

Report of Analysis

Client:	Sciacca General Contractors, LLC	Date Collected:	08/04/25
Project:	92 Clifford Street, Newark	Date Received:	08/04/25
Client Sample ID:	TP-2MSD	SDG No.:	Q2758
Lab Sample ID:	Q2763-01MSD	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	81.8
Sample Wt/Vol:	50.05 Units: g	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	SVOC-PAH
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
BF143328.D	1	08/05/25 08:50	08/06/25 14:18	PB169119

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
91-20-3	Naphthalene	1000		16.6	120	ug/Kg
208-96-8	Acenaphthylene	1100		21.2	120	ug/Kg
83-32-9	Acenaphthene	1100		15.6	120	ug/Kg
86-73-7	Fluorene	1100		18.5	120	ug/Kg
85-01-8	Phenanthrene	1300		15.3	120	ug/Kg
120-12-7	Anthracene	1100		24.4	120	ug/Kg
206-44-0	Fluoranthene	1700		22.0	120	ug/Kg
129-00-0	Pyrene	1000		26.4	120	ug/Kg
56-55-3	Benzo(a)anthracene	1300		16.9	120	ug/Kg
218-01-9	Chrysene	1200		14.6	120	ug/Kg
205-99-2	Benzo(b)fluoranthene	1700		13.9	120	ug/Kg
207-08-9	Benzo(k)fluoranthene	1300		16.4	120	ug/Kg
50-32-8	Benzo(a)pyrene	1300		21.6	120	ug/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	920		21.3	120	ug/Kg
53-70-3	Dibenzo(a,h)anthracene	810		20.1	120	ug/Kg
191-24-2	Benzo(g,h,i)perylene	950		18.8	120	ug/Kg
SURROGATES						
4165-60-0	Nitrobenzene-d5	50.2		18 - 107	50%	SPK: 100
321-60-8	2-Fluorobiphenyl	44.7		20 - 109	45%	SPK: 100
1718-51-0	Terphenyl-d14	35.0		10 - 105	35%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	77300		6.963		
1146-65-2	Naphthalene-d8	265000		8.245		
15067-26-2	Acenaphthene-d10	133000		9.998		
1517-22-2	Phenanthrene-d10	243000		11.486		
1719-03-5	Chrysene-d12	227000		14.133		
1520-96-3	Perylene-d12	163000		15.645		

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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