Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	04/28/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	04/29/25
Client Sample ID:	SS-8MEDL	SDG No.:	Q1914
Lab Sample ID:	Q1914-09MEDL	Matrix:	SOIL
Analytical Method:	8260D	% Solid:	92.4
Sample Wt/Vol:	8 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	100 uL	Test:	VOCMS Group3
GC Column:	RXI-624 ID : 0.25	Level :	MED
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VN086496.D	20		05/05/25 19:14	VN050525

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
71-43-2	Benzene	3400	UD	530	3400	ug/Kg
108-88-3	Toluene	1100	JD	530	3400	ug/Kg
100-41-4	Ethyl Benzene	2700	JD	450	3400	ug/Kg
1330-20-7	Total Xylenes	15200	D	1390	10200	ug/Kg
98-82-8	Isopropylbenzene	1000	JD	530	3400	ug/Kg
103-65-1	n-propylbenzene	2900	JD	490	3400	ug/Kg
108-67-8	1,3,5-Trimethylbenzene	5200	D	550	3400	ug/Kg
98-06-6	tert-Butylbenzene	3400	UD	450	3400	ug/Kg
95-63-6	1,2,4-Trimethylbenzene	18900	D	430	3400	ug/Kg
135-98-8	sec-Butylbenzene	1300	JD	450	3400	ug/Kg
99-87-6	p-Isopropyltoluene	900	JD	420	3400	ug/Kg
104-51-8	n-Butylbenzene	2100	JD	980	3400	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	47.5		63 - 155	95%	SPK: 50
1868-53-7	Dibromofluoromethane	59.2		70 - 134	118%	SPK: 50
2037-26-5	Toluene-d8	51.5		74 - 123	103%	SPK: 50
460-00-4	4-Bromofluorobenzene	53.6		38 - 136	107%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	168000	8.224			
540-36-3	1,4-Difluorobenzene	313000	9.1			
3114-55-4	Chlorobenzene-d5	300000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	141000	13.788			

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