

Report of Analysis

Client:	LaBella Associates P.C.	Date Collected:	
Project:	Mackenna Parcels	Date Received:	
Client Sample ID:	VD0717SBS01	SDG No.:	O3645
Lab Sample ID:	VD0717SBS01	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:	RTX-VMS ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VD076747.D	1		07/17/23 11:46	VD071723

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	22.5		1.60	5.00	ug/Kg
74-87-3	Chloromethane	19.5		0.91	5.00	ug/Kg
75-01-4	Vinyl Chloride	19.7		0.93	5.00	ug/Kg
74-83-9	Bromomethane	19.1		1.20	5.00	ug/Kg
75-00-3	Chloroethane	19.1		0.88	5.00	ug/Kg
75-69-4	Trichlorofluoromethane	21.9		1.10	5.00	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	21.9		0.72	5.00	ug/Kg
75-35-4	1,1-Dichloroethene	20.5		0.79	5.00	ug/Kg
67-64-1	Acetone	100		9.40	25.0	ug/Kg
75-15-0	Carbon Disulfide	20.5		2.20	5.00	ug/Kg
1634-04-4	Methyl tert-butyl Ether	20.0		0.65	5.00	ug/Kg
79-20-9	Methyl Acetate	20.5		1.60	5.00	ug/Kg
75-09-2	Methylene Chloride	17.3		6.10	10.0	ug/Kg
156-60-5	trans-1,2-Dichloroethene	20.2		0.73	5.00	ug/Kg
75-34-3	1,1-Dichloroethane	20.1		0.72	5.00	ug/Kg
110-82-7	Cyclohexane	20.4		0.70	5.00	ug/Kg
78-93-3	2-Butanone	100		7.30	25.0	ug/Kg
56-23-5	Carbon Tetrachloride	21.4		0.78	5.00	ug/Kg
156-59-2	cis-1,2-Dichloroethene	20.8		0.64	5.00	ug/Kg
74-97-5	Bromochloromethane	23.7		2.40	5.00	ug/Kg
67-66-3	Chloroform	20.6		1.30	5.00	ug/Kg
71-55-6	1,1,1-Trichloroethane	22.5		0.76	5.00	ug/Kg
108-87-2	Methylcyclohexane	20.8		3.40	5.00	ug/Kg
71-43-2	Benzene	20.6		0.66	5.00	ug/Kg
107-06-2	1,2-Dichloroethane	21.4		0.72	5.00	ug/Kg
79-01-6	Trichloroethene	21.1		0.66	5.00	ug/Kg
78-87-5	1,2-Dichloropropane	20.7		0.59	5.00	ug/Kg
75-27-4	Bromodichloromethane	20.5		0.70	5.00	ug/Kg
108-10-1	4-Methyl-2-Pentanone	99.9		4.50	25.0	ug/Kg
108-88-3	Toluene	20.4		0.65	5.00	ug/Kg
10061-02-6	t-1,3-Dichloropropene	19.8		0.77	5.00	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	19.8		0.74	5.00	ug/Kg

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79-00-5	1,1,2-Trichloroethane	21.0		0.86	5.00	ug/Kg
591-78-6	2-Hexanone	99.6		5.30	25.0	ug/Kg
124-48-1	Dibromochloromethane	21.4		0.85	5.00	ug/Kg
106-93-4	1,2-Dibromoethane	20.6		0.79	5.00	ug/Kg
127-18-4	Tetrachloroethene	23.3		0.77	5.00	ug/Kg
108-90-7	Chlorobenzene	21.1		0.63	5.00	ug/Kg
100-41-4	Ethyl Benzene	20.7		0.67	5.00	ug/Kg
179601-23-1	m/p-Xylenes	41.3		1.40	10.0	ug/Kg
1330-20-7	Total Xylenes	61.8		2.17	15.0	ug/Kg
95-47-6	o-Xylene	20.5		0.77	5.00	ug/Kg
100-42-5	Styrene	20.7		0.69	5.00	ug/Kg
75-25-2	Bromoform	22.4		0.95	5.00	ug/Kg
98-82-8	Isopropylbenzene	21.1		0.71	5.00	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	20.5		1.10	5.00	ug/Kg
103-65-1	n-propylbenzene	21.1		0.70	5.00	ug/Kg
108-67-8	1,3,5-Trimethylbenzene	21.0		0.64	5.00	ug/Kg
98-06-6	tert-Butylbenzene	21.3		0.77	5.00	ug/Kg
95-63-6	1,2,4-Trimethylbenzene	20.3		0.67	5.00	ug/Kg
135-98-8	sec-Butylbenzene	21.4		0.77	5.00	ug/Kg
99-87-6	p-Isopropyltoluene	21.1		0.74	5.00	ug/Kg
541-73-1	1,3-Dichlorobenzene	21.5		0.68	5.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	21.3		0.60	5.00	ug/Kg
104-51-8	n-Butylbenzene	21.2		0.68	5.00	ug/Kg
95-50-1	1,2-Dichlorobenzene	21.6		0.60	5.00	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	20.1		1.20	5.00	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	20.8		0.61	5.00	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	21.3		0.63	5.00	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	49.1		50 - 163	98%	SPK: 50
1868-53-7	Dibromofluoromethane	49.0		54 - 147	98%	SPK: 50
2037-26-5	Toluene-d8	47.3		58 - 134	95%	SPK: 50
460-00-4	4-Bromofluorobenzene	48.4		30 - 143	97%	SPK: 50

INTERNAL STANDARDS

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363-72-4	Pentafluorobenzene	204000	7.875			
540-36-3	1,4-Difluorobenzene	374000	8.775			
3114-55-4	Chlorobenzene-d5	340000	11.581			
3855-82-1	1,4-Dichlorobenzene-d4	159000	13.516			

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