

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU040120\
 Data File : VU037536.D
 Acq On : 01 Apr 2020 21:40
 Operator : JC/MD
 Sample : L2065-15
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DBF11

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\SOMUTR040120WMA.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	1.369	3	9	20	rBV	147164	141866	38.66%	3.959%
2	1.485	39	45	49	rBV	31189	31647	8.62%	0.883%
3	1.610	73	84	98	rBV3	84336	131042	35.71%	3.657%
4	1.749	120	127	136	rVB3	5747	7764	2.12%	0.217%
5	1.929	173	183	194	rVB2	31475	44460	12.12%	1.241%
6	2.012	201	209	218	rBV	12689	17414	4.75%	0.486%
7	2.173	250	259	267	rBV	19506	26515	7.23%	0.740%
8	2.221	267	274	283	rVB	19742	26860	7.32%	0.750%
9	2.430	332	339	350	rVB7	5392	10463	2.85%	0.292%
10	2.591	379	389	402	rVB	49933	80900	22.04%	2.257%
11	2.665	404	412	425	rVB2	3104	5395	1.47%	0.151%
12	2.819	448	460	475	rVB	19336	32313	8.81%	0.902%
13	3.006	503	518	531	rBV5	3674	9317	2.54%	0.260%
14	3.109	545	550	565	rVB4	2517	4313	1.18%	0.120%
15	3.395	628	639	652	rVB5	2735	5821	1.59%	0.162%
16	3.626	696	711	724	rVB6	5863	12655	3.45%	0.353%
17	3.742	733	747	759	rVB6	2953	6790	1.85%	0.189%
18	4.671	1021	1036	1057	rBV	61495	143682	39.15%	4.009%
19	5.105	1155	1171	1193	rVB2	52978	125105	34.09%	3.491%
20	5.761	1354	1375	1392	rBV2	99865	256644	69.93%	7.161%
21	6.282	1524	1537	1556	rVB	113503	210017	57.23%	5.860%
22	6.723	1661	1674	1688	rBV	68083	128663	35.06%	3.590%
23	7.591	1933	1944	1959	rBV2	35029	61952	16.88%	1.729%
24	7.922	2034	2047	2061	rBV	120537	204003	55.59%	5.693%
25	7.996	2061	2070	2080	rVB2	3643	6065	1.65%	0.169%
26	8.202	2125	2134	2147	rBV2	23258	37753	10.29%	1.053%
27	8.658	2263	2276	2300	rBV	211484	366982	100.00%	10.240%
28	9.436	2505	2518	2533	rBV	163809	266479	72.61%	7.436%
29	10.774	2921	2934	2946	rBV	56198	89501	24.39%	2.497%
30	10.903	2966	2974	2986	rVB2	6823	10562	2.88%	0.295%
31	11.587	3181	3187	3195	rVB7	2818	3728	1.02%	0.104%
32	11.648	3195	3206	3215	rBV8	3458	6001	1.64%	0.167%
33	11.825	3250	3261	3273	rVB	174127	267850	72.99%	7.474%
34	11.983	3298	3310	3321	rVB	3497	8995	2.45%	0.251%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

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Stop Thrs : 0

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Max Peaks: 100

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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 Title : TRACE VOA SOM01.0

35	12.044	3321	3329	3333	rBV5	3291	4562	1.24%	0.127%
36	12.076	3333	3339	3346	rVV3	3528	4947	1.35%	0.138%
37	12.156	3356	3364	3368	rBV6	4644	6030	1.64%	0.168%
38	12.205	3368	3379	3395	rVB	144403	243030	66.22%	6.782%
39	12.398	3429	3439	3441	rBV6	2996	4033	1.10%	0.113%
40	12.726	3535	3541	3546	rBV9	3220	4928	1.34%	0.138%
41	12.767	3547	3554	3557	rBV6	4386	5431	1.48%	0.152%
42	12.854	3570	3581	3594	rVV5	23957	40458	11.02%	1.129%
43	12.912	3594	3599	3607	rVB7	3140	3756	1.02%	0.105%
44	12.986	3612	3622	3629	rVB9	2956	4548	1.24%	0.127%
45	13.070	3637	3648	3657	rBV5	16842	28479	7.76%	0.795%
46	13.150	3661	3673	3681	rBV5	12303	24862	6.77%	0.694%
47	13.330	3721	3729	3736	rVB10	5042	7851	2.14%	0.219%
48	13.401	3744	3751	3762	rVB7	6352	10525	2.87%	0.294%
49	13.475	3762	3774	3785	rBV10	10005	18945	5.16%	0.529%
50	13.555	3789	3799	3809	rBV10	5423	9974	2.72%	0.278%
51	13.632	3813	3823	3834	rBV10	6159	12996	3.54%	0.363%
52	13.758	3854	3862	3867	rBV10	2626	4577	1.25%	0.128%
53	13.873	3889	3898	3909	rBV10	2496	4970	1.35%	0.139%
54	14.115	3961	3973	3983	rBV	62013	99722	27.17%	2.783%
55	15.259	4317	4329	4340	rBV	21372	34254	9.33%	0.956%
56	15.449	4378	4388	4403	rVB	11214	17146	4.67%	0.478%
57	16.005	4552	4561	4569	rBV4	9264	14066	3.83%	0.392%
58	16.201	4615	4622	4629	rBV4	3498	5171	1.41%	0.144%
59	16.314	4649	4657	4666	rVB6	5746	8259	2.25%	0.230%
60	16.462	4693	4703	4711	rBV7	5727	10085	2.75%	0.281%
61	16.504	4711	4716	4724	rVB7	3407	4763	1.30%	0.133%
62	16.944	4844	4853	4862	rVB5	7919	12596	3.43%	0.351%
63	17.159	4907	4920	4932	rBV	49146	87173	23.75%	2.432%
64	17.478	5008	5019	5035	rVB	28382	56061	15.28%	1.564%

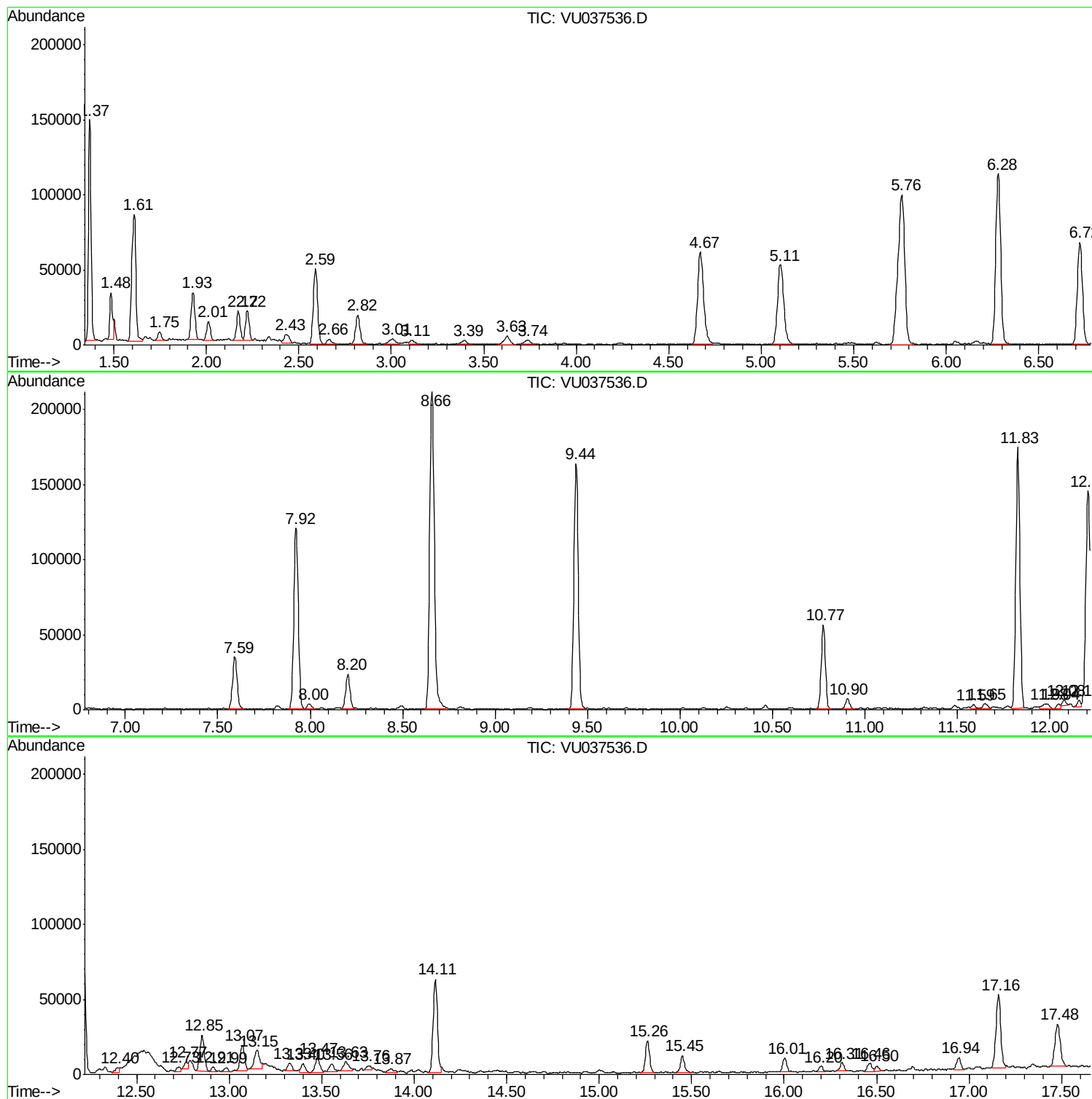
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ClientSampled :
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Quant Title : TRACE VOA SOM01.0

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



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ClientSampleId :
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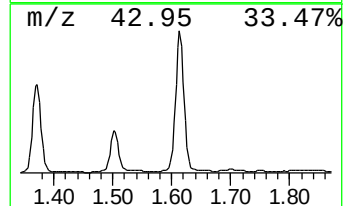
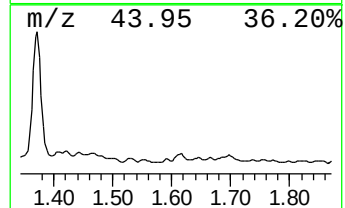
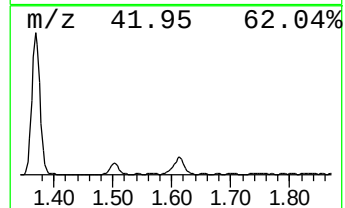
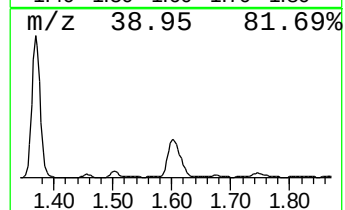
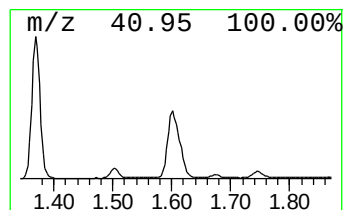
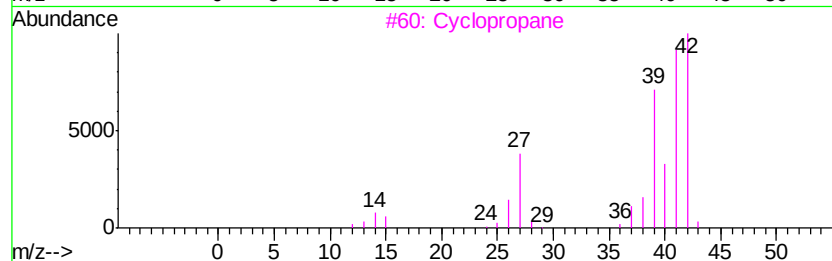
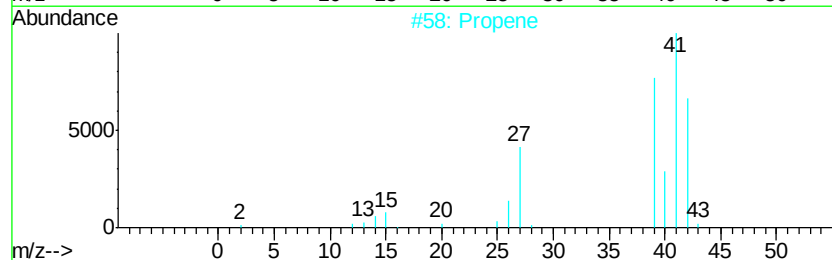
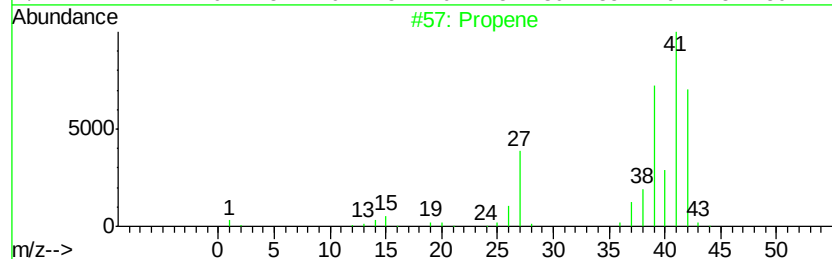
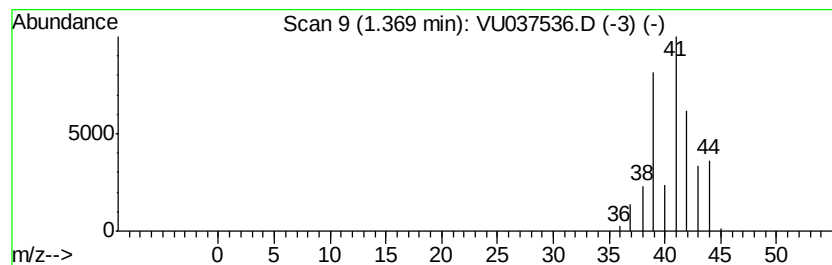
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Cyclic1.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.37	3.38 ug/L	141866	1,4-Difluorobenzene	6.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene			42	C3H6	000115-07-1	90
2	Propene			42	C3H6	000115-07-1	42
3	Cyclopropane			42	C3H6	000075-19-4	9
4	Cyclopropene			40	C3H4	002781-85-3	9
5	Cyclopropane			42	C3H6	000075-19-4	7



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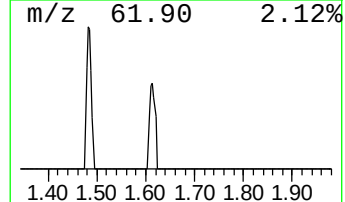
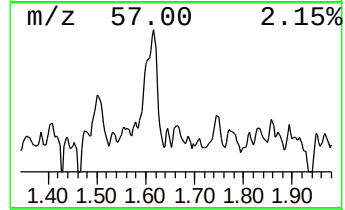
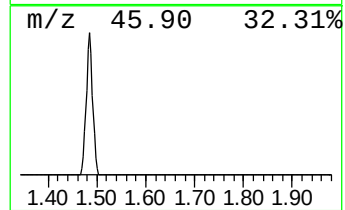
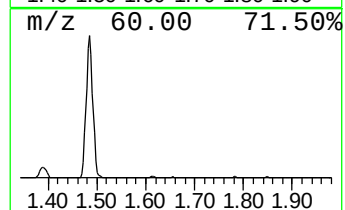
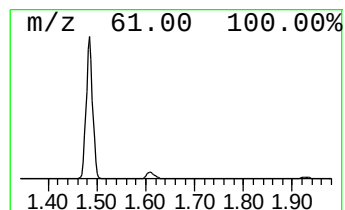
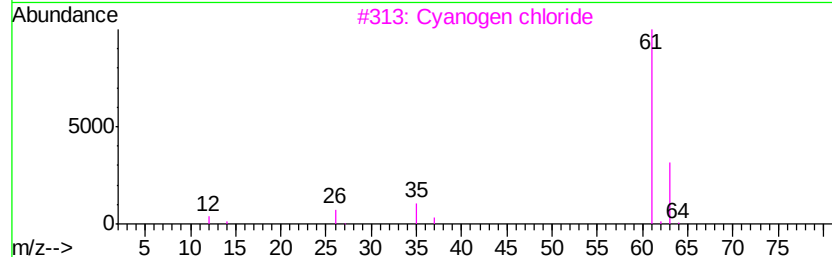
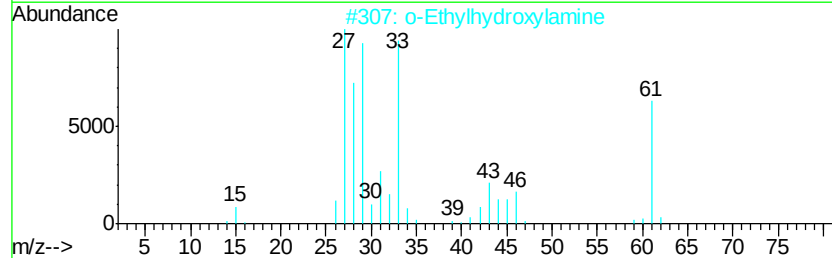
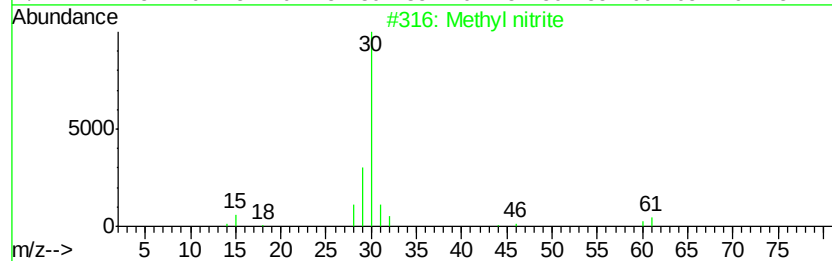
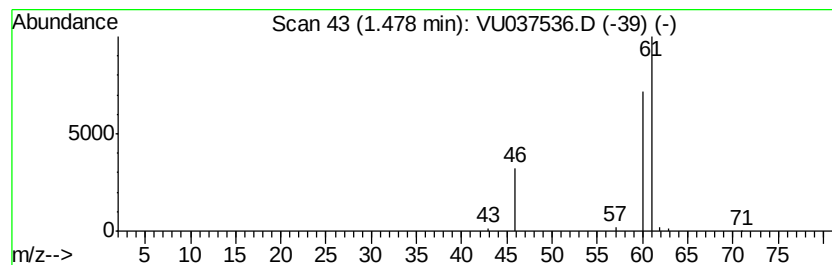
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR040120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.48	0.75 ug/L	31647	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl nitrite	61	CH3NO2	000624-91-9	5
2		o-Ethylhydroxylamine	61	C2H7NO	1000322-42-5	4
3		Cyanogen chloride	61	CCLN	000506-77-4	4
4		Trifluoroquandine	113	CH2F3N3	037950-72-4	4
5		Cyanogen chloride	61	CCLN	000506-77-4	4



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ALS Vial : 1 Sample Multiplier: 1

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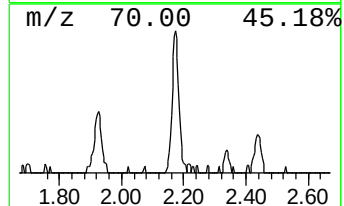
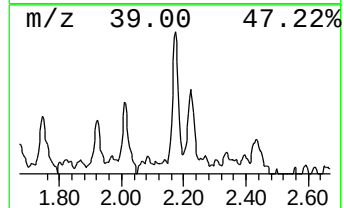
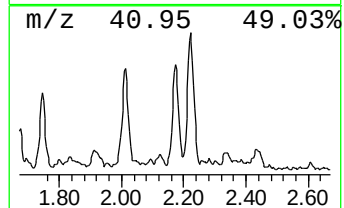
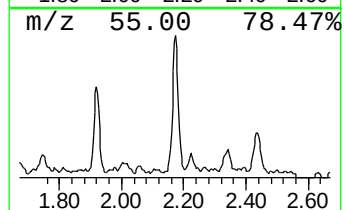
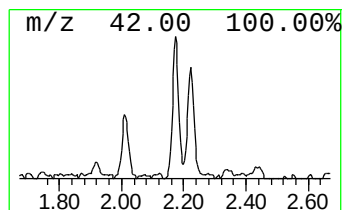
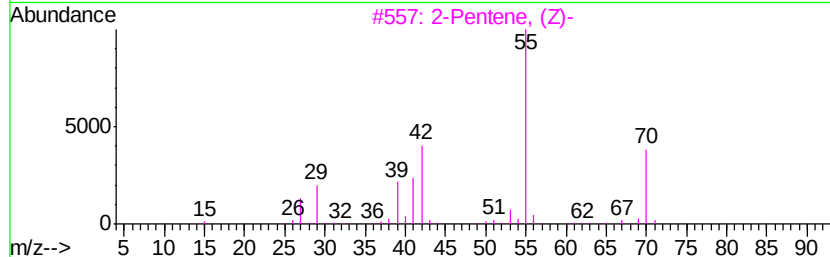
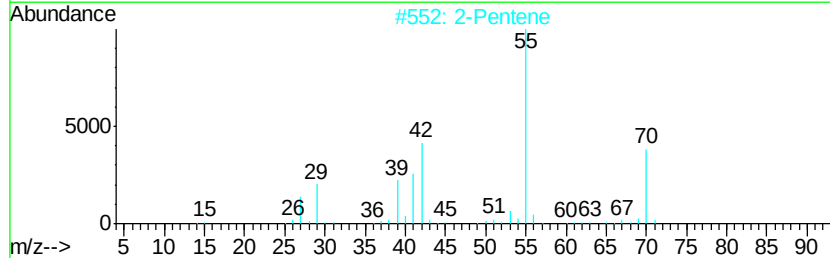
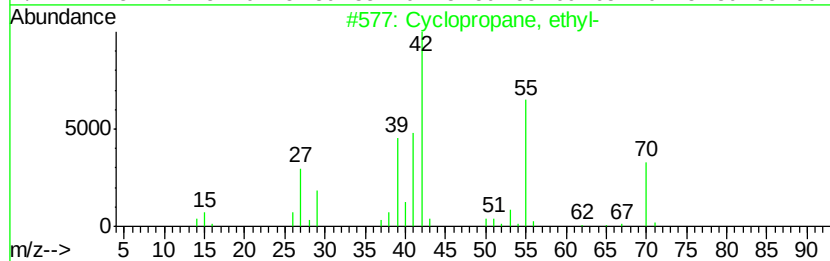
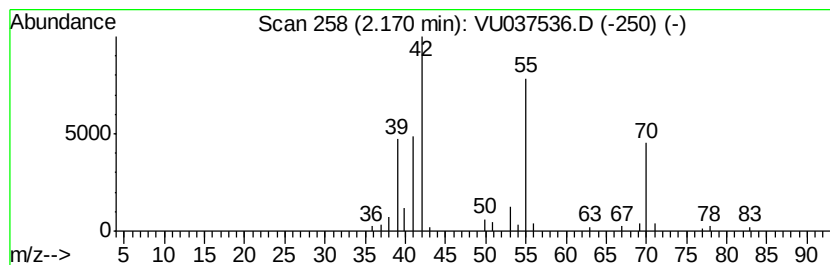
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 (DEL) Alkane: Cyclic2.17 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.17	0.63 ug/L	26515	1,4-Difluorobenzene	6.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, ethyl-	70	C5H10	001191-96-4	87
2			2-Pentene	70	C5H10	000109-68-2	83
3			2-Pentene, (Z)-	70	C5H10	000627-20-3	83
4			Cyclopropane, ethyl-	70	C5H10	001191-96-4	78
5			1-Pentene	70	C5H10	000109-67-1	72



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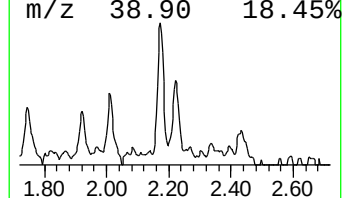
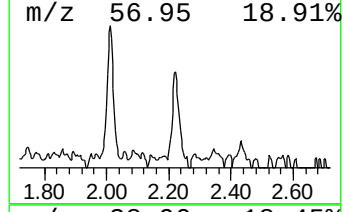
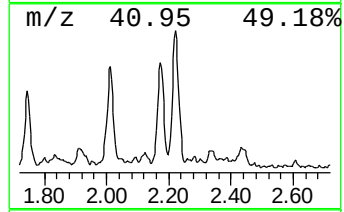
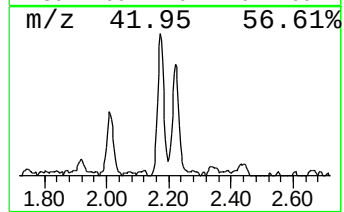
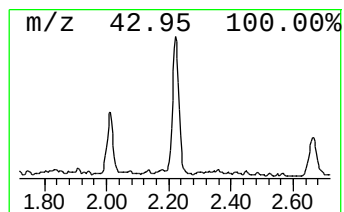
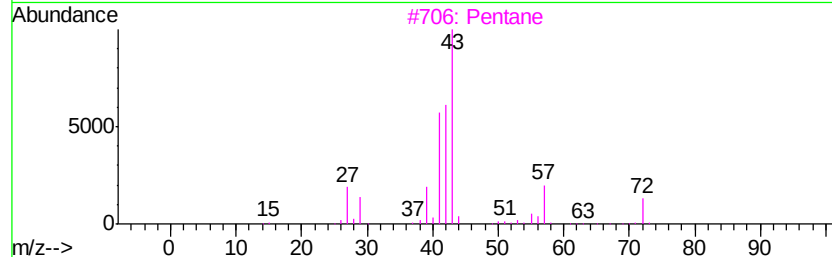
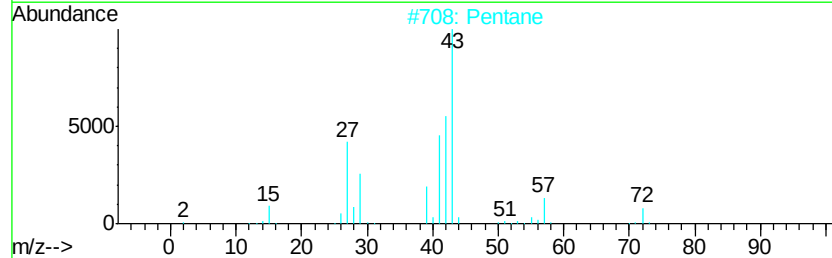
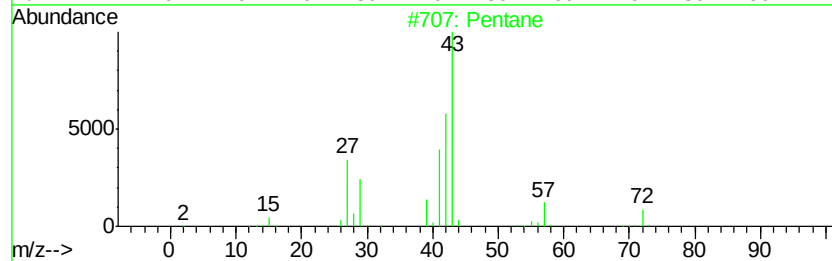
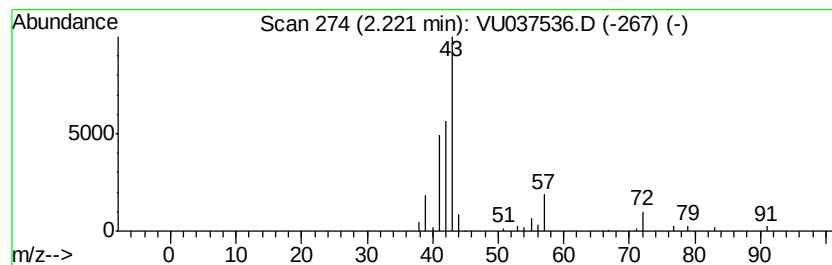
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR040120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.22	0.64 ug/L	26860	1,4-Difluorobenzene	6.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane		72	C5H12	000109-66-0	86
2	Pentane		72	C5H12	000109-66-0	86
3	Pentane		72	C5H12	000109-66-0	86
4	Pentane, 2-bromo-		150	C5H11Br	000107-81-3	39
5	Butane, 2-methyl-		72	C5H12	000078-78-4	33



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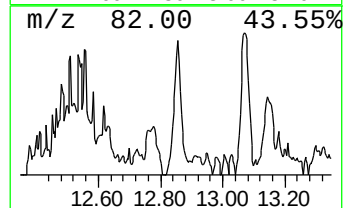
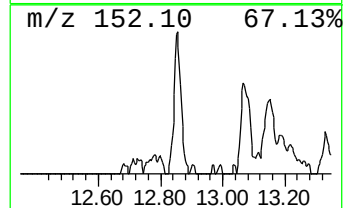
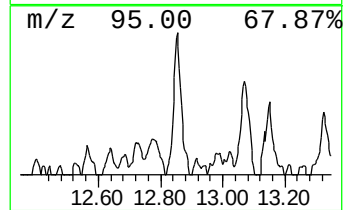
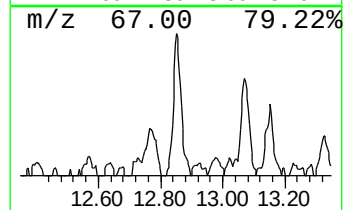
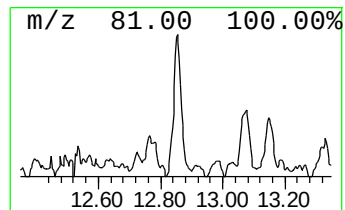
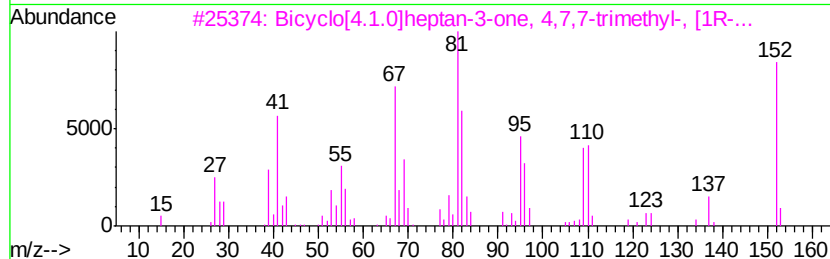
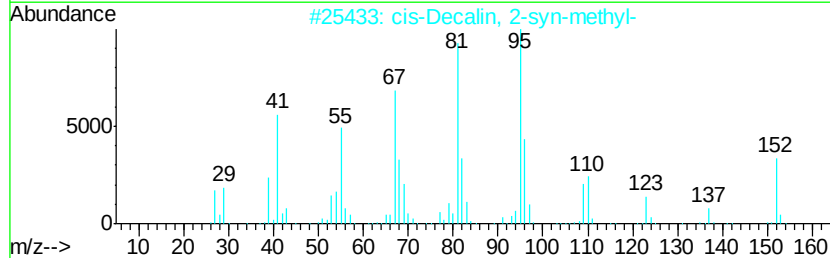
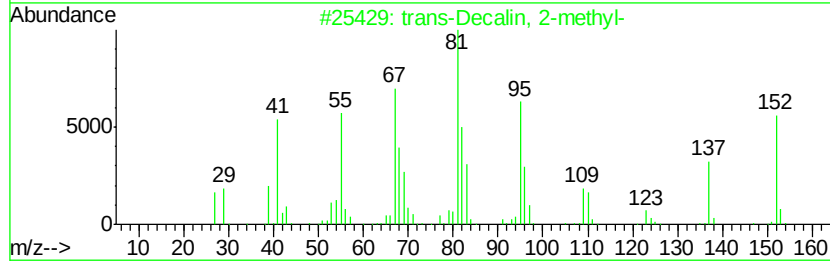
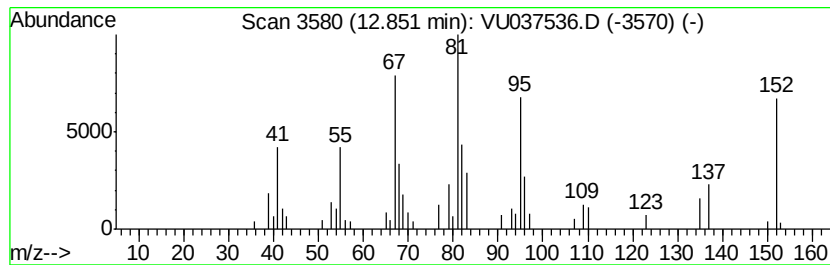
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MSVOA_U
ClientSampled :
DBF11

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR040120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 trans-Decalin, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.85	0.76 ug/L	40458	1,4-Dichlorobenzene-d4	11.83		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-Decalin, 2-methyl-		152	C11H20	1000152-47-3	87
2	cis-Decalin, 2-syn-methyl-		152	C11H20	1000155-85-6	80
3	Bicyclo[4.1.0]heptan-3-one, 4,7,...		152	C10H16O	004176-04-9	74
4	Cyclohexanone, 5-methyl-2-(1-met...		152	C10H16O	015932-80-6	72
5	5-Hexadecyne		222	C16H30	071899-37-1	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040120\
Data File : VU037536.D
Acq On : 01 Apr 2020 21:40
Operator : JC/MD
Sample : L2065-15
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF11

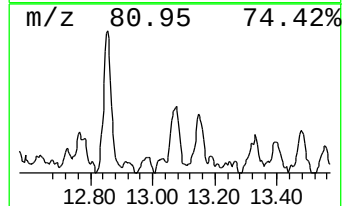
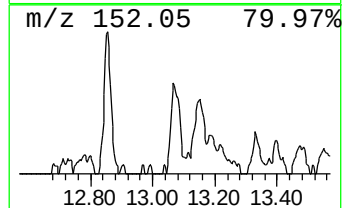
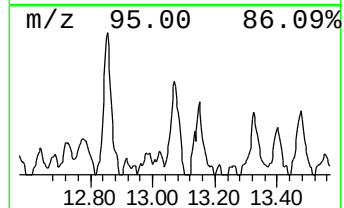
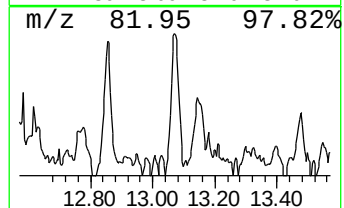
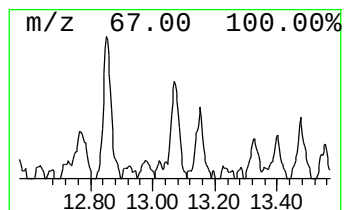
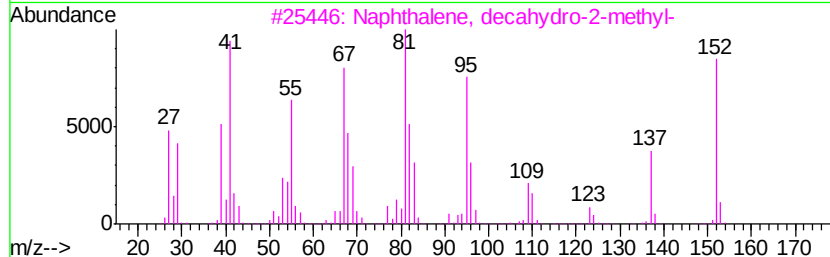
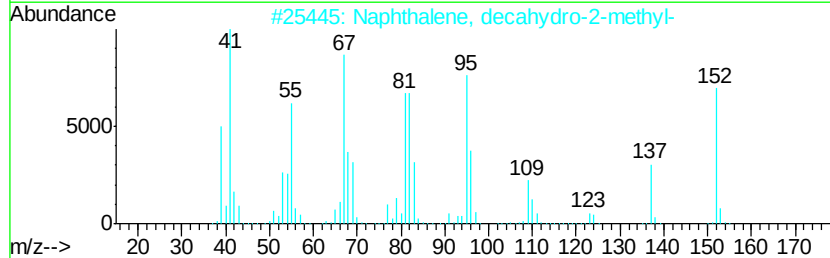
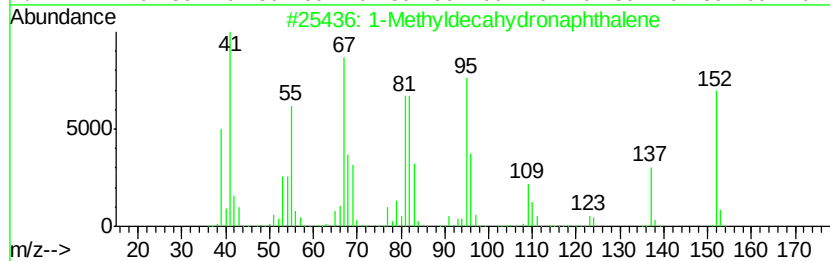
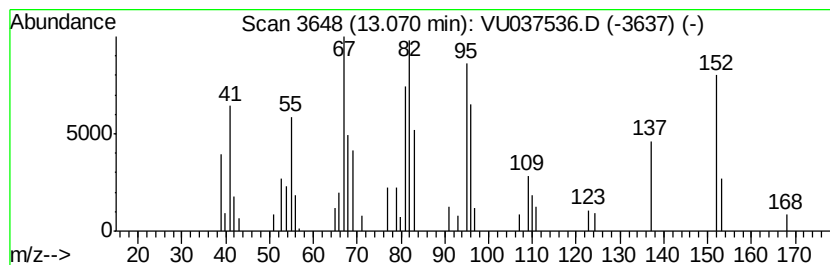
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 1-Methyldecahydronaphthalene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.07	0.53 ug/L	28479	1,4-Dichlorobenzene-d4	11.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	93
2		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	93
3		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	78
4		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	78
5		Bicyclo[4.1.0]heptane, 7-butyl-	152	C11H20	018645-10-8	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040120\
Data File : VU037536.D
Acq On : 01 Apr 2020 21:40
Operator : JC/MD
Sample : L2065-15
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF11

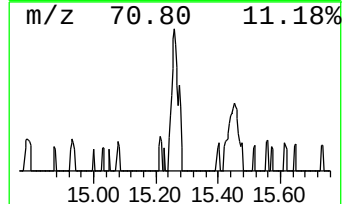
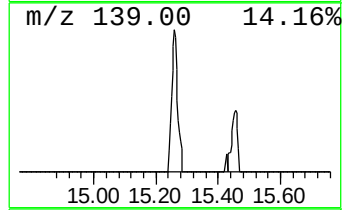
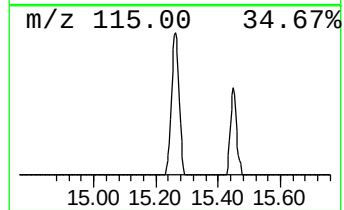
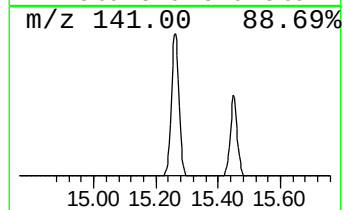
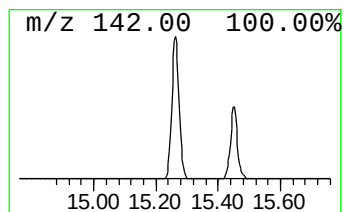
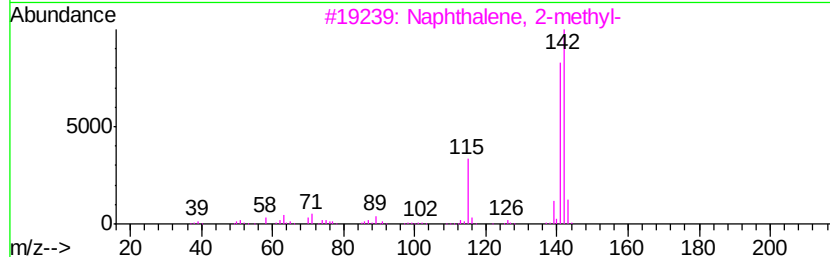
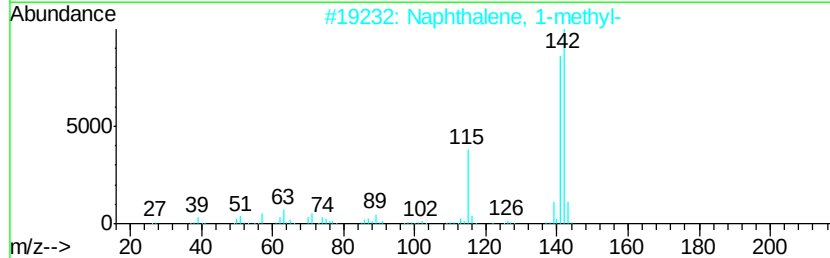
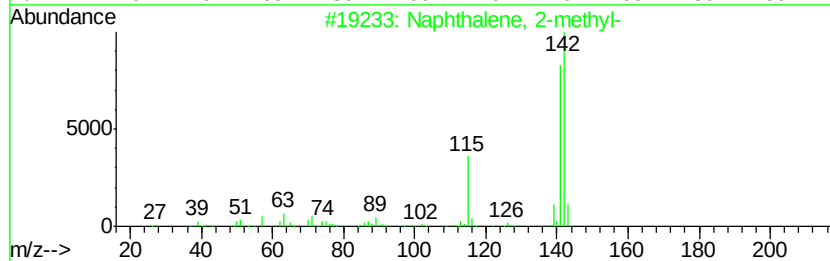
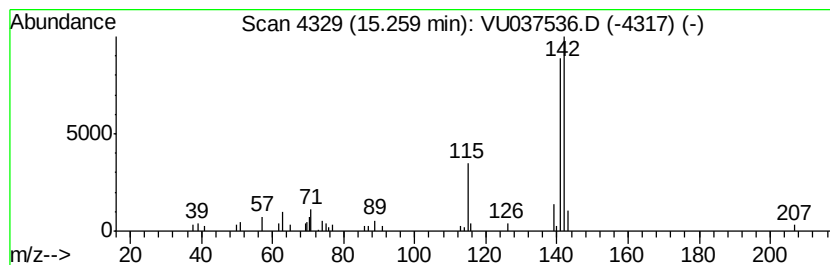
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	0.64 ug/L	34254	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU040120\
Data File : VU037536.D
Acq On : 01 Apr 2020 21:40
Operator : JC/MD
Sample : L2065-15
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF11

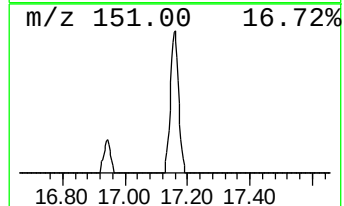
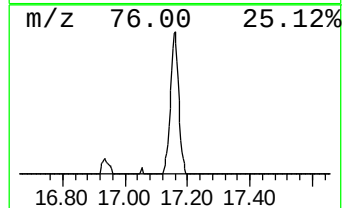
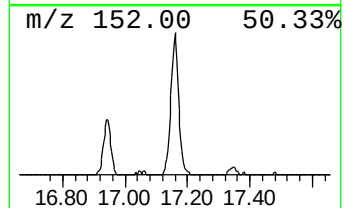
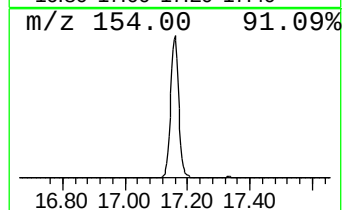
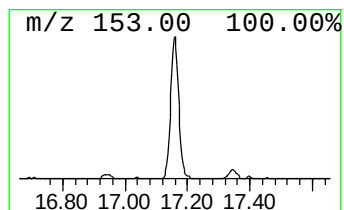
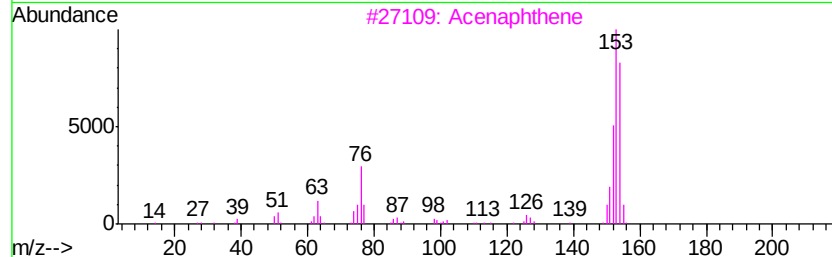
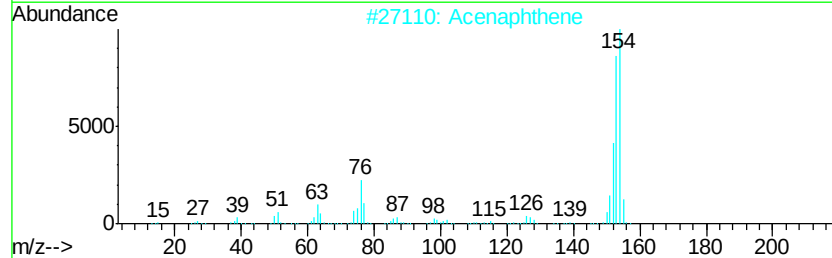
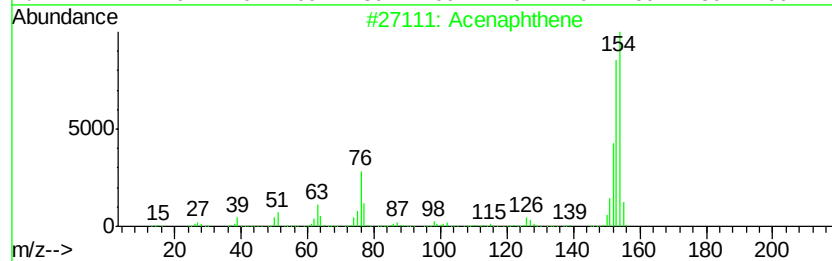
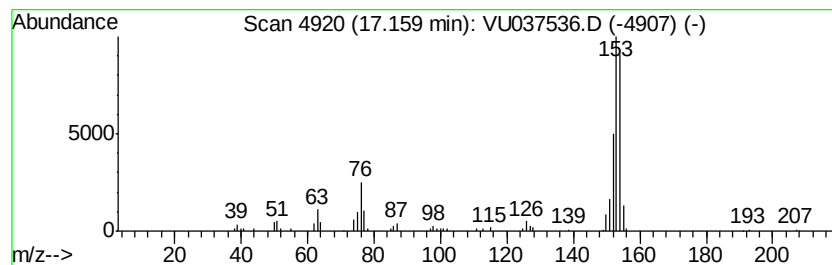
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Acenaphthene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.16	1.63 ug/L	87173	1,4-Dichlorobenzene-d4	11.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthene	154	C12H10	000083-32-9	96	
2	Acenaphthene	154	C12H10	000083-32-9	96	
3	Acenaphthene	154	C12H10	000083-32-9	96	
4	Acenaphthene	154	C12H10	000083-32-9	76	
5	1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	72	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040120\
Data File : VU037536.D
Acq On : 01 Apr 2020 21:40
Operator : JC/MD
Sample : L2065-15
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF11

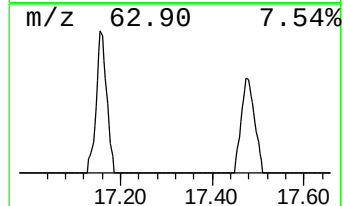
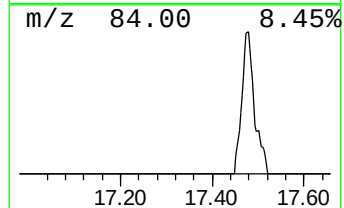
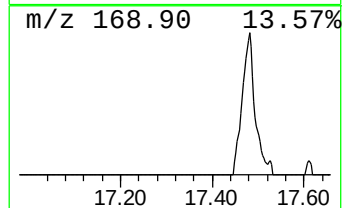
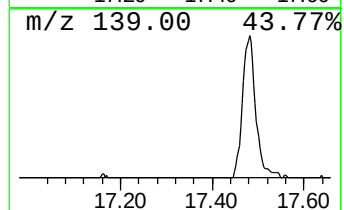
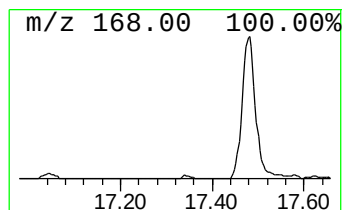
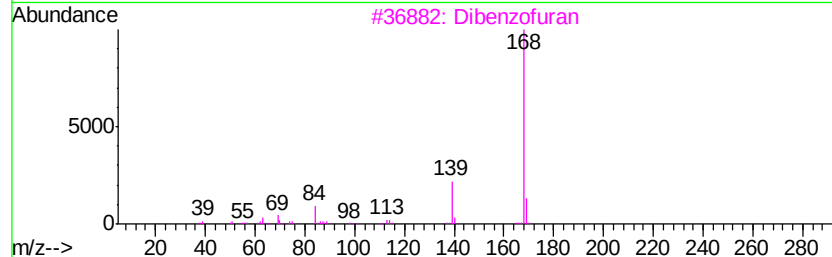
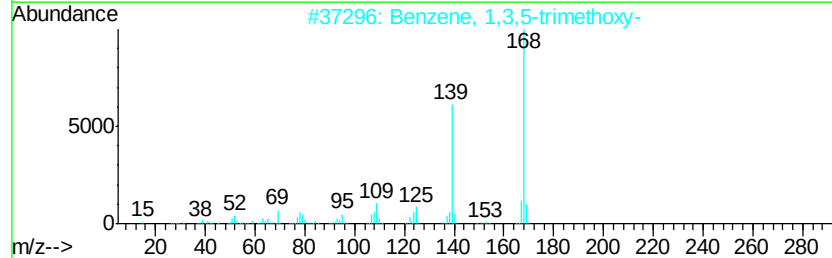
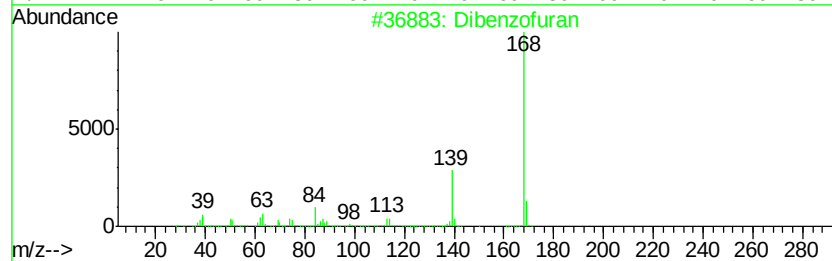
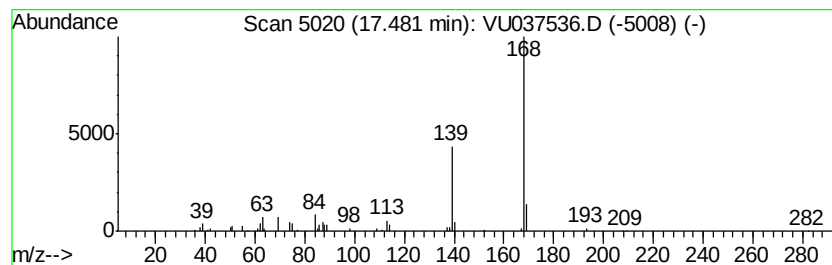
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR040120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Dibenzofuran Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.48	1.05 ug/L	56061	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzofuran	168	C12H8O	000132-64-9	94
2		Benzene, 1,3,5-trimethoxy-	168	C9H12O3	000621-23-8	64
3		Dibenzofuran	168	C12H8O	000132-64-9	64
4		9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	53
5		1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	53



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU040120\
Data File : VU037536.D
Acq On : 01 Apr 2020 21:40
Operator : JC/MD
Sample : L2065-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR040120WMA.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
(DEL) Alkane: Cyc...	1.37	3.4	ug/L	141866	1	6.28	210017	5.0
unknown-01	1.48	0.8	ug/L	31647	1	6.28	210017	5.0
(DEL) Alkane: Cyc...	2.17	0.6	ug/L	26515	1	6.28	210017	5.0
(DEL) Alkane: Str...	2.22	0.6	ug/L	26860	1	6.28	210017	5.0
trans-Decalin, 2-...	12.85	0.8	ug/L	40458	3	11.83	267850	5.0
1-Methyldecahydro...	13.07	0.5	ug/L	28479	3	11.83	267850	5.0
Naphthalene, 1-me...	15.26	0.6	ug/L	34254	3	11.83	267850	5.0
Acenaphthene	17.16	1.6	ug/L	87173	3	11.83	267850	5.0
Dibenzofuran	17.48	1.1	ug/L	56061	3	11.83	267850	5.0