

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU053020\
 Data File : VU038386.D
 Acq On : 29 May 2020 19:15
 Operator : JC/MD
 Sample : L2749-14
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BE4X8

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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_U\METHOD\SOMUTR050720WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.510	43	53	70	rBV	59544	186627	6.19%	1.290%
2	1.613	77	85	95	rBV	141455	175913	5.83%	1.216%
3	1.932	177	184	203	rVB	127031	173342	5.75%	1.199%
4	2.401	322	330	344	rVB	95251	145141	4.81%	1.004%
5	2.594	376	390	405	rBV	238488	388100	12.87%	2.684%
6	3.395	625	639	655	rVB5	15568	34857	1.16%	0.241%
7	4.034	821	838	858	rVB2	50655	133939	4.44%	0.926%
8	4.674	1019	1037	1067	rBV2	206252	613012	20.32%	4.239%
9	5.102	1151	1170	1189	rBV	243102	542981	18.00%	3.755%
10	5.761	1354	1375	1404	rBV2	445846	1176924	39.01%	8.138%
11	6.282	1521	1537	1555	rBV	364295	697935	23.14%	4.826%
12	6.722	1659	1674	1700	rBV	279513	558101	18.50%	3.859%
13	7.591	1931	1944	1961	rBV	142133	249139	8.26%	1.723%
14	7.922	2032	2047	2061	rBV	530137	955439	31.67%	6.607%
15	7.992	2061	2069	2083	rVB2	48056	89042	2.95%	0.616%
16	8.201	2120	2134	2149	rBV	91469	152355	5.05%	1.053%
17	8.655	2262	2275	2314	rBV	805725	1478871	49.02%	10.226%
18	9.436	2505	2518	2522	rBV	529096	842423	27.93%	5.825%
19	9.780	2614	2625	2638	rBV3	27888	47401	1.57%	0.328%
20	10.497	2838	2848	2860	rBV	50482	79394	2.63%	0.549%
21	10.770	2922	2933	2944	rBV	221095	358863	11.90%	2.481%
22	11.304	3087	3099	3117	rVB2	15611	31341	1.04%	0.217%
23	11.825	3248	3261	3281	rBV2	579722	1088004	36.07%	7.523%
24	12.028	3316	3324	3335	rVB2	21930	32192	1.07%	0.223%
25	12.105	3335	3348	3364	rVB	100356	159164	5.28%	1.101%
26	12.204	3364	3379	3397	rVB	600886	994789	32.98%	6.879%
27	16.966	4847	4860	4881	rBV	1913599	3016657	100.00%	20.859%
28	17.162	4910	4921	4932	rBV	31859	59874	1.98%	0.414%

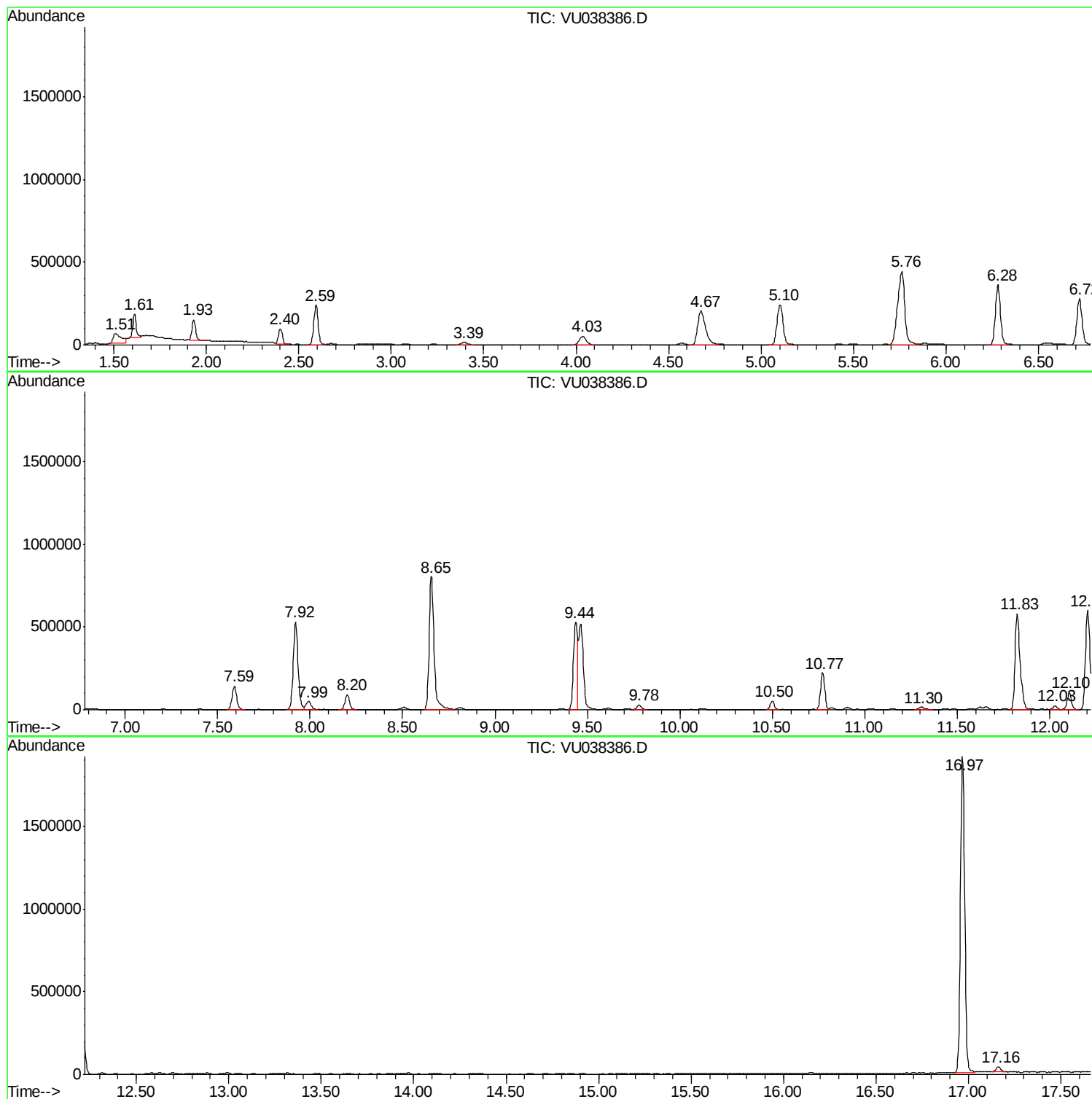
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ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4X8

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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



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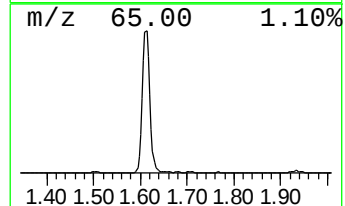
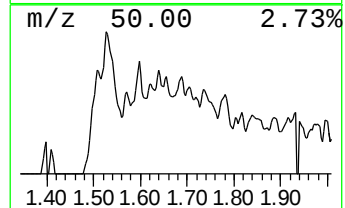
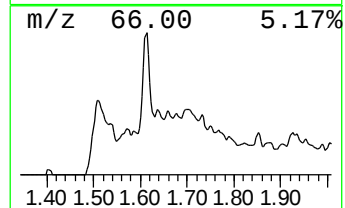
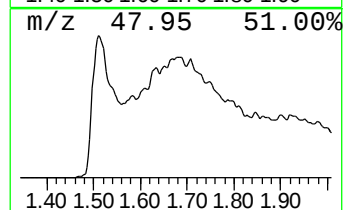
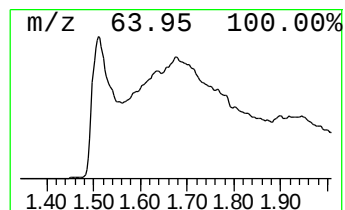
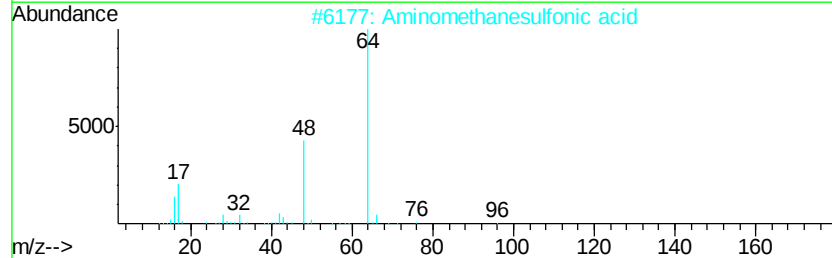
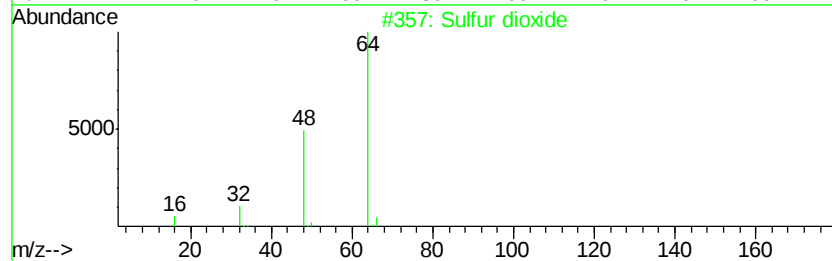
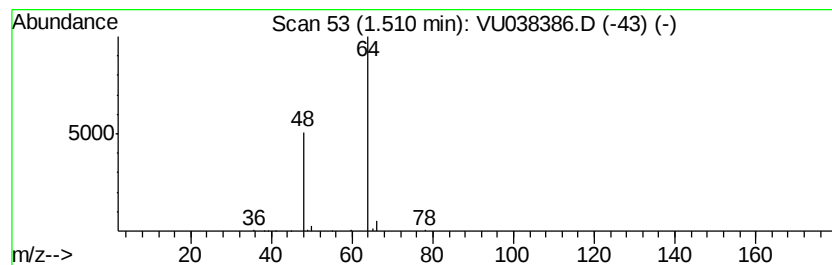
Instrument :
MSVOA_U
ClientSampleId :
BE4X8

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Sulfur dioxide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.51	1.34 ug/L	186627	1,4-Difluorobenzene	6.28
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Sulfur dioxide	64 02S	007446-09-5	83
2	Aminomethanesulfonic acid	111 CH5N03S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169 C3H7N05S	000498-40-8	9
4	Aminomethanesulfonic acid	111 CH5N03S	013881-91-9	9
5	Aminomethanesulfonic acid	111 CH5N03S	013881-91-9	9



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Instrument :
MSVOA_U
ClientSampled :
BE4X8

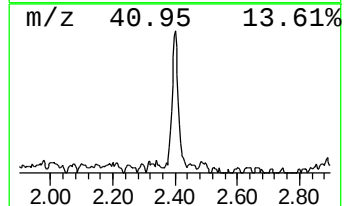
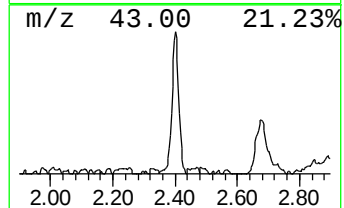
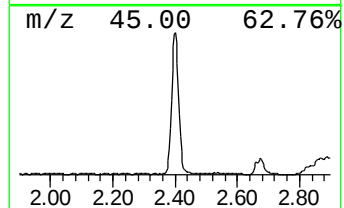
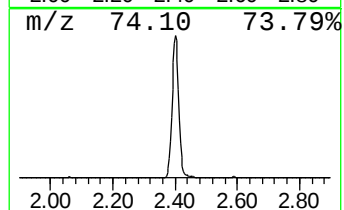
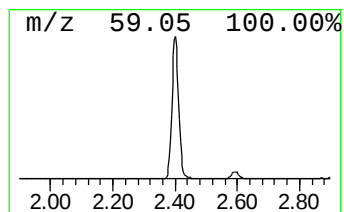
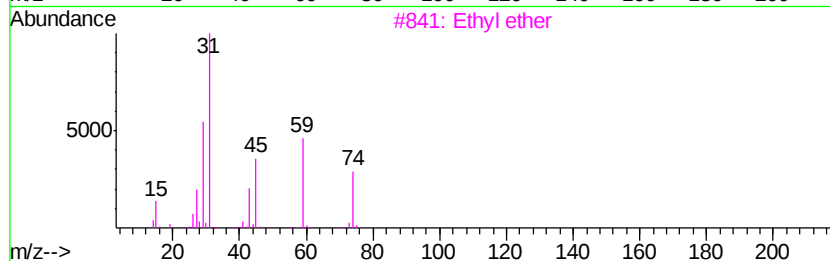
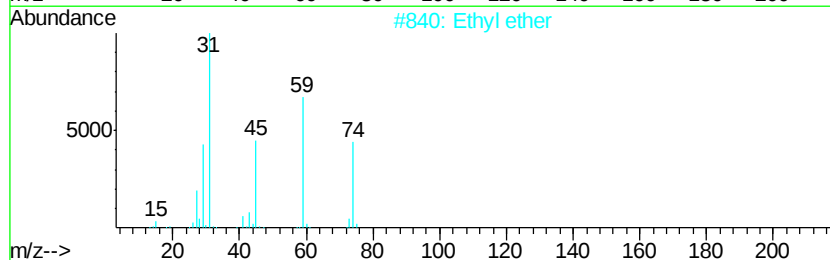
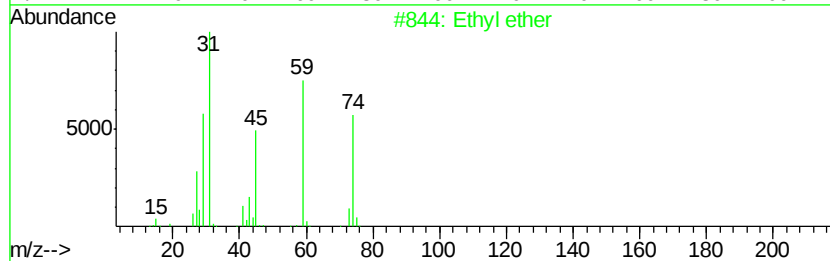
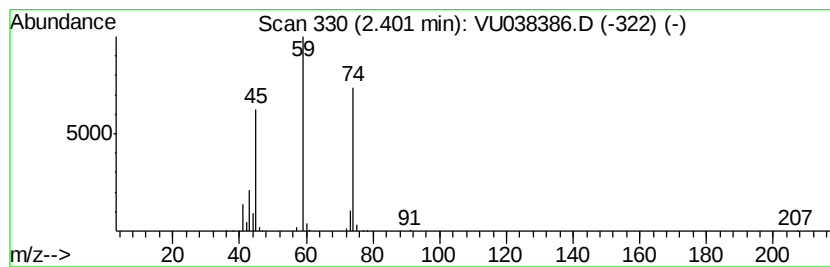
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Ethyl ether Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.40	1.04 ug/L	145141	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethyl ether	74	C4H10O	000060-29-7	90
2		Ethyl ether	74	C4H10O	000060-29-7	86
3		Ethyl ether	74	C4H10O	000060-29-7	78
4		Ethane, 1,2-diethoxy-	118	C6H14O2	000629-14-1	74
5		Ethane, 1,2-diethoxy-	118	C6H14O2	000629-14-1	74



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ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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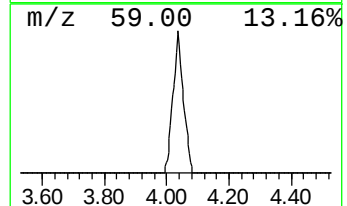
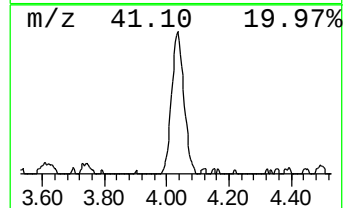
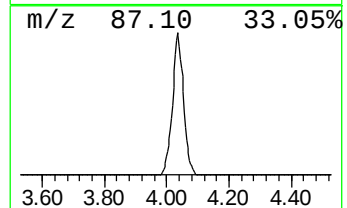
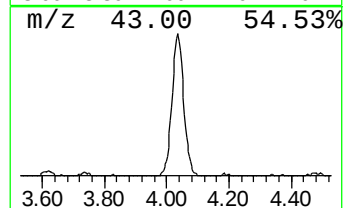
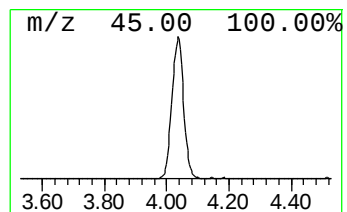
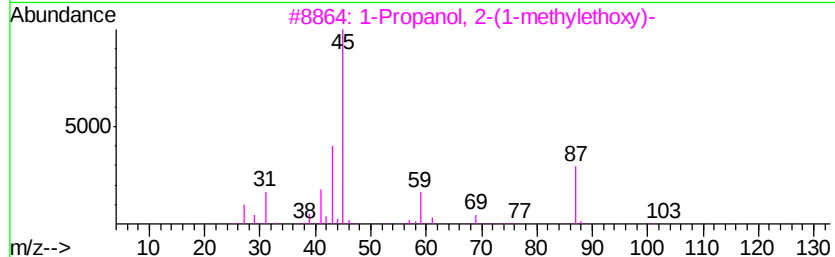
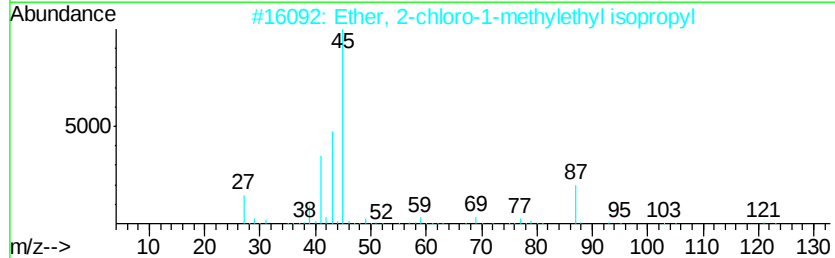
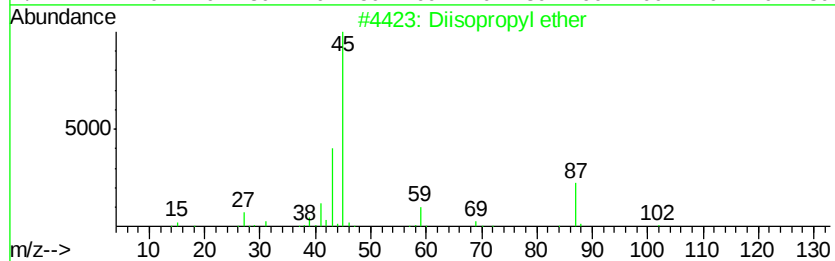
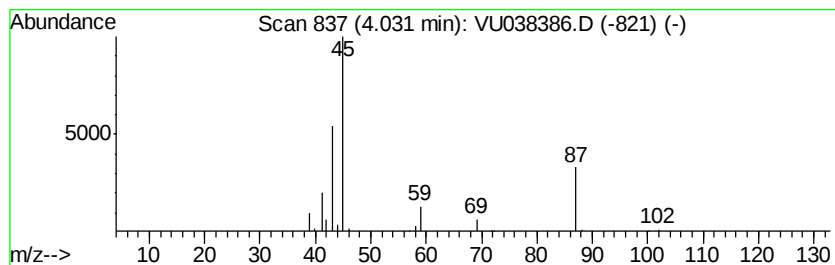
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Diisopropyl ether Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.03	0.96 ug/L	133939	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diisopropyl ether	102	C6H14O	000108-20-3	83
2		Ether, 2-chloro-1-methylethyl is...	136	C6H13ClO	098277-76-0	78
3		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	78
4		Diisopropyl ether	102	C6H14O	000108-20-3	72
5		Diisopropyl ether	102	C6H14O	000108-20-3	64



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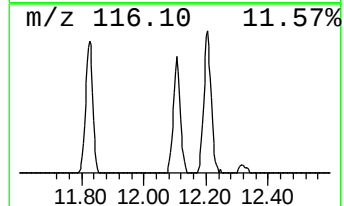
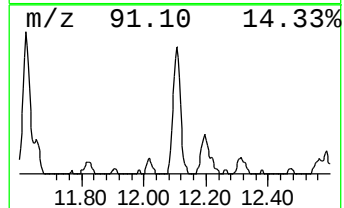
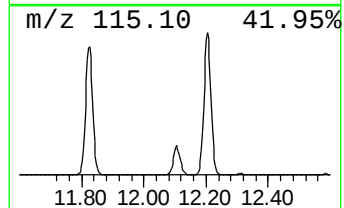
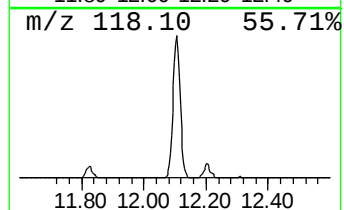
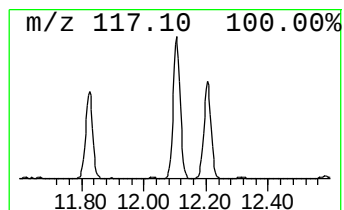
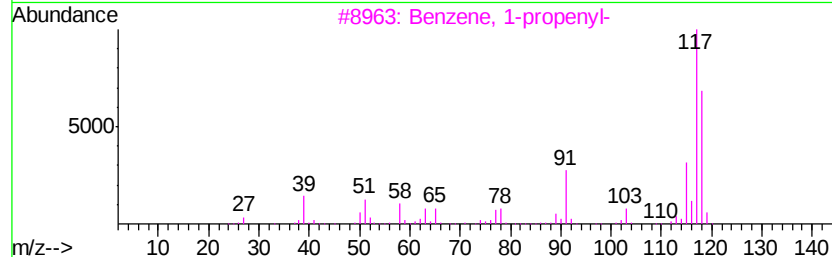
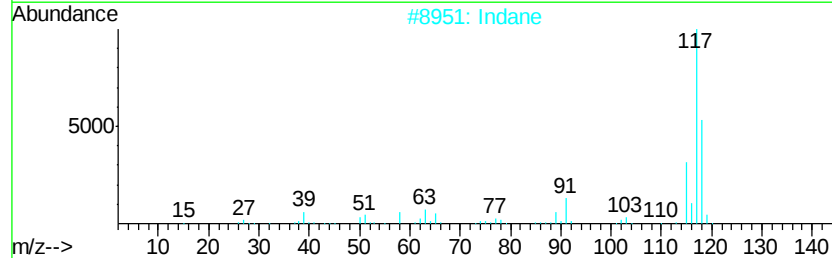
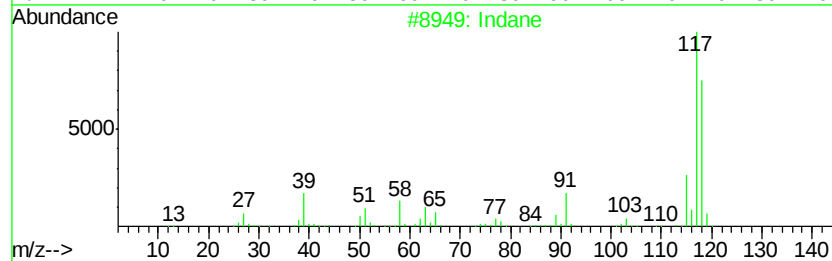
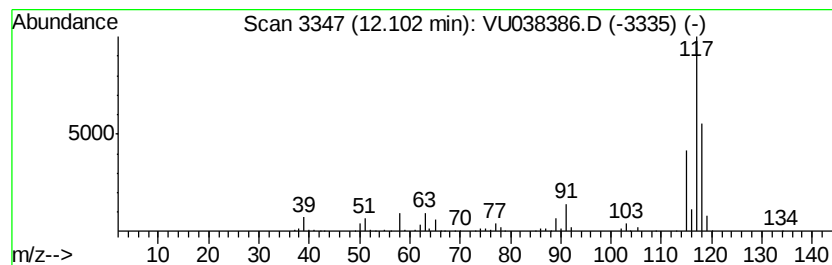
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	0.73 ug/L	159164	1,4-Dichlorobenzene-d4	11.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	96
2	Indane		118	C9H10	000496-11-7	94
3	Benzene, 1-propenyl-		118	C9H10	000637-50-3	76
4	Indane		118	C9H10	000496-11-7	76
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-	...	118	C9H10	1000191-13-7	64



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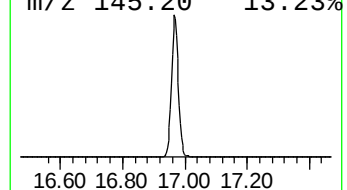
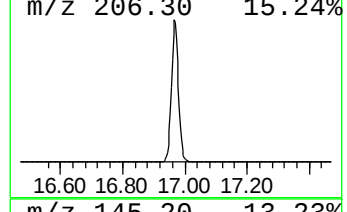
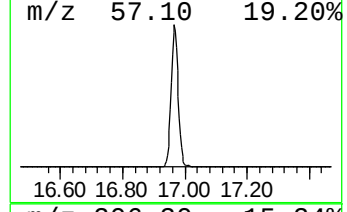
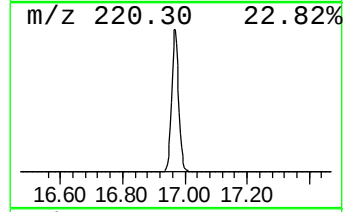
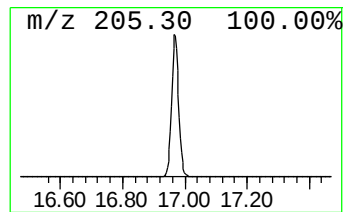
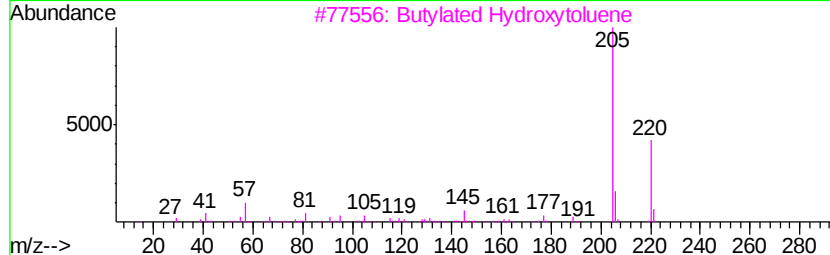
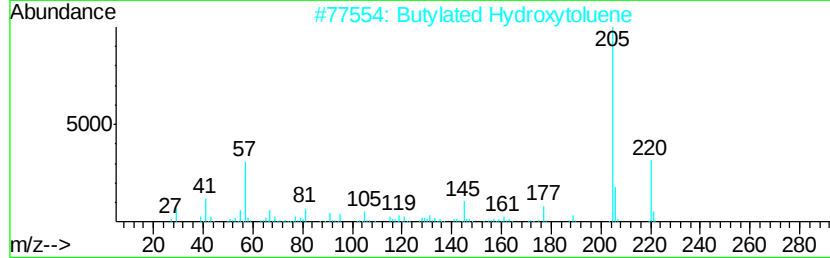
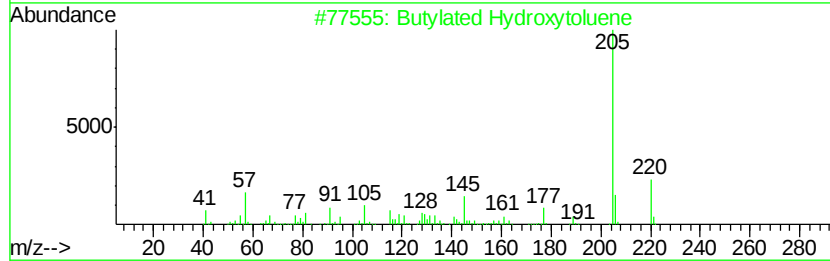
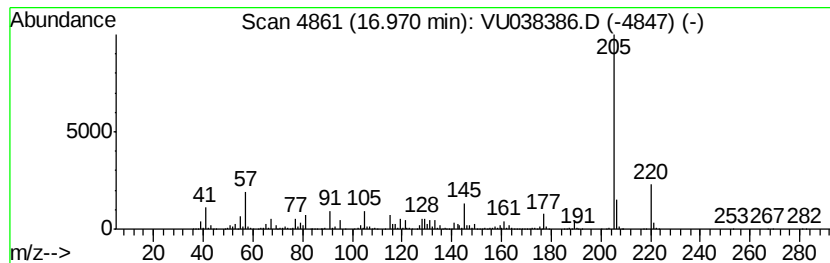
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 Butylated Hydroxytoluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.97	13.86 ug/L	3016660	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	99
2		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	96
3		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	94
4		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	93
5		Phenol, 2,4,6-tris(1-methylethyl)-	220	C15H24O	002934-07-8	91



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

TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Sulfur dioxide	1.51	1.3	ug/L	186627	1	6.28	697935	5.0
Ethyl ether	2.40	1.0	ug/L	145141	1	6.28	697935	5.0
Diisopropyl ether	4.03	1.0	ug/L	133939	1	6.28	697935	5.0
Indane	12.10	0.7	ug/L	159164	3	11.83	1088000	5.0
Butylated Hydroxy...	16.97	13.9	ug/L	3016660	3	11.83	1088000	5.0