

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD050919\
Data File : VD062197.D
Acq On : 9 May 2019 19:28
Operator : FY/SY
Sample : K2717-03
Misc : 5.66g/5mL/MSVOA D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P026-SS001-0612-01

Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D050719S.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.244	79	83	86	rVB3	5375	13895	1.11%	0.172%
2	3.799	235	241	247	rBV	64337	167712	13.36%	2.076%
3	6.899	550	556	563	rBV2	290177	841352	67.03%	10.416%
4	7.017	563	568	584	rVB	226660	662051	52.74%	8.196%
5	7.430	604	610	621	rBV	234136	617294	49.18%	7.642%
6	8.198	682	688	703	rVB	396650	977333	77.86%	12.100%
7	8.542	719	723	732	rBV2	49296	132922	10.59%	1.646%
8	10.392	905	911	918	rBV	577860	1255224	100.00%	15.540%
9	11.219	990	995	1000	rBV	9368	25971	2.07%	0.322%
10	12.351	1106	1110	1121	rVB	763331	1223457	97.47%	15.147%
11	13.542	1228	1231	1235	rBV	694673	1056350	84.16%	13.078%
12	13.709	1245	1248	1251	rVB	9851	15743	1.25%	0.195%
13	14.506	1326	1329	1335	rVB	853636	1087981	86.68%	13.470%

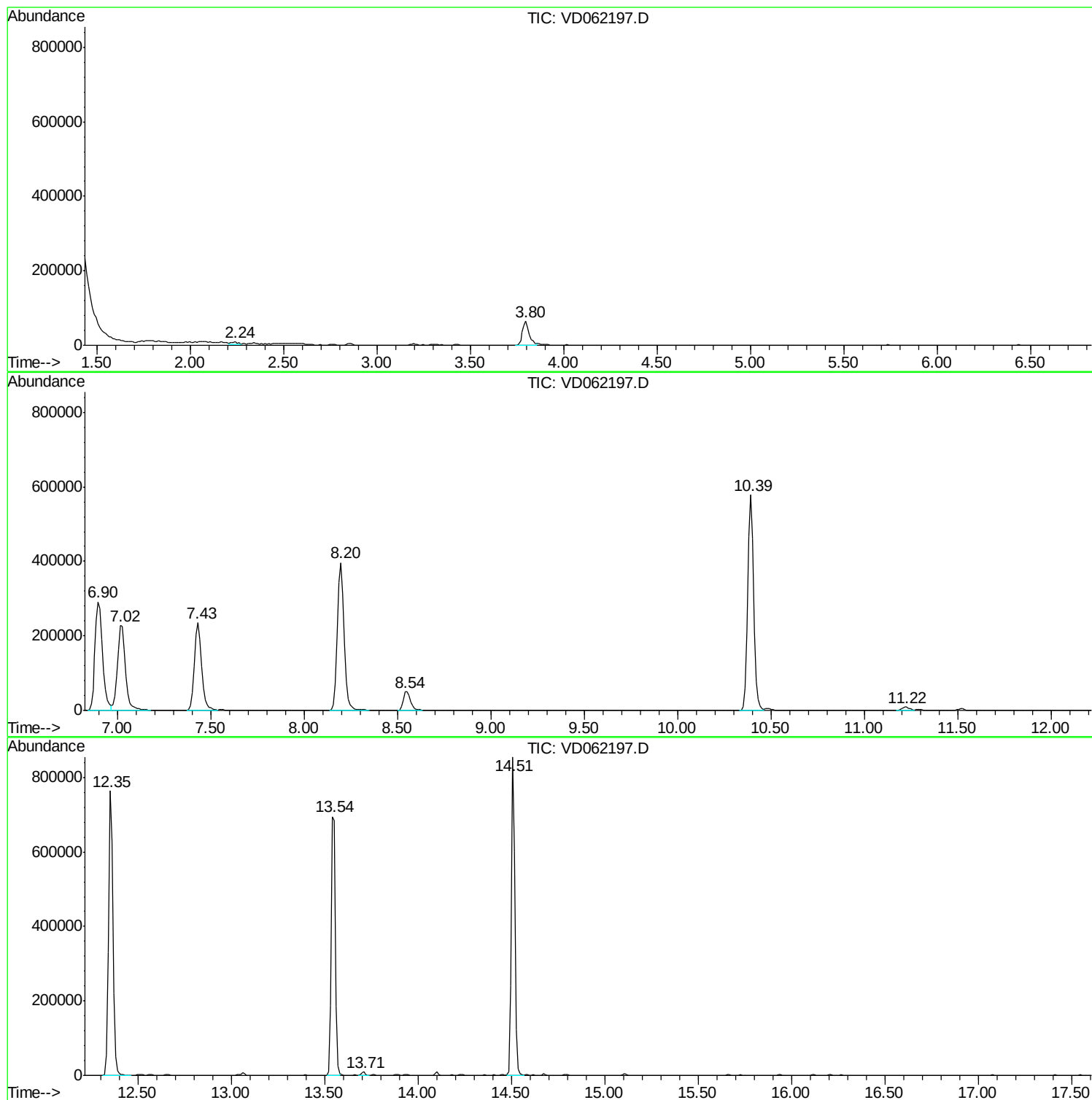
Sum of corrected areas: 8077285

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D050719S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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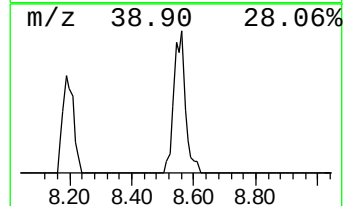
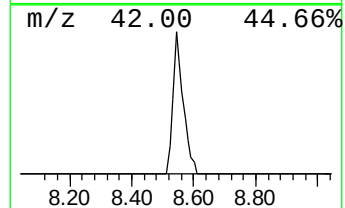
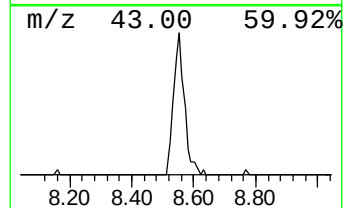
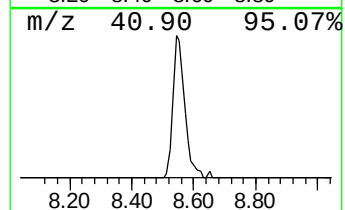
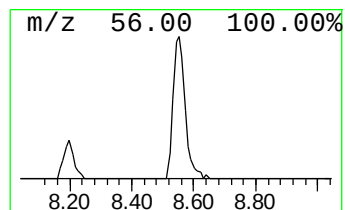
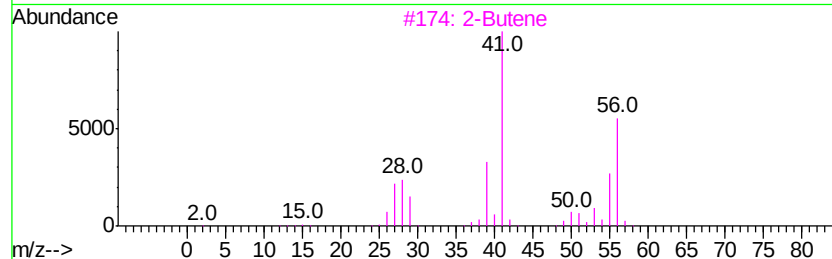
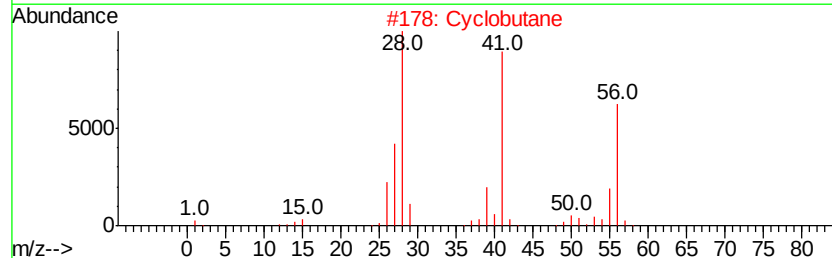
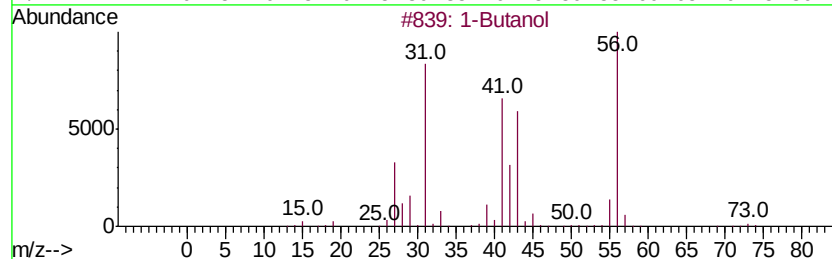
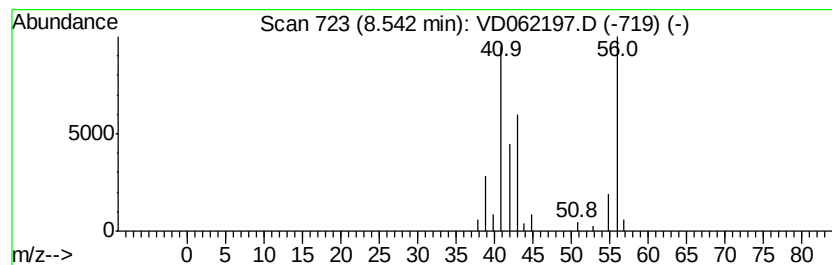
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Butanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.54	6.80 ug/l	132922	1,4-Difluorobenzene	8.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butanol	74	C4H10O	000071-36-3	83
2		Cyclobutane	56	C4H8	000287-23-0	7
3		2-Butene	56	C4H8	000107-01-7	5
4		2-Butene, (E)-	56	C4H8	000624-64-6	5
5		2-Propenal	56	C3H4O	000107-02-8	5



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Butanol	8.54	6.8	ug/l	132922	2	8.20	977333	50.0