

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV091318\
 Data File : VV007525.D
 Acq On : 13 Sep 2018 22:25
 Operator : SY/MD
 Sample : J4864-04
 Misc : 25.0 mL/MSVOA_V/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0D10

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_V\METHOD\SOMVTR091318WMA.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	1.312	60	69	86	rVB3	83925	106560	5.05%	0.821%
2	1.422	97	103	113	rVB	44596	49416	2.34%	0.381%
3	1.576	145	151	165	rVB	45413	52630	2.49%	0.405%
4	2.129	314	323	338	rBV3	112075	172859	8.19%	1.332%
5	2.544	444	452	465	rVB	49263	79969	3.79%	0.616%
6	3.978	883	898	912	rBV	284501	683677	32.39%	5.267%
7	4.409	1021	1032	1048	rVB	73227	164226	7.78%	1.265%
8	4.656	1098	1109	1124	rVB	58088	132231	6.26%	1.019%
9	5.093	1228	1245	1267	rVB2	153646	332047	15.73%	2.558%
10	5.669	1410	1424	1442	rBV	154215	303902	14.40%	2.341%
11	5.965	1498	1516	1539	rBV	1096715	2110978	100.00%	16.263%
12	6.061	1539	1546	1557	rVV	29308	48762	2.31%	0.376%
13	6.132	1558	1568	1586	rVB	79248	163388	7.74%	1.259%
14	7.055	1843	1855	1870	rBV	45489	82345	3.90%	0.634%
15	7.380	1944	1956	1970	rBV	186994	320532	15.18%	2.469%
16	7.685	2041	2051	2070	rVB	28562	49719	2.36%	0.383%
17	8.161	2182	2199	2228	rBV	173289	346009	16.39%	2.666%
18	8.907	2417	2431	2448	rBV	219504	360536	17.08%	2.778%
19	9.068	2471	2481	2490	rBV	18589	28164	1.33%	0.217%
20	9.190	2507	2519	2538	rVB	66413	106474	5.04%	0.820%
21	9.598	2637	2646	2657	rVB	21838	35158	1.67%	0.271%
22	10.193	2816	2831	2836	rBV4	15924	28043	1.33%	0.216%
23	10.235	2836	2844	2850	rVV	61452	105173	4.98%	0.810%
24	10.270	2850	2855	2870	rVB	66210	103659	4.91%	0.799%
25	10.662	2967	2977	2992	rVB7	13070	33158	1.57%	0.255%
26	11.293	3161	3173	3194	rVB3	1139376	1799026	85.22%	13.860%
27	11.582	3252	3263	3276	rVV3	30571	53311	2.53%	0.411%
28	11.678	3280	3293	3313	rVB	203096	335667	15.90%	2.586%
29	11.794	3319	3329	3339	rBV2	17572	26416	1.25%	0.204%
30	12.167	3431	3445	3452	rBV	20236	33641	1.59%	0.259%
31	12.350	3495	3502	3513	rVB2	19095	32224	1.53%	0.248%
32	12.437	3519	3529	3533	rBV3	33450	53198	2.52%	0.410%
33	12.463	3533	3537	3547	rVV	37339	53894	2.55%	0.415%
34	12.524	3547	3556	3566	rVB7	11491	23786	1.13%	0.183%

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 Title : TRACE VOA SOM01.0

35	12.601	3571	3580	3592	rVB3	24850	36926	1.75%	0.284%
36	12.736	3603	3622	3631	rBV3	41736	90021	4.26%	0.694%
37	12.810	3640	3645	3653	rVB	17238	22142	1.05%	0.171%
38	12.936	3677	3684	3696	rVB5	17391	25289	1.20%	0.195%
39	13.010	3696	3707	3716	rVB	75891	119060	5.64%	0.917%
40	13.090	3716	3732	3736	rBV9	11137	27628	1.31%	0.213%
41	13.219	3758	3772	3781	rBV	104396	186289	8.82%	1.435%
42	13.267	3781	3787	3795	rVV2	32463	52767	2.50%	0.407%
43	13.321	3796	3804	3810	rVV2	49969	73861	3.50%	0.569%
44	13.360	3811	3816	3819	rVV2	34253	42415	2.01%	0.327%
45	13.389	3820	3825	3858	rVB2	70723	164460	7.79%	1.267%
46	13.553	3859	3876	3885	rBV	73942	155541	7.37%	1.198%
47	13.662	3902	3910	3923	rVB	155184	241746	11.45%	1.862%
48	13.759	3924	3940	3952	rBV2	163120	285866	13.54%	2.202%
49	13.826	3953	3961	3971	rVB5	23988	41067	1.95%	0.316%
50	13.961	3992	4003	4018	rBV5	26521	59999	2.84%	0.462%
51	14.067	4028	4036	4042	rBV8	14151	26518	1.26%	0.204%
52	14.141	4050	4059	4073	rBV2	98993	219604	10.40%	1.692%
53	14.283	4094	4103	4112	rBV2	74312	116040	5.50%	0.894%
54	14.350	4114	4124	4136	rVB5	76613	155046	7.34%	1.194%
55	14.415	4136	4144	4157	rVB3	33529	55789	2.64%	0.430%
56	14.501	4160	4171	4180	rBV2	67266	124219	5.88%	0.957%
57	14.627	4203	4210	4214	rBV4	18938	26842	1.27%	0.207%
58	14.669	4215	4223	4238	rVV4	108785	255349	12.10%	1.967%
59	14.743	4238	4246	4259	rVB3	96336	178909	8.48%	1.378%
60	14.820	4259	4270	4276	rBV3	34518	60010	2.84%	0.462%
61	14.874	4279	4287	4298	rVB7	39738	78430	3.72%	0.604%
62	14.936	4298	4306	4311	rBV2	40555	68771	3.26%	0.530%
63	15.061	4335	4345	4355	rBV5	24165	50364	2.39%	0.388%
64	15.228	4384	4397	4407	rBV9	23948	59088	2.80%	0.455%
65	15.395	4434	4449	4464	rVB2	97109	247726	11.74%	1.908%
66	15.476	4465	4474	4489	rBV6	46393	92010	4.36%	0.709%
67	15.617	4509	4518	4519	rBV6	34067	39014	1.85%	0.301%
68	15.643	4521	4526	4536	rVB2	65465	99659	4.72%	0.768%
69	15.730	4539	4553	4561	rBV10	39814	86737	4.11%	0.668%
70	15.871	4589	4597	4603	rBV7	42510	67751	3.21%	0.522%
71	15.977	4626	4630	4643	rVB4	12936	21184	1.00%	0.163%

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72	16.064	4650	4657	4663	rBV3	48639	74007	3.51%	0.570%
73	16.103	4663	4669	4685	rVB4	59715	108596	5.14%	0.837%
74	16.238	4705	4711	4720	rVV2	27187	38989	1.85%	0.300%
75	16.286	4721	4726	4734	rVB5	19434	27031	1.28%	0.208%
76	16.408	4758	4764	4777	rVB7	23495	44865	2.13%	0.346%
77	16.498	4785	4792	4795	rBV2	22785	29613	1.40%	0.228%
78	16.540	4797	4805	4816	rVB3	72603	136686	6.48%	1.053%
79	16.601	4818	4824	4830	rBV2	16691	23108	1.09%	0.178%
80	16.842	4892	4899	4909	rVB2	32329	46639	2.21%	0.359%
81	16.971	4932	4939	4950	rVB	86610	130532	6.18%	1.006%
82	17.135	4984	4990	4996	rBV3	19439	31385	1.49%	0.242%
83	17.312	5041	5045	5056	rVB4	23518	35498	1.68%	0.273%
84	17.386	5061	5068	5078	rVV3	23582	39205	1.86%	0.302%
85	17.450	5082	5088	5098	rVB2	31255	49176	2.33%	0.379%

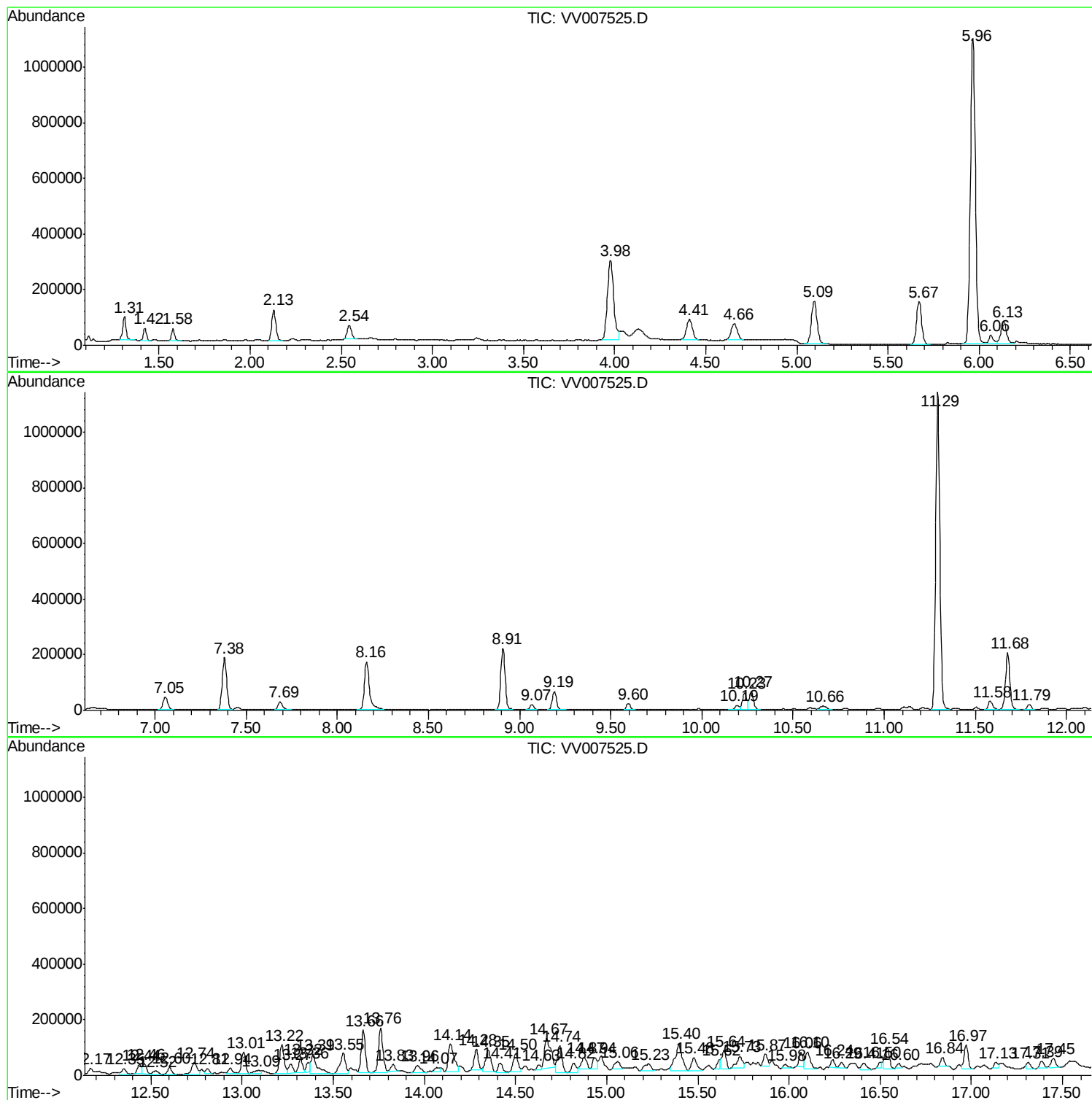
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR091318WMA.M
Quant Title : TRACE VOA SOM01.0

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



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Instrument :
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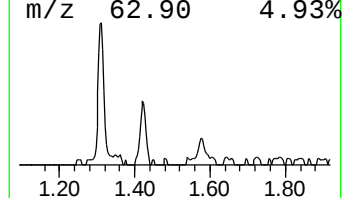
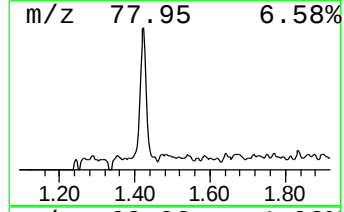
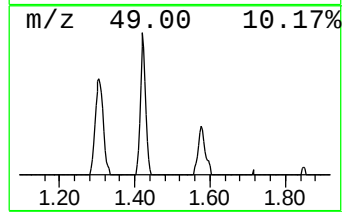
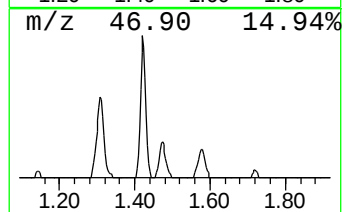
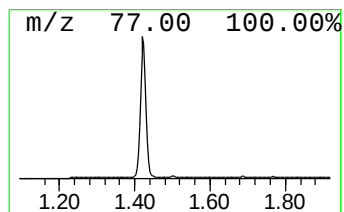
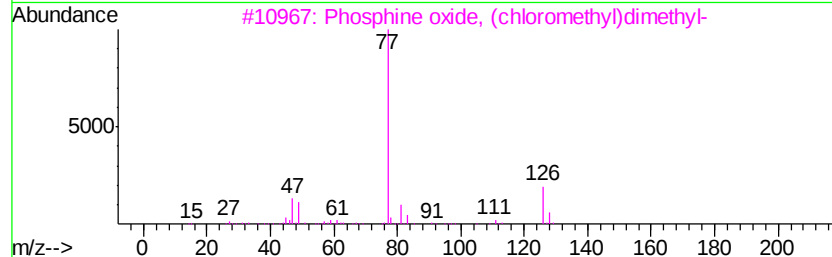
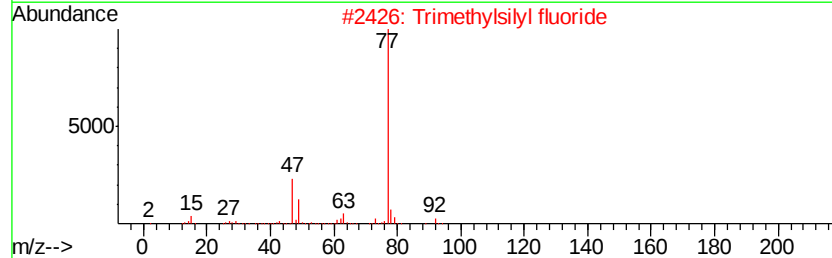
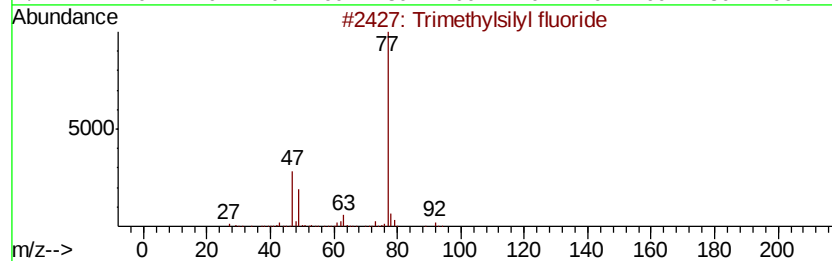
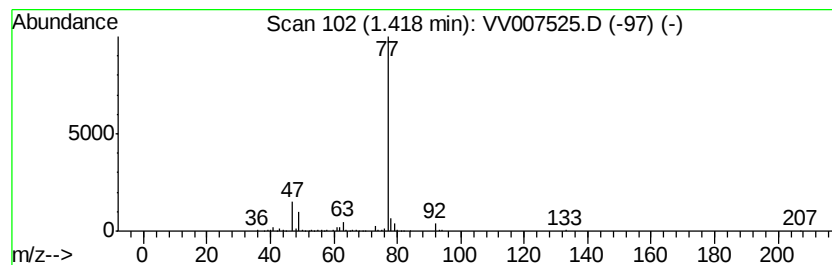
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR091318WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Trimethylsilyl fluoride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.42	0.81 ug/L	49416	1,4-Difluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trimethylsilyl fluoride	92	C3H9FSi	000420-56-4	91
2			Trimethylsilyl fluoride	92	C3H9FSi	000420-56-4	91
3			Phosphine oxide, (chloromethyl)d...	126	C3H8ClOP	001638-75-1	39
4			Propane, 1,2-dichloro-2-methyl-	126	C4H8Cl2	000594-37-6	9
5			Propane, 1,2-dichloro-2-methyl-	126	C4H8Cl2	000594-37-6	9



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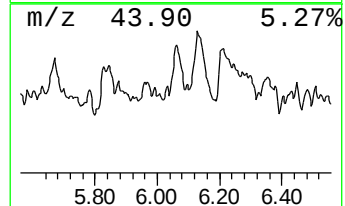
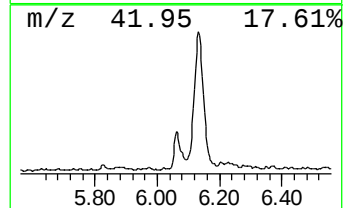
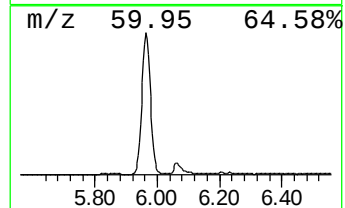
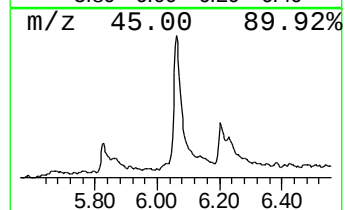
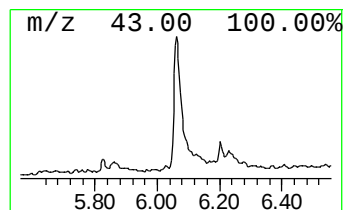
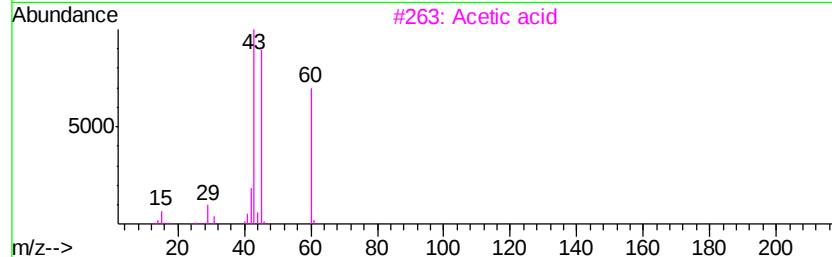
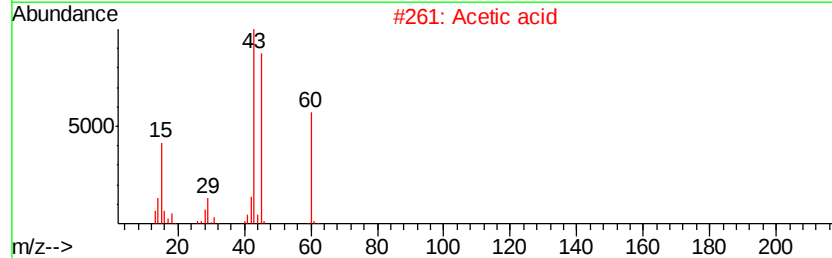
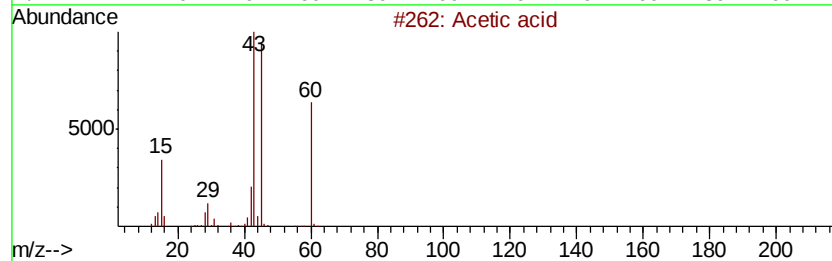
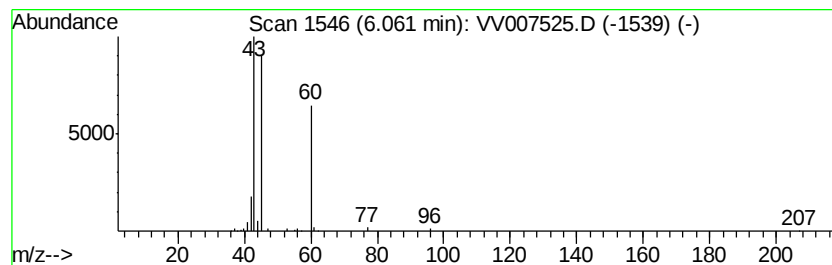
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Acetic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.06	0.80 ug/L	48762	1,4-Difluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetic acid	60	C2H4O2	000064-19-7	90
2		Acetic acid	60	C2H4O2	000064-19-7	90
3		Acetic acid	60	C2H4O2	000064-19-7	90
4		Acetic acid	60	C2H4O2	000064-19-7	78
5		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9



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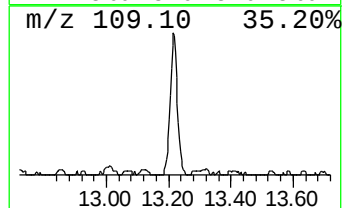
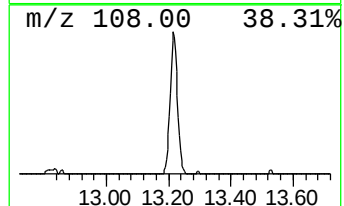
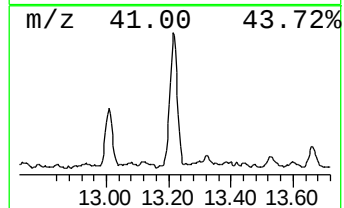
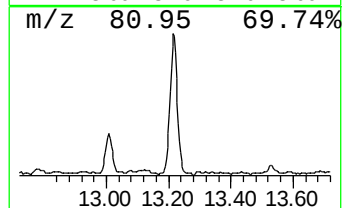
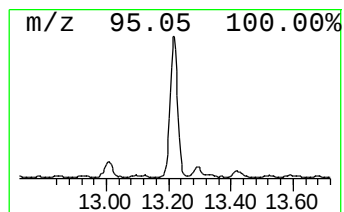
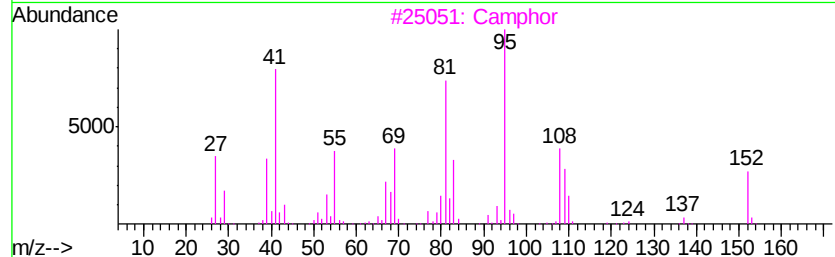
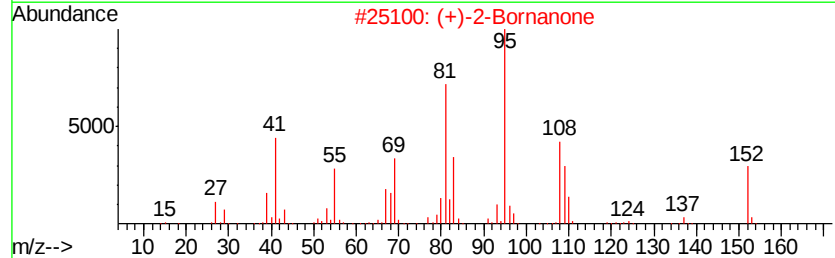
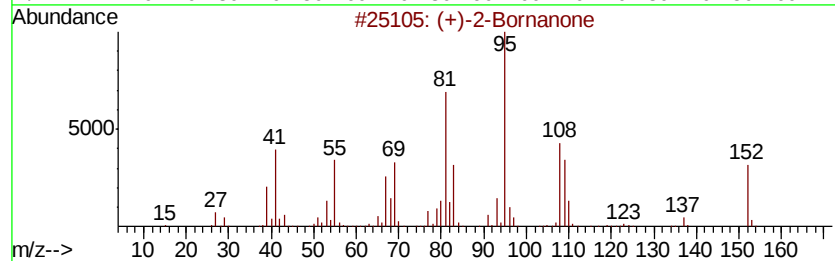
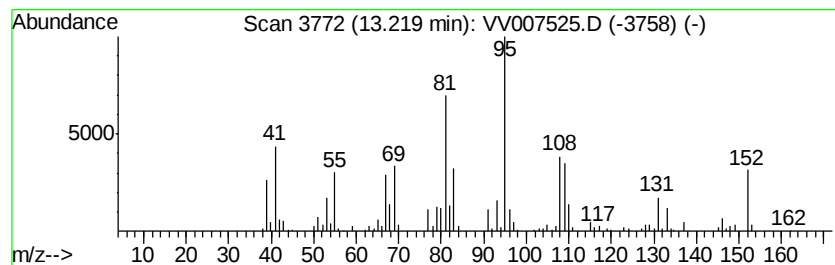
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (+)-2-Bornanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.22	0.52 ug/L	186289	1,4-Dichlorobenzene-d4	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
3	Camphor	152	C10H16O	000076-22-2	96
4	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96
5	Camphor	152	C10H16O	000076-22-2	93



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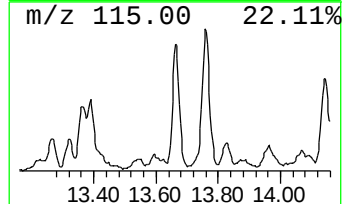
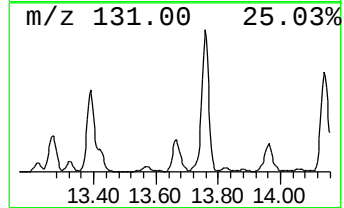
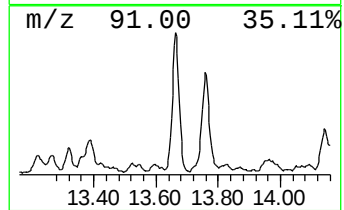
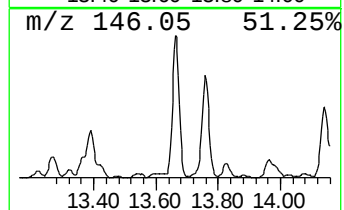
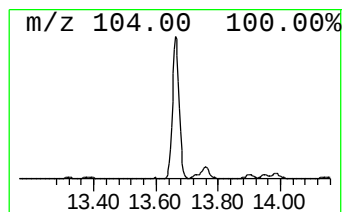
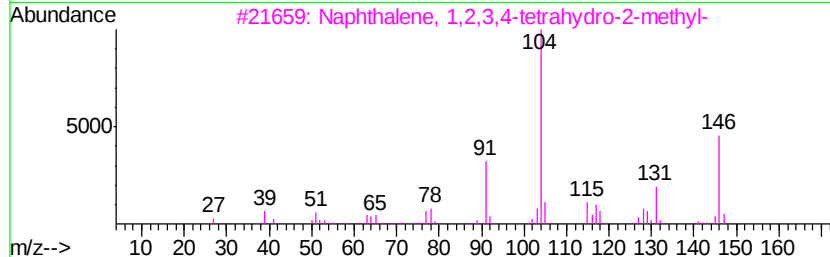
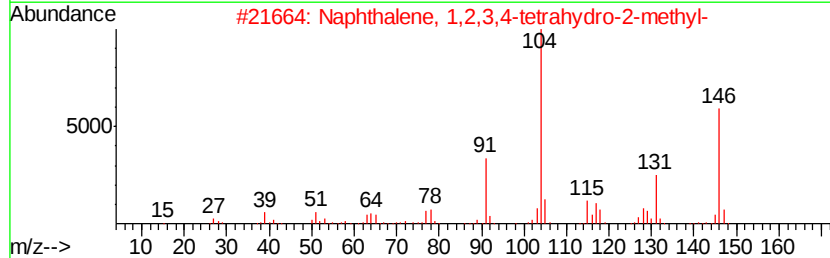
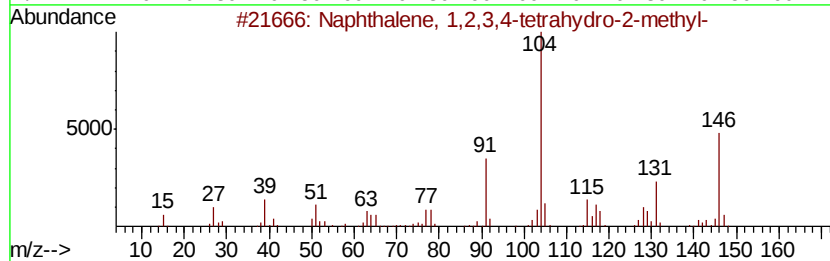
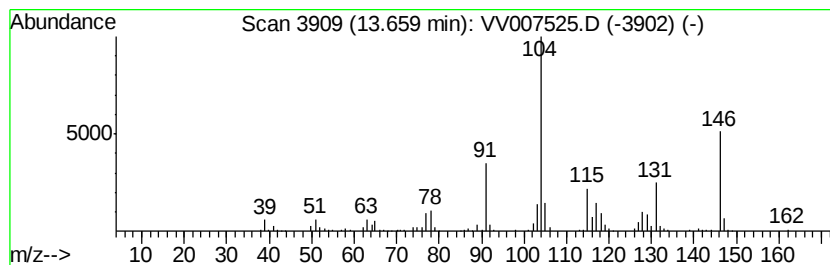
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	0.67 ug/L	241746	1,4-Dichlorobenzene-d4	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	93
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	90
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	90
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	81
5		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV091318\
Data File : VV007525.D
Acq On : 13 Sep 2018 22:25
Operator : SY/MD
Sample : J4864-04
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

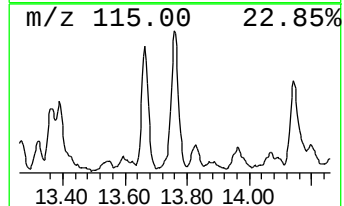
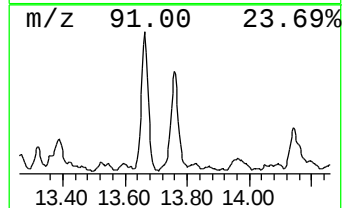
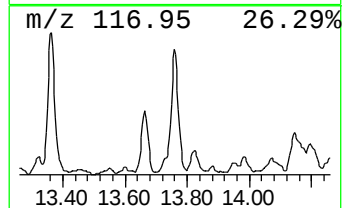
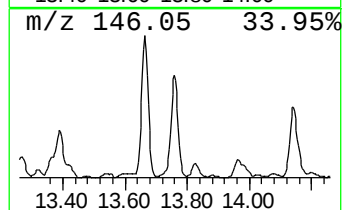
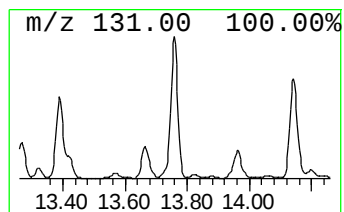
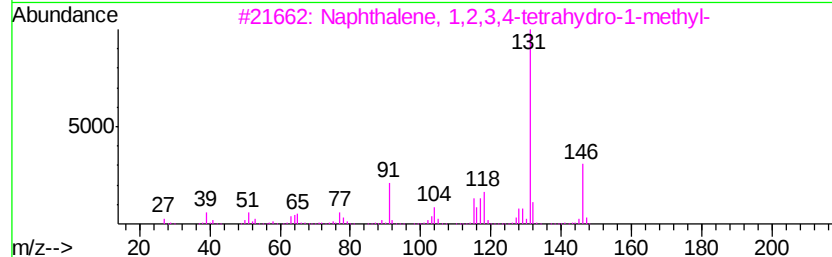
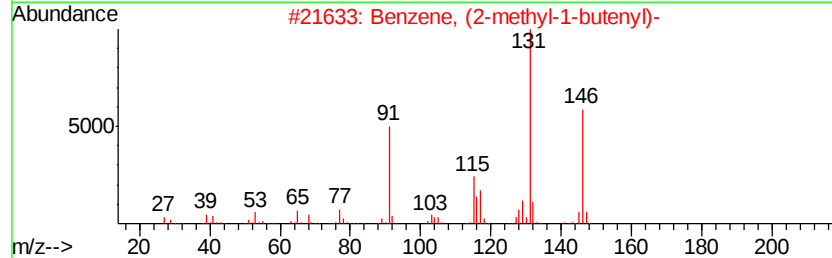
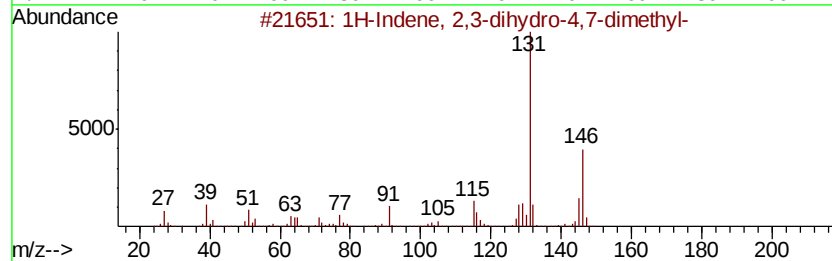
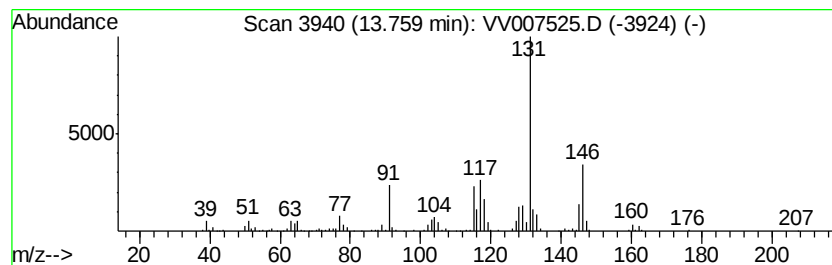
Instrument :
MSVOA_V
ClientSampleId :
C0D10

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR091318WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	0.79 ug/L	285866	1,4-Dichlorobenzene-d4	11.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	94
2	Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6	94
3	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	93
4	Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3	90
5	1H-Inden-1-one, 2,3-dihydro-2-me...	146 C10H10O	017496-14-9	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV091318\
Data File : VV007525.D
Acq On : 13 Sep 2018 22:25
Operator : SY/MD
Sample : J4864-04
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0D10

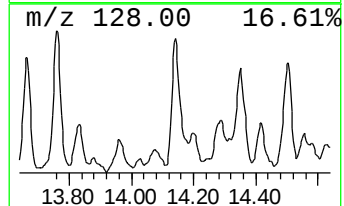
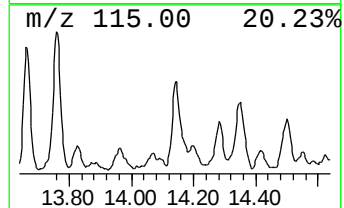
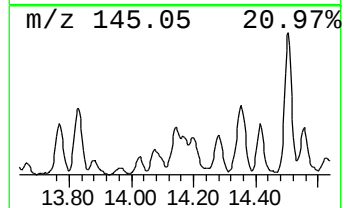
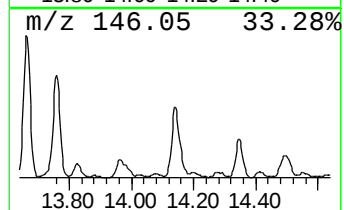
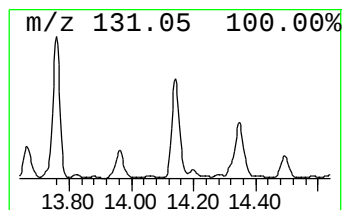
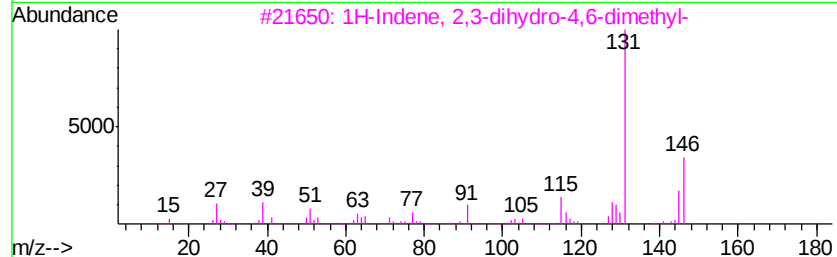
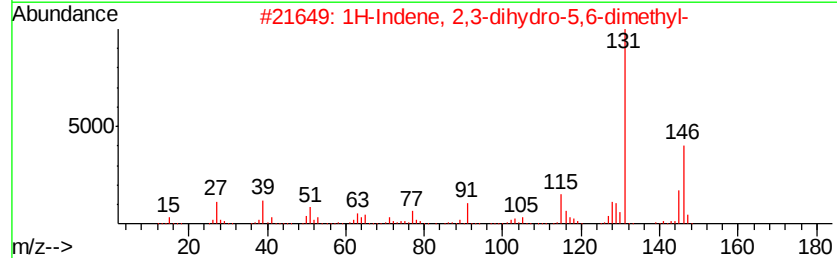
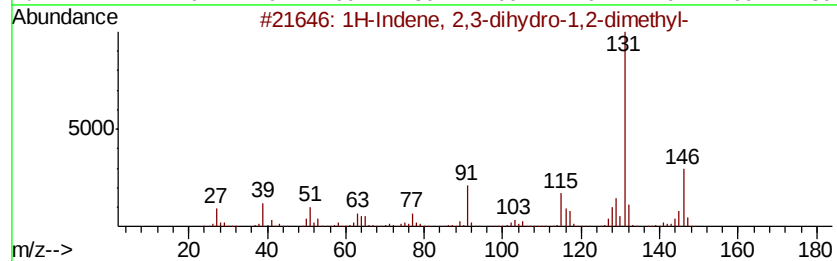
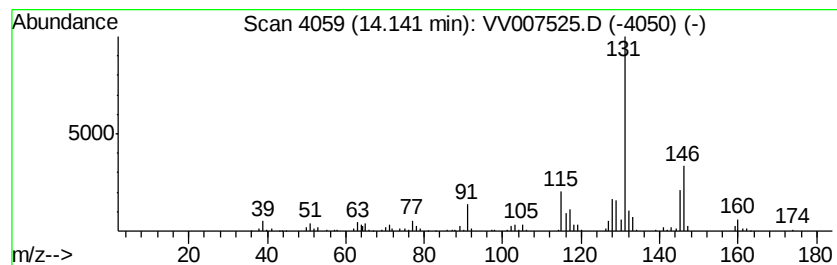
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	0.61 ug/L	219604	1,4-Dichlorobenzene-d4	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14		017057-82-8	90
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14		001075-22-5	90
3	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14		001685-82-1	87
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	87
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV091318\
Data File : VV007525.D
Acq On : 13 Sep 2018 22:25
Operator : SY/MD
Sample : J4864-04
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

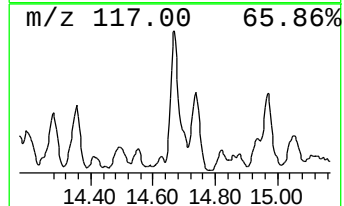
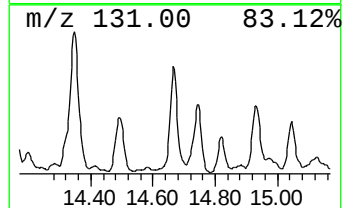
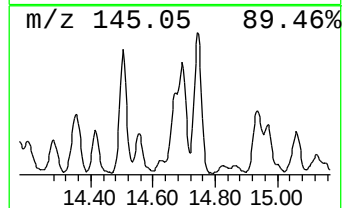
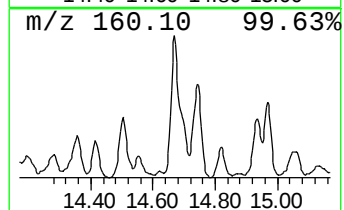
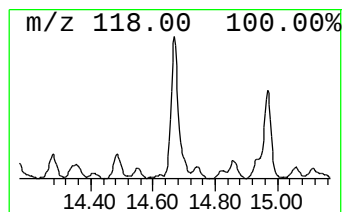
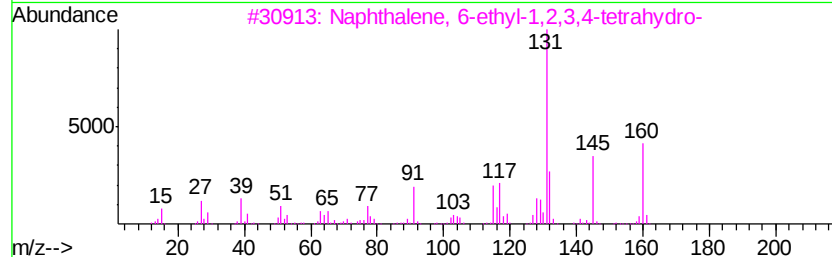
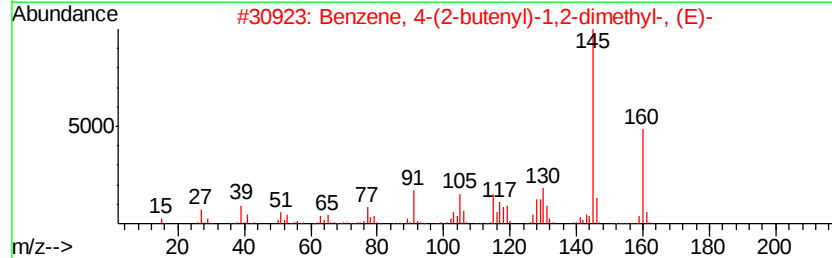
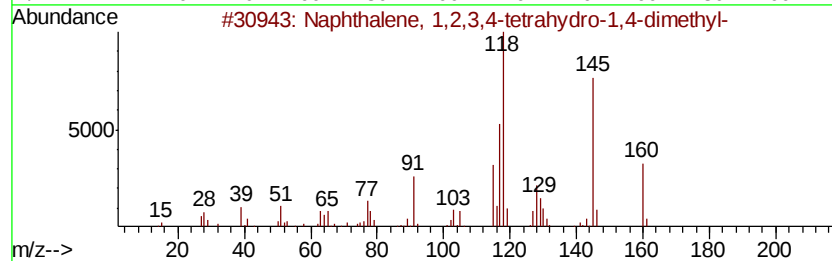
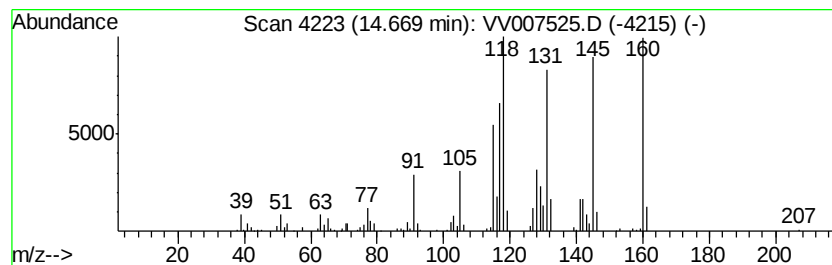
Instrument :
MSVOA_V
ClientSampleId :
C0D10

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR091318WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	0.71 ug/L	255349	1,4-Dichlorobenzene-d4	11.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6	76
2	Benzene, 4-(2-butenyl)-1,2-dimet...	160 C12H16	054340-86-2	60
3	Naphthalene, 6-ethyl-1,2,3,4-tet...	160 C12H16	022531-20-0	53
4	Naphthalene, 5-ethyl-1,2,3,4-tet...	160 C12H16	042775-75-7	49
5	Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	020027-77-4	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV091318\
Data File : VV007525.D
Acq On : 13 Sep 2018 22:25
Operator : SY/MD
Sample : J4864-04
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0D10

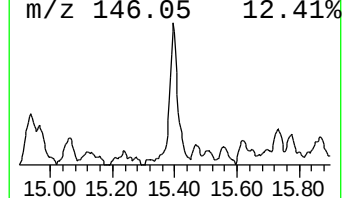
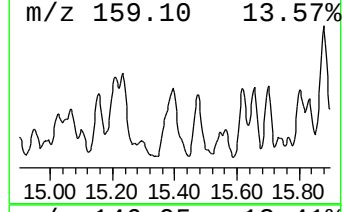
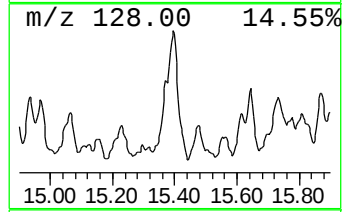
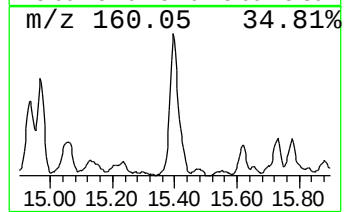
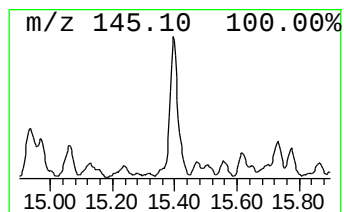
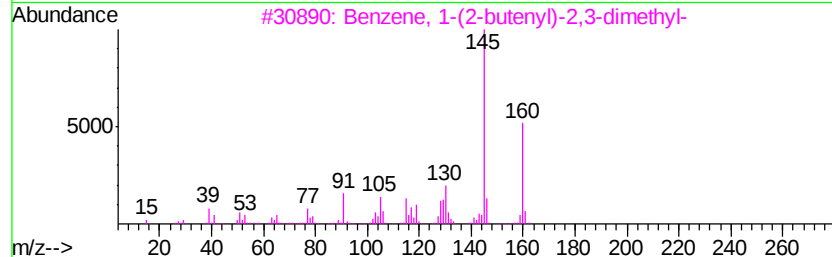
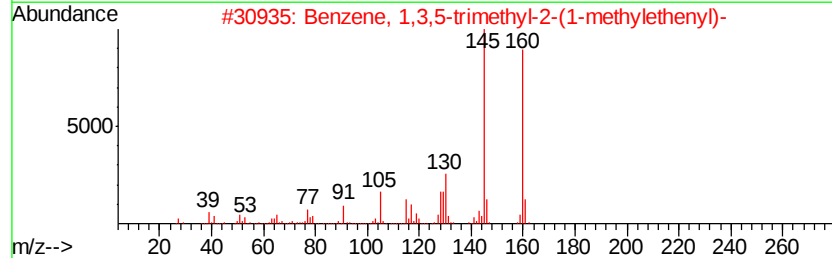
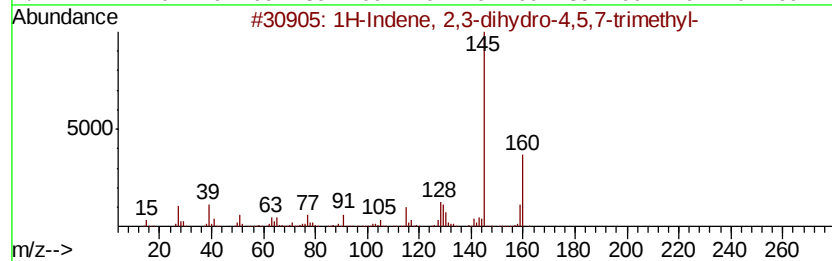
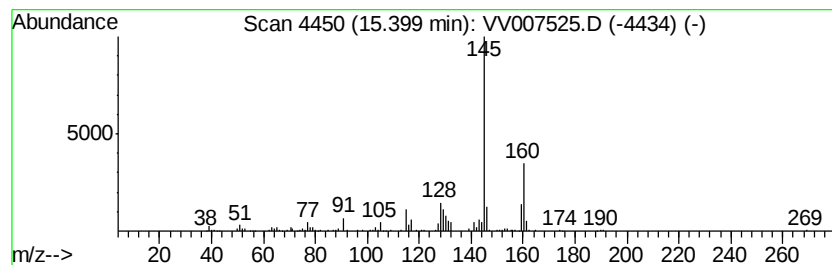
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 1H-Indene, 2,3-dihydro-4,5,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	0.69 ug/L	247726	1,4-Dichlorobenzene-d4	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,5,7-tri...	160	C12H16		006682-06-0	95
2	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16		014679-13-1	94
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16		054340-85-1	93
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16		020027-77-4	91
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16		021564-91-0	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV091318\
Data File : VV007525.D
Acq On : 13 Sep 2018 22:25
Operator : SY/MD
Sample : J4864-04
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0D10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR091318WMA.M
Quant Title : TRACE VOA SOM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Trimethylsilyl fl...	1.42	0.8	ug/L	49416	1	5.67	303902	5.0
Acetic acid	6.06	0.8	ug/L	48762	1	5.67	303902	5.0
(+)-2-Bornanone	13.22	0.5	ug/L	186289	3	11.30	1799030	5.0
Naphthalene, 1,2,...	13.66	0.7	ug/L	241746	3	11.30	1799030	5.0
1H-Indene, 2,3-di...	13.76	0.8	ug/L	285866	3	11.30	1799030	5.0
1H-Indene, 2,3-di...	14.14	0.6	ug/L	219604	3	11.30	1799030	5.0
Naphthalene, 1,2,...	14.67	0.7	ug/L	255349	3	11.30	1799030	5.0
1H-Indene, 2,3-di...	15.40	0.7	ug/L	247726	3	11.30	1799030	5.0