/22/25 10:26	PB168929

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units(Dry Weight)
TARGETS							
91-20-3	Naphthalene	1.00		0.016	0.092	0.12	mg/Kg
91-57-6	2-Methylnaphthalene	1.00		0.018	0.092	0.12	mg/Kg
208-96-8	Acenaphthylene	1.00		0.020	0.092	0.12	mg/Kg
83-32-9	Acenaphthene	1.10		0.015	0.092	0.12	mg/Kg
86-73-7	Fluorene	1.00		0.018	0.092	0.12	mg/Kg
85-01-8	Phenanthrene	1.30		0.015	0.092	0.12	mg/Kg
120-12-7	Anthracene	1.10		0.024	0.092	0.12	mg/Kg
206-44-0	Fluoranthene	1.50		0.021	0.092	0.12	mg/Kg
129-00-0	Pyrene	1.40		0.025	0.092	0.12	mg/Kg
56-55-3	Benzo(a)anthracene	1.30		0.016	0.092	0.12	mg/Kg
218-01-9	Chrysene	1.30		0.014	0.092	0.12	mg/Kg
205-99-2	Benzo(b)fluoranthene	1.60		0.013	0.092	0.12	mg/Kg
207-08-9	Benzo(k)fluoranthene	1.20		0.016	0.092	0.12	mg/Kg
50-32-8	Benzo(a)pyrene	1.30		0.021	0.092	0.12	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.84		0.021	0.092	0.12	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.73		0.019	0.092	0.12	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.85		0.018	0.092	0.12	mg/Kg
SURROGATES							
4165-60-0	Nitrobenzene-d5	56.5		37 - 122		56%	SPK: 100
321-60-8	2-Fluorobiphenyl	52.0		44 - 115		52%	SPK: 100
1718-51-0	Terphenyl-d14	44.2	*	54 - 127		44%	SPK: 100
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	123000		6.963			
1146-65-2	Naphthalene-d8	483000		8.245			
15067-26-2	Acenaphthene-d10	259000		10.004			
1517-22-2	Phenanthrene-d10	421000		11.492			
1719-03-5	Chrysene-d12	278000		14.133			
1520-96-3	Perylene-d12	251000		15.633			

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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