

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW062318\
 Data File : VW003485.D
 Acq On : 24 Jun 2018 03:38
 Operator : SY/AP
 Sample : J3670-05
 Misc : 6.62G/5ML/MSVOA W/SOIL
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WP021SB487-17-19-180621

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	1.806	28	36	58	rVB	2124236	4527146	99.85%	31.577%
2	2.154	87	93	104	rBV	24884	72493	1.60%	0.506%
3	7.171	903	916	931	rBV	291749	766804	16.91%	5.349%
4	7.891	1023	1034	1038	rBV	120908	286374	6.32%	1.997%
5	7.952	1038	1044	1055	rVB	206343	461815	10.19%	3.221%
6	8.311	1093	1103	1115	rBV	136298	301511	6.65%	2.103%
7	8.848	1182	1191	1206	rBV	314477	608254	13.42%	4.243%
8	10.329	1427	1434	1441	rBV	455281	784560	17.30%	5.472%
9	11.664	1639	1653	1661	rBV2	2272557	4533937	100.00%	31.624%
10	12.341	1755	1764	1768	rBV4	16167	51246	1.13%	0.357%
11	12.628	1804	1811	1817	rBV	372751	620377	13.68%	4.327%
12	12.951	1858	1864	1869	rBV6	22828	46432	1.02%	0.324%
13	13.567	1959	1965	1975	rVB2	490554	876142	19.32%	6.111%
14	13.878	2010	2016	2024	rBV	192281	329501	7.27%	2.298%
15	14.085	2044	2050	2067	rVB4	22434	70210	1.55%	0.490%

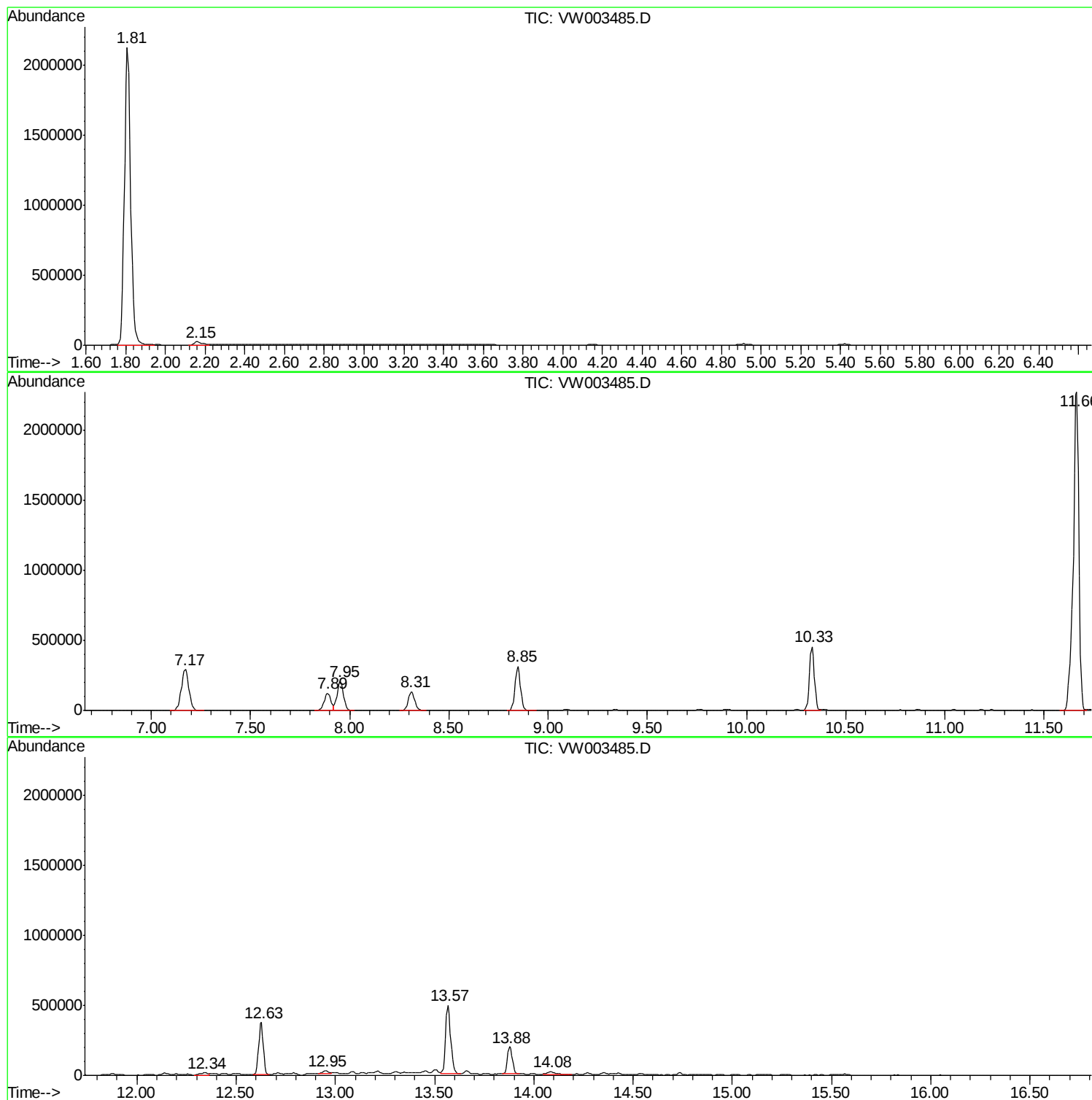
Sum of corrected areas: 14336802

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Quant Method : Z:\V0ASRV\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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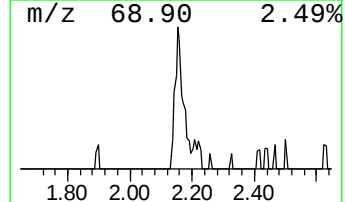
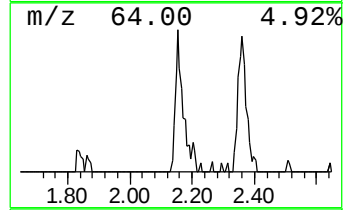
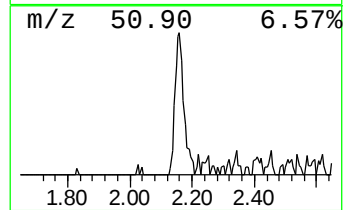
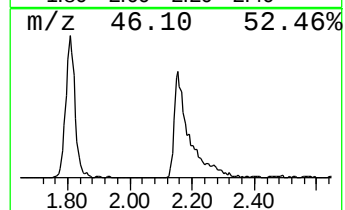
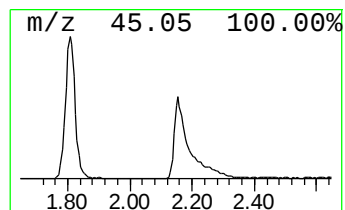
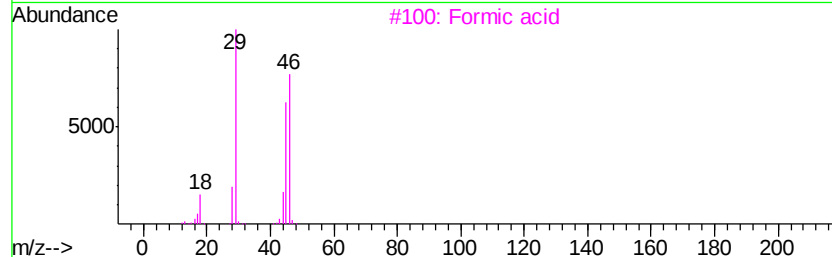
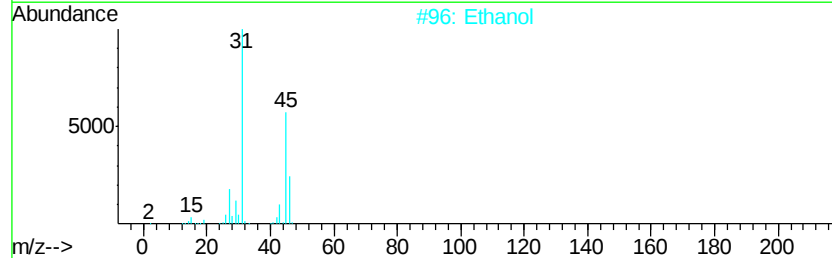
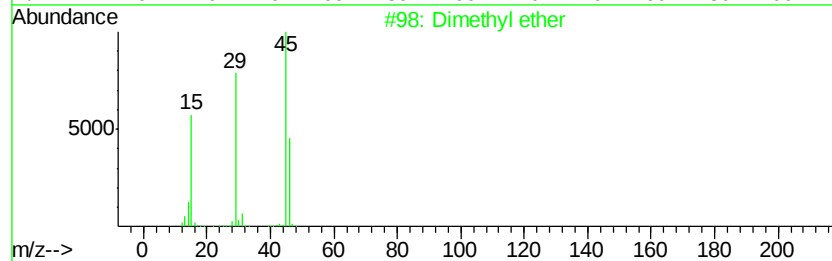
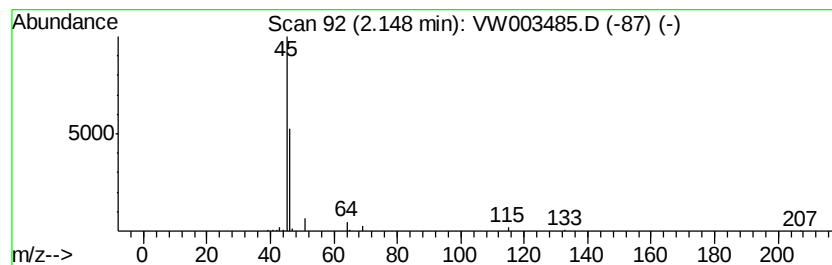
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Dimethyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	7.85 ug/l	72493	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl ether	46	C2H6O	000115-10-6	80
2		Ethanol	46	C2H6O	000064-17-5	9
3		Formic acid	46	CH2O2	000064-18-6	5
4		Propanedioic acid, dihydroxy-	136	C3H4O6	000560-27-0	4
5		Oxalic acid	90	C2H2O4	000144-62-7	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dimethyl ether	2.15	7.8	ug/l	72493	1	7.95	461815	50.0