

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062318\
 Data File : VW003467.D
 Acq On : 23 Jun 2018 18:11
 Operator : SY/AP
 Sample : J3664-10
 Misc : 6.83G/5ML/MSVOA_W/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WP021SB476-31-33-180619

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\82W061318S.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	1.807	28	36	58	rVB	2269803	4764172	100.00%	39.081%
2	2.154	87	93	105	rBV	24870	77825	1.63%	0.638%
3	2.355	120	126	135	rBV	52556	93291	1.96%	0.765%
4	5.422	616	629	643	rBV	108065	330705	6.94%	2.713%
5	7.171	902	916	932	rBV	989161	2572532	54.00%	21.103%
6	7.885	1023	1033	1038	rBV	99100	234439	4.92%	1.923%
7	7.952	1038	1044	1057	rVB	222236	490740	10.30%	4.026%
8	8.312	1088	1103	1113	rBV	105241	242767	5.10%	1.991%
9	8.848	1180	1191	1201	rBV	328641	633737	13.30%	5.199%
10	9.092	1220	1231	1240	rBV	84832	172407	3.62%	1.414%
11	10.329	1426	1434	1443	rBV	363528	629245	13.21%	5.162%
12	11.634	1640	1648	1659	rBV	437392	759829	15.95%	6.233%
13	12.628	1802	1811	1819	rBV	260903	437447	9.18%	3.588%
14	13.567	1958	1965	1976	rVB	470332	751380	15.77%	6.164%

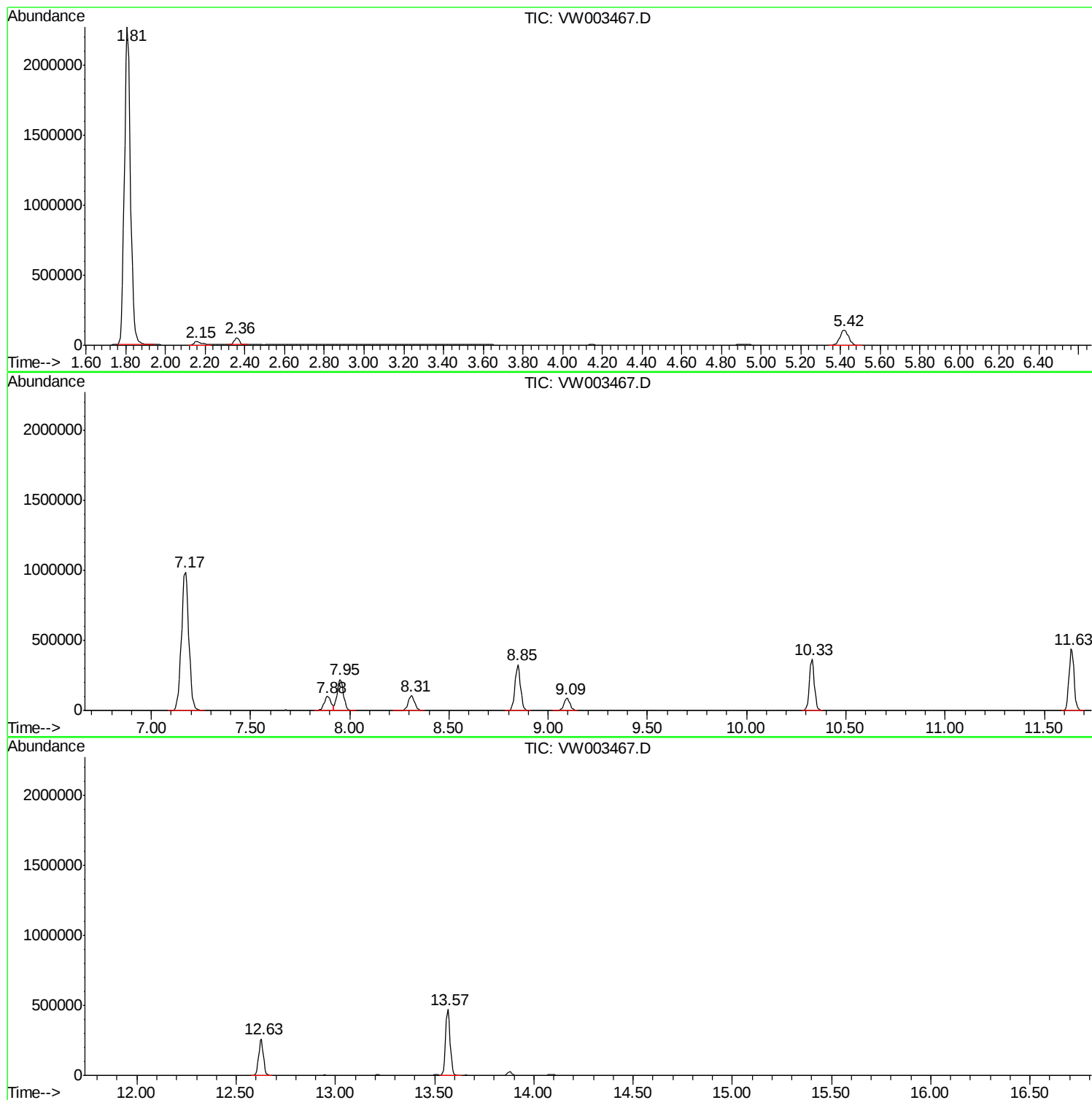
Sum of corrected areas: 12190516

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W061318S.M
Quant Title : SW846 8260

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



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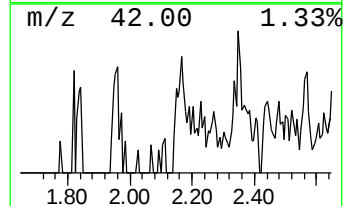
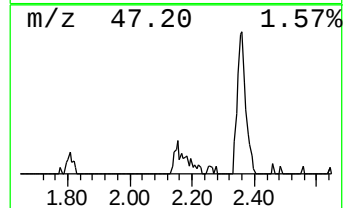
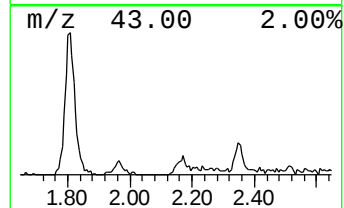
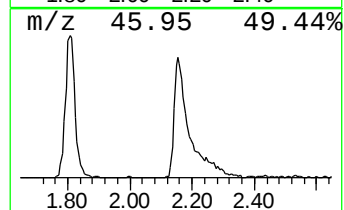
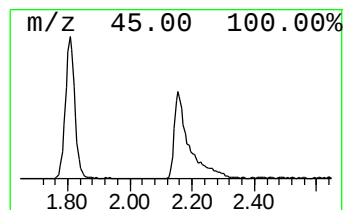
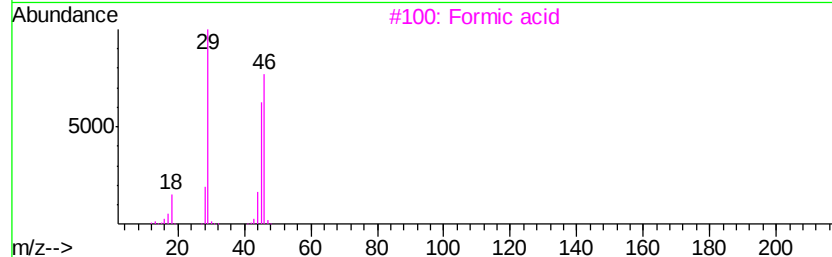
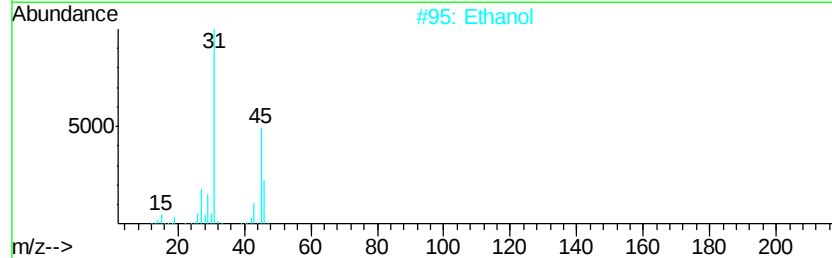
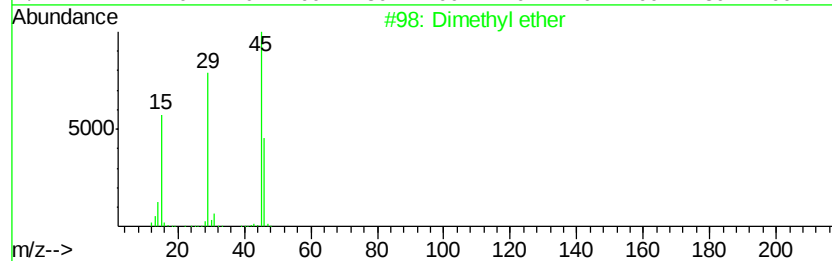
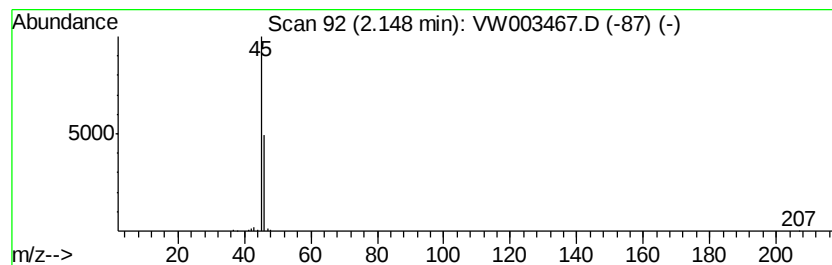
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dimethyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	7.93 ug/l	77825	Pentafluorobenzene	7.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Dimethyl ether	46 C2H6O	000115-10-6 74
2		Ethanol	46 C2H6O	000064-17-5 7
3		Formic acid	46 CH2O2	000064-18-6 5
4		Propanedioic acid, dihydroxy-	136 C3H4O6	000560-27-0 4
5		Methane, nitroso-	45 CH3NO	000865-40-7 3



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dimethyl ether	2.15	7.9	ug/l	77825	1	7.95	490740	50.0