

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU040120\
 Data File : VU037535.D
 Acq On : 01 Apr 2020 21:15
 Operator : JC/MD
 Sample : L2065-14
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DBF10

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	1.366	3	8	20	rVB	91724	86118	1.03%	0.485%
2	1.601	71	81	90	rBV2	55687	83847	1.00%	0.473%
3	4.752	1033	1061	1082	rBV2	35321	123598	1.47%	0.697%
4	5.128	1163	1178	1196	rBV3	55907	135723	1.62%	0.765%
5	5.774	1361	1379	1404	rBV3	108432	258083	3.08%	1.455%
6	6.295	1529	1541	1557	rBV	106025	196305	2.34%	1.107%
7	6.735	1670	1678	1696	rVB	68566	134606	1.61%	0.759%
8	7.928	2038	2049	2063	rBV	120056	199574	2.38%	1.125%
9	8.671	2264	2280	2308	rBV	193838	353555	4.22%	1.993%
10	9.440	2504	2519	2535	rBV	150788	243668	2.91%	1.374%
11	10.774	2922	2934	2946	rBV2	54946	88103	1.05%	0.497%
12	11.825	3251	3261	3276	rVB	163797	271130	3.23%	1.528%
13	12.108	3326	3349	3362	rBV	331363	562575	6.71%	3.171%
14	12.205	3368	3379	3393	rVB2	138509	225892	2.69%	1.273%
15	12.327	3402	3417	3431	rVB	555941	872590	10.41%	4.919%
16	13.047	3630	3641	3662	rBV2	52969	113348	1.35%	0.639%
17	14.118	3956	3974	3990	rBV	5185425	8383487	100.00%	47.257%
18	14.237	3999	4011	4022	rBV	114376	181750	2.17%	1.025%
19	15.262	4318	4330	4354	rVB	933631	1520721	18.14%	8.572%
20	15.452	4376	4389	4405	rBV	480127	791184	9.44%	4.460%
21	16.002	4542	4560	4573	rBV2	208505	328273	3.92%	1.850%
22	16.198	4603	4621	4630	rBV	112120	196868	2.35%	1.110%
23	16.314	4644	4657	4677	rVB	141065	254102	3.03%	1.432%
24	16.462	4691	4703	4710	rBV	143510	248479	2.96%	1.401%
25	16.500	4710	4715	4727	rVB	84157	127618	1.52%	0.719%
26	16.693	4758	4775	4786	rBV3	38525	94360	1.13%	0.532%
27	16.944	4839	4853	4865	rBV2	61470	103905	1.24%	0.586%
28	17.156	4906	4919	4934	rBV	707594	1253777	14.96%	7.067%
29	17.475	5003	5018	5034	rVB	159104	306981	3.66%	1.730%

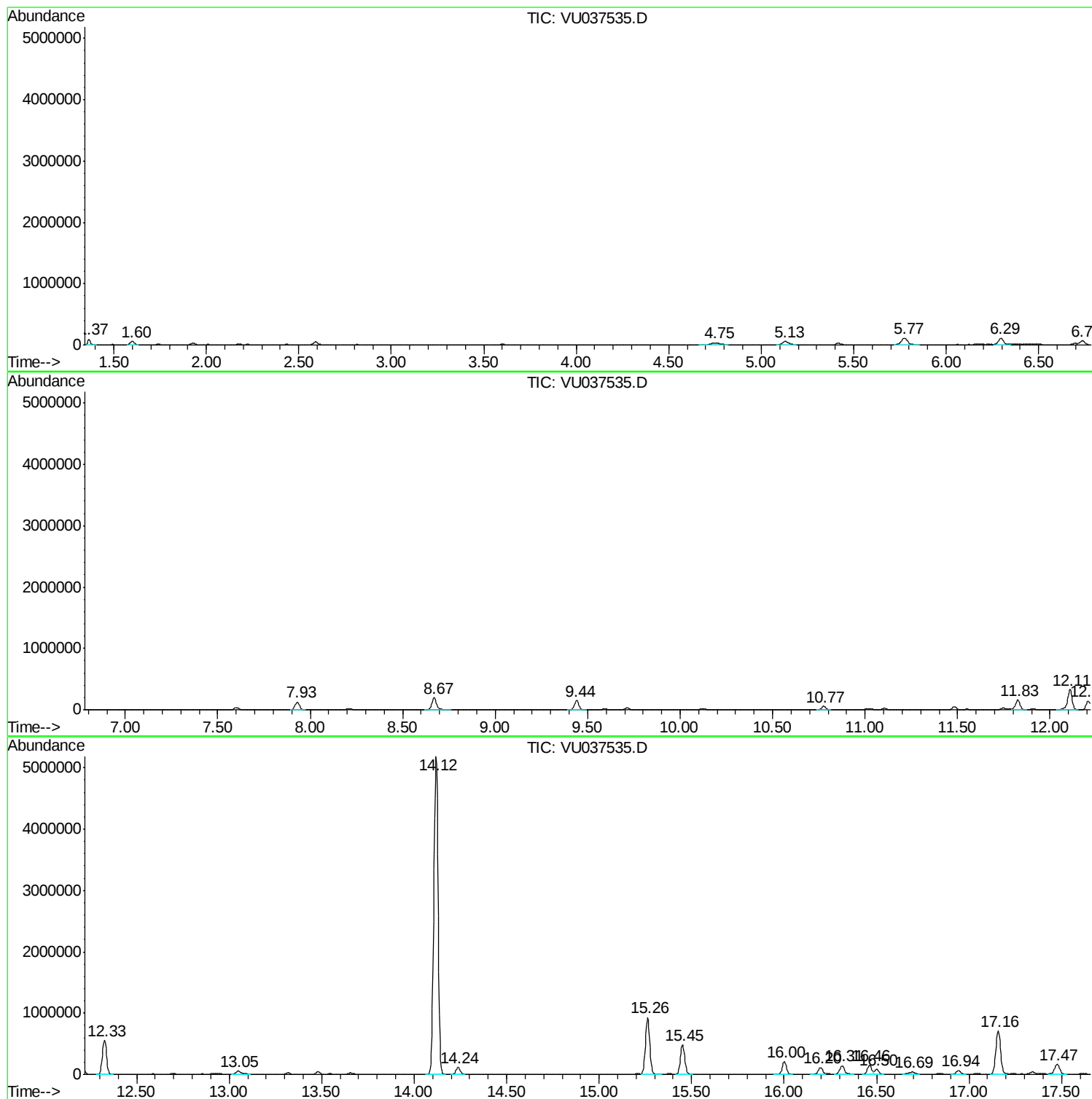
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF10

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



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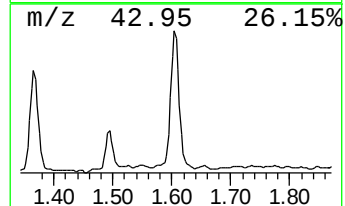
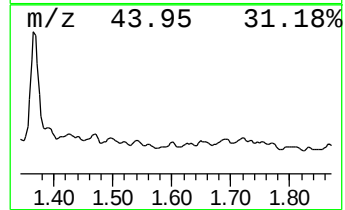
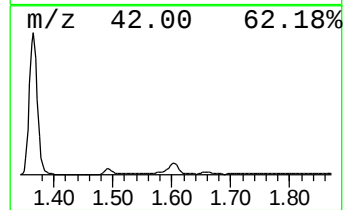
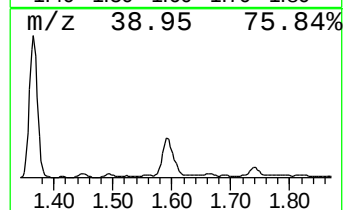
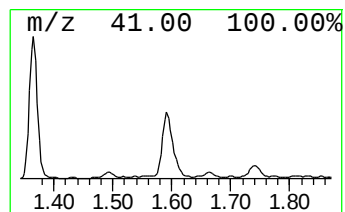
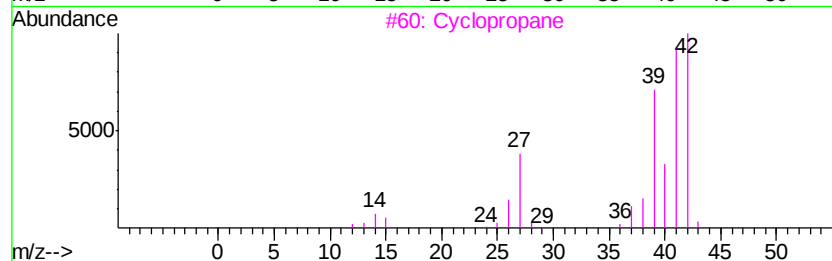
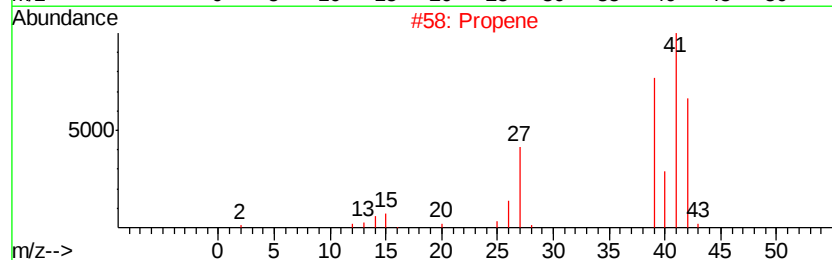
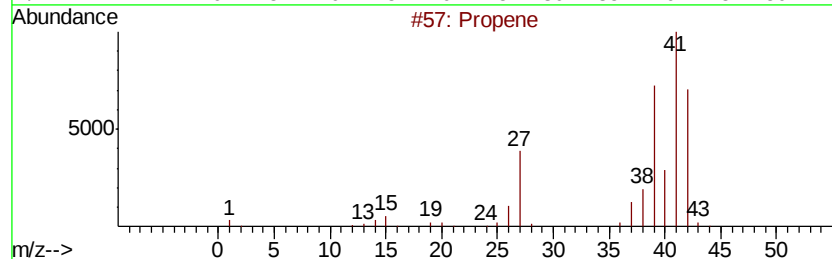
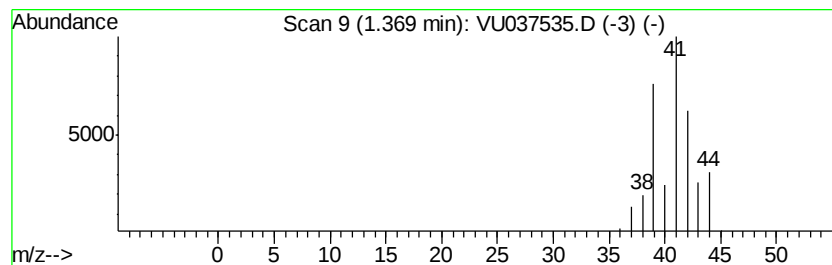
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Cyclic1.37 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.37	2.19 ug/L	86118	1,4-Difluorobenzene	6.29

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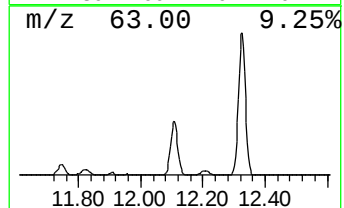
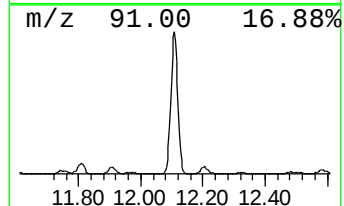
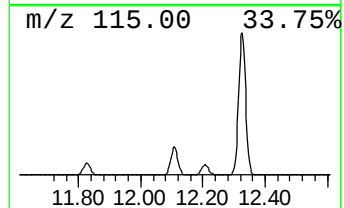
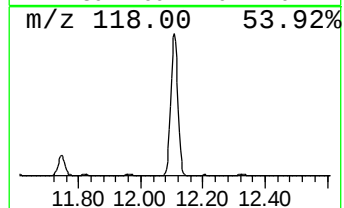
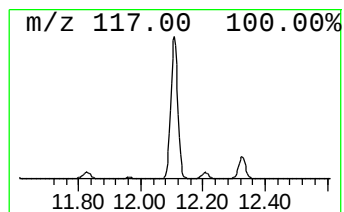
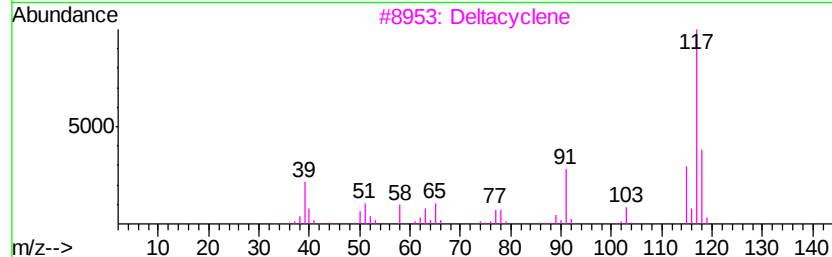
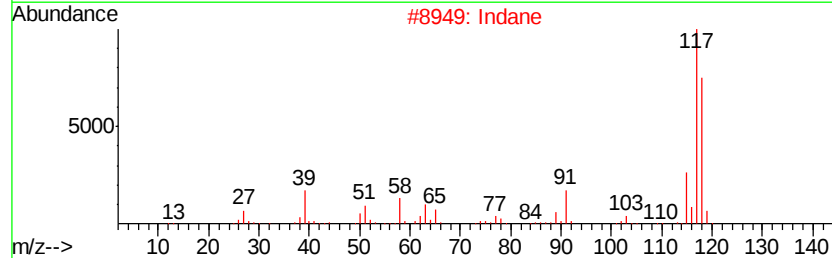
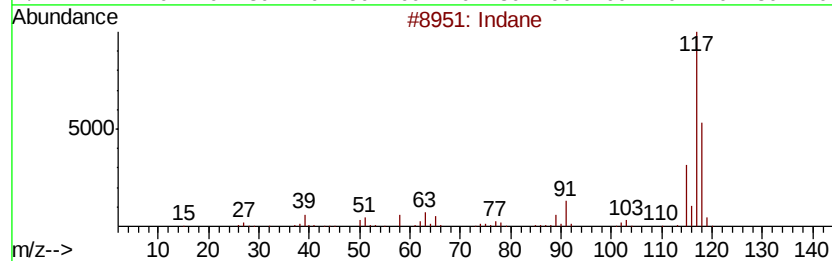
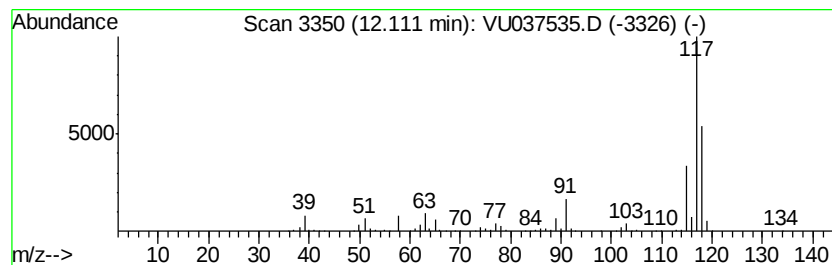
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR040120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	10.37 ug/L	562575	1,4-Dichlorobenzene-d4	11.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Indane		118	C9H10	000496-11-7	93
3	Deltacyclene		118	C9H10	007785-10-6	68
4	Benzene, 1-ethenyl-4-methyl-		118	C9H10	000622-97-9	64
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...		118	C9H10	1000191-13-7	64



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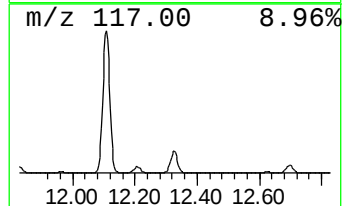
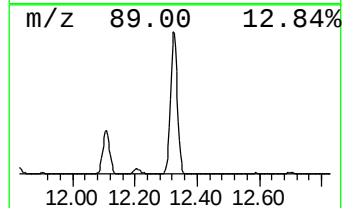
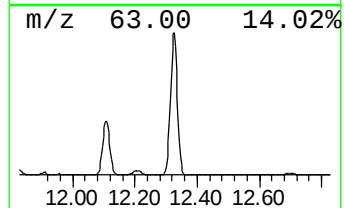
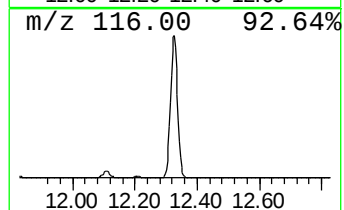
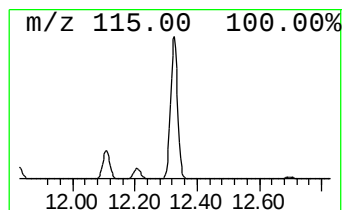
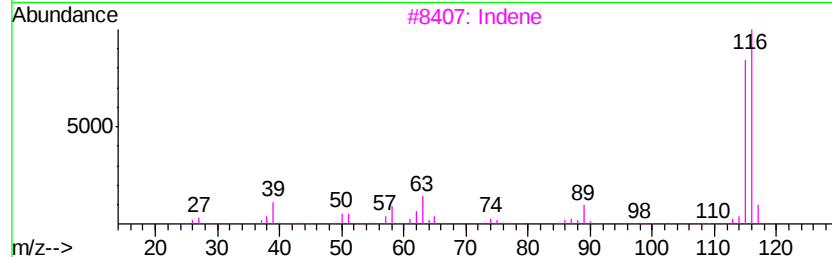
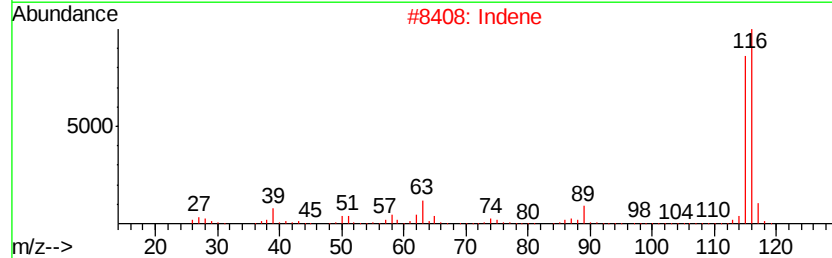
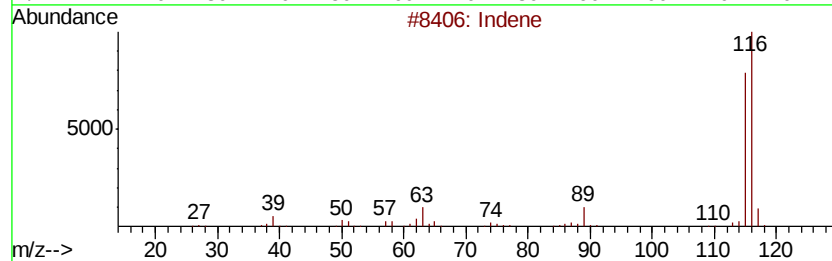
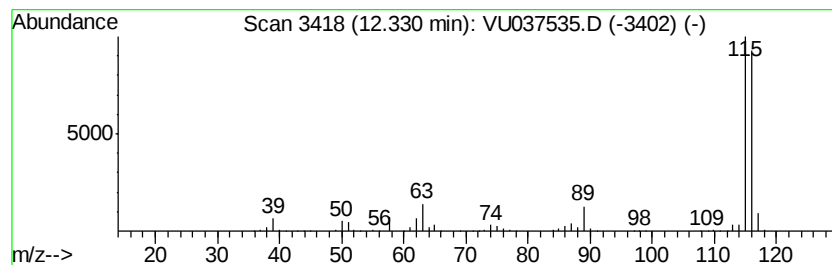
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	16.09 ug/L	872590	1,4-Dichlorobenzene-d4	11.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene		116	C9H8	000095-13-6	95
2	Indene		116	C9H8	000095-13-6	94
3	Indene		116	C9H8	000095-13-6	94
4	Benzene, 1,2-propadienyl-		116	C9H8	002327-99-3	91
5	3-Methylphenylacetylene		116	C9H8	000766-82-5	91



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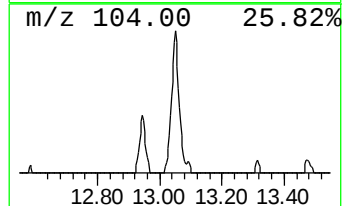
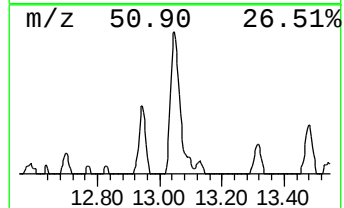
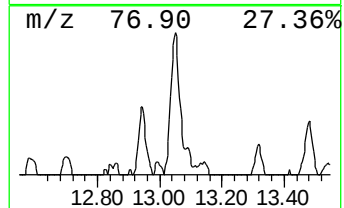
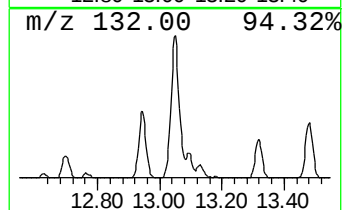
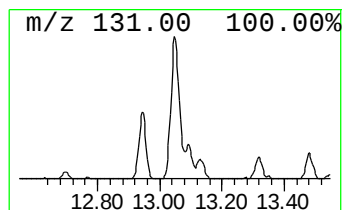
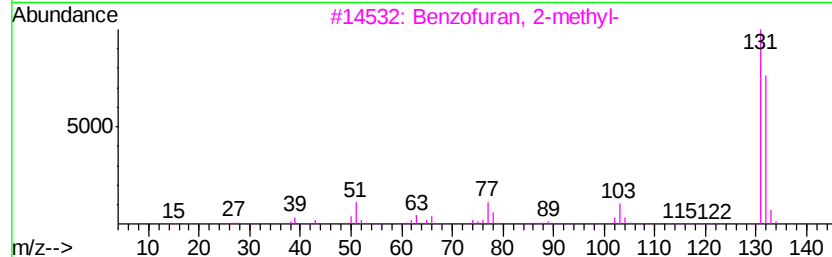
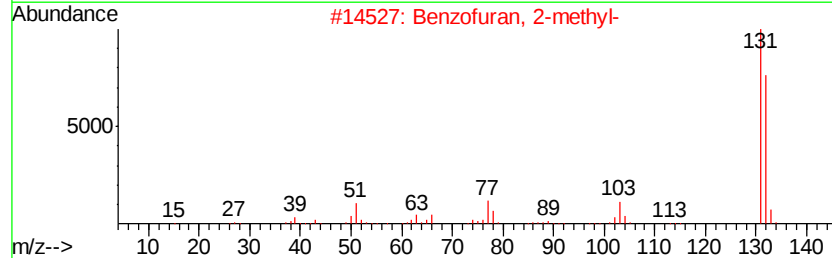
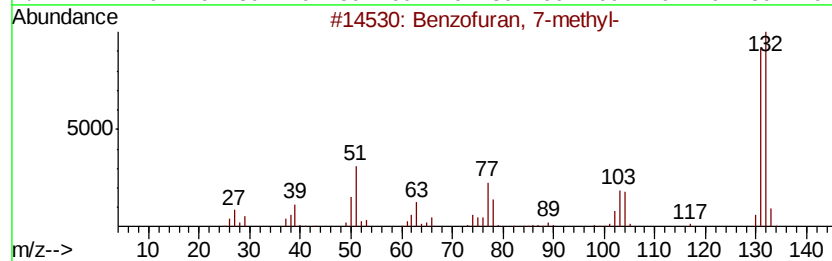
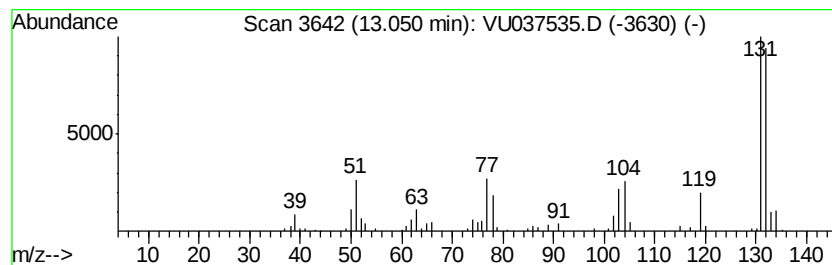
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzofuran, 7-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.05	2.09 ug/L	113348	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
2		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90
5		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	87



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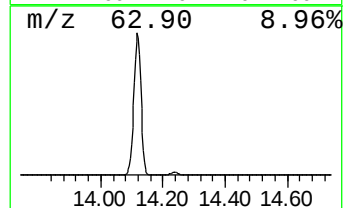
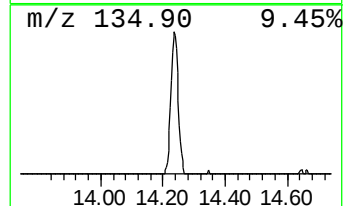
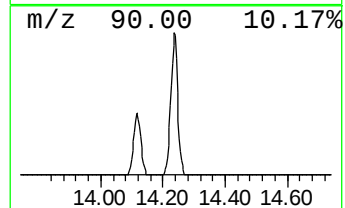
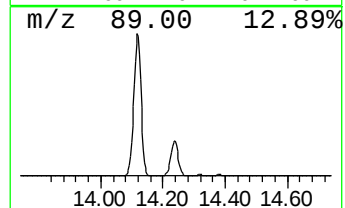
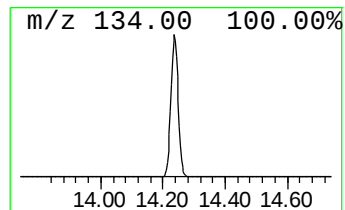
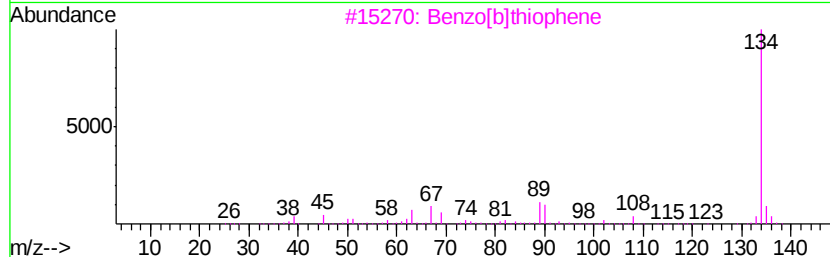
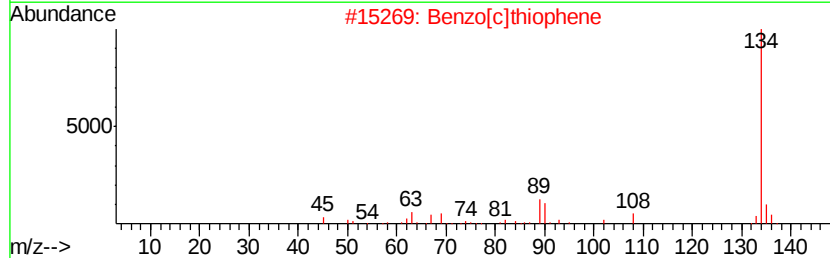
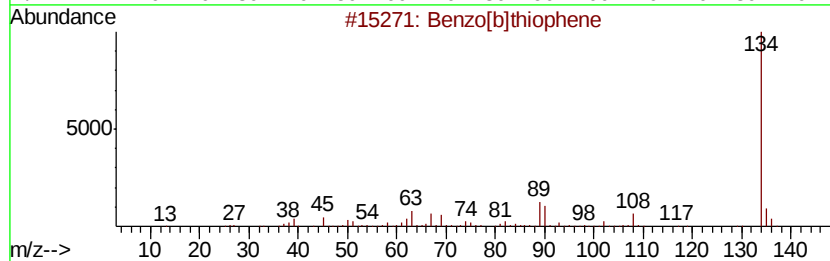
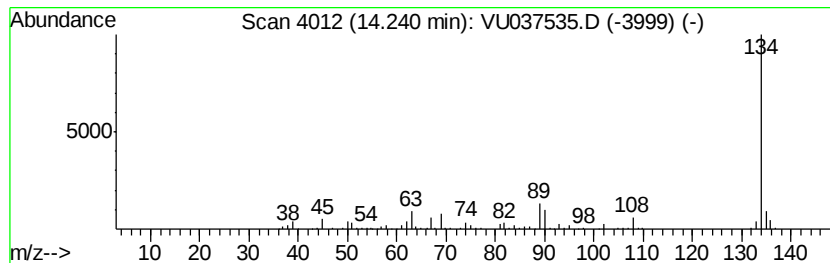
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 Benzo[b]thiophene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	3.35 ug/L	181750	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene	134	C8H6S	000095-15-8	95
2		Benzo[c]thiophene	134	C8H6S	000270-82-6	95
3		Benzo[b]thiophene	134	C8H6S	000095-15-8	94
4		Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5		Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	90



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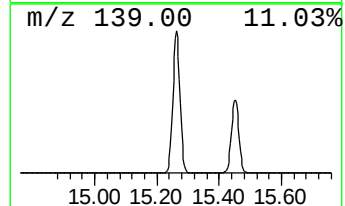
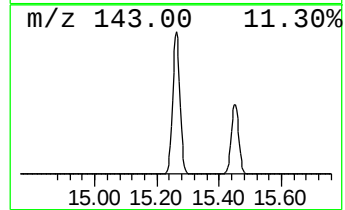
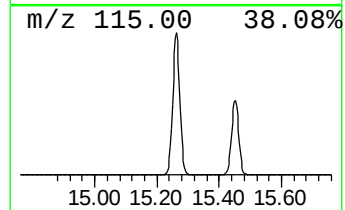
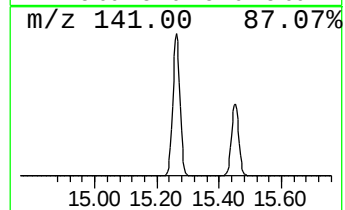
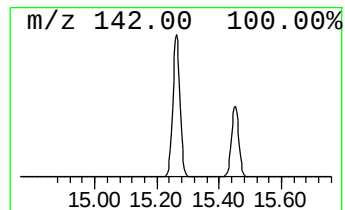
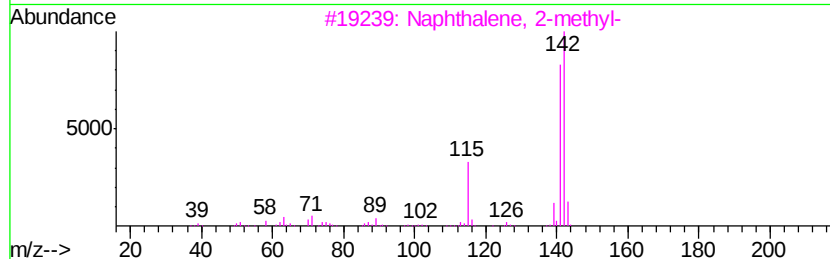
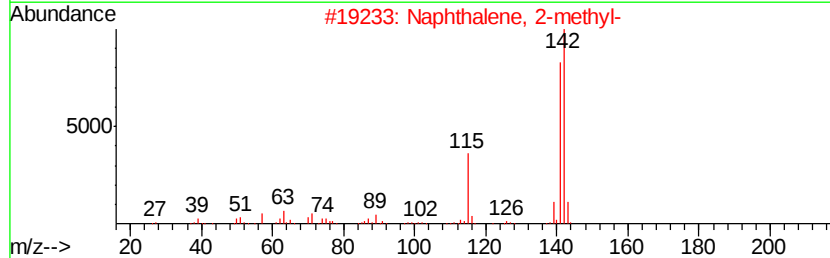
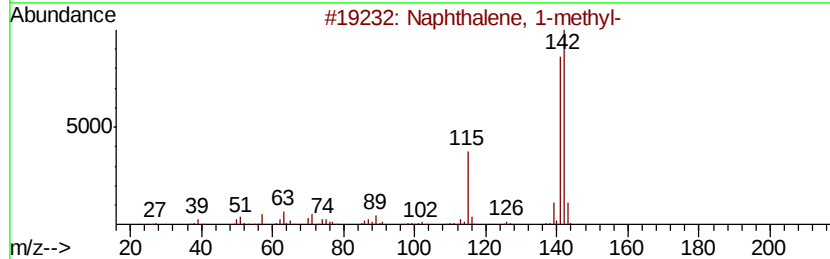
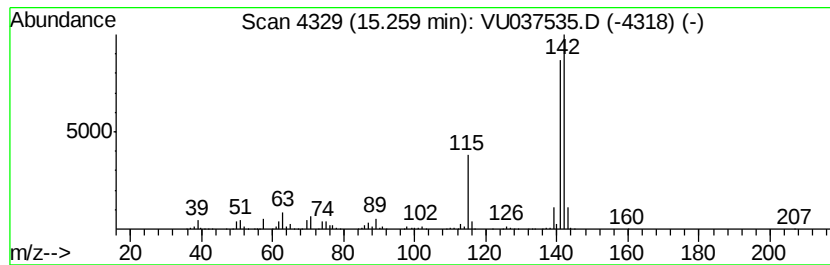
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	28.04 ug/L	1520720	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU040120\
Data File : VU037535.D
Acq On : 01 Apr 2020 21:15
Operator : JC/MD
Sample : L2065-14
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DBF10

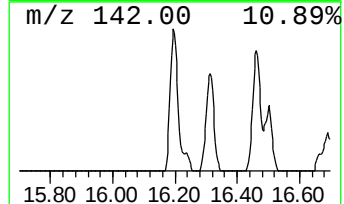
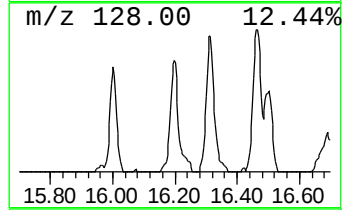
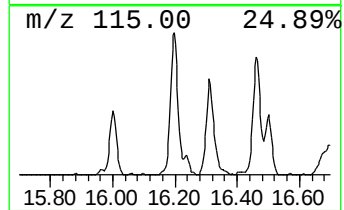
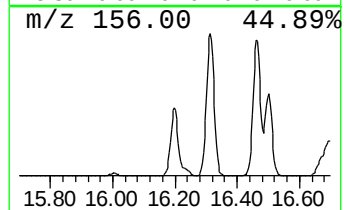
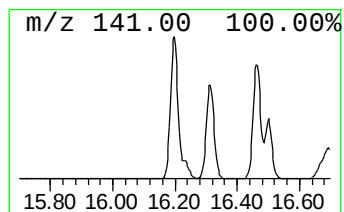
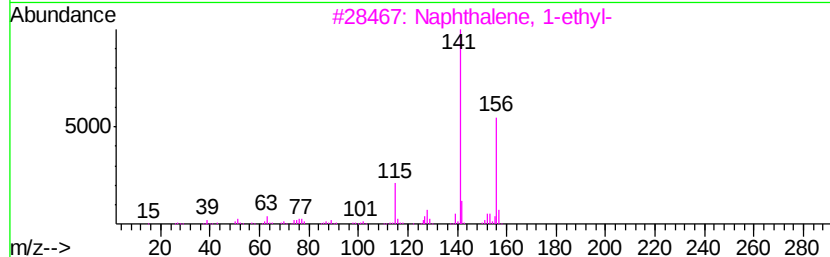
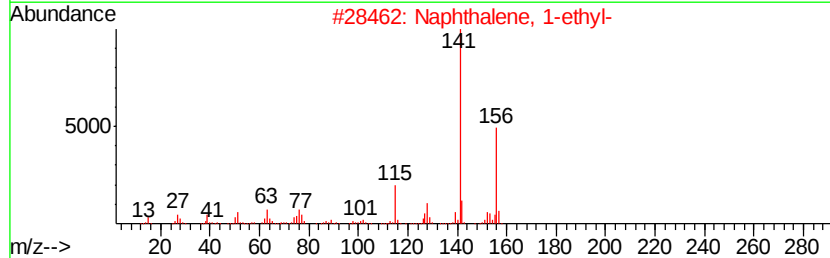
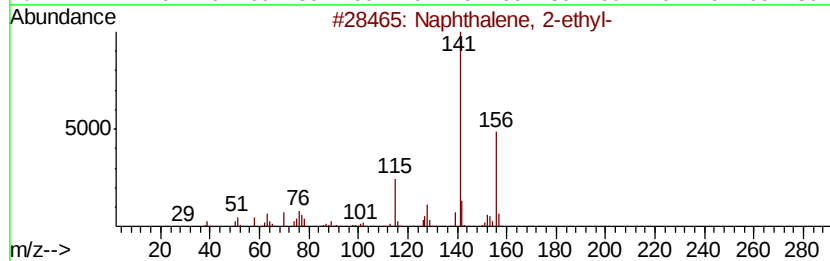
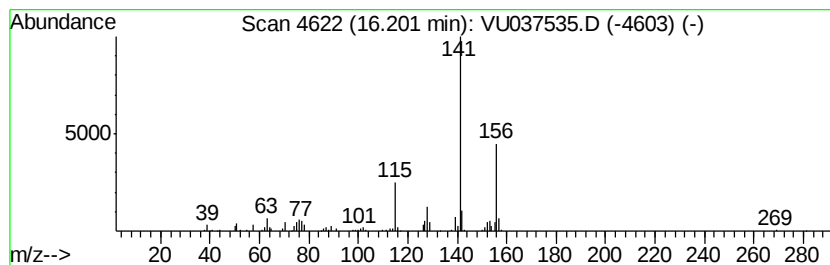
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Naphthalene, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	3.63 ug/L	196868	1,4-Dichlorobenzene-d4	11.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
2			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
3			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
4			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	93
5			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU040120\
Data File : VU037535.D
Acq On : 01 Apr 2020 21:15
Operator : JC/MD
Sample : L2065-14
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

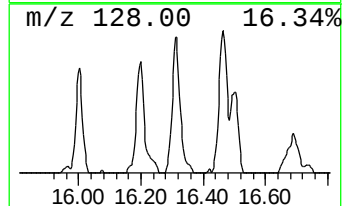
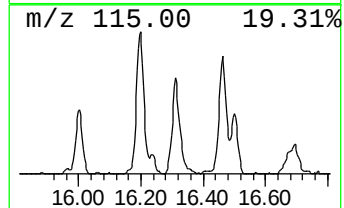
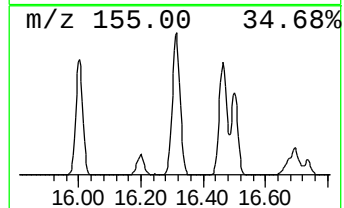
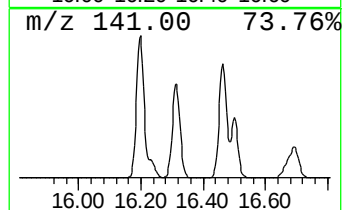
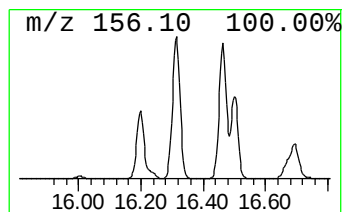
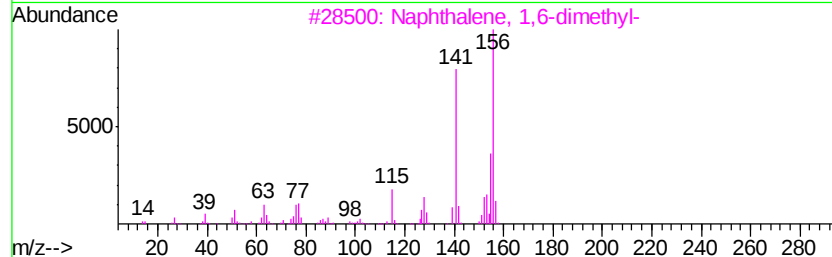
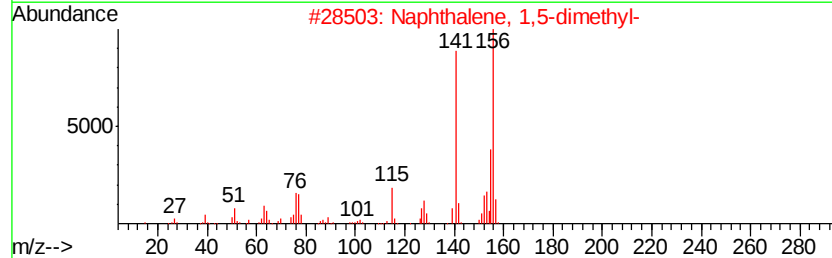
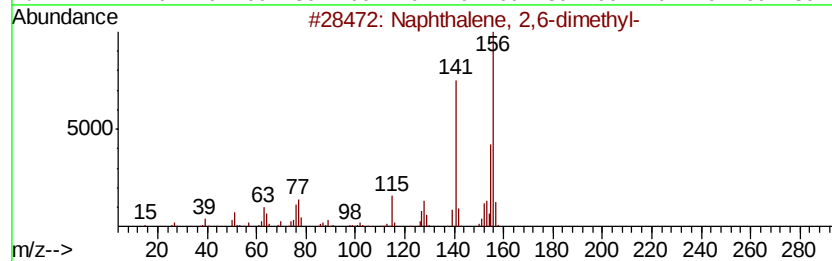
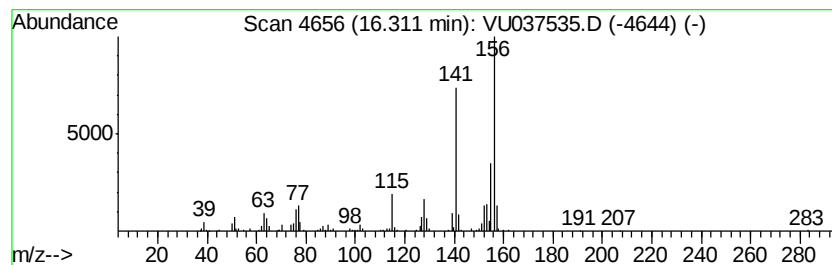
Instrument :
MSVOA_U
ClientSampleId :
DBF10

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

Peak Number 11 Naphthalene, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.31	4.69 ug/L	254102	1,4-Dichlorobenzene-d4	11.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 98
2		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 98
3		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 98
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 98
5		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 98



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU040120\
Data File : VU037535.D
Acq On : 01 Apr 2020 21:15
Operator : JC/MD
Sample : L2065-14
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

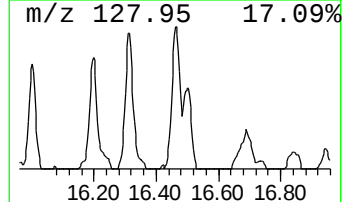
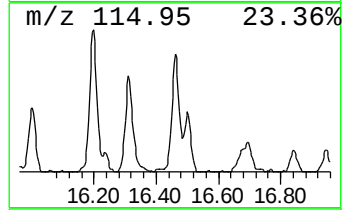
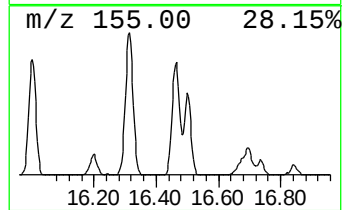
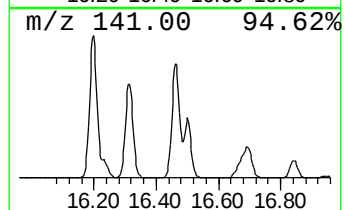
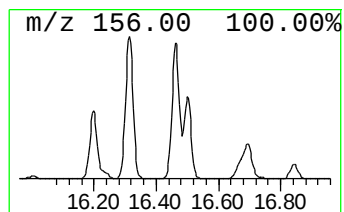
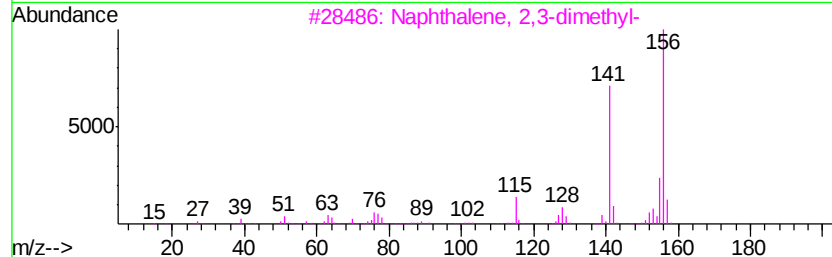
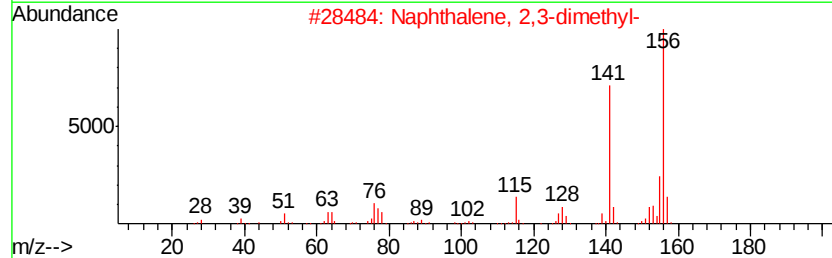
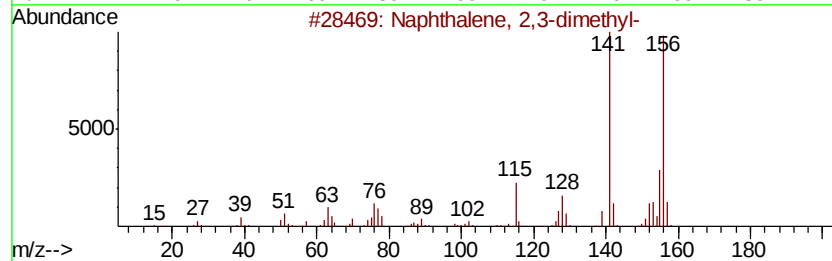
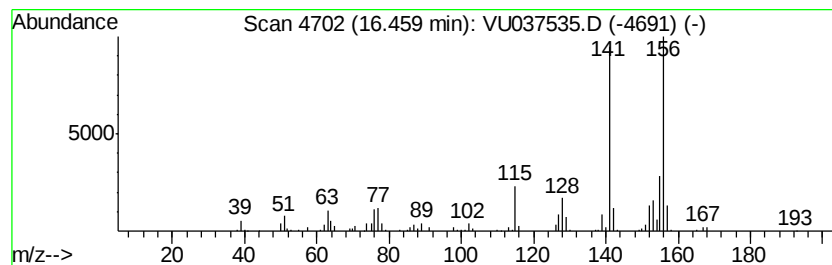
Instrument :
MSVOA_U
ClientSampleId :
DBF10

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

Peak Number 12 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	4.58 ug/L	248479	1,4-Dichlorobenzene-d4	11.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 98
2		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
3		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 96
4		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 96
5		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 96



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU040120\
 Data File : VU037535.D
 Acq On : 01 Apr 2020 21:15
 Operator : JC/MD
 Sample : L2065-14
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DBF10

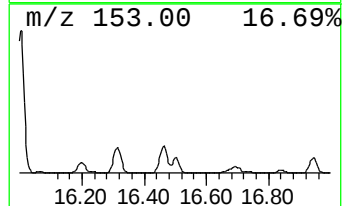
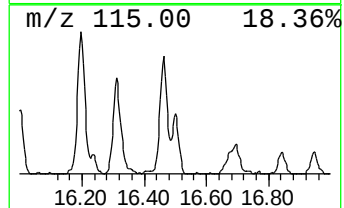
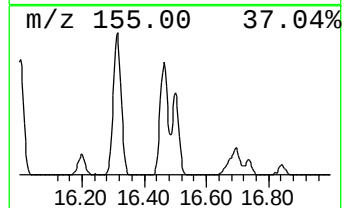
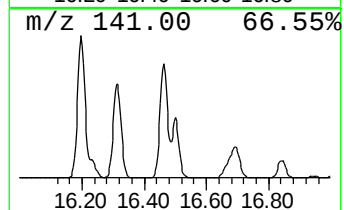
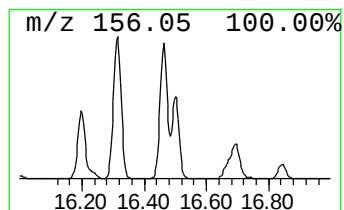
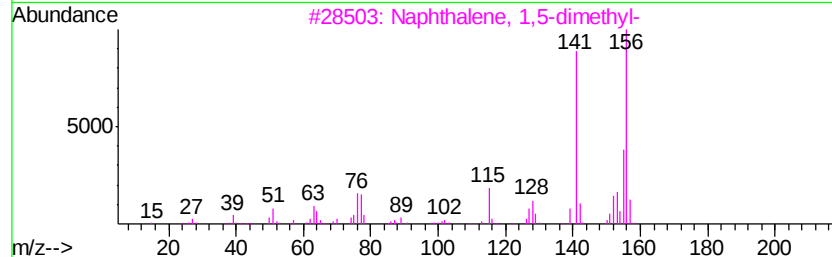
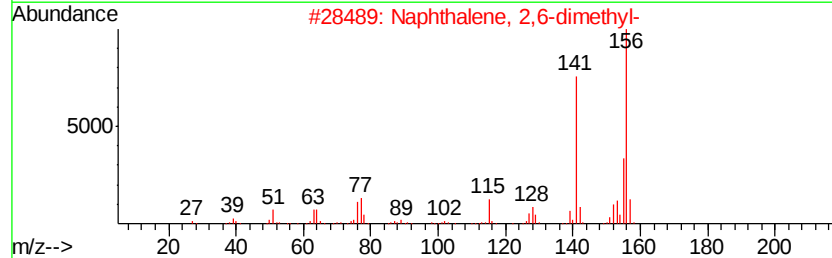
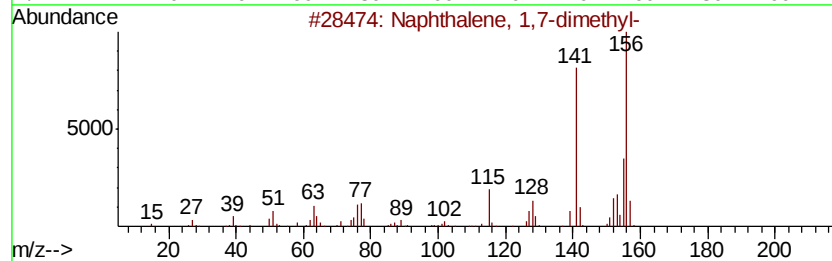
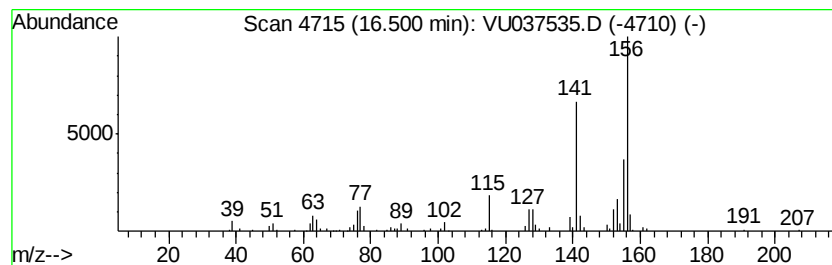
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 Quant Title : TRACE VOA SOM01.0

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

 Peak Number 13 Naphthalene, 1,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	2.35 ug/L	127618	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	93
4		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	93
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU040120\
Data File : VU037535.D
Acq On : 01 Apr 2020 21:15
Operator : JC/MD
Sample : L2065-14
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

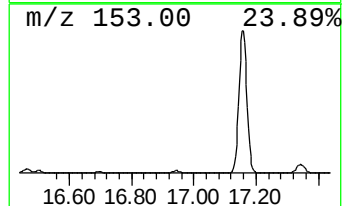
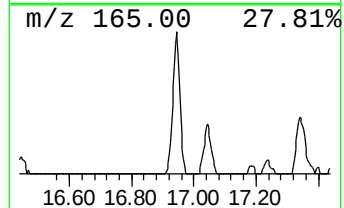
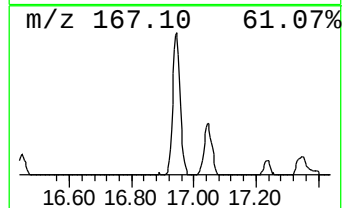
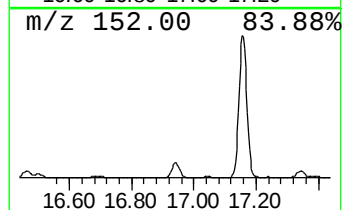
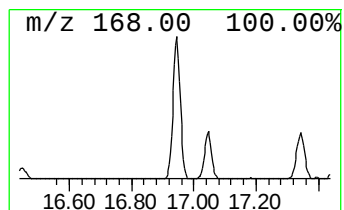
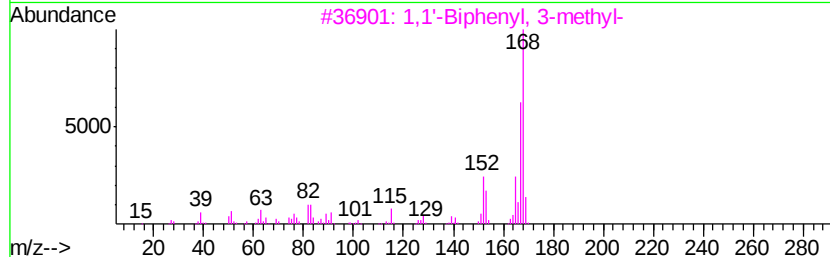
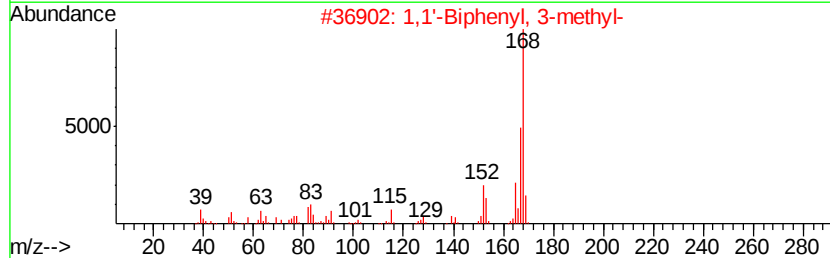
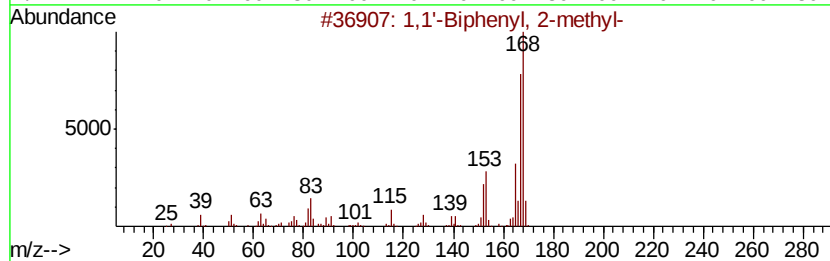
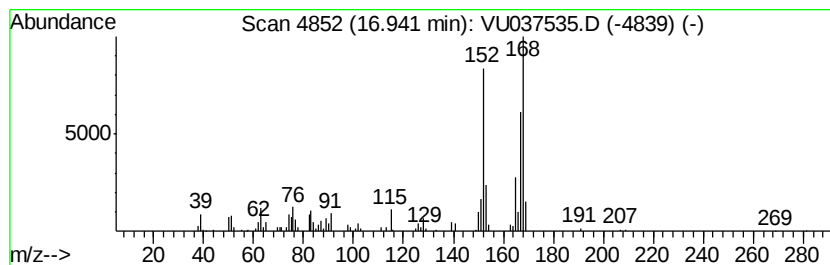
Instrument :
MSVOA_U
ClientSampleId :
DBF10

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

Peak Number 15 1,1'-Biphenyl, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.94	1.92 ug/L	103905	1,4-Dichlorobenzene-d4	11.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	92
2	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	64
3	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	60
4	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	60
5	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040120\
 Data File : VU037535.D
 Acq On : 01 Apr 2020 21:15
 Operator : JC/MD
 Sample : L2065-14
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DBF10

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	2.2	ug/L	86118	1	6.29	196305	5.0
Indane	12.11	10.4	ug/L	562575	3	11.83	271130	5.0
Indene	12.33	16.1	ug/L	872590	3	11.83	271130	5.0
Benzofuran, 7-met...	13.05	2.1	ug/L	113348	3	11.83	271130	5.0
Benzo[b]thiophene	14.24	3.4	ug/L	181750	3	11.83	271130	5.0
Naphthalene, 1-me...	15.26	28.0	ug/L	1520720	3	11.83	271130	5.0
Naphthalene, 2-et...	16.20	3.6	ug/L	196868	3	11.83	271130	5.0
Naphthalene, 2,6-...	16.31	4.7	ug/L	254102	3	11.83	271130	5.0
Naphthalene, 2,3-...	16.46	4.6	ug/L	248479	3	11.83	271130	5.0
Naphthalene, 1,7-...	16.50	2.4	ug/L	127618	3	11.83	271130	5.0
1,1'-Biphenyl, 2-...	16.94	1.9	ug/L	103905	3	11.83	271130	5.0