

Report of Analysis

Client:	Nobis Group		Date Collected:		
Project:	Raymark Superfund Site		Date Received:		
Client Sample ID:	PB166776BS		SDG No.:	Q1383	
Lab Sample ID:	PB166776BS		Matrix:	SOIL	
Analytical Method:	SW8270		% Solid:	100	
Sample Wt/Vol:	30.02	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
BF141699.D	1	02/19/25 10:05	02/20/25 13:03	PB166776

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units(Dry Weight)
TARGETS							
91-20-3	Naphthalene	1.40		0.083	0.13	0.17	mg/Kg
91-57-6	2-Methylnaphthalene	1.40		0.082	0.13	0.17	mg/Kg
208-96-8	Acenaphthylene	1.50		0.086	0.13	0.17	mg/Kg
83-32-9	Acenaphthene	1.40		0.081	0.13	0.17	mg/Kg
86-73-7	Fluorene	1.40		0.085	0.13	0.17	mg/Kg
85-01-8	Phenanthrene	1.50		0.084	0.13	0.17	mg/Kg
120-12-7	Anthracene	1.60		0.084	0.13	0.17	mg/Kg
206-44-0	Fluoranthene	1.40		0.082	0.13	0.17	mg/Kg
129-00-0	Pyrene	1.40		0.083	0.13	0.17	mg/Kg
56-55-3	Benzo(a)anthracene	1.50		0.081	0.13	0.17	mg/Kg
218-01-9	Chrysene	1.50		0.079	0.13	0.17	mg/Kg
205-99-2	Benzo(b)fluoranthene	1.30		0.081	0.13	0.17	mg/Kg
207-08-9	Benzo(k)fluoranthene	1.60		0.083	0.13	0.17	mg/Kg
50-32-8	Benzo(a)pyrene	1.60		0.093	0.13	0.17	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	1.60		0.078	0.13	0.17	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	1.60		0.081	0.13	0.17	mg/Kg
191-24-2	Benzo(g,h,i)perylene	1.50		0.080	0.13	0.17	mg/Kg
SURROGATES							
4165-60-0	Nitrobenzene-d5	88.5		37 - 122		88%	SPK: 100
321-60-8	2-Fluorobiphenyl	94.1		44 - 115		94%	SPK: 100
1718-51-0	Terphenyl-d14	96.1		54 - 127		96%	SPK: 100
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	87200	6.781				
1146-65-2	Naphthalene-d8	333000	8.063				
15067-26-2	Acenaphthene-d10	174000	9.816				
1517-22-2	Phenanthrene-d10	270000	11.298				
1719-03-5	Chrysene-d12	174000	13.939				
1520-96-3	Perylene-d12	177000	15.38				

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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