

Report of Analysis

Client:	Kleinfelder	Date Collected:	10/13/25
Project:	Webster School	Date Received:	10/13/25
Client Sample ID:	LAYDOWN-YARD-1MS	SDG No.:	Q3341
Lab Sample ID:	Q3337-01MS	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	86.9
Sample Wt/Vol:	50.03 g	Test:	SVOCMS Group1
Prep Method :	SW3541		
	Level : LOW		
	Final Vol: 1000 uL		
	Prep Date: 10/15/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS									
78-59-1	Isophorone	800		2	45.3	230	ug/Kg	10/15/25 22:49	PB170106
91-20-3	Naphthalene	800		2	31.3	230	ug/Kg	10/15/25 22:49	PB170106
86-73-7	Fluorene	840		2	34.9	230	ug/Kg	10/15/25 22:49	PB170106
85-01-8	Phenanthrene	960		2	28.8	230	ug/Kg	10/15/25 22:49	PB170106
120-12-7	Anthracene	860		2	46.0	230	ug/Kg	10/15/25 22:49	PB170106
129-00-0	Pyrene	880		2	49.7	230	ug/Kg	10/15/25 22:49	PB170106
56-55-3	Benzo(a)anthracene	930		2	31.7	230	ug/Kg	10/15/25 22:49	PB170106
218-01-9	Chrysene	890		2	27.5	230	ug/Kg	10/15/25 22:49	PB170106
205-99-2	Benzo(b)fluoranthene	1000		2	26.2	230	ug/Kg	10/15/25 22:49	PB170106
50-32-8	Benzo(a)pyrene	940		2	40.7	230	ug/Kg	10/15/25 22:49	PB170106
191-24-2	Benzo(g,h,i)perylene	720		2	35.5	230	ug/Kg	10/15/25 22:49	PB170106
SURROGATES									
4165-60-0	Nitrobenzene-d5	51.2			18 - 107	51%	SPK: 100		
321-60-8	2-Fluorobiphenyl	52.1			20 - 109	52%	SPK: 100		
1718-51-0	Terphenyl-d14	40.0			10 - 105	40%	SPK: 100		
INTERNAL STANDARDS									
		Area Count							
3855-82-1	1,4-Dichlorobenzene-d4	182000							
1146-65-2	Naphthalene-d8	654000							
15067-26-2	Acenaphthene-d10	309000							
1517-22-2	Phenanthrene-d10	475000							
1719-03-5	Chrysene-d12	360000							
1520-96-3	Perylene-d12	304000							

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