

Report of Analysis

Client:	LaBella Associates P.C.	Date Collected:	
Project:	613 Pine Avenue	Date Received:	
Client Sample ID:	VY0717SBS01	SDG No.:	O3631
Lab Sample ID:	VY0717SBS01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	100
Sample Wt/Vol:	5 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group1
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY014665.D	1		07/17/23 12:23	VY071723

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	23.5		1.60	5.00	ug/Kg
74-87-3	Chloromethane	25.4		0.91	5.00	ug/Kg
75-01-4	Vinyl Chloride	24.9		0.93	5.00	ug/Kg
74-83-9	Bromomethane	25.0		1.20	5.00	ug/Kg
75-00-3	Chloroethane	24.9		0.88	5.00	ug/Kg
75-69-4	Trichlorofluoromethane	22.2		1.10	5.00	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	21.7		0.72	5.00	ug/Kg
75-35-4	1,1-Dichloroethene	22.1		0.79	5.00	ug/Kg
67-64-1	Acetone	120		9.40	25.0	ug/Kg
75-15-0	Carbon Disulfide	22.2		2.20	5.00	ug/Kg
1634-04-4	Methyl tert-butyl Ether	23.4		0.65	5.00	ug/Kg
79-20-9	Methyl Acetate	25.4		1.60	5.00	ug/Kg
75-09-2	Methylene Chloride	23.3		6.10	10.0	ug/Kg
156-60-5	trans-1,2-Dichloroethene	22.4		0.73	5.00	ug/Kg
75-34-3	1,1-Dichloroethane	22.1		0.72	5.00	ug/Kg
110-82-7	Cyclohexane	20.5		0.70	5.00	ug/Kg
78-93-3	2-Butanone	130		7.30	25.0	ug/Kg
56-23-5	Carbon Tetrachloride	21.5		0.78	5.00	ug/Kg
156-59-2	cis-1,2-Dichloroethene	22.9		0.64	5.00	ug/Kg
74-97-5	Bromochloromethane	23.6		2.40	5.00	ug/Kg
67-66-3	Chloroform	22.4		1.30	5.00	ug/Kg
71-55-6	1,1,1-Trichloroethane	21.7		0.76	5.00	ug/Kg
108-87-2	Methylcyclohexane	20.7		3.40	5.00	ug/Kg
71-43-2	Benzene	22.2		0.66	5.00	ug/Kg
107-06-2	1,2-Dichloroethane	22.5		0.72	5.00	ug/Kg
79-01-6	Trichloroethene	22.2		0.66	5.00	ug/Kg
78-87-5	1,2-Dichloropropane	21.8		0.59	5.00	ug/Kg
75-27-4	Bromodichloromethane	21.8		0.70	5.00	ug/Kg
108-10-1	4-Methyl-2-Pentanone	120		4.50	25.0	ug/Kg
108-88-3	Toluene	22.1		0.65	5.00	ug/Kg

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10061-02-6	t-1,3-Dichloropropene	22.1		0.77	5.00	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	21.8		0.74	5.00	ug/Kg
79-00-5	1,1,2-Trichloroethane	22.9		0.86	5.00	ug/Kg
591-78-6	2-Hexanone	120		5.30	25.0	ug/Kg
124-48-1	Dibromochloromethane	22.8		0.85	5.00	ug/Kg
106-93-4	1,2-Dibromoethane	23.4		0.79	5.00	ug/Kg
127-18-4	Tetrachloroethene	22.9		0.77	5.00	ug/Kg
108-90-7	Chlorobenzene	22.0		0.63	5.00	ug/Kg
100-41-4	Ethyl Benzene	22.1		0.67	5.00	ug/Kg
179601-23-1	m/p-Xylenes	43.0		1.40	10.0	ug/Kg
1330-20-7	Total Xylenes	65.0		2.17	15.0	ug/Kg
95-47-6	o-Xylene	22.0		0.77	5.00	ug/Kg
100-42-5	Styrene	22.2		0.69	5.00	ug/Kg
75-25-2	Bromoform	23.6		0.95	5.00	ug/Kg
98-82-8	Isopropylbenzene	21.0		0.71	5.00	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	22.5		1.10	5.00	ug/Kg
103-65-1	n-propylbenzene	20.1		0.70	5.00	ug/Kg
108-67-8	1,3,5-Trimethylbenzene	20.9		0.64	5.00	ug/Kg
98-06-6	tert-Butylbenzene	20.3		0.77	5.00	ug/Kg
95-63-6	1,2,4-Trimethylbenzene	21.3		0.67	5.00	ug/Kg
135-98-8	sec-Butylbenzene	20.2		0.77	5.00	ug/Kg
99-87-6	p-Isopropyltoluene	20.5		0.74	5.00	ug/Kg
541-73-1	1,3-Dichlorobenzene	21.6		0.68	5.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	21.5		0.60	5.00	ug/Kg
104-51-8	n-Butylbenzene	20.0		0.68	5.00	ug/Kg
95-50-1	1,2-Dichlorobenzene	21.7		0.60	5.00	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	23.7		1.20	5.00	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	22.0		0.61	5.00	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	21.6		0.63	5.00	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	51.3		50 - 163	103%	SPK: 50

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1868-53-7	Dibromofluoromethane	50.0		54 - 147	100%	SPK: 50
2037-26-5	Toluene-d8	49.6		58 - 134	99%	SPK: 50
460-00-4	4-Bromofluorobenzene	49.6		30 - 143	99%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	291000	7.783			
540-36-3	1,4-Difluorobenzene	487000	8.685			
3114-55-4	Chlorobenzene-d5	432000	11.489			
3855-82-1	1,4-Dichlorobenzene-d4	217000	13.422			

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