

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
 Data File : VU061507.D
 Acq On : 06 Nov 2024 17:22
 Operator : MD/SY
 Sample : P4714-03DL 4X
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AF5DL

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: VU061507.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	382	395	414	rBV	109822	179607	1.78%	0.647%
2	4.624	1024	1037	1070	rBV	76236	202240	2.00%	0.729%
3	5.052	1156	1170	1197	rBV	101593	227583	2.25%	0.820%
4	5.718	1358	1377	1389	rBV2	211260	551284	5.45%	1.987%
5	6.242	1527	1540	1561	rBV	167650	326026	3.23%	1.175%
6	6.682	1664	1677	1697	rBV	125867	252094	2.49%	0.909%
7	7.891	2040	2053	2069	rBV	254556	443078	4.38%	1.597%
8	8.627	2271	2282	2322	rBV2	272328	531657	5.26%	1.916%
9	9.412	2514	2526	2553	rBV	237344	412607	4.08%	1.487%
10	9.563	2561	2573	2595	rBV	341389	552959	5.47%	1.993%
11	9.688	2596	2612	2632	rVB	167906	295594	2.92%	1.065%
12	10.094	2728	2738	2762	rVB	135946	218037	2.16%	0.786%
13	10.476	2844	2857	2881	rBV	88952	144592	1.43%	0.521%
14	10.746	2929	2941	2959	rBV	98246	159063	1.57%	0.573%
15	11.081	3036	3045	3061	rVB	68119	103956	1.03%	0.375%
16	11.283	3097	3108	3127	rBV	78702	129913	1.29%	0.468%
17	11.460	3150	3163	3178	rBV	332398	513119	5.08%	1.850%
18	11.804	3257	3270	3286	rVV	254697	429477	4.25%	1.548%
19	11.888	3286	3296	3306	rVB	67784	103165	1.02%	0.372%
20	12.087	3344	3358	3377	rBV	6547775	10108719	100.00%	36.437%
21	12.183	3377	3388	3406	rVB	294044	489562	4.84%	1.765%
22	12.563	3495	3506	3513	rBV	73286	112732	1.12%	0.406%
23	12.672	3528	3540	3560	rVB	314237	508680	5.03%	1.834%
24	12.917	3605	3616	3625	rBV	357713	559422	5.53%	2.016%
25	13.019	3638	3648	3659	rBV	466641	903068	8.93%	3.255%
26	13.100	3667	3673	3689	rVB	132137	213601	2.11%	0.770%
27	13.283	3720	3730	3747	rVB	298309	464615	4.60%	1.675%
28	13.447	3759	3781	3793	rBV2	649389	1042538	10.31%	3.758%
29	14.077	3959	3977	3995	rBV	3626761	5937207	58.73%	21.401%
30	14.196	3996	4014	4033	rVV	569025	1050665	10.39%	3.787%
31	15.216	4322	4331	4356	rVB	99131	186789	1.85%	0.673%
32	15.402	4377	4389	4419	rVB	232066	389546	3.85%	1.404%

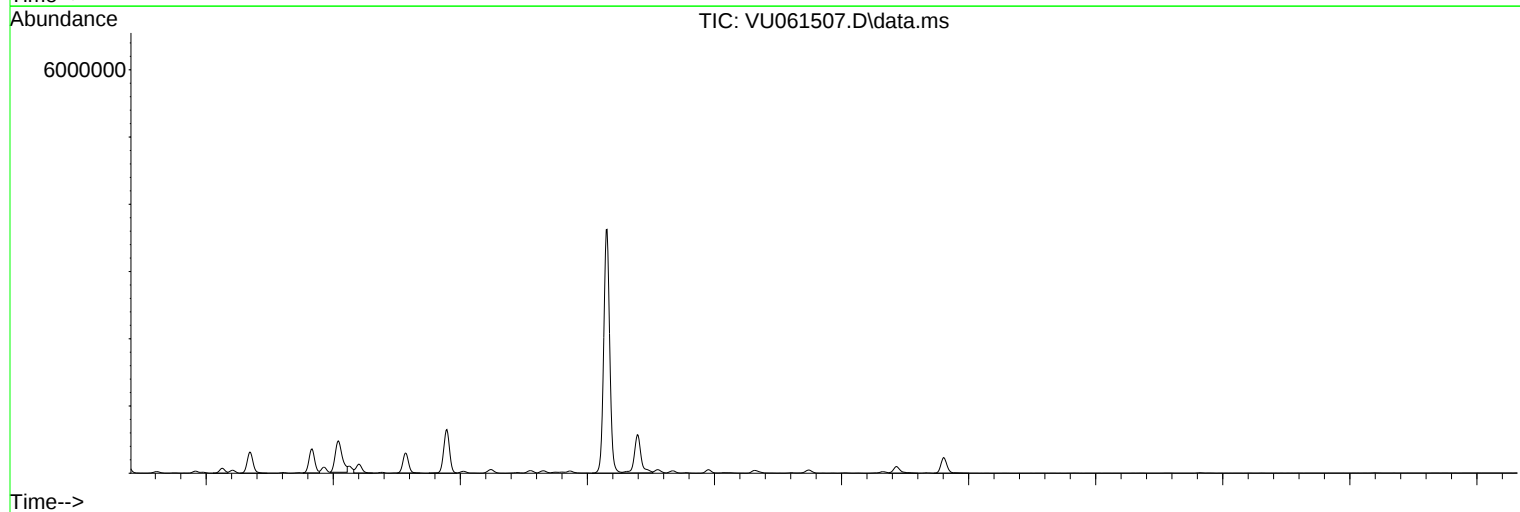
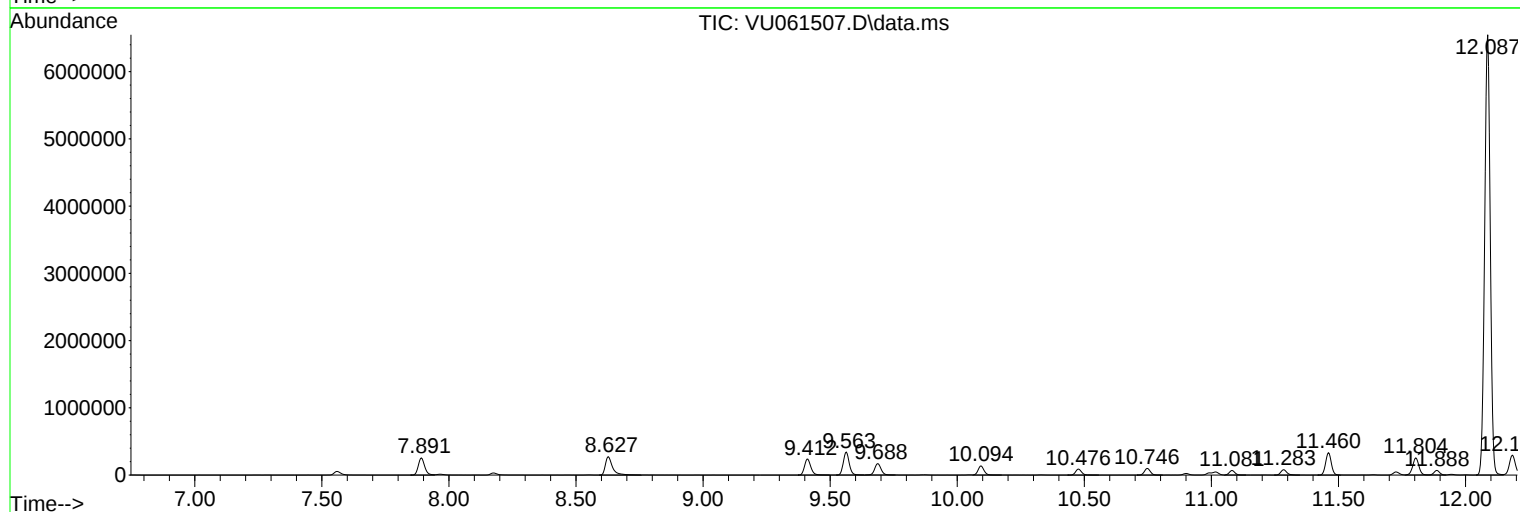
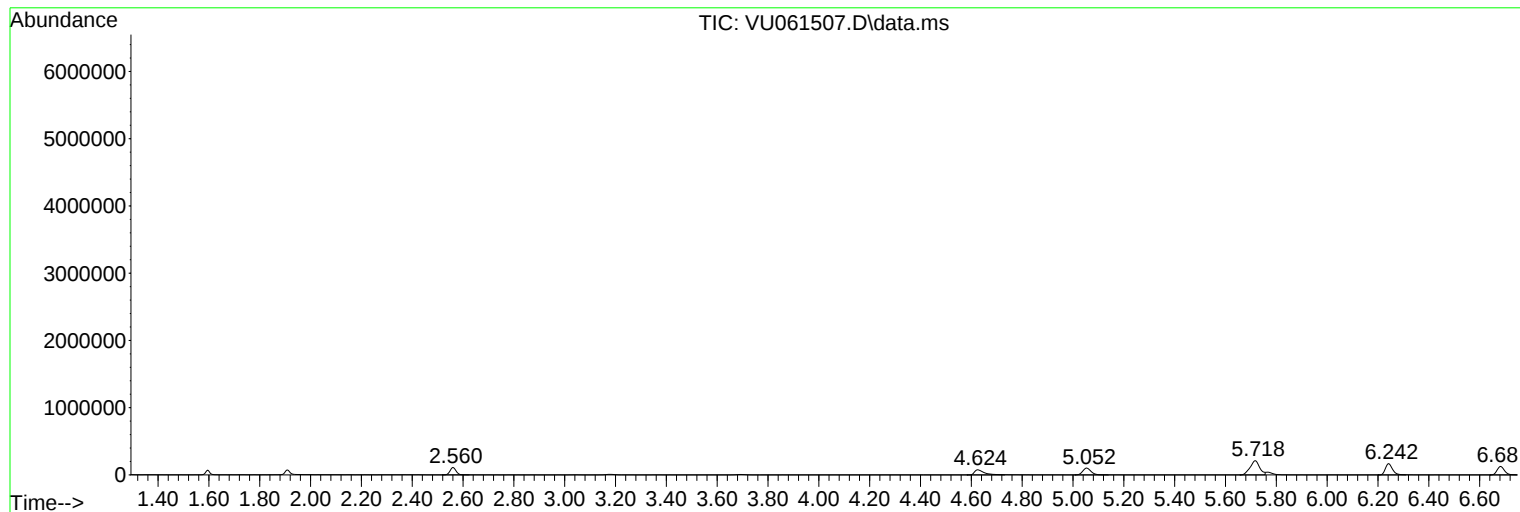
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0AF5DL

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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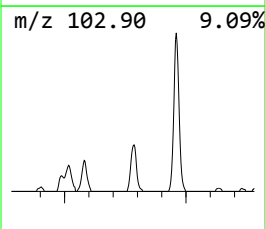
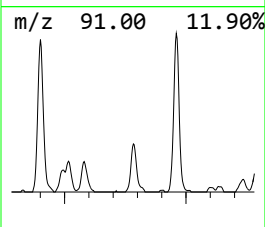
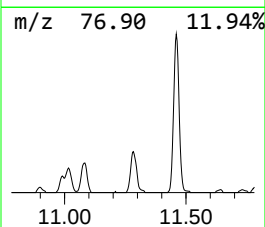
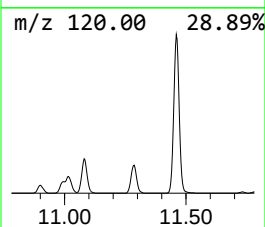
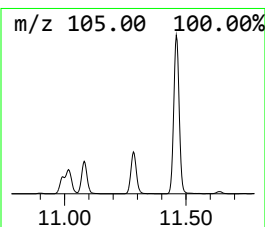
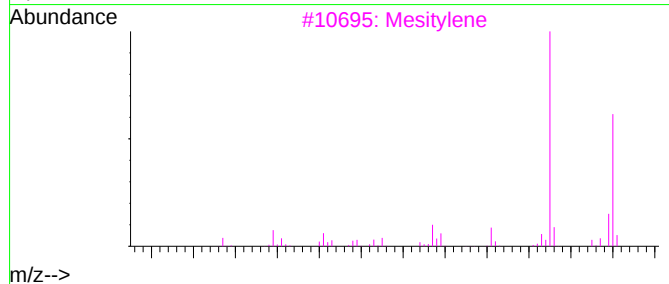
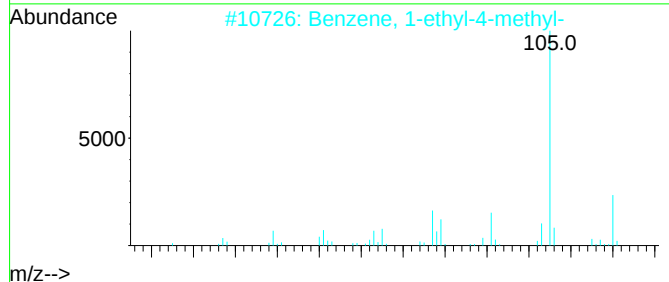
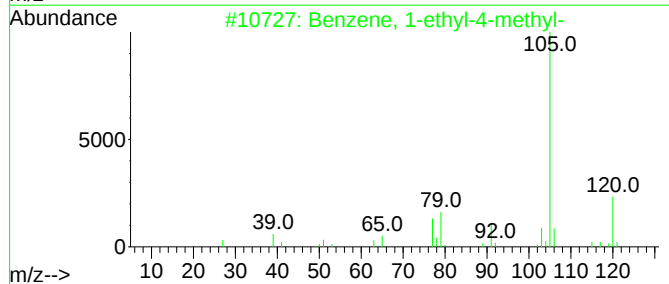
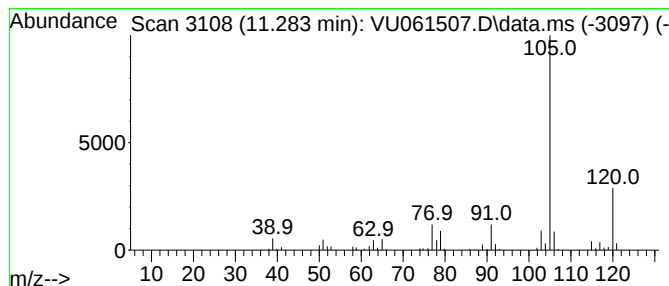
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.283	1.51 ug/L	129913	1,4-Dichlorobenzene-d4		11.804	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-		120	C9H12	000622-96-8	94
2	Benzene, 1-ethyl-4-methyl-		120	C9H12	000622-96-8	91
3	Mesitylene		120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-		120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-		120	C9H12	000620-14-4	91



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Instrument :
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ClientSampleId :
C0AF5DL

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

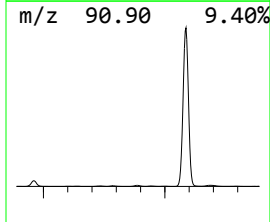
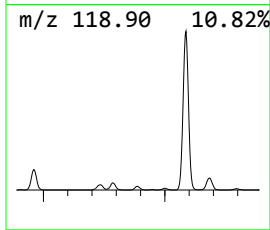
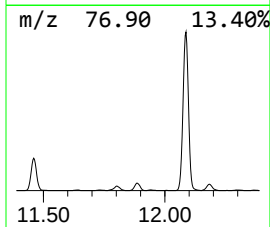
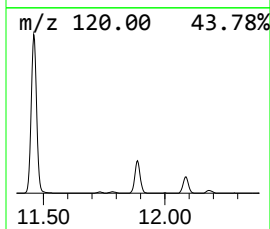
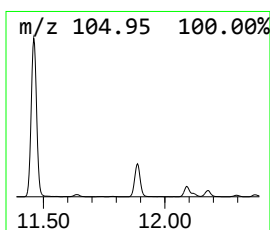
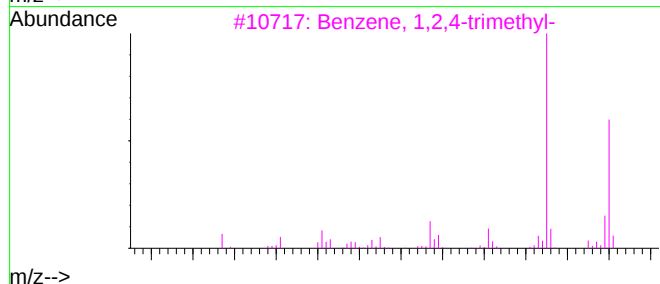
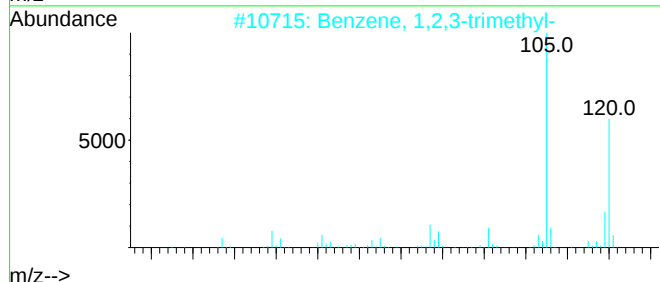
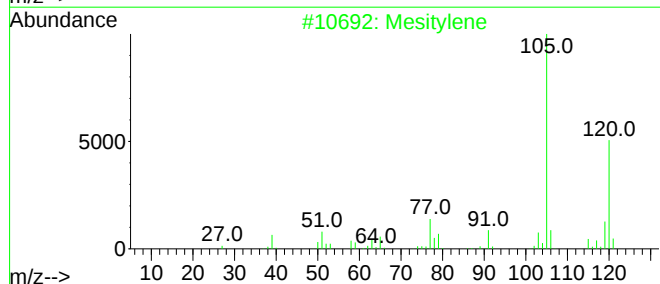
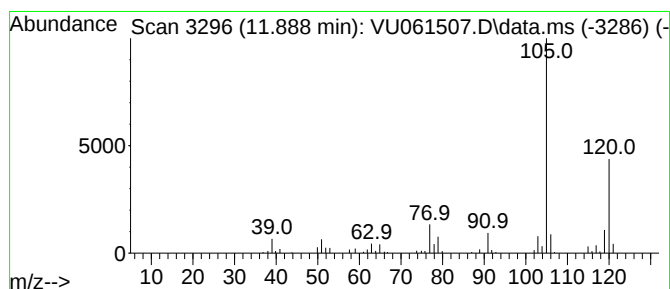
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 2 Mesitylene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.888	1.20 ug/L	103165	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Mesitylene	120	C9H12	000108-67-8	93



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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AF5DL

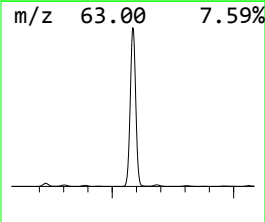
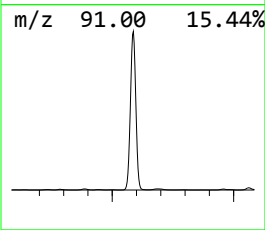
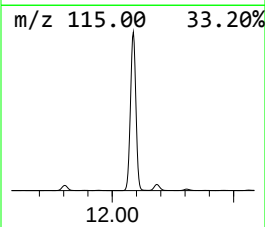
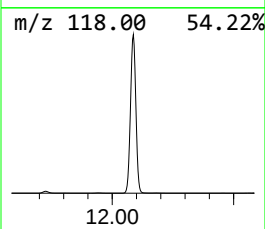
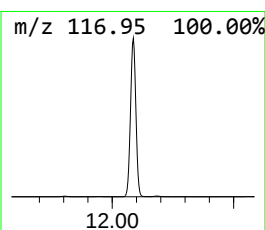
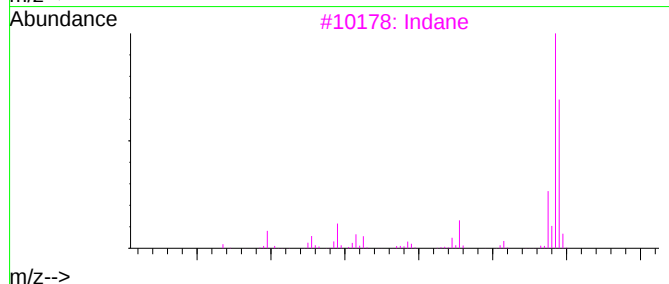
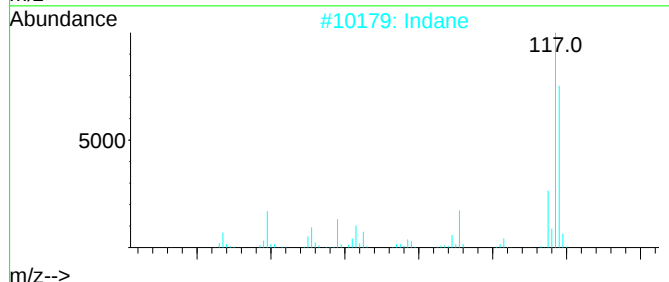
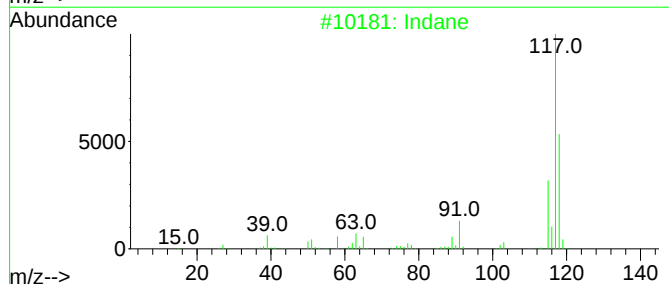
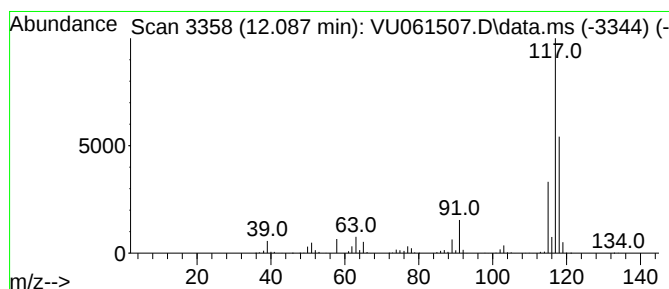
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.087	117.69 ug/L	10108700	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	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
5	Deltacyclene	118	C9H10	007785-10-6	72



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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

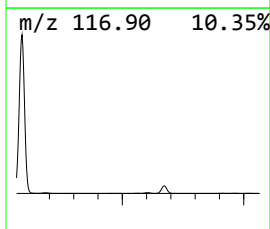
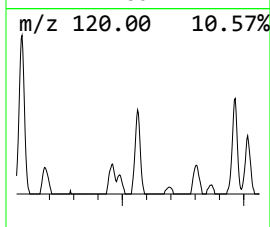
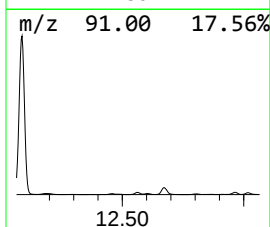
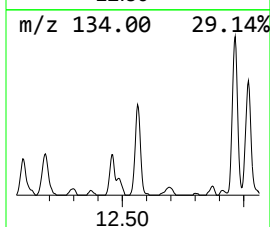
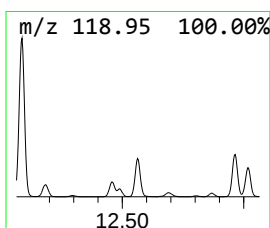
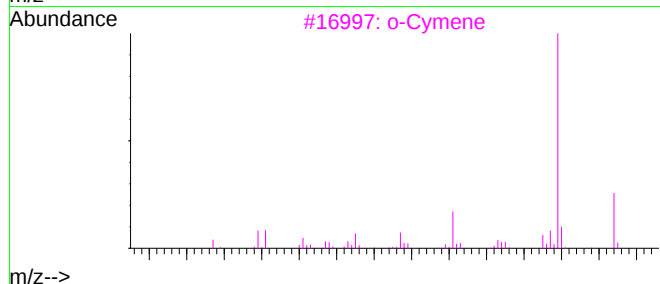
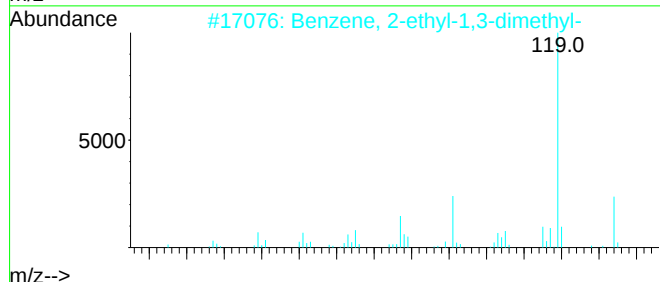
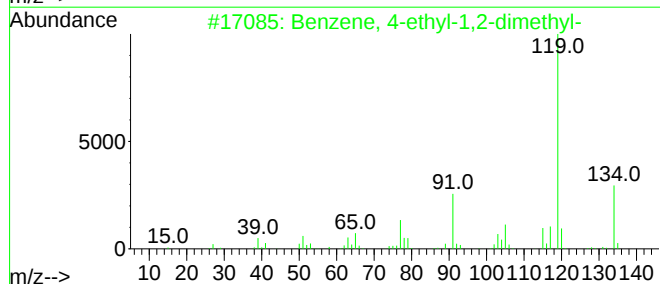
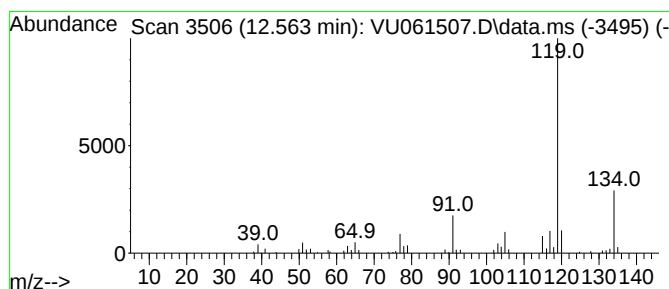
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.563	1.31 ug/L	112732	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



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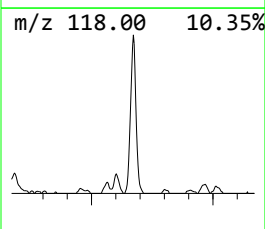
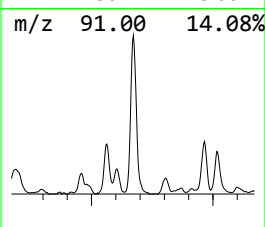
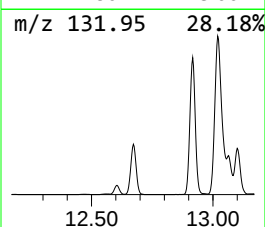
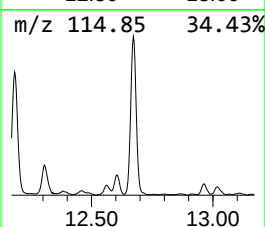
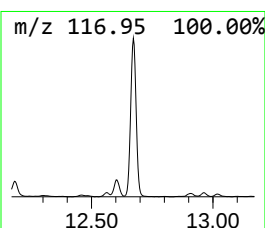
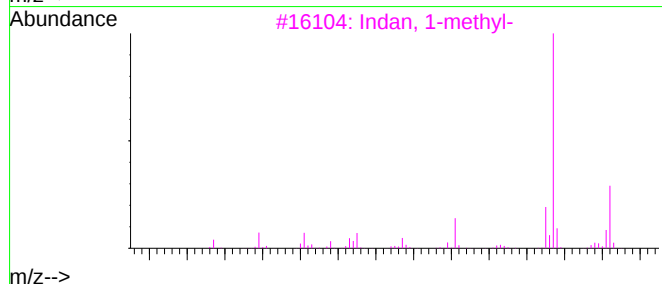
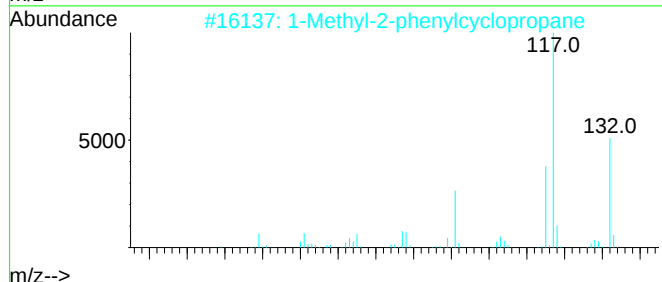
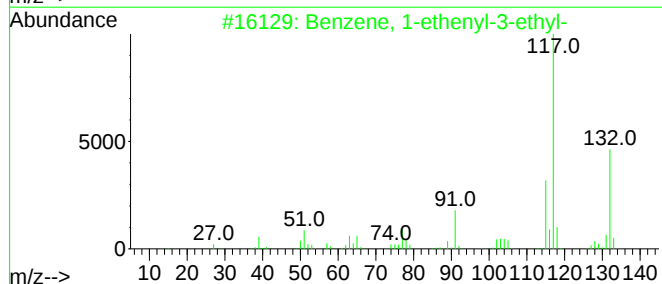
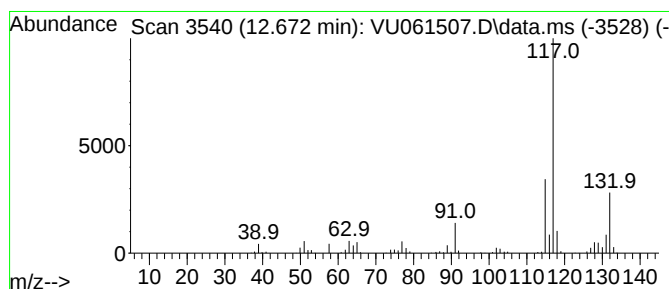
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.672	5.92 ug/L	508680	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	1-Methyl-2-phenylcyclopropane		132	C10H12	003145-76-4	90
3	Indan, 1-methyl-		132	C10H12	000767-58-8	90
4	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	87
5	Indan, 1-methyl-		132	C10H12	000767-58-8	81



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ClientSampleId :
C0AF5DL

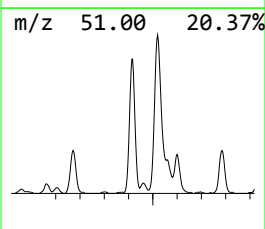
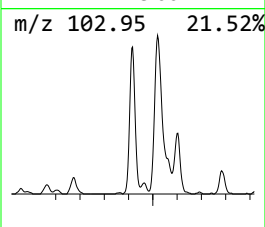
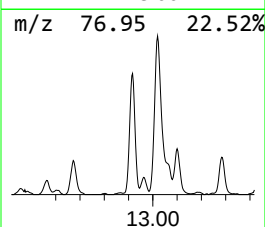
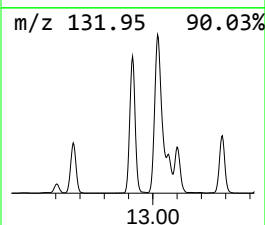
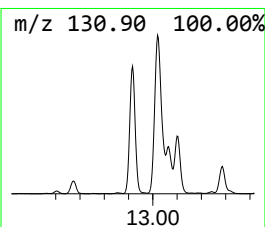
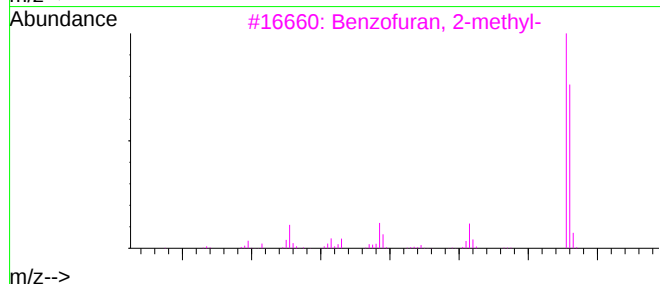
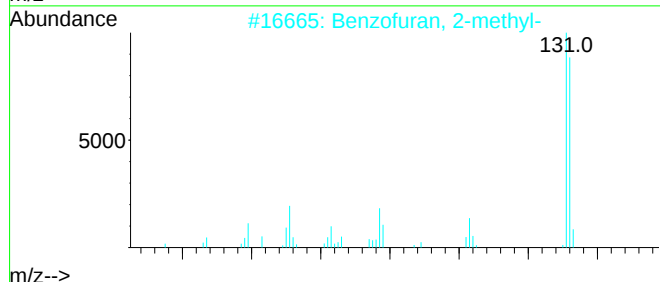
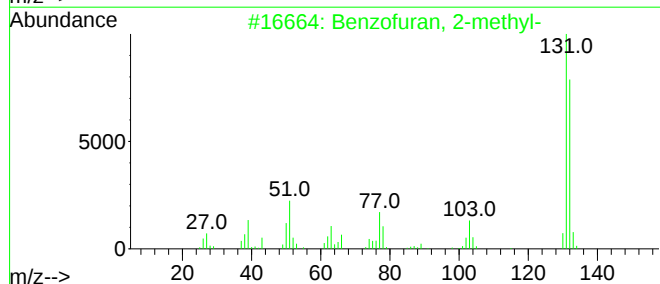
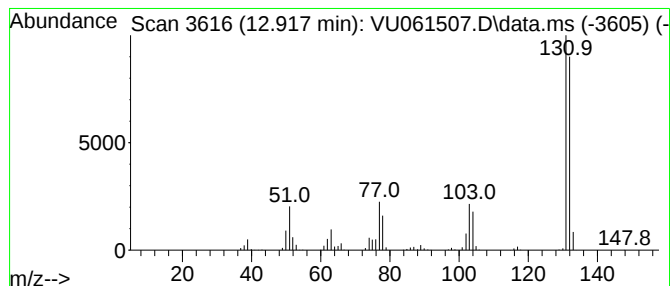
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.917	6.51 ug/L	559422	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	94
4	Benzofuran, 2-methyl-		132	C9H8O	004265-25-2	91
5	3-Phenyl-2-propyn-1-ol		132	C9H8O	001504-58-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
Data File : VU061507.D
Acq On : 06 Nov 2024 17:22
Operator : MD/SY
Sample : P4714-03DL 4X
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AF5DL

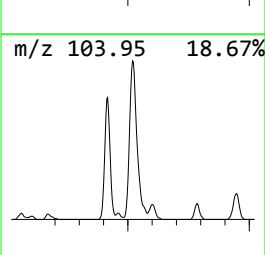
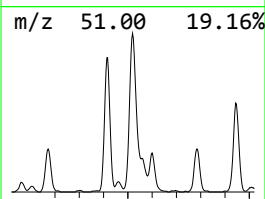
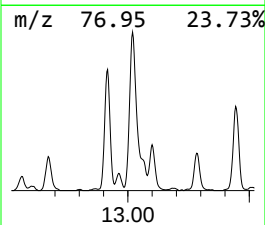
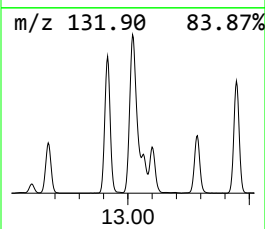
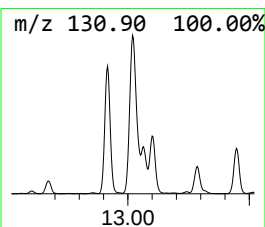
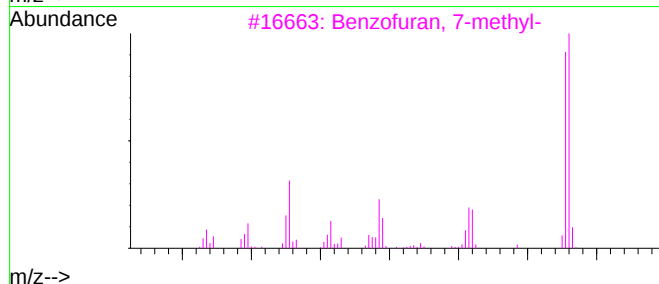
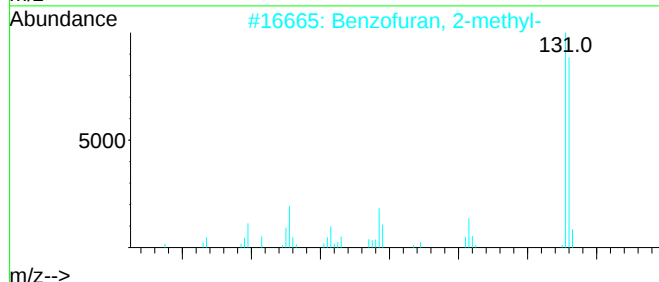
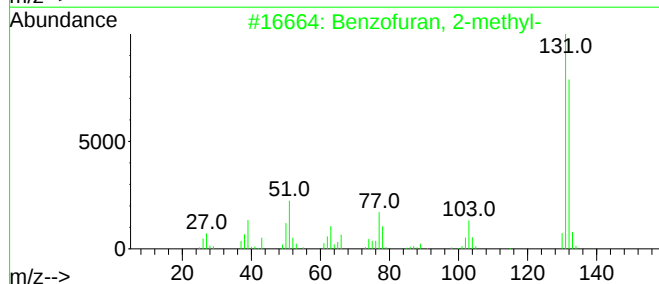
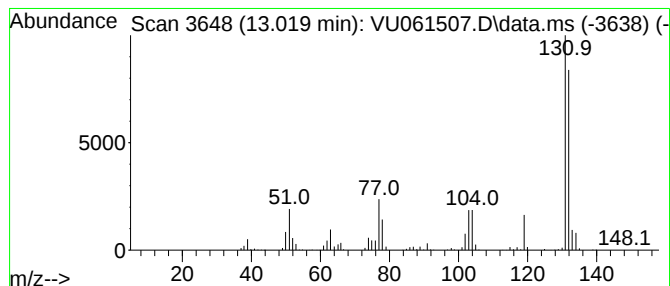
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzofuran, 7-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.020	10.51 ug/L	903068	1,4-Dichlorobenzene-d4		11.804	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-		132	C9H8O	004265-25-2	93
2	Benzofuran, 2-methyl-		132	C9H8O	004265-25-2	93
3	Benzofuran, 7-methyl-		132	C9H8O	017059-52-8	93
4	Cinnamaldehyde, (E)-		132	C9H8O	014371-10-9	90
5	Benzofuran, 2-methyl-		132	C9H8O	004265-25-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
Data File : VU061507.D
Acq On : 06 Nov 2024 17:22
Operator : MD/SY
Sample : P4714-03DL 4X
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

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

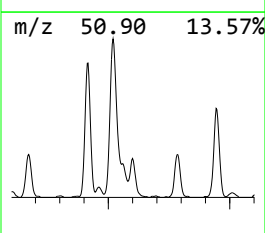
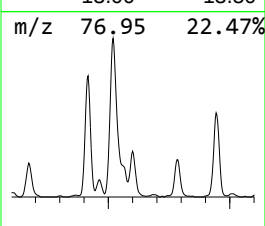
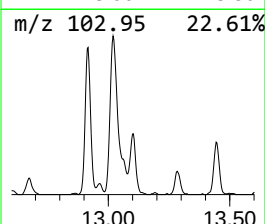
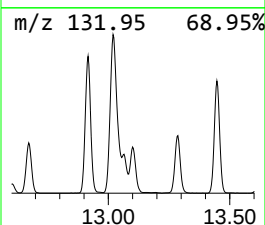
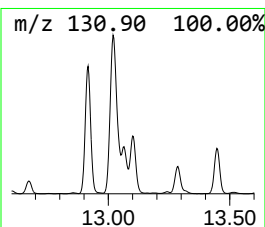
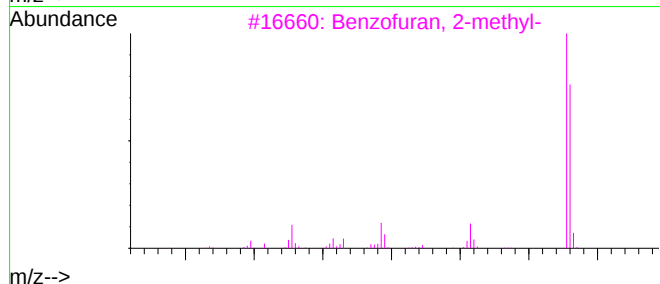
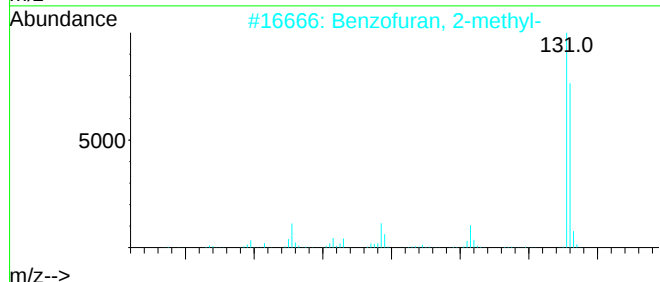
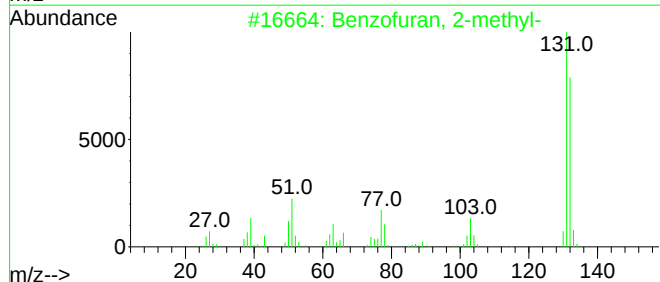
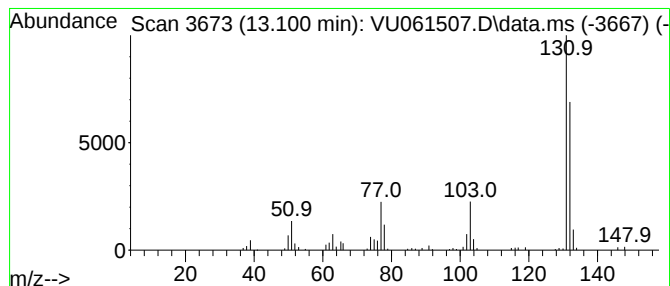
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 Cinnamaldehyde, (E)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	2.49 ug/L	213601	1,4-Dichlorobenzene-d4	11.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	78
5	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
Data File : VU061507.D
Acq On : 06 Nov 2024 17:22
Operator : MD/SY
Sample : P4714-03DL 4X
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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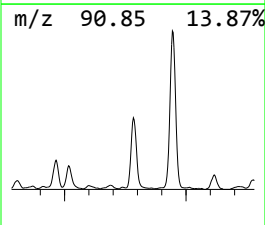
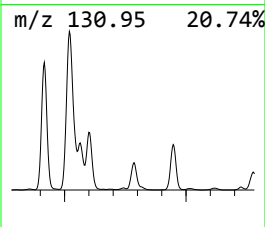
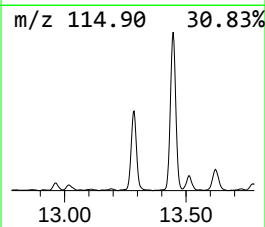
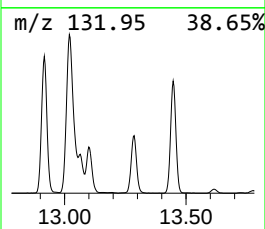
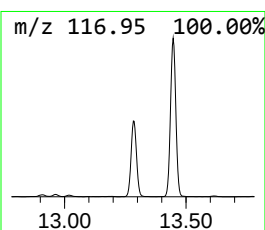
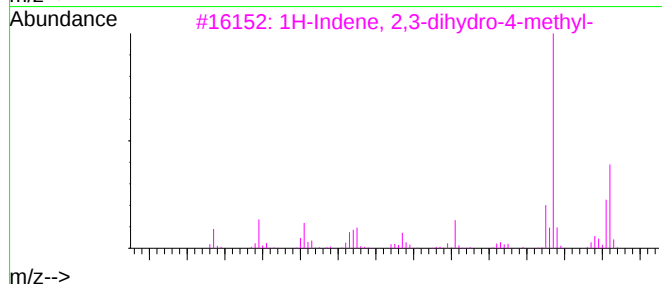
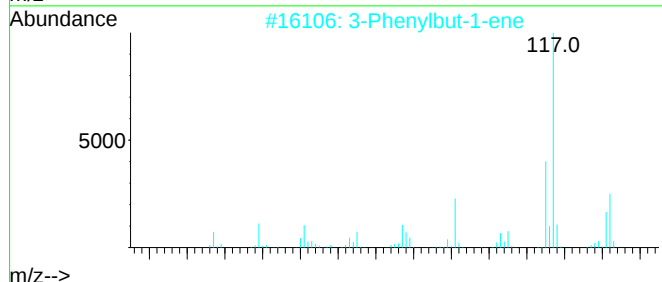
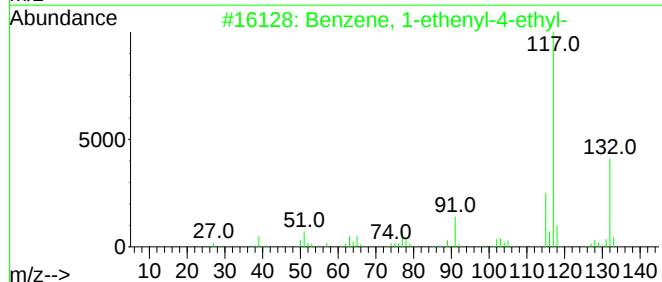
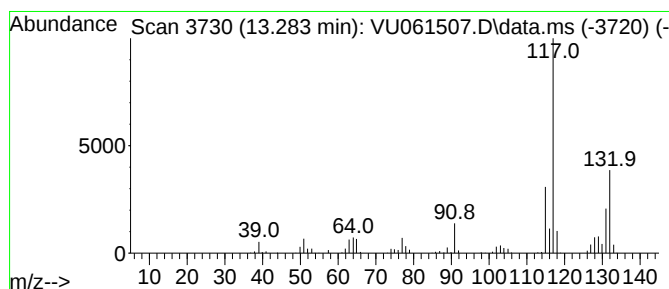
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.283	5.41 ug/L	464615	1,4-Dichlorobenzene-d4		11.804	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	94
2	3-Phenylbut-1-ene		132	C10H12	000934-10-1	91
3	1H-Indene, 2,3-dihydro-4-methyl-		132	C10H12	000824-22-6	90
4	1H-Indene, 2,3-dihydro-5-methyl-		132	C10H12	000874-35-1	90
5	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
Data File : VU061507.D
Acq On : 06 Nov 2024 17:22
Operator : MD/SY
Sample : P4714-03DL 4X
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

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

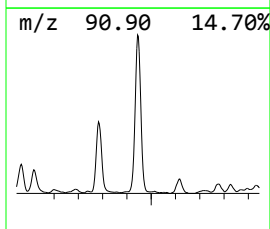
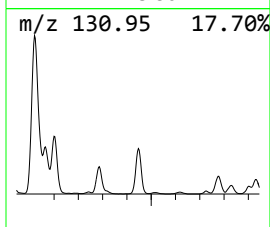
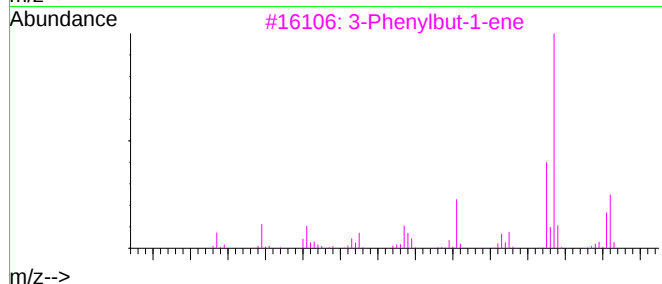
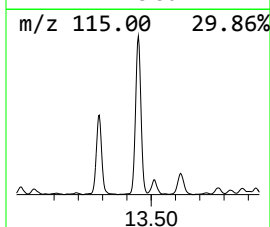
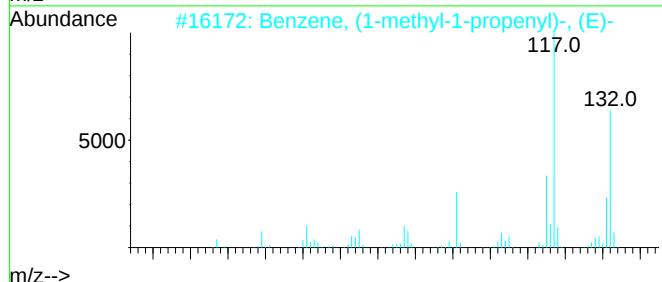
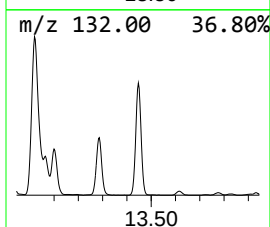
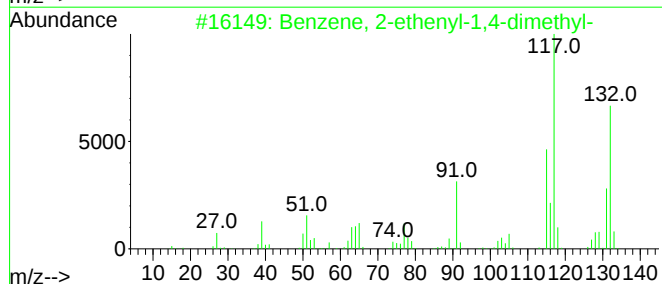
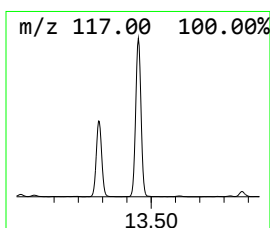
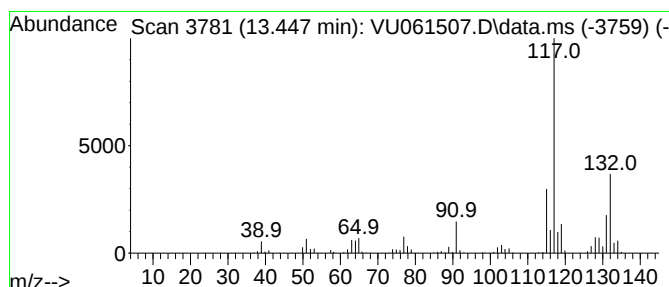
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.447	12.14 ug/L	1042540	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	96
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
Data File : VU061507.D
Acq On : 06 Nov 2024 17:22
Operator : MD/SY
Sample : P4714-03DL 4X
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

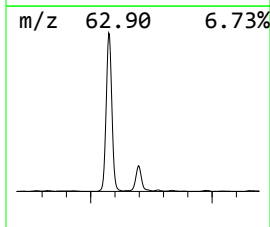
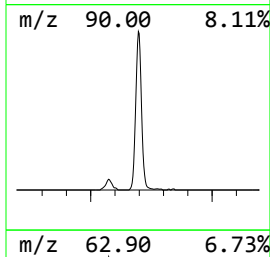
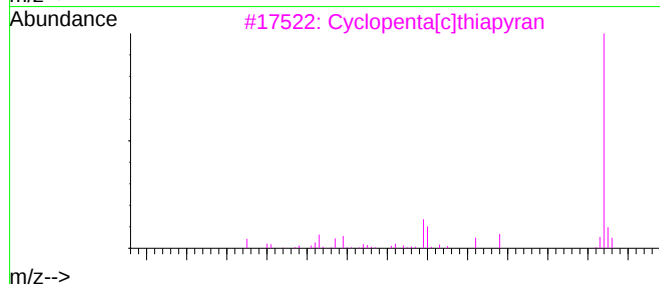
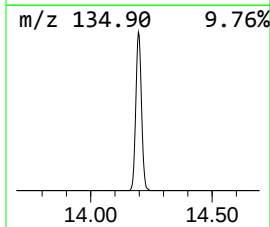
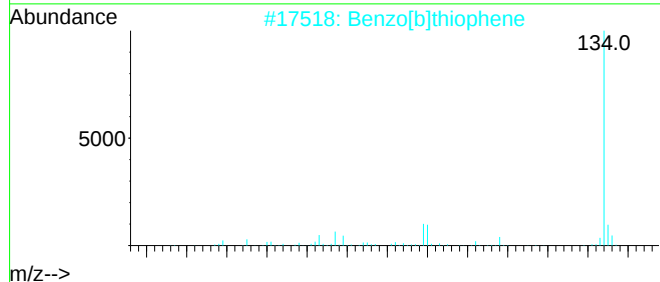
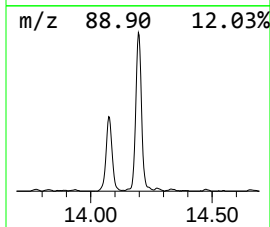
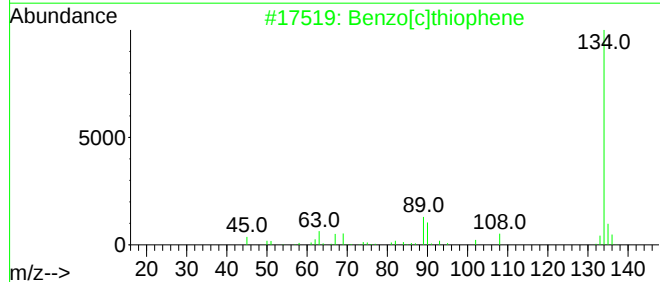
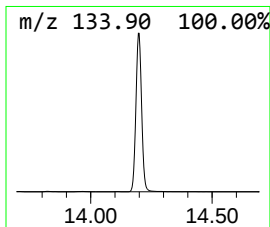
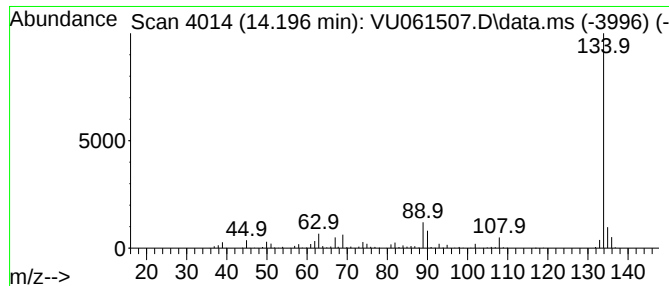
Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Benzo[c]thiophene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.196	12.23 ug/L	1050670	1,4-Dichlorobenzene-d4		11.804	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzo[c]thiophene		134	C8H6S	000270-82-6	95
2	Benzo[b]thiophene		134	C8H6S	000095-15-8	94
3	Cyclopenta[c]thiapyran		134	C8H6S	000270-63-3	94
4	Benzo[b]thiophene		134	C8H6S	000095-15-8	94
5	Benzo[b]thiophene		134	C8H6S	000095-15-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110624\
Data File : VU061507.D
Acq On : 06 Nov 2024 17:22
Operator : MD/SY
Sample : P4714-03DL 4X
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AF5DL

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
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Mesitylene	11.888	1.2	ug/L	103165	3	11.804	429477	5.0
Indane	12.087	117.7	ug/L	10108700	3	11.804	429477	5.0
Benzene, 4-ethy...	12.563	1.3	ug/L	112732	3	11.804	429477	5.0
Benzene, 1-ethe...	12.672	5.9	ug/L	508680	3	11.804	429477	5.0
Benzofuran, 2-m...	12.917	6.5	ug/L	559422	3	11.804	429477	5.0
Benzofuran, 7-m...	13.020	10.5	ug/L	903068	3	11.804	429477	5.0
Cinnamaldehyde,...	13.100	2.5	ug/L	213601	3	11.804	429477	5.0
Benzene, 1-ethe...	13.283	5.4	ug/L	464615	3	11.804	429477	5.0
Benzene, 2-ethe...	13.447	12.1	ug/L	1042540	3	11.804	429477	5.0
Benzo[c]thiophene	14.196	12.2	ug/L	1050670	3	11.804	429477	5.0