

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW083119\
 Data File : VW012425.D
 Acq On : 01 Sep 2019 03:39
 Operator : SY/VA
 Sample : K4516-02
 Misc : 6.65G/5ML/MSVOA W/SOIL
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 T1-S-B2-(2.5)

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_W\METHOD\82W082319S.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	2.160	35	42	49	rBV	15624	29007	7.07%	1.422%
2	3.678	286	291	298	rBV7	1842	4802	1.17%	0.235%
3	4.909	484	493	504	rBV2	9920	31327	7.64%	1.536%
4	7.128	852	857	865	rBV4	1534	4126	1.01%	0.202%
5	7.531	918	923	929	rBV6	2123	4514	1.10%	0.221%
6	7.884	972	981	986	rBV	52933	130402	31.79%	6.393%
7	7.945	986	991	1002	rVB2	102046	226955	55.33%	11.127%
8	8.305	1039	1050	1058	rBV	60416	134364	32.75%	6.587%
9	8.842	1128	1138	1146	rBV	151492	300544	73.26%	14.735%
10	10.323	1370	1381	1389	rBV	233276	410216	100.00%	20.112%
11	10.378	1389	1390	1397	rVB4	3754	5872	1.43%	0.288%
12	10.707	1438	1444	1449	rBV2	3488	6031	1.47%	0.296%
13	11.634	1585	1596	1607	rBV	182970	328032	79.97%	16.082%
14	12.621	1750	1758	1767	rBV	124575	218240	53.20%	10.700%
15	13.560	1904	1912	1921	rBV	116537	205253	50.04%	10.063%

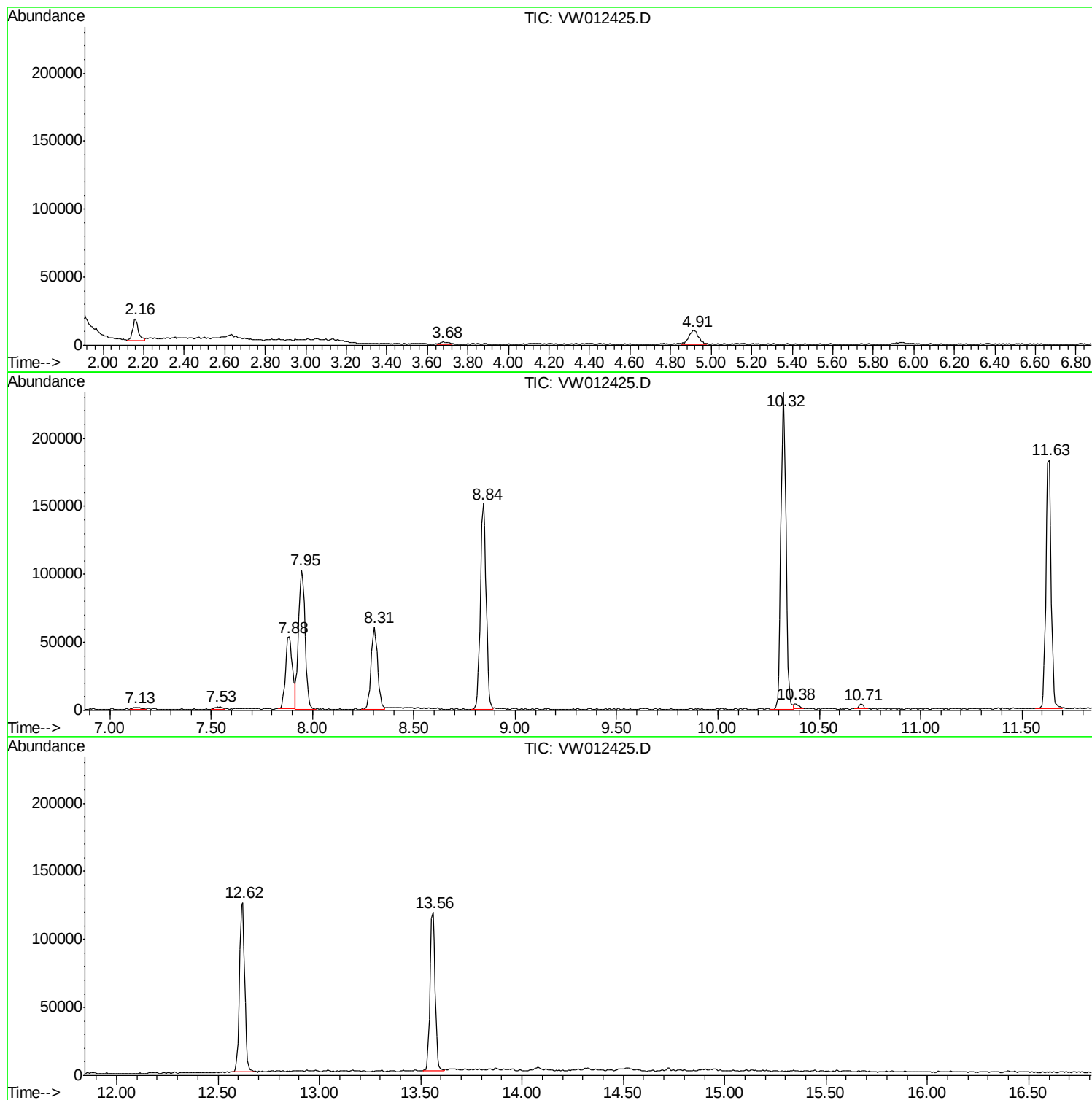
Sum of corrected areas: 2039685

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082319S.M
Quant Title : SW846 8260

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



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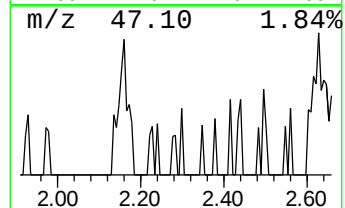
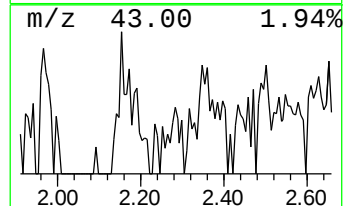
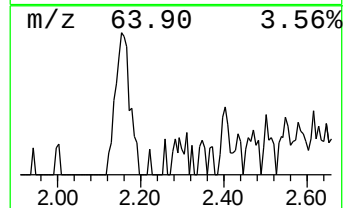
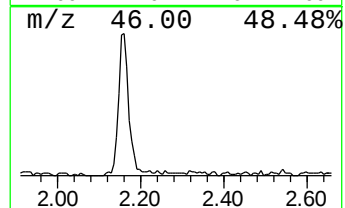
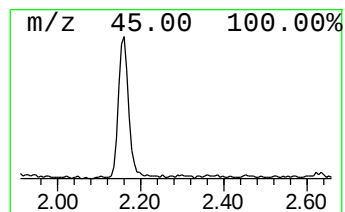
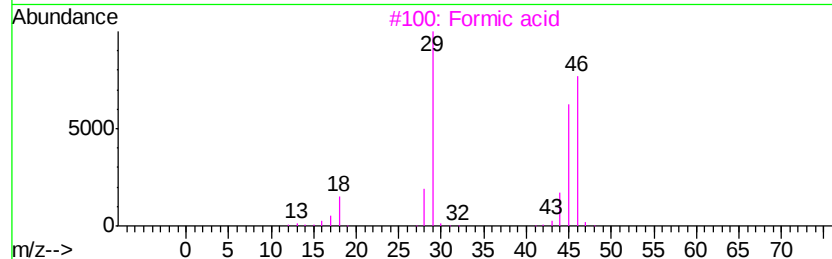
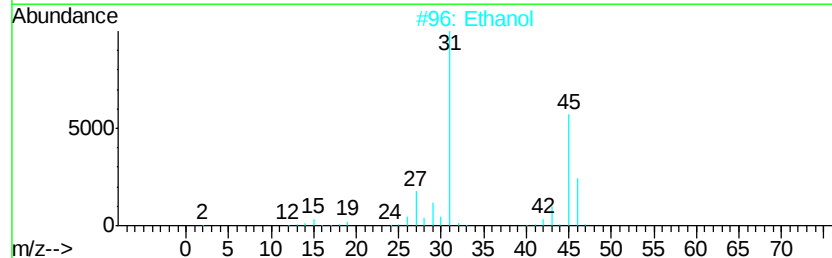
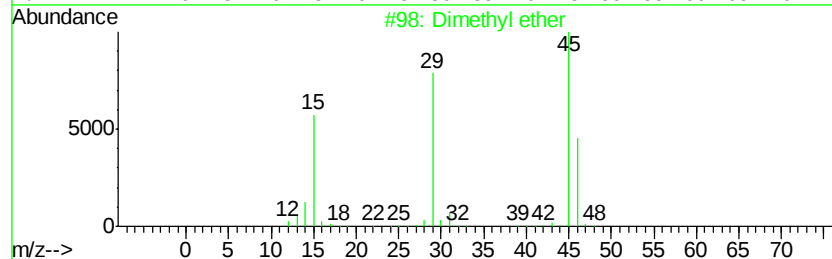
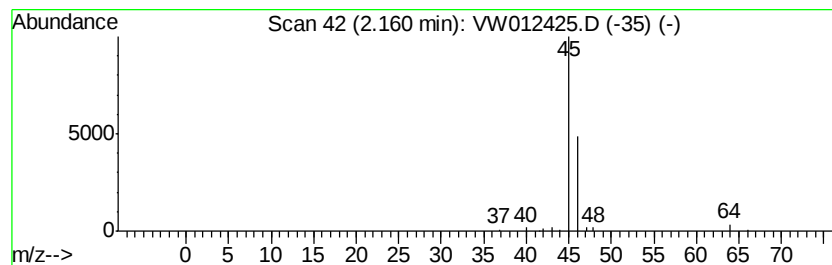
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Peak Number 1 Dimethyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	6.39 ug/l	29007	Pentafluorobenzene	7.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dimethyl ether		46	C2H6O	000115-10-6	64
2	Ethanol		46	C2H6O	000064-17-5	7
3	Formic acid		46	CH2O2	000064-18-6	5
4	Methane, nitroso-		45	CH3NO	000865-40-7	3
5	Formamide		45	CH3NO	000075-12-7	3



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dimethyl ether	2.16	6.4	ug/l	29007	1	7.95	226955	50.0