

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW062318\
 Data File : VW003461.D
 Acq On : 23 Jun 2018 15:36
 Operator : SY/AP
 Sample : J3664-04
 Misc : 6.46G/5ML/MSVOA W/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WP021SB482-17-19-180619

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	56	rVV	1604512	3400325	100.00%	38.139%
2	1.953	56	60	70	rVB2	25328	40780	1.20%	0.457%
3	2.154	87	93	105	rBV	35360	102241	3.01%	1.147%
4	4.148	411	420	438	rVB2	22513	78090	2.30%	0.876%
5	4.904	536	544	558	rBV3	14484	52774	1.55%	0.592%
6	7.891	1024	1034	1038	rBV	91991	219894	6.47%	2.466%
7	7.952	1038	1044	1053	rVB	277376	609736	17.93%	6.839%
8	8.318	1090	1104	1115	rBV3	161216	422458	12.42%	4.738%
9	8.848	1182	1191	1207	rBV	417806	829794	24.40%	9.307%
10	10.330	1425	1434	1440	rBV	370325	644569	18.96%	7.230%
11	10.391	1440	1444	1456	rVV	20669	49853	1.47%	0.559%
12	11.634	1639	1648	1659	rBV	565148	1014132	29.82%	11.375%
13	12.476	1777	1786	1790	rBV2	20346	38329	1.13%	0.430%
14	12.530	1790	1795	1801	rVV3	28489	50389	1.48%	0.565%
15	12.628	1804	1811	1821	rVB	270468	457369	13.45%	5.130%
16	13.567	1958	1965	1974	rVB	567919	904935	26.61%	10.150%

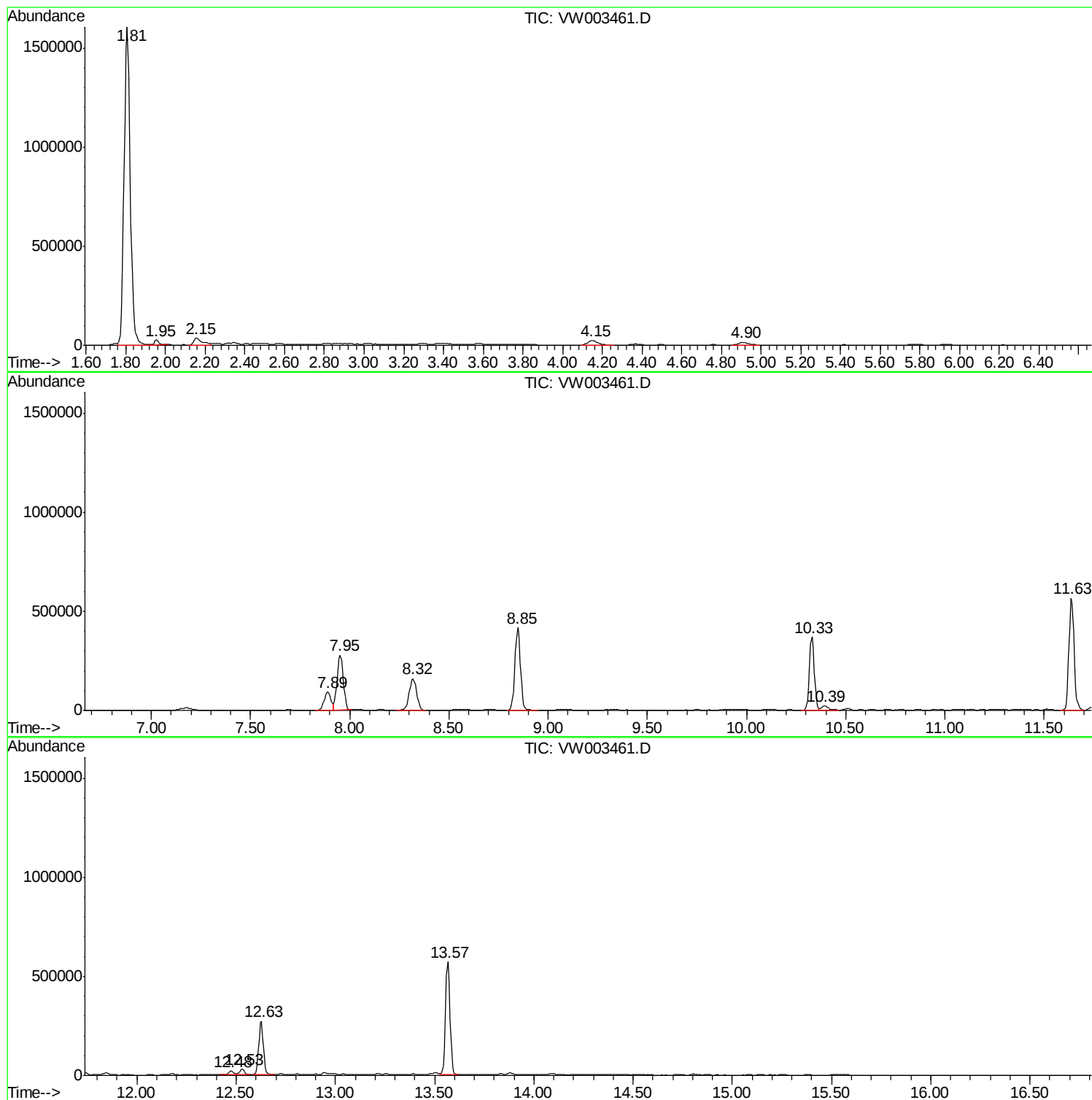
Sum of corrected areas: 8915668

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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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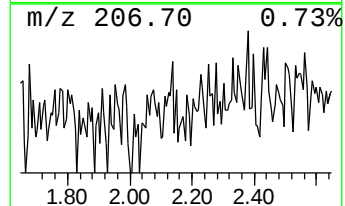
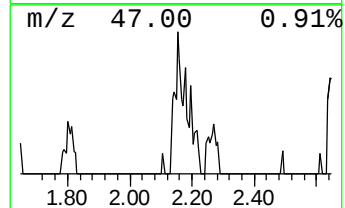
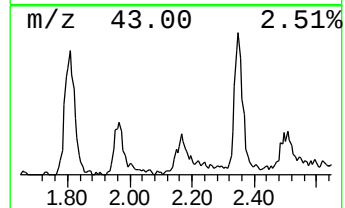
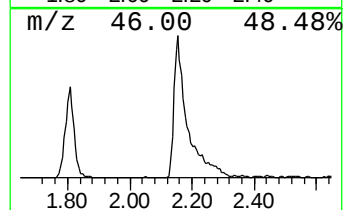
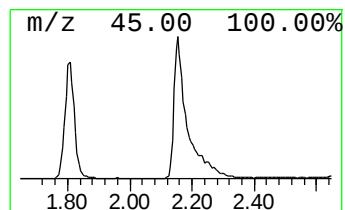
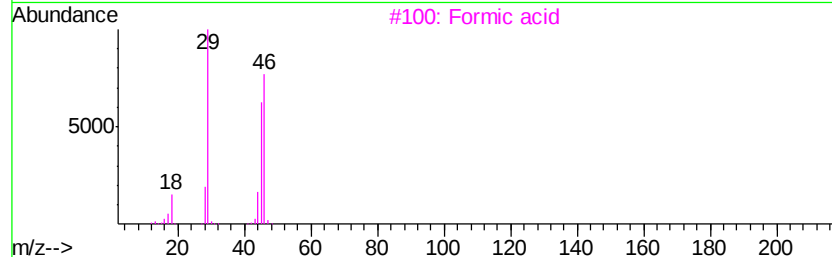
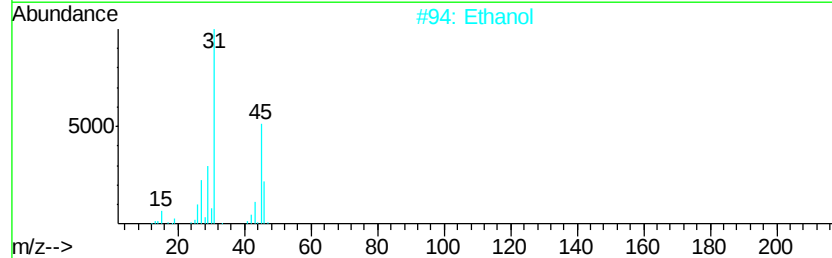
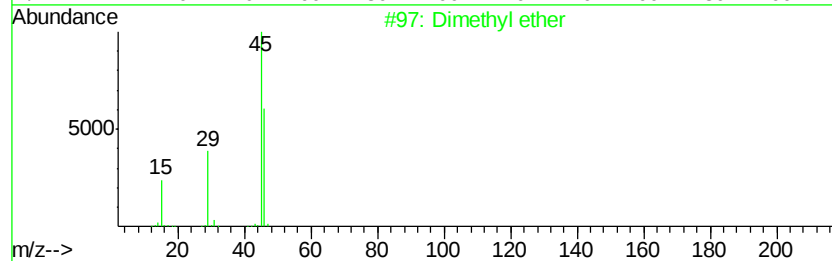
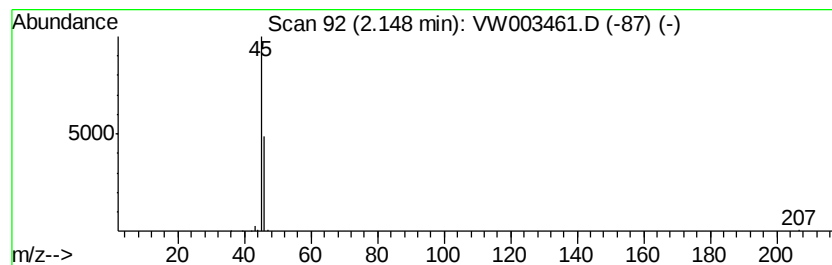
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Dimethyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	8.38 ug/l	102241	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		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	8.4	ug/l	102241	1	7.95	609736	50.0