

Report of Analysis

Client:	Alliance Technical Group, LLC - Newark		Date Collected:	04/21/25	
Project:	NJ Soil PT		Date Received:	04/24/25	
Client Sample ID:	HW0425-PT-BNA-SOIL		SDG No.:	Q1872	
Lab Sample ID:	Q1872-11		Matrix:	SOIL	
Analytical Method:	SW8270ESIM		% Solid:	100	
Sample Wt/Vol:	30.21	Units: g	Final Vol:	1000	uL
Soil Aliquot Vol:		uL	Test:	SVOCMS Group2	
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
BN036994.D	1	05/08/25 10:03	05/13/25 13:23	PB167915

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
62-75-9	n-Nitrosodimethylamine	1.70	U	1.70	6.60	ug/Kg
87-68-3	Hexachlorobutadiene	4000	E	1.10	3.30	ug/Kg
534-52-1	4,6-Dinitro-2-methylphenol	4.20	U	4.20	6.60	ug/Kg
118-74-1	Hexachlorobenzene	0.97	U	0.97	3.30	ug/Kg
87-86-5	Pentachlorophenol	3600	E	3.20	6.60	ug/Kg
123-91-1	1,4-Dioxane	1.60	U	1.60	6.60	ug/Kg
SURROGATES						
7297-45-2	2-Methylnaphthalene-d10	0	*	17 - 161	0%	SPK: 100
93951-69-0	Fluoranthene-d10	0	*	23 - 138	0%	SPK: 100
367-12-4	2-Fluorophenol	76.9		28 - 122	51%	SPK: 150
13127-88-3	Phenol-d6	87.7		25 - 125	58%	SPK: 150
4165-60-0	Nitrobenzene-d5	61.0		33 - 121	61%	SPK: 100
321-60-8	2-Fluorobiphenyl	48.1		32 - 121	48%	SPK: 100
118-79-6	2,4,6-Tribromophenol	111		19 - 110	74%	SPK: 150
1718-51-0	Terphenyl-d14	55.8		21 - 130	56%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	1720		7.618		
1146-65-2	Naphthalene-d8	4560		10.404		
15067-26-2	Acenaphthene-d10	2710		14.267		
1517-22-2	Phenanthrene-d10	5810		17.009		
1719-03-5	Chrysene-d12	6220		21.215		
1520-96-3	Perylene-d12	6500		23.415		

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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