

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110124\
 Data File : VU061388.D
 Acq On : 02 Nov 2024 02:29
 Operator : MD/SY
 Sample : P4668-01
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AD1

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.1

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VU061388.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.560	383	395	412	rBV	87997	142927	1.13%	0.693%
2	4.624	1028	1037	1062	rBV	63677	165701	1.31%	0.803%
3	5.052	1156	1170	1196	rBV	90155	203095	1.61%	0.985%
4	5.717	1355	1377	1410	rBV3	182353	501695	3.97%	2.432%
5	6.241	1528	1540	1569	rBV	193443	383935	3.04%	1.861%
6	6.682	1661	1677	1695	rBV2	114092	226367	1.79%	1.097%
7	7.891	2033	2053	2071	rBV	227440	400933	3.17%	1.944%
8	8.627	2271	2282	2321	rBV2	245082	493538	3.91%	2.393%
9	9.409	2513	2525	2546	rBV	274177	476556	3.77%	2.310%
10	10.746	2930	2941	2961	rBV	88728	144436	1.14%	0.700%
11	11.804	3256	3270	3288	rBV	284430	463265	3.67%	2.246%
12	12.087	3342	3358	3376	rBV	8061167	12634648	100.00%	61.252%
13	12.183	3377	3388	3406	rVB	253196	433366	3.43%	2.101%
14	12.601	3509	3518	3529	rVB	102776	157456	1.25%	0.763%
15	12.672	3529	3540	3559	rVB	653571	1038832	8.22%	5.036%
16	12.916	3605	3616	3626	rBV	170361	270642	2.14%	1.312%
17	13.103	3665	3674	3688	rVB	182934	304370	2.41%	1.476%
18	13.283	3722	3730	3738	rBV	132328	203291	1.61%	0.986%
19	13.444	3767	3780	3792	rBV3	316820	520229	4.12%	2.522%
20	13.617	3814	3834	3848	rVB3	124119	215304	1.70%	1.044%
21	13.929	3926	3931	3943	rVB	103524	155721	1.23%	0.755%
22	14.077	3958	3977	3994	rBV3	349970	940078	7.44%	4.557%
23	14.871	4209	4224	4236	rVB3	82694	150984	1.19%	0.732%

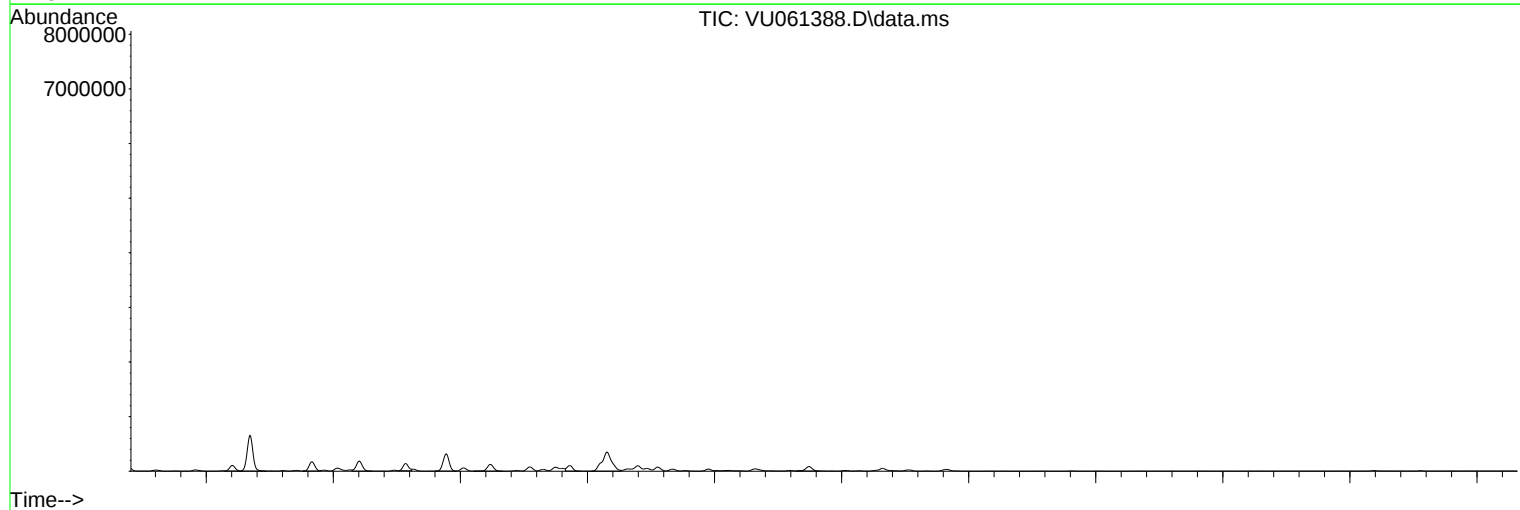
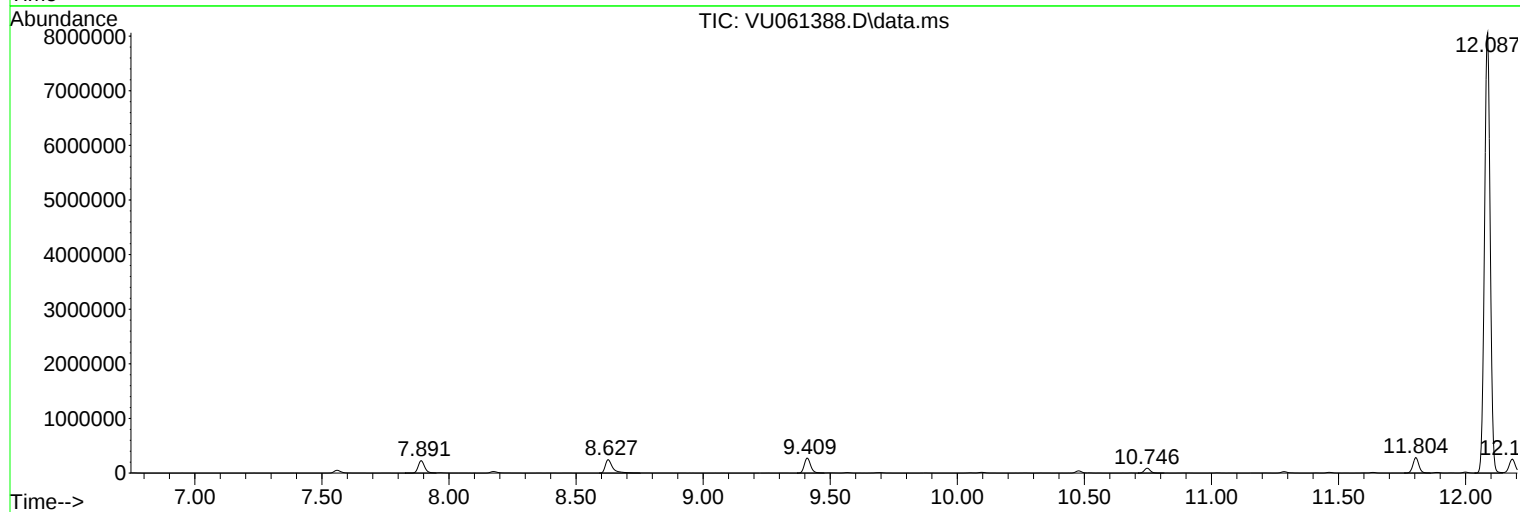
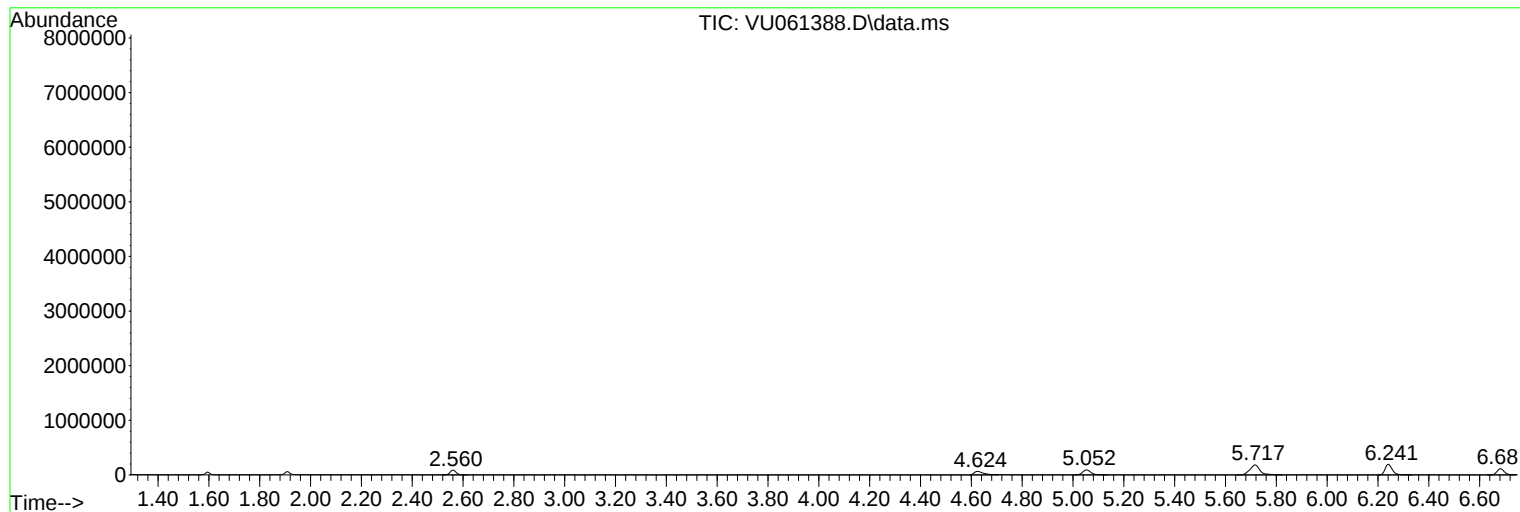
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ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0AD1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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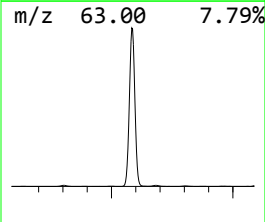
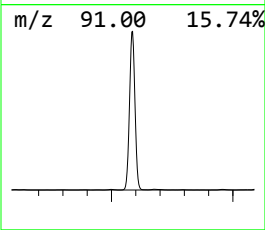
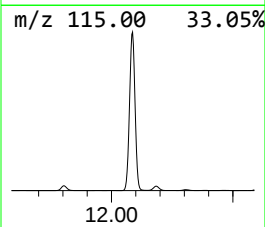
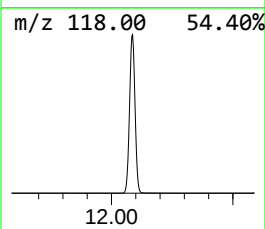
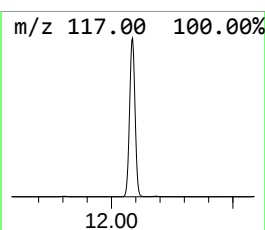
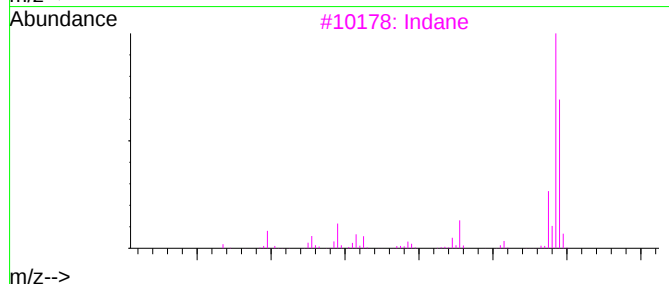
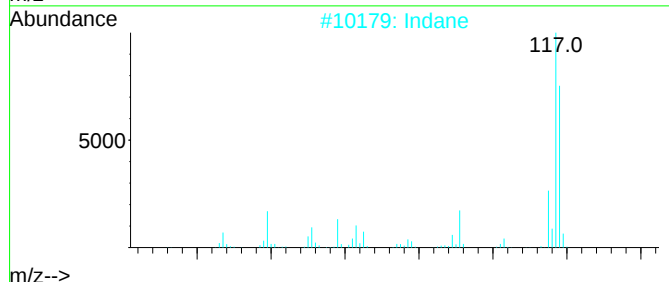
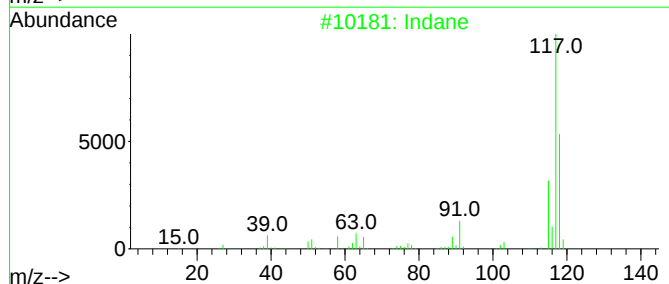
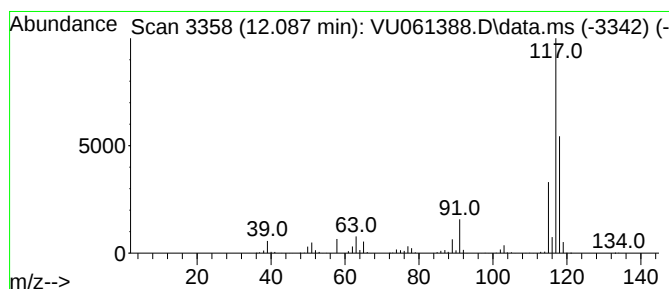
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.087	136.37 ug/L	12634600	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	83
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
5	Deltacyclene	118	C9H10	007785-10-6	72



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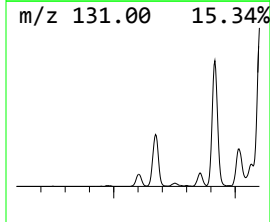
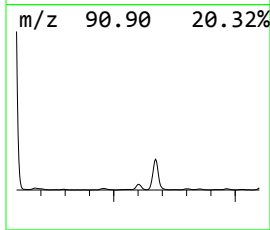
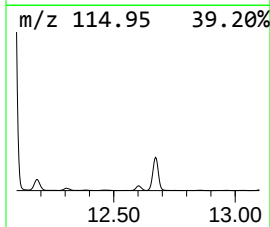
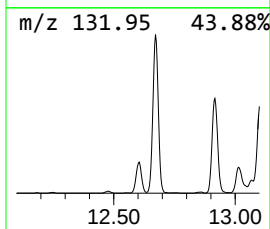
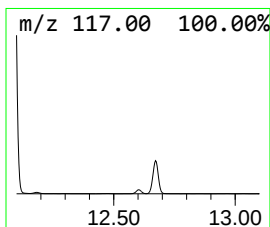
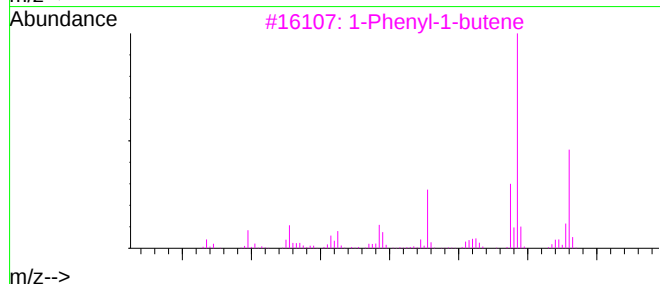
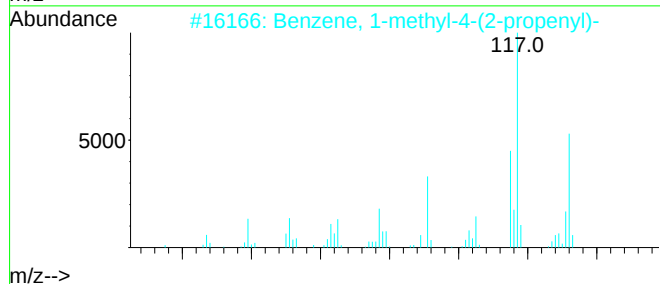
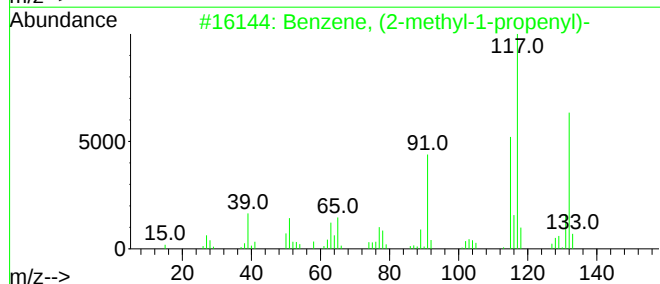
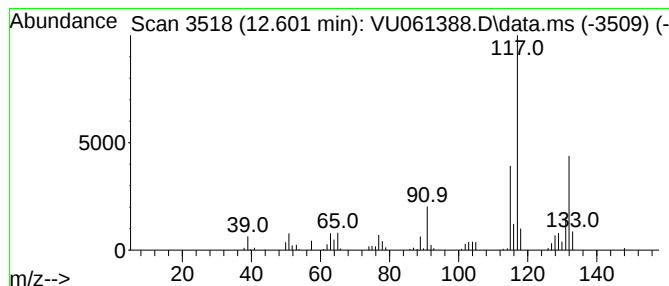
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.601	1.70 ug/L	157456	1,4-Dichlorobenzene-d4	11.804	
Hit# of 5	Tentative ID		MW MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-		132 C10H12	000768-49-0	94
2	Benzene, 1-methyl-4-(2-propenyl)-		132 C10H12	003333-13-9	94
3	1-Phenyl-1-butene		132 C10H12	000824-90-8	94
4	Benzene, (2-methyl-2-propenyl)-		132 C10H12	003290-53-7	94
5	Benzene, 1-methyl-2-(2-propenyl)-		132 C10H12	001587-04-8	91



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ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

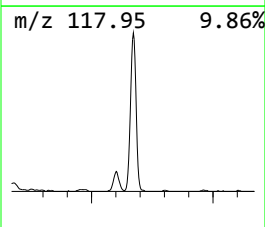
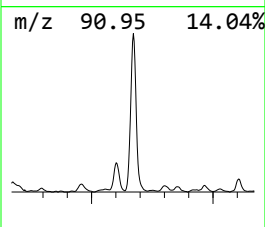
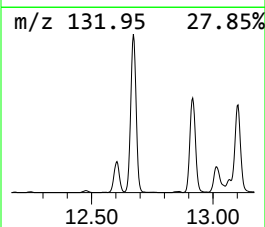
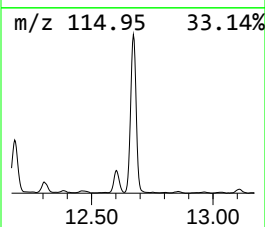
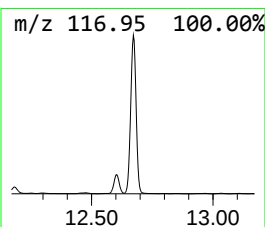
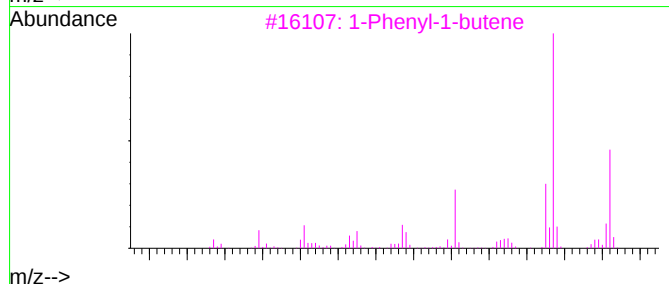
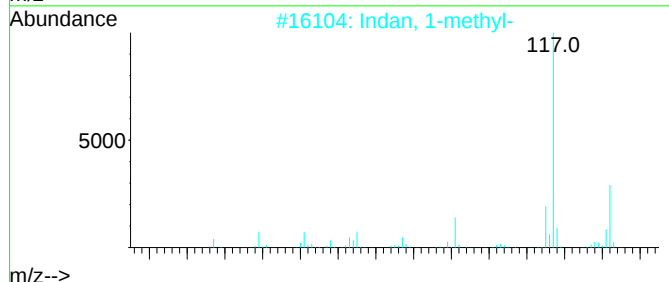
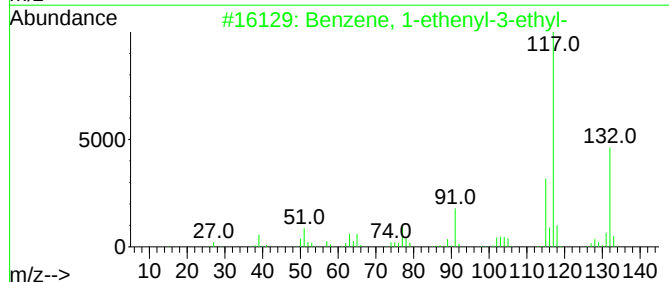
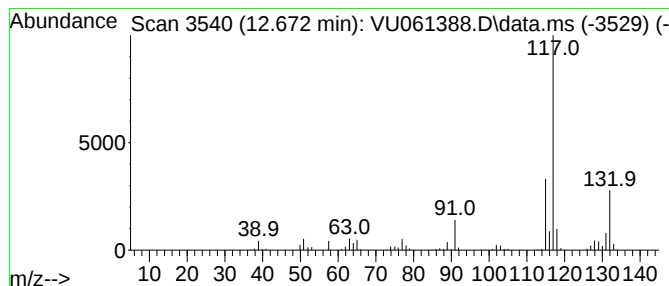
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.672	11.21 ug/L	1038830	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86



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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

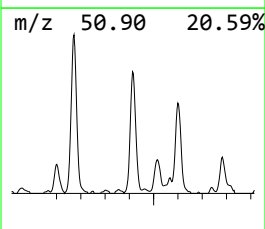
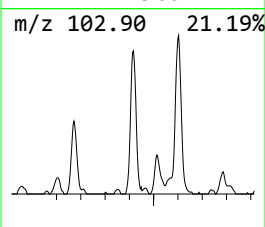
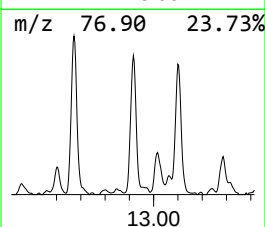
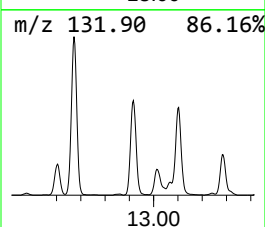
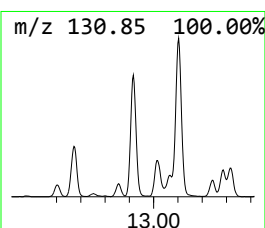
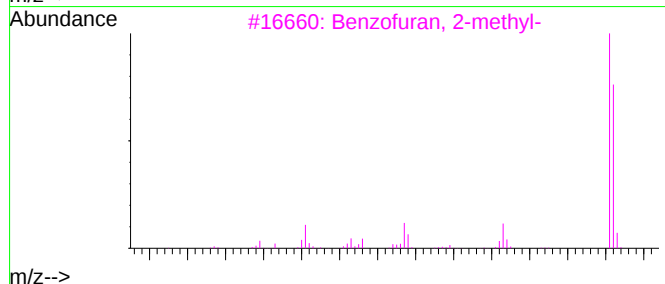
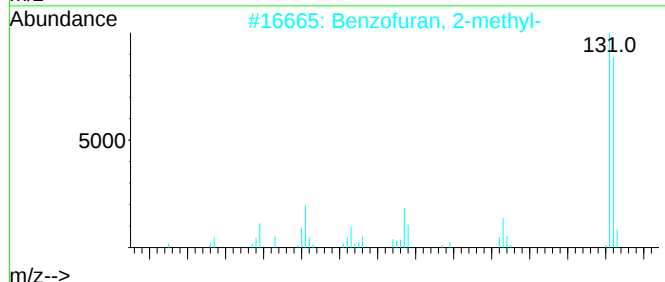
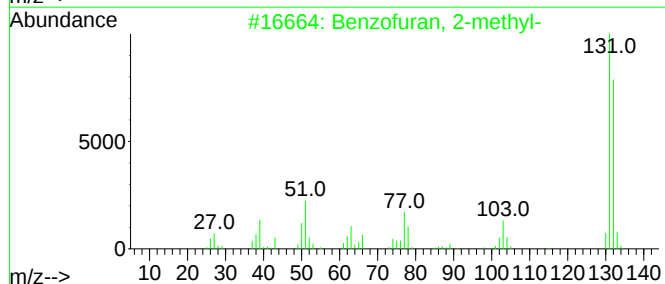
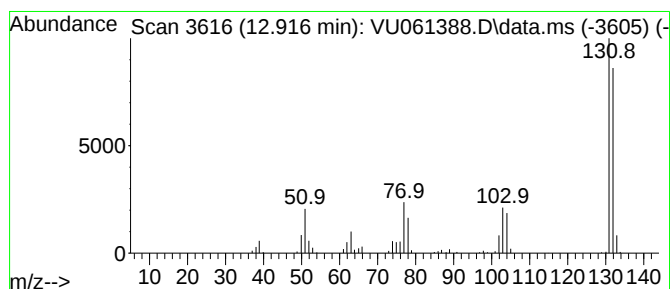
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzofuran, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.916	2.92 ug/L	270642	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91



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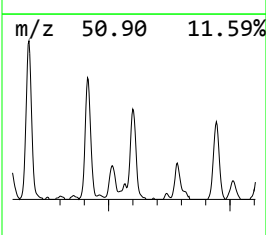
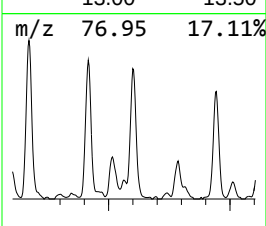
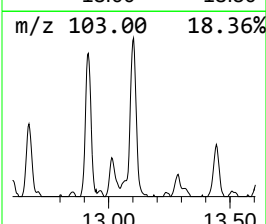
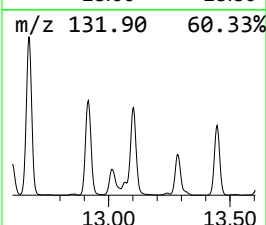
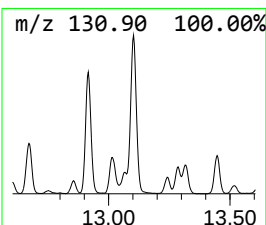
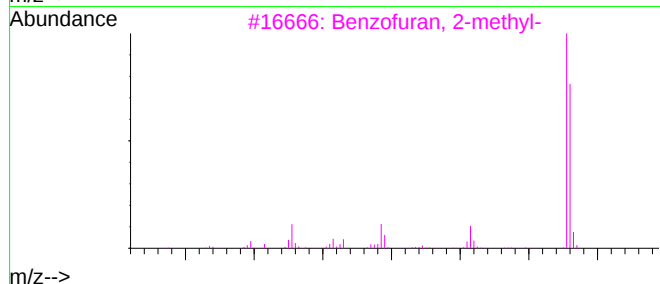
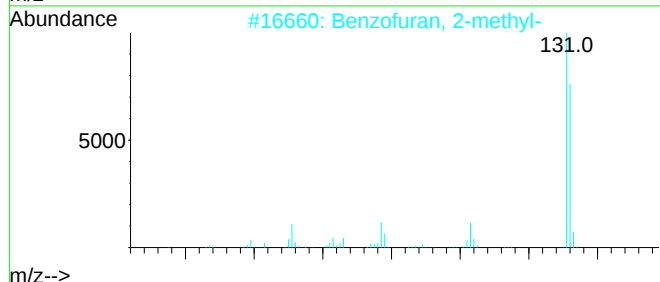
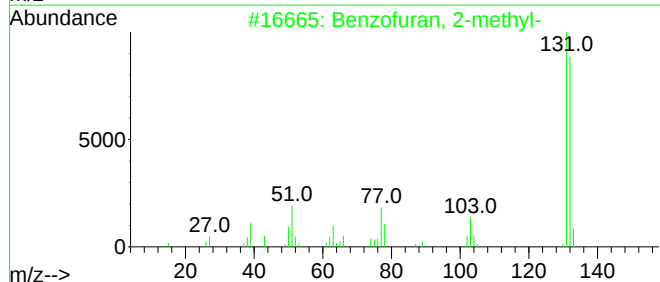
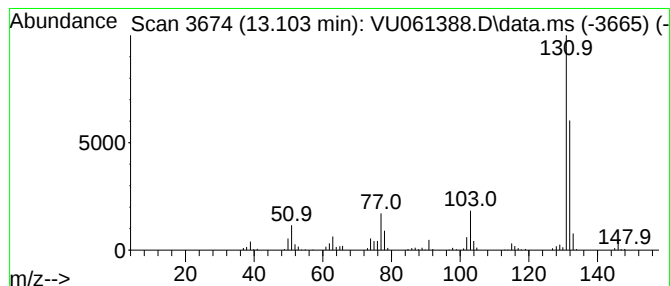
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Propenal, 3-phenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.103	3.29 ug/L	304370	1,4-Dichlorobenzene-d4		11.804	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-		132	C9H8O	004265-25-2	94
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	87
5	Cinnamaldehyde, (E)-		132	C9H8O	014371-10-9	87



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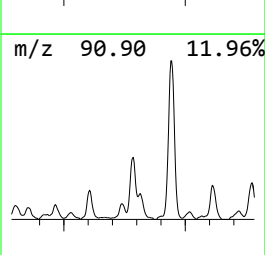
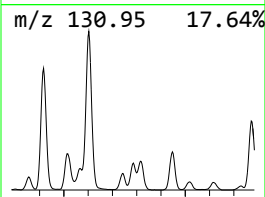
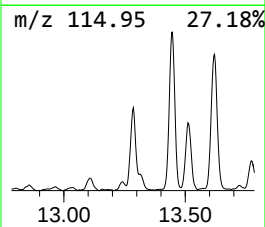
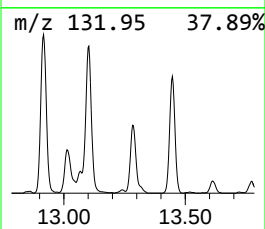
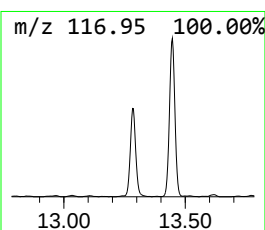
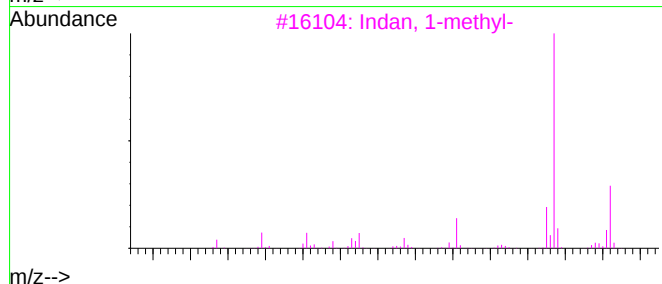
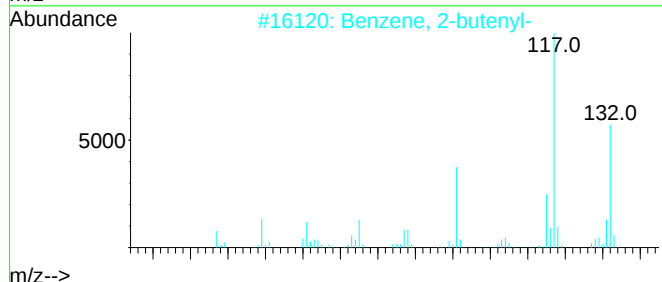
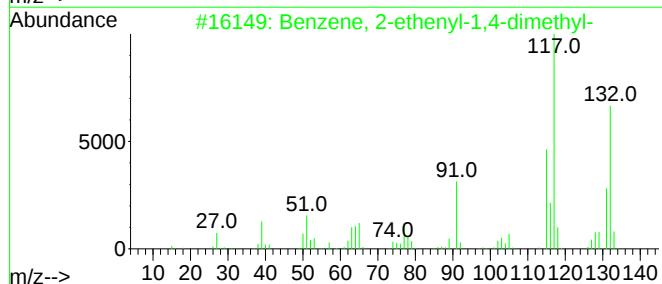
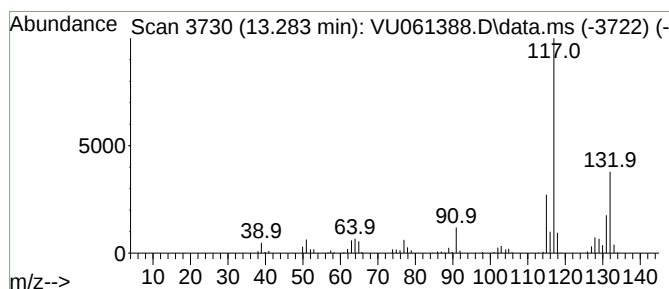
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.283	2.19 ug/L	203291	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
3	Indan, 1-methyl-	132	C10H12	000767-58-8	90
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110124\
Data File : VU061388.D
Acq On : 02 Nov 2024 02:29
Operator : MD/SY
Sample : P4668-01
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AD1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

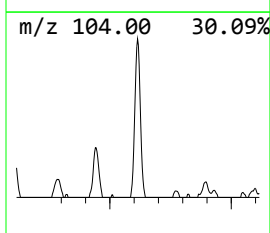
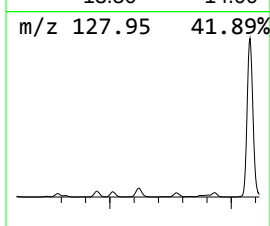
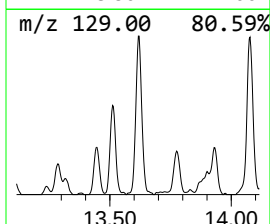
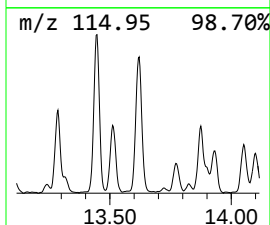
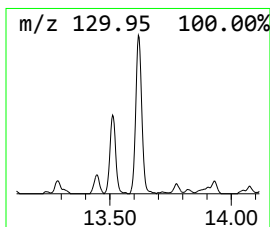
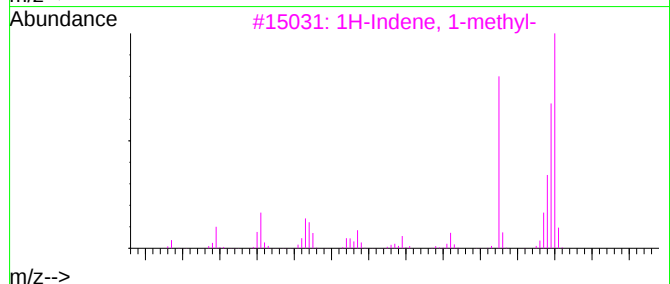
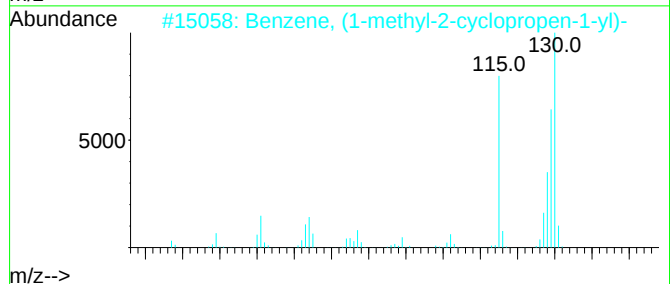
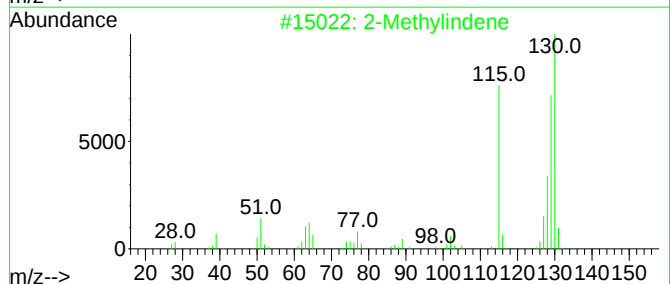
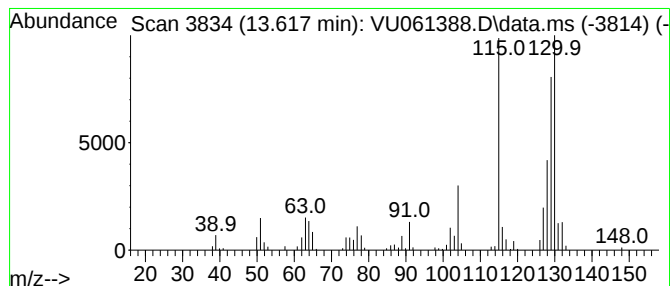
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 2-Methylindene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.617	2.32 ug/L	215304	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene	130	C10H10	002177-47-1	96
2	Benzene, (1-methyl-2-cyclopropen-1-yl)-	130	C10H10	065051-83-4	96
3	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
4	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	95
5	Tetracyclo[5.3.0.0<2,6>.0<3,10>]...	130	C10H10	034324-40-8	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110124\
Data File : VU061388.D
Acq On : 02 Nov 2024 02:29
Operator : MD/SY
Sample : P4668-01
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AD1

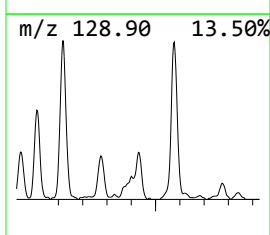
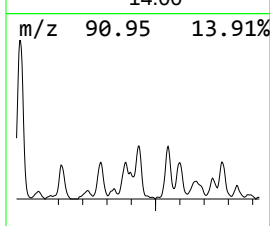
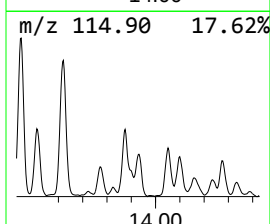
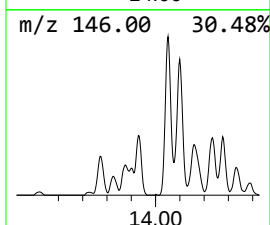
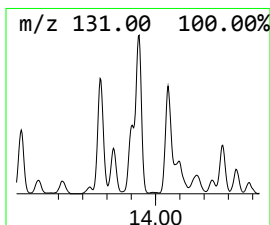
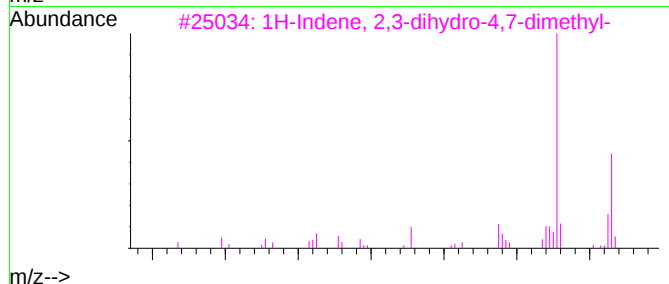
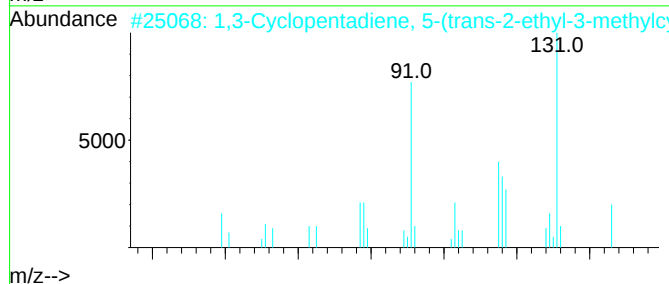
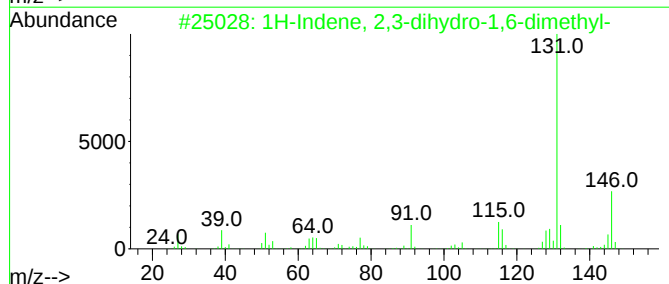
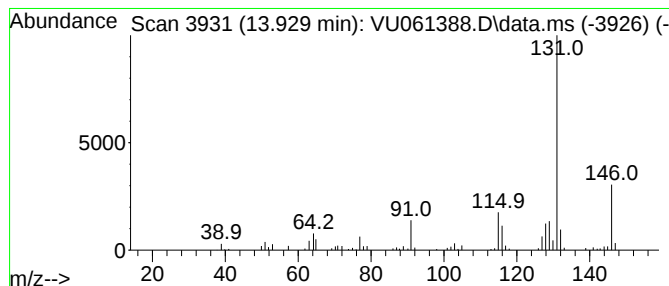
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.929	1.68 ug/L	155721	1,4-Dichlorobenzene-d4	11.804	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2	1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	91
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110124\
Data File : VU061388.D
Acq On : 02 Nov 2024 02:29
Operator : MD/SY
Sample : P4668-01
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AD1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

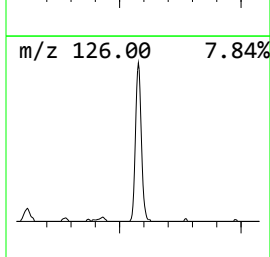
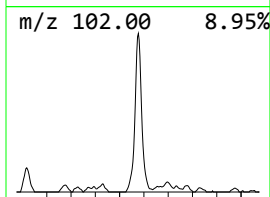
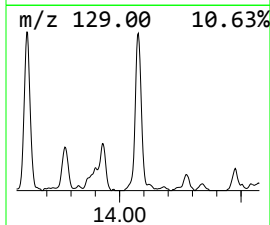
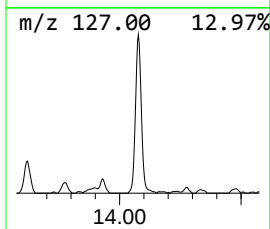
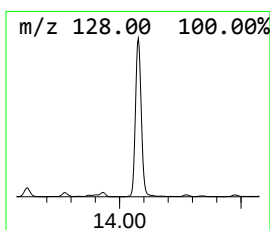
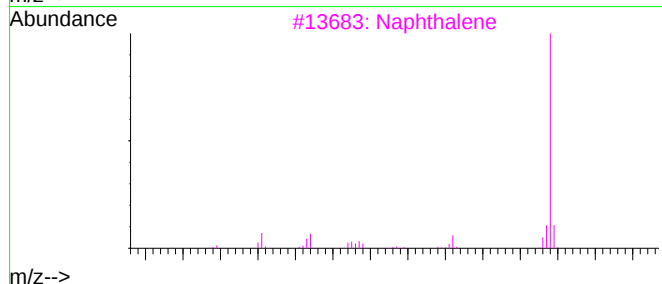
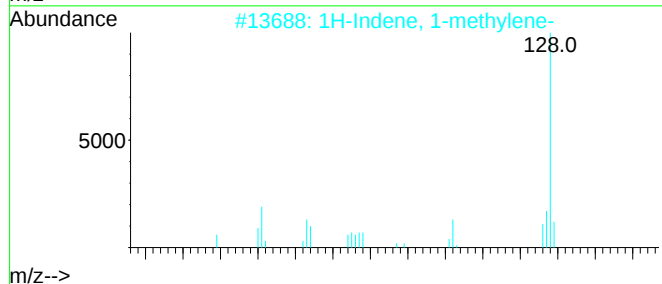
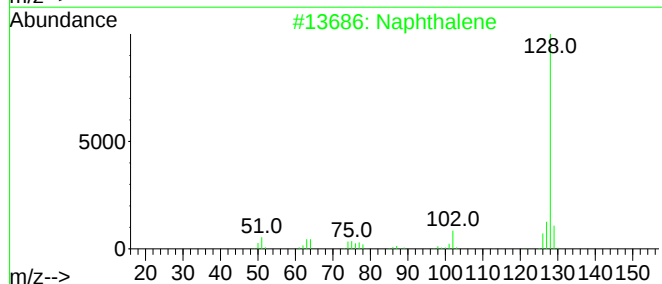
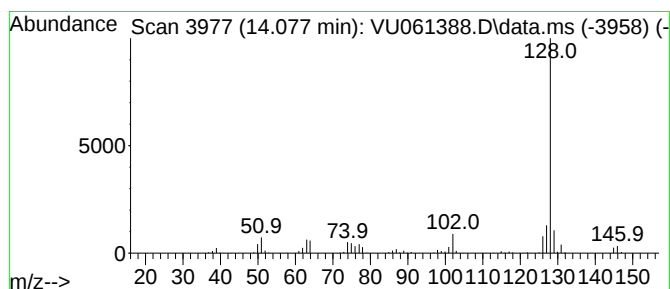
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 1-methylene- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.077	10.15 ug/L	940078	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	94
2	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	93
3	Naphthalene	128	C10H8	000091-20-3	90
4	Azulene	128	C10H8	000275-51-4	90
5	Azulene	128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110124\
Data File : VU061388.D
Acq On : 02 Nov 2024 02:29
Operator : MD/SY
Sample : P4668-01
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AD1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

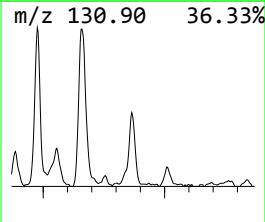
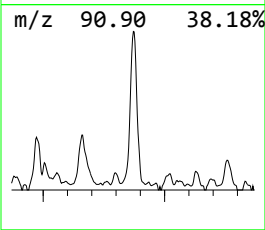
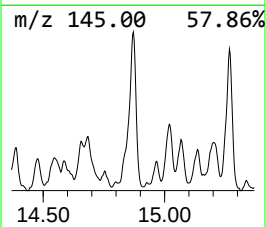
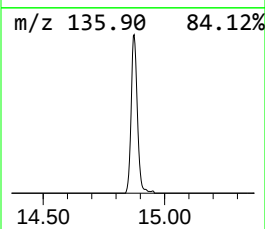
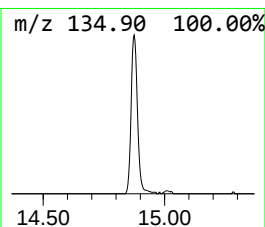
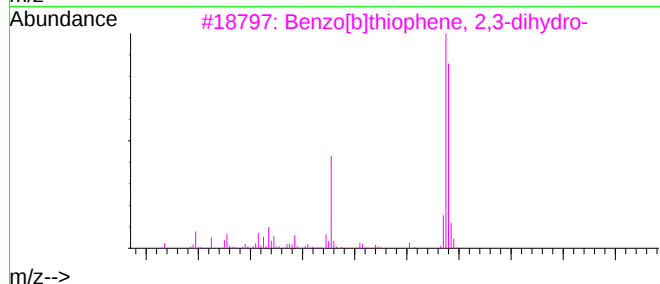
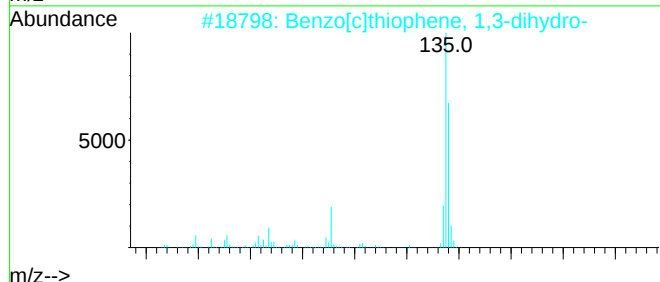
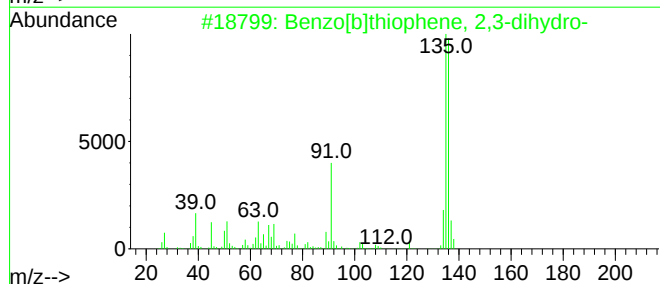
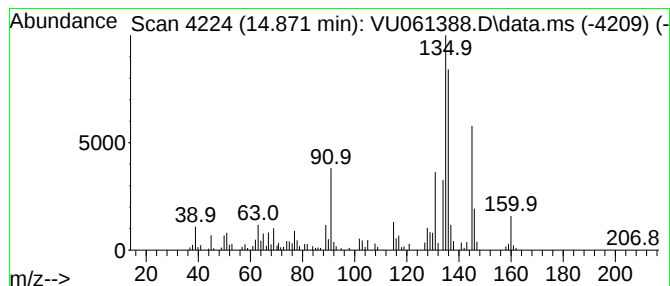
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.871	1.63 ug/L	150984	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	92
2	Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	60
3	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	53
4	Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	53
5	Benzaldehyde, 3-methoxy-	136	C8H8O2	000591-31-1	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110124\
Data File : VU061388.D
Acq On : 02 Nov 2024 02:29
Operator : MD/SY
Sample : P4668-01
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AD1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	12.087	136.4	ug/L	12634600	3	11.804	463265	5.0
Benzene, (2-met...	12.601	1.7	ug/L	157456	3	11.804	463265	5.0
Benzene, 1-ethe...	12.672	11.2	ug/L	1038830	3	11.804	463265	5.0
Benzofuran, 2-m...	12.916	2.9	ug/L	270642	3	11.804	463265	5.0
2-Propenal, 3-p...	13.103	3.3	ug/L	304370	3	11.804	463265	5.0
Benzene, 2-ethe...	13.283	2.2	ug/L	203291	3	11.804	463265	5.0
2-Methylindene	13.617	2.3	ug/L	215304	3	11.804	463265	5.0
1H-Indene, 2,3-...	13.929	1.7	ug/L	155721	3	11.804	463265	5.0
1H-Indene, 1-me...	14.077	10.2	ug/L	940078	3	11.804	463265	5.0
Benzo[b]thiophe...	14.871	1.6	ug/L	150984	3	11.804	463265	5.0