

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW091718\
 Data File : VW005472.D
 Acq On : 18 Sep 2018 10:06
 Operator : SY/AP
 Sample : J4936-39
 Misc : 5.49G/5ML/MSVOA W/SOIL
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 26KV

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W091618S.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	1.959	5	9	17	rVB	23008	33727	12.25%	2.104%
2	2.172	39	44	48	rBV3	5945	9043	3.29%	0.564%
3	2.324	64	69	71	rVV2	10317	16935	6.15%	1.057%
4	2.349	71	73	78	rVB2	11489	15252	5.54%	0.952%
5	3.019	180	183	193	rVB6	3819	9539	3.47%	0.595%
6	3.318	227	232	236	rVB6	1534	3086	1.12%	0.193%
7	3.397	239	245	255	rVB4	5641	14466	5.26%	0.903%
8	4.690	447	457	461	rBV5	1416	3903	1.42%	0.244%
9	4.921	485	495	510	rBV3	14318	54130	19.67%	3.377%
10	5.946	652	663	674	rBV4	6047	18624	6.77%	1.162%
11	7.147	853	860	865	rBV2	4247	11286	4.10%	0.704%
12	7.890	971	982	987	rBV2	46187	115505	41.97%	7.207%
13	7.957	987	993	1003	rVB	119796	271768	98.74%	16.957%
14	8.171	1020	1028	1035	rVB4	2378	5689	2.07%	0.355%
15	8.311	1042	1051	1061	rBV	46726	106908	38.84%	6.671%
16	8.585	1090	1096	1102	rBV5	1360	2794	1.02%	0.174%
17	8.701	1105	1115	1121	rBV5	4263	9294	3.38%	0.580%
18	8.848	1128	1139	1151	rBV	137255	275225	100.00%	17.173%
19	9.341	1214	1220	1228	rVB7	3570	8122	2.95%	0.507%
20	9.896	1303	1311	1316	rBV6	2160	5057	1.84%	0.316%
21	9.963	1316	1322	1326	rBV5	2070	4471	1.62%	0.279%
22	10.110	1339	1346	1348	rBV3	1639	2992	1.09%	0.187%
23	10.329	1375	1382	1389	rVV	116983	209147	75.99%	13.050%
24	10.390	1390	1392	1396	rVV3	2679	4002	1.45%	0.250%
25	10.512	1405	1412	1417	rBV3	5888	10531	3.83%	0.657%
26	10.713	1440	1445	1451	rBV3	10035	21273	7.73%	1.327%
27	10.865	1465	1470	1475	rBV6	3735	7349	2.67%	0.459%
28	11.067	1498	1503	1511	rBV9	1634	3938	1.43%	0.246%
29	11.182	1516	1522	1527	rBV7	1739	3844	1.40%	0.240%
30	11.243	1527	1532	1537	rVB8	1647	3129	1.14%	0.195%
31	11.347	1543	1549	1554	rBV10	1604	3884	1.41%	0.242%
32	11.512	1573	1576	1581	rBV5	1711	2903	1.05%	0.181%
33	11.634	1590	1596	1608	rBV	80253	144022	52.33%	8.986%
34	11.847	1627	1631	1634	rVV6	3925	5311	1.93%	0.331%

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35	12.152	1677	1681	1687	rVB8	1750	2961	1.08%	0.185%
36	12.621	1751	1758	1767	rBV2	48778	92109	33.47%	5.747%
37	13.566	1908	1913	1919	rBV	38179	65015	23.62%	4.057%
38	14.084	1994	1998	2003	rBV2	14646	25434	9.24%	1.587%

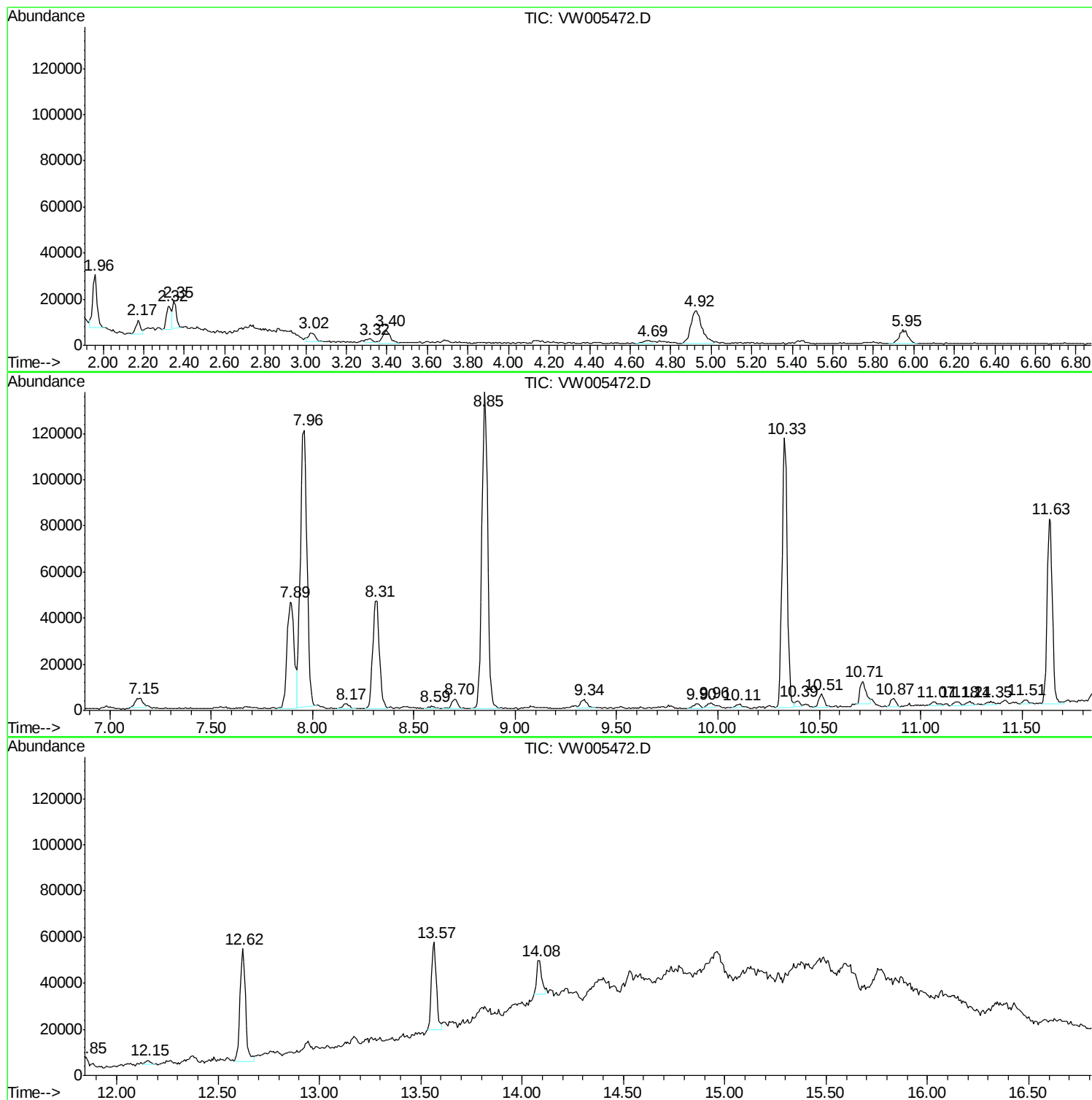
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W091618S.M
Quant Title : SW846 8260

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



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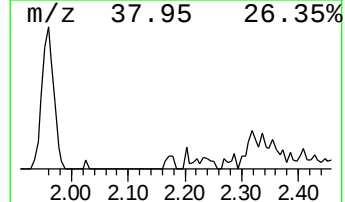
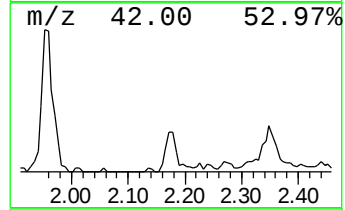
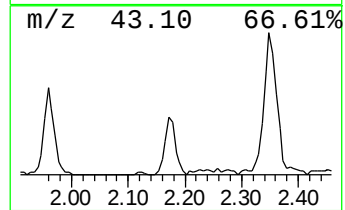
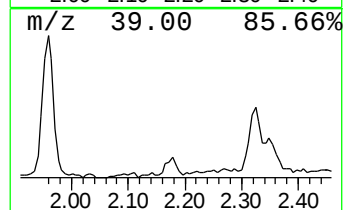
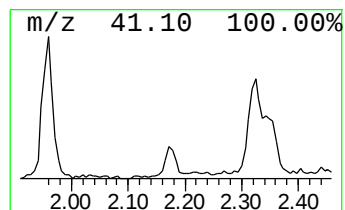
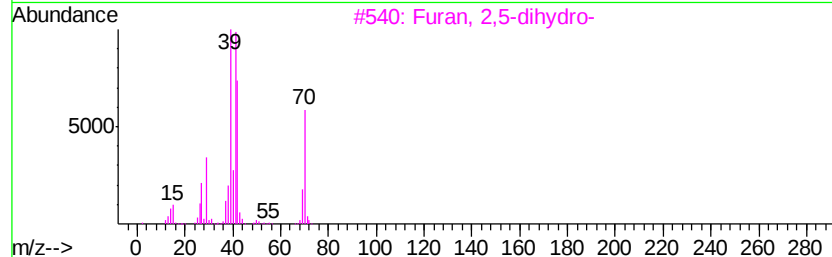
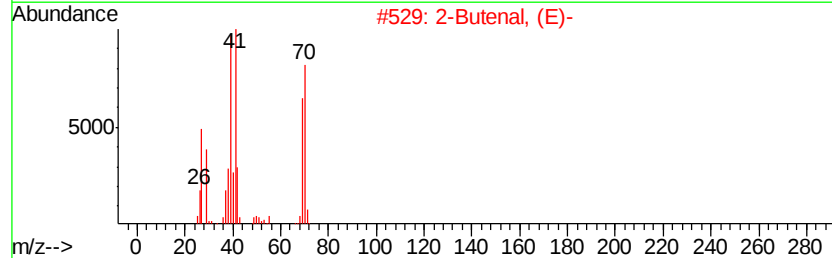
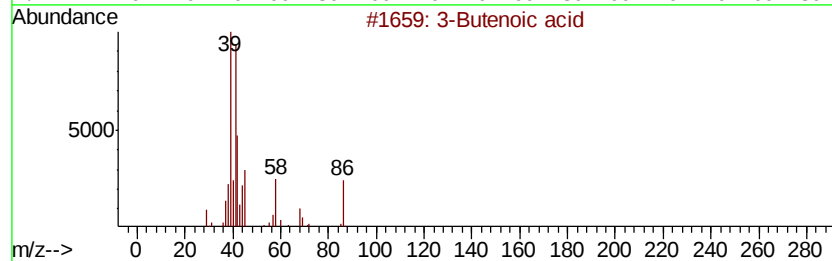
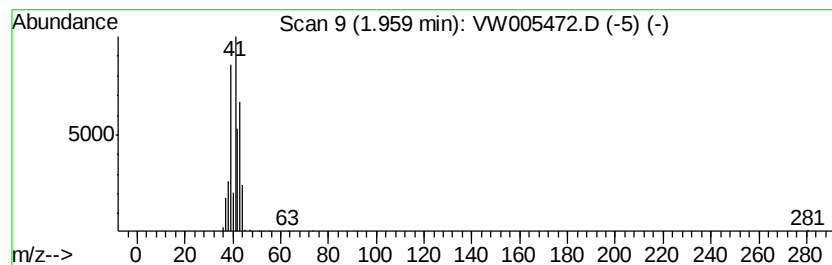
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Peak Number 1 3-Butenoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	6.21 ug/l	33727	Pentafluorobenzene	7.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Butenoic acid	86	C4H6O2	000625-38-7	83
2			2-Butenal, (E)-	70	C4H6O	000123-73-9	78
3			Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	50
4			Bis(3-methylbutyl) fluorene-2,7-...	466	C23H30O6S2	253664-95-8	43
5			Propene	42	C3H6	000115-07-1	40



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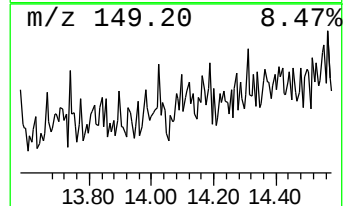
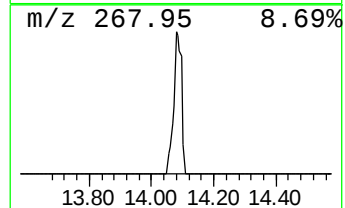
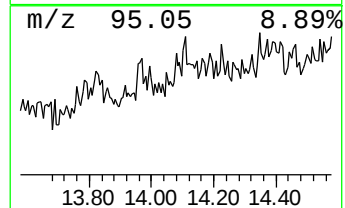
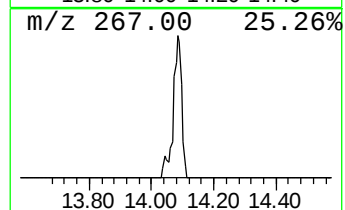
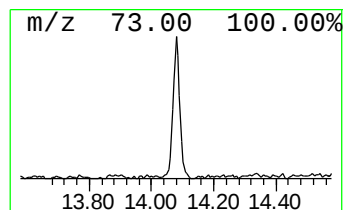
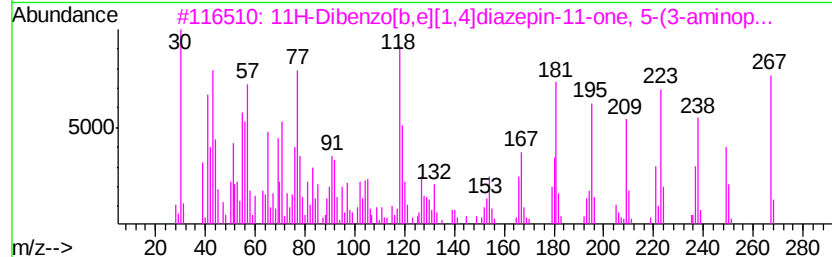
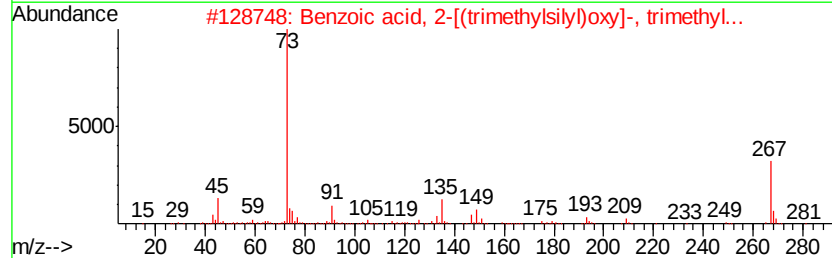
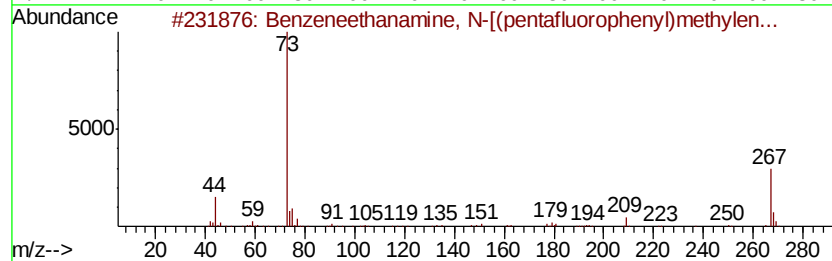
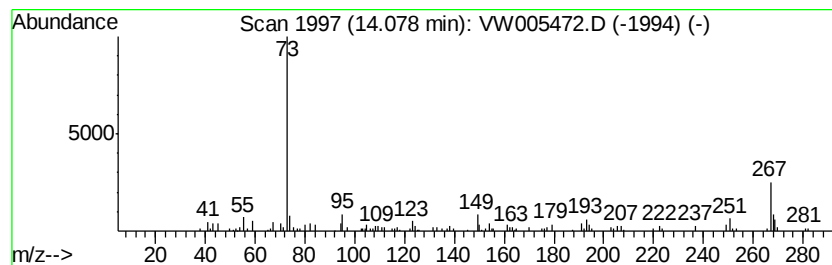
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Peak Number 4 unknown14.08 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.08	19.56 ug/l	25434	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzeneethanamine, N-[(pentafluo...	475 C21H26F5N02Si2	055429-85-1 42
2		Benzoic acid, 2-[(trimethylsilyl...]	282 C13H22O3Si2	003789-85-3 40
3		11H-Dibenzo[b,e][1,4]diazepin-11...	267 C16H17N3O	013450-73-2 30
4		3,5-Dimethoxyphenylacetic acid, ...	268 C13H20O4Si	1000071-82-4 10
5		Monobenzylidene-d-glucose	268 C13H16O6	1000127-04-3 9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Butenoic acid	1.96	6.2	ug/l	33727	1	7.96	271768	50.0
unknown14.08	14.08	19.6	ug/l	25434	4	13.56	65015	50.0