

Report of Analysis

Client: Kleinfelder
Project: Webster School
Client Sample ID: C-3
Lab Sample ID: Q3341-03
Analytical Method: 8270E
Sample Wt/Vol: 30.05 g
Prep Method: SW3541

Level: LOW
Final Vol: 1000 uL
Prep Date: 10/15/25

Date Collected: 10/13/25
Date Received: 10/14/25
SDG No.: Q3341
Matrix: SOIL
% Solid: 80.1
Test: SVOCMS Group1

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS									
78-59-1	Isophorone	40.9	U	1	40.9	210	ug/Kg	10/15/25 18:56	PB170106
91-20-3	Naphthalene	28.3	U	1	28.3	210	ug/Kg	10/15/25 18:56	PB170106
86-73-7	Fluorene	31.5	U	1	31.5	210	ug/Kg	10/15/25 18:56	PB170106
85-01-8	Phenanthrene	26.0	U	1	26.0	210	ug/Kg	10/15/25 18:56	PB170106
120-12-7	Anthracene	41.5	U	1	41.5	210	ug/Kg	10/15/25 18:56	PB170106
129-00-0	Pyrene	44.9	U	1	44.9	210	ug/Kg	10/15/25 18:56	PB170106
56-55-3	Benzo(a)anthracene	28.7	U	1	28.7	210	ug/Kg	10/15/25 18:56	PB170106
218-01-9	Chrysene	24.8	U	1	24.8	210	ug/Kg	10/15/25 18:56	PB170106
205-99-2	Benzo(b)fluoranthene	23.7	U	1	23.7	210	ug/Kg	10/15/25 18:56	PB170106
50-32-8	Benzo(a)pyrene	36.8	U	1	36.8	210	ug/Kg	10/15/25 18:56	PB170106
191-24-2	Benzo(g,h,i)perylene	32.0	U	1	32.0	210	ug/Kg	10/15/25 18:56	PB170106
SURROGATES									
4165-60-0	Nitrobenzene-d5	55.8			18 - 107	56%	SPK: 100		
321-60-8	2-Fluorobiphenyl	51.5			20 - 109	52%	SPK: 100		
1718-51-0	Terphenyl-d14	59.1			10 - 105	59%	SPK: 100		
INTERNAL STANDARDS									
		Area Count							
3855-82-1	1,4-Dichlorobenzene-d4	110000							
1146-65-2	Naphthalene-d8	434000							
15067-26-2	Acenaphthene-d10	242000							
1517-22-2	Phenanthrene-d10	411000							
1719-03-5	Chrysene-d12	222000							
1520-96-3	Perylene-d12	220000							

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