

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
 Data File : VU060060.D
 Acq On : 19 Jul 2024 15:08
 Operator : MD/SY
 Sample : P3244-03 100X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 YDZD4

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\SFAMUTR071824WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU060060.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.602	85	97	113	rBV2	355721	428433	10.32%	2.953%
2	1.911	185	193	205	rBV	42682	57045	1.37%	0.393%
3	2.563	384	396	410	rBV	72270	119037	2.87%	0.820%
4	4.657	1027	1047	1094	rBV	1315629	3141079	75.70%	21.648%
5	5.055	1155	1171	1194	rBV2	79725	178365	4.30%	1.229%
6	5.718	1359	1377	1412	rBV2	159297	441023	10.63%	3.040%
7	6.242	1527	1540	1561	rBV	150004	292824	7.06%	2.018%
8	6.682	1663	1677	1693	rBV	97919	192686	4.64%	1.328%
9	7.560	1939	1950	1972	rBV	44295	79692	1.92%	0.549%
10	7.891	2040	2053	2063	rBV	184362	316711	7.63%	2.183%
11	7.962	2063	2075	2116	rVB	2437418	4149597	100.00%	28.599%
12	8.174	2132	2141	2157	rVB2	27419	46556	1.12%	0.321%
13	8.627	2270	2282	2313	rBV	217664	423494	10.21%	2.919%
14	9.412	2513	2526	2550	rBV	221488	409289	9.86%	2.821%
15	9.566	2561	2574	2590	rVB	62265	102618	2.47%	0.707%
16	9.685	2597	2611	2632	rBV	305791	514761	12.41%	3.548%
17	10.094	2725	2738	2757	rBV	189071	313588	7.56%	2.161%
18	10.750	2931	2942	2954	rBV	85284	135007	3.25%	0.930%
19	10.901	2974	2989	3002	rBV2	23906	46308	1.12%	0.319%
20	10.991	3006	3017	3036	rVV	103410	238338	5.74%	1.643%
21	11.081	3036	3045	3062	rVB	84124	134827	3.25%	0.929%
22	11.287	3097	3109	3126	rVB	81585	127691	3.08%	0.880%
23	11.463	3152	3164	3179	rBV	212588	335205	8.08%	2.310%
24	11.631	3197	3216	3230	rBV9	30005	64010	1.54%	0.441%
25	11.804	3256	3270	3286	rVV2	250486	487359	11.74%	3.359%
26	11.888	3286	3296	3318	rVB	167709	296272	7.14%	2.042%
27	12.090	3347	3359	3365	rBV3	34357	58823	1.42%	0.405%
28	12.187	3377	3389	3411	rVB2	251859	517930	12.48%	3.570%
29	12.679	3533	3542	3561	rVB5	17698	45259	1.09%	0.312%
30	12.939	3613	3623	3640	rVV3	35980	85193	2.05%	0.587%
31	13.020	3640	3648	3663	rVB	30099	47419	1.14%	0.327%
32	13.447	3772	3781	3792	rVB5	50340	84753	2.04%	0.584%
33	13.512	3792	3801	3810	rVB	32732	49790	1.20%	0.343%
34	13.730	3855	3869	3880	rBV3	41779	72456	1.75%	0.499%
35	14.084	3963	3979	3999	rBV	48377	84636	2.04%	0.583%
36	15.219	4312	4332	4356	rBV	130895	223672	5.39%	1.542%

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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\SFAMUTR071824WMA.M
Title : TRACE VOA SFAM1.0

37	15.405	4376	4390	4402	rBV	103149	167819	4.04%	1.157%
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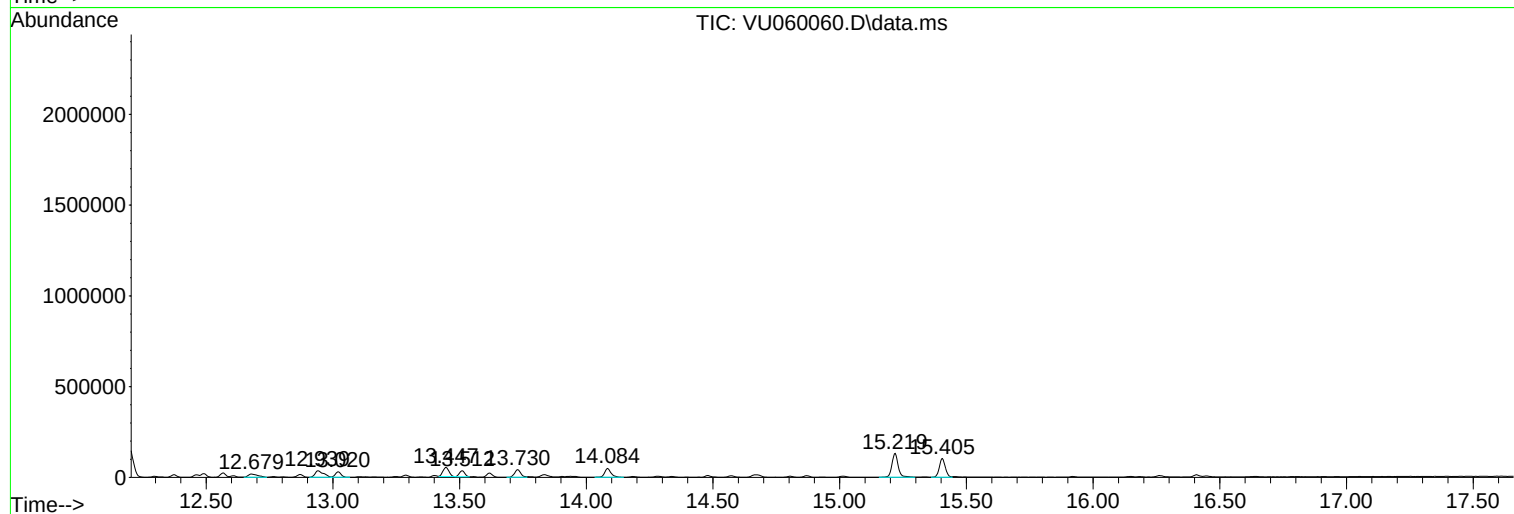
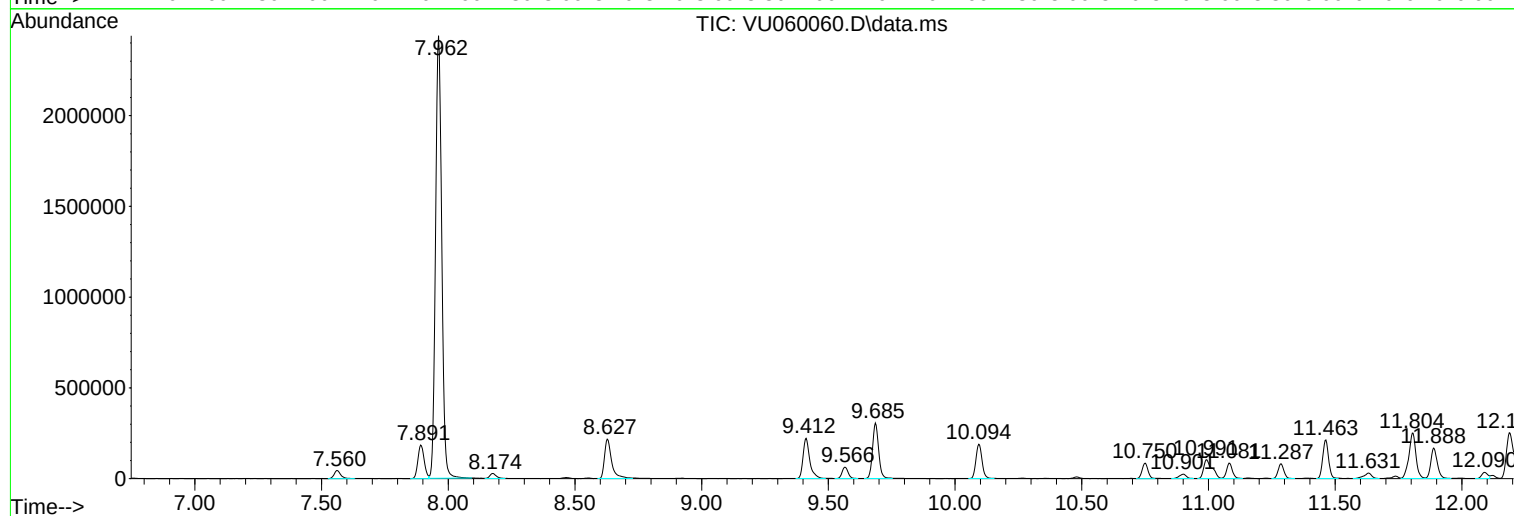
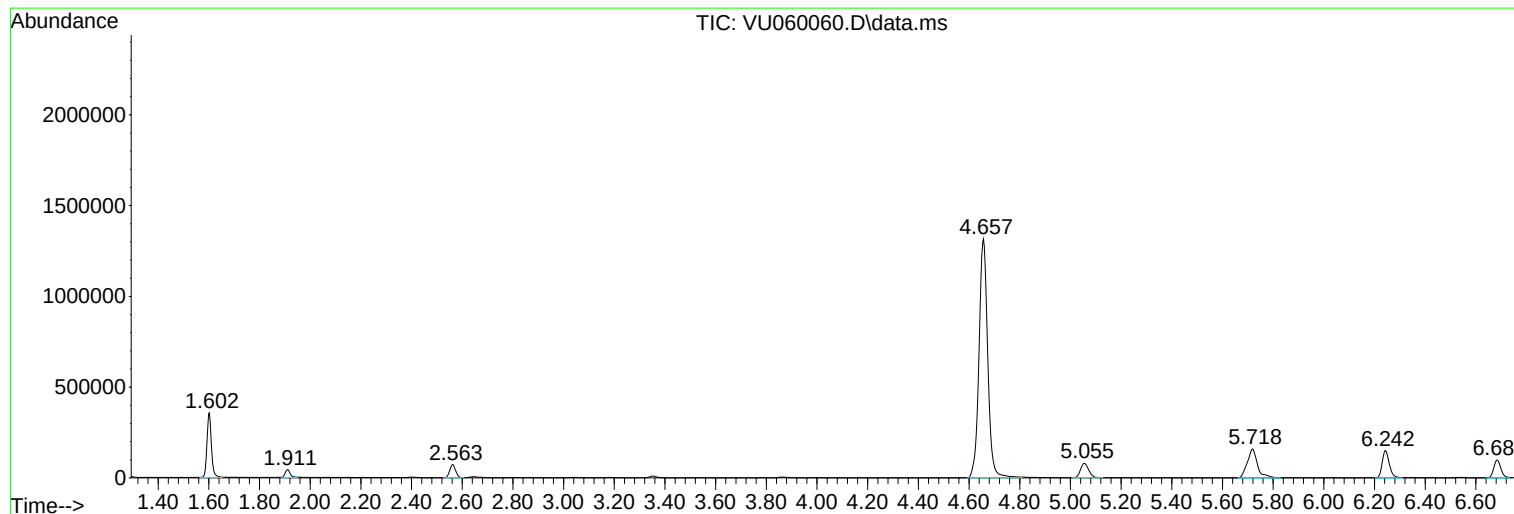
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Misc : 25.0mL/MSVOA_U/WATER
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Instrument :
MSVOA_U
ClientSampleId :
YDZD4

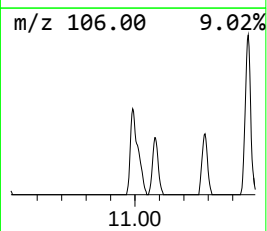
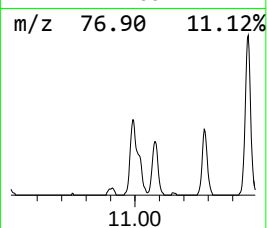
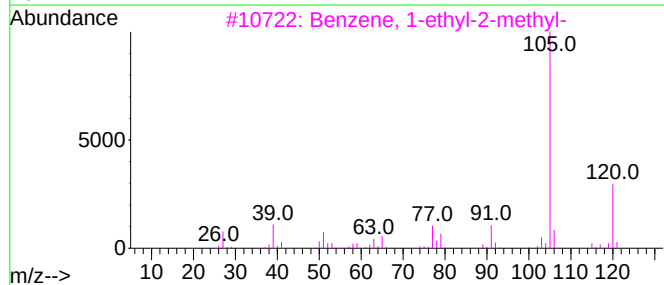
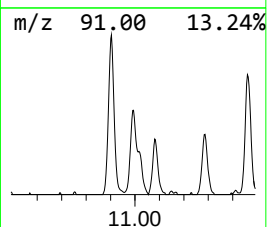
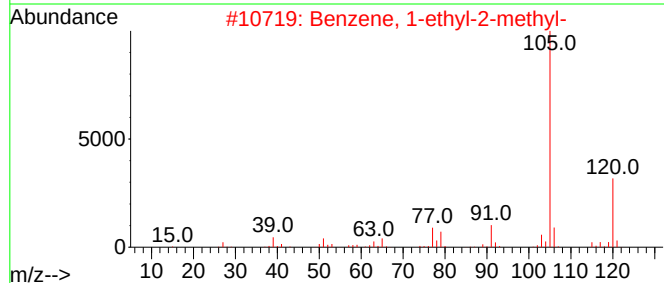
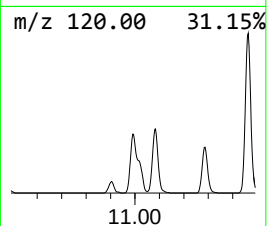
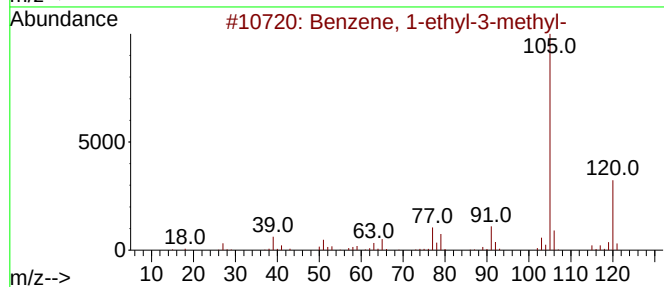
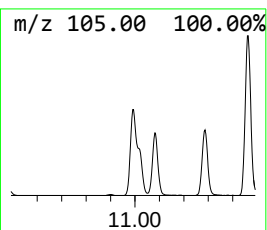
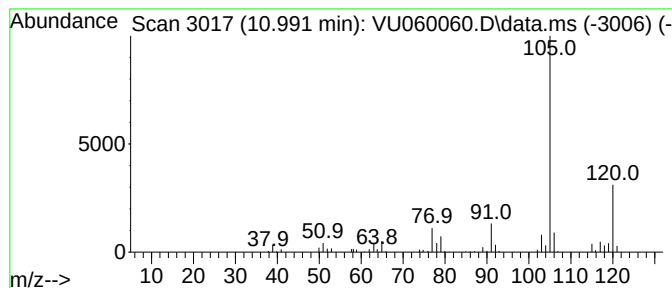
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.991	2.45 ug/L	238338	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

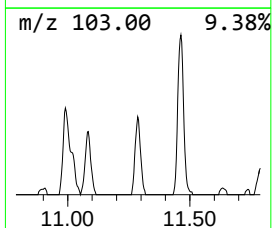
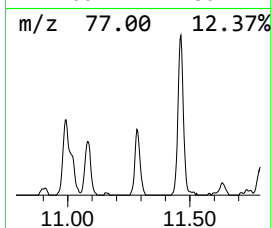
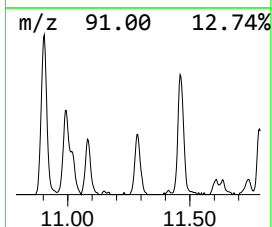
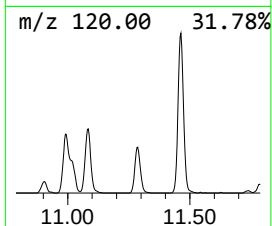
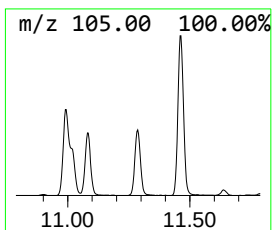
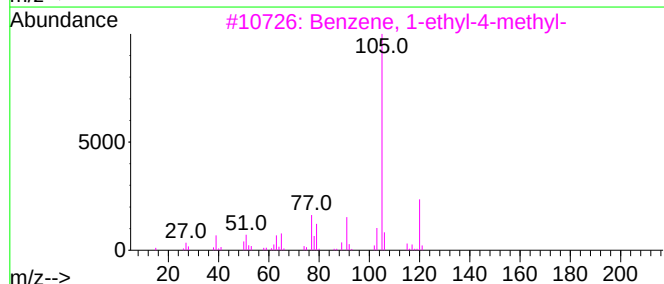
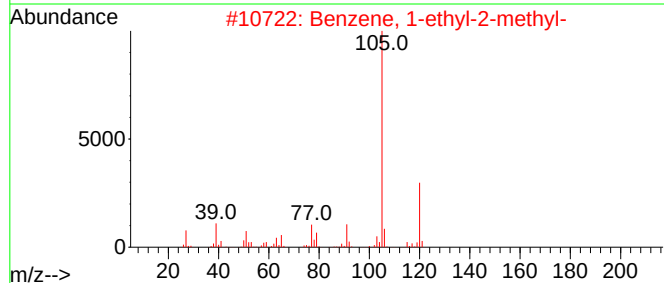
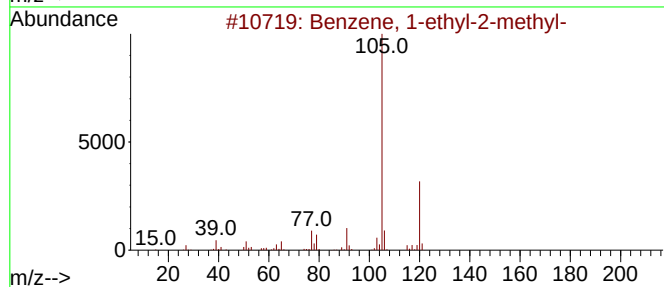
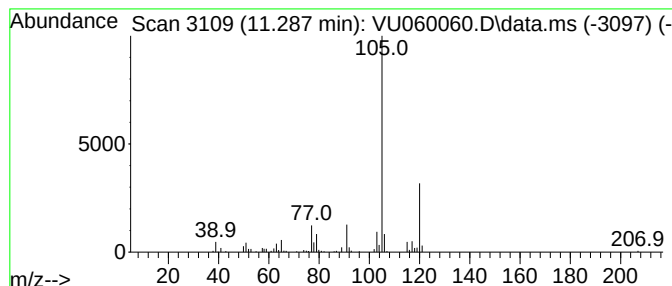
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.287	1.31 ug/L	127691	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

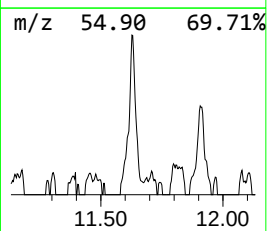
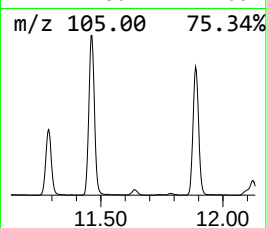
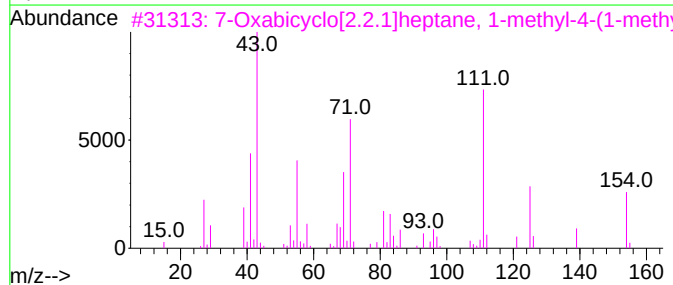
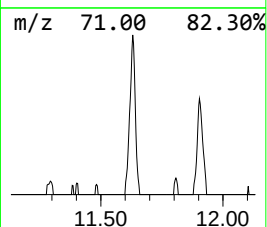
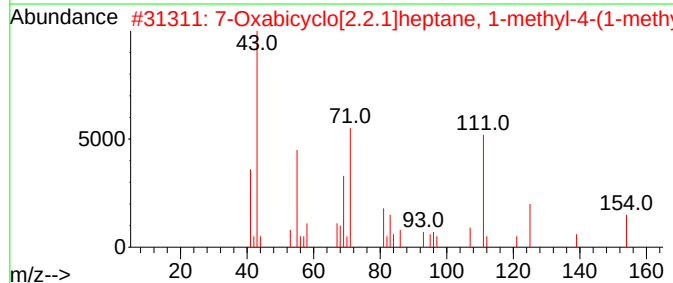
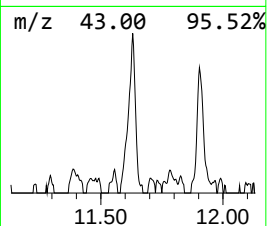
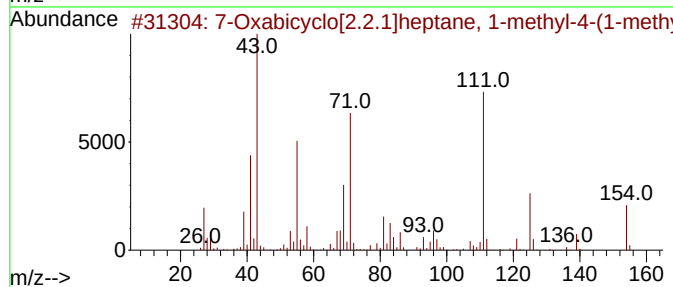
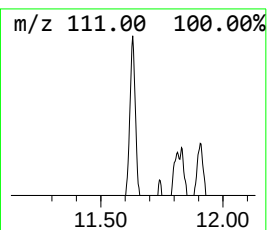
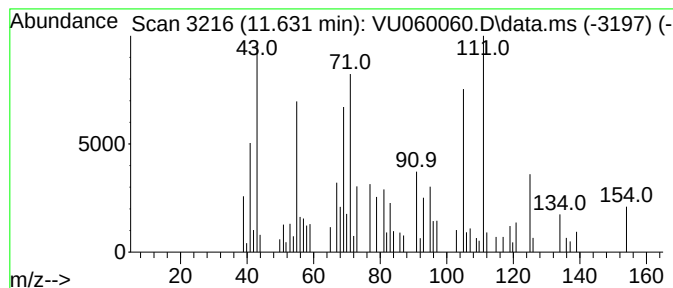
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 7-Oxabicyclo[2.2.1]heptane,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.631	0.66 ug/L	64010	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	89
2			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	81
3			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	76
4			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	64
5			Tetrahydrocarvone	154	C10H18O	000499-70-7	42



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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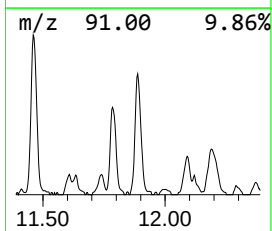
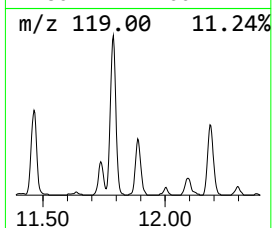
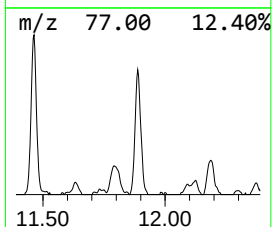
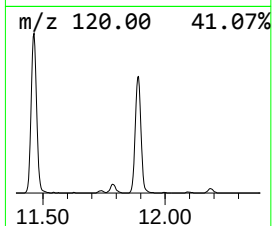
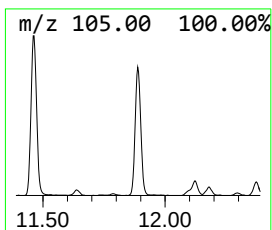
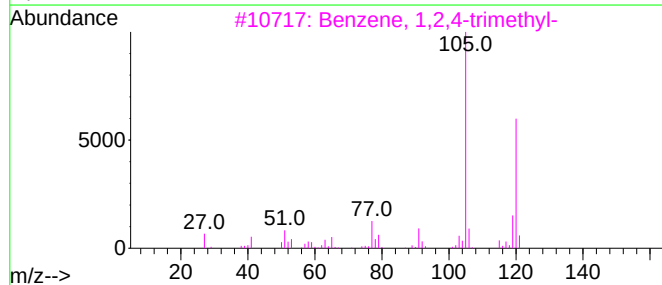
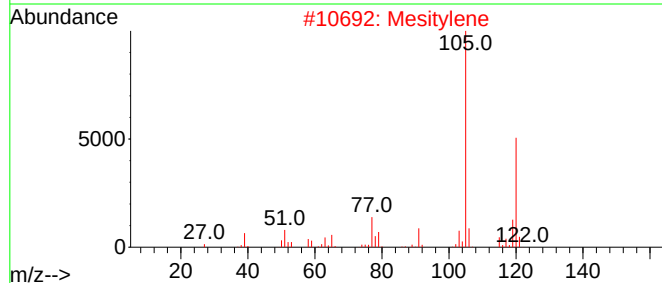
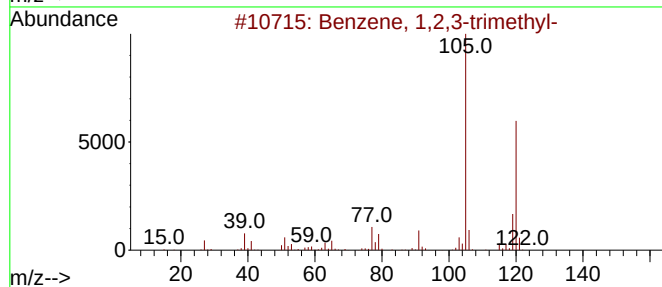
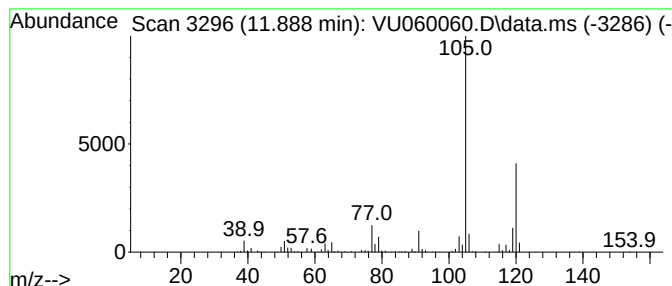
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.888	3.04 ug/L	296272	1,4-Dichlorobenzene-d4	11.807

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



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YDZD4

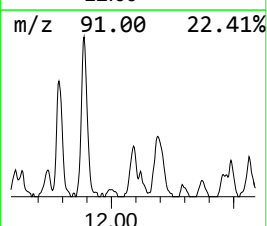
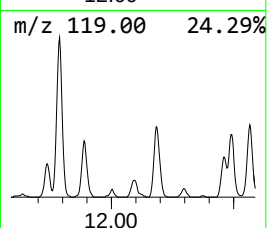
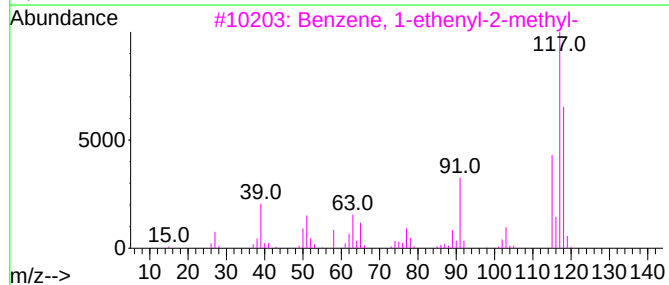
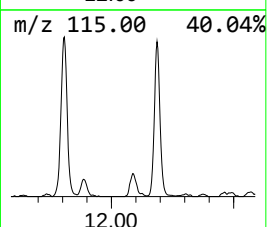
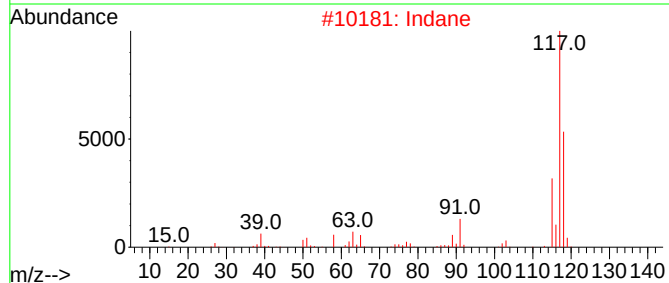
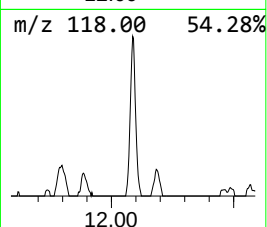
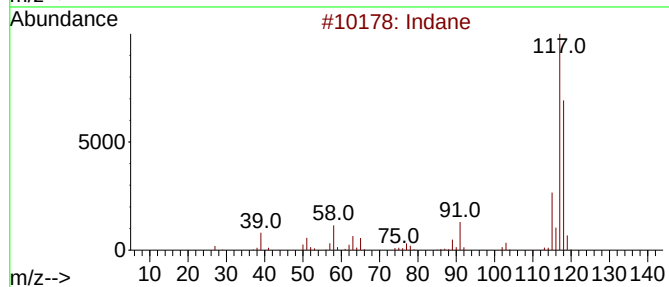
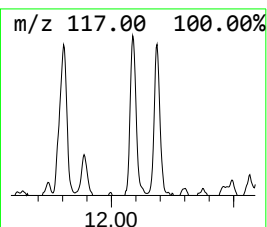
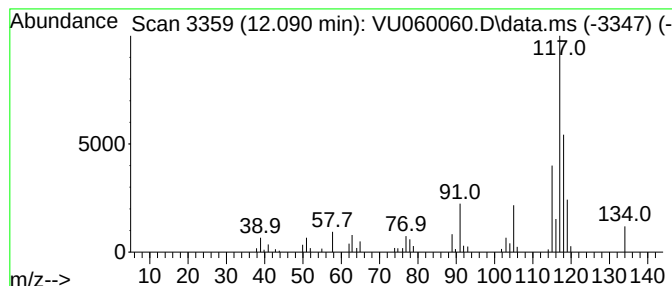
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Indane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.090	0.60 ug/L	58823	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	60
2			Indane	118	C9H10	000496-11-7	60
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
4			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	60
5			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060060.D
Acq On : 19 Jul 2024 15:08
Operator : MD/SY
Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

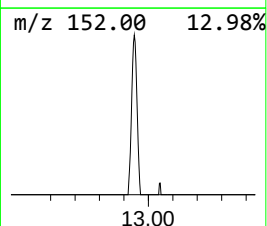
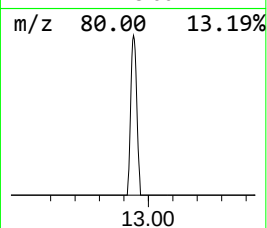
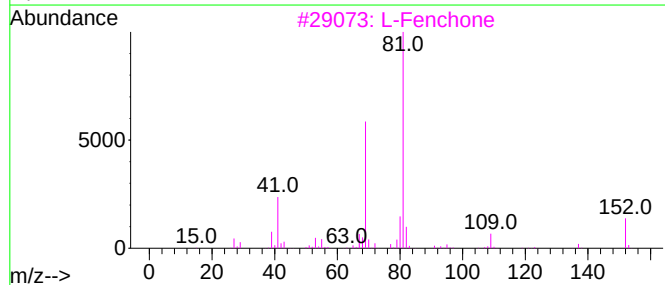
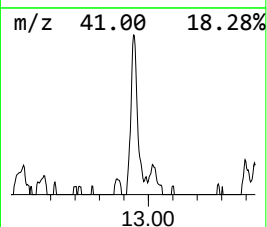
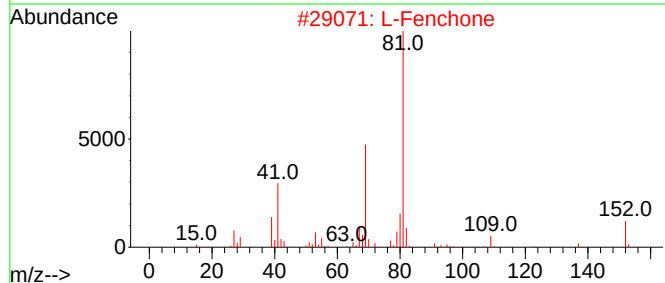
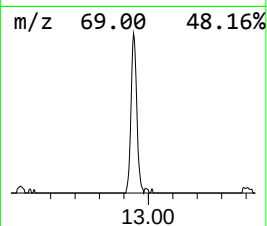
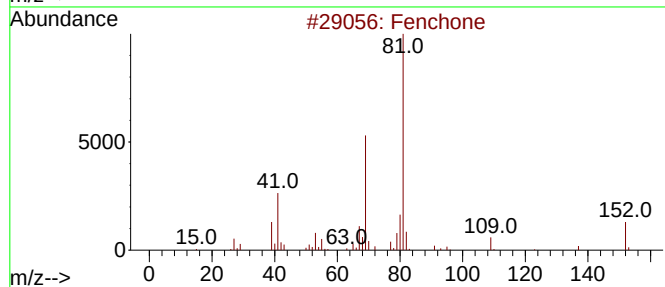
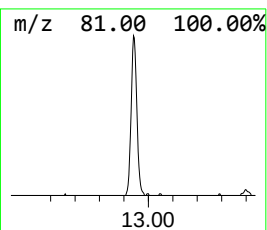
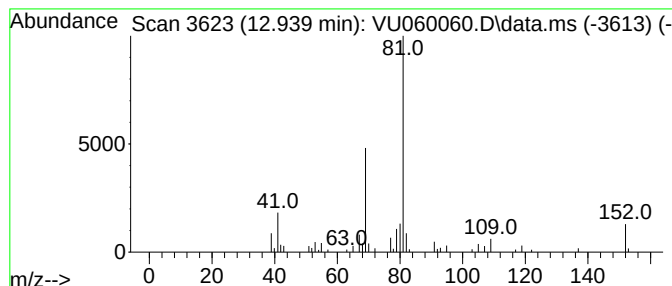
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Fenchone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.939	0.87 ug/L	85193	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchone	152	C10H16O	001195-79-5	91
2			L-Fenchone	152	C10H16O	007787-20-4	91
3			L-Fenchone	152	C10H16O	007787-20-4	90
4			D-Fenchone	152	C10H16O	004695-62-9	90
5			Fenchone	152	C10H16O	001195-79-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060060.D
Acq On : 19 Jul 2024 15:08
Operator : MD/SY
Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

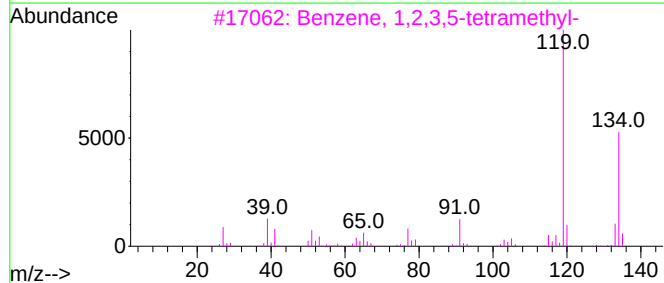
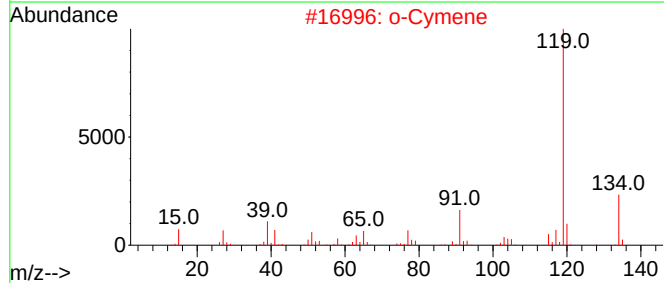
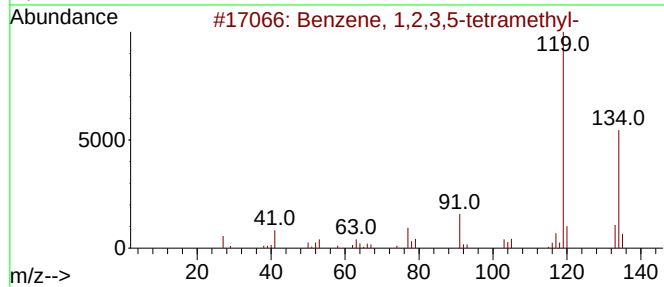
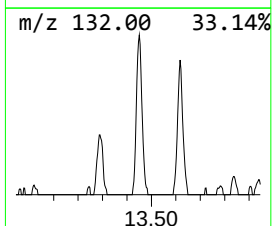
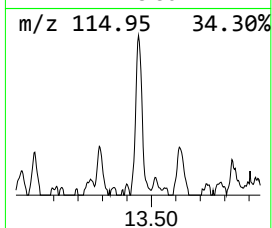
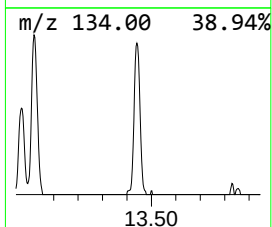
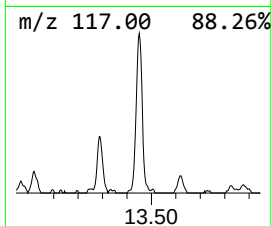
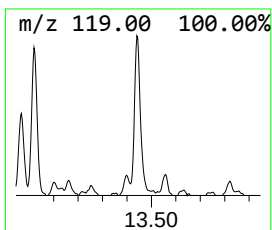
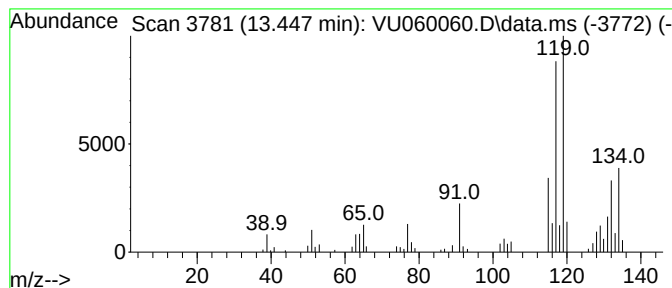
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.447	0.87 ug/L	84753	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	50
2			o-Cymene	134	C10H14	000527-84-4	50
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	50
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060060.D
Acq On : 19 Jul 2024 15:08
Operator : MD/SY
Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

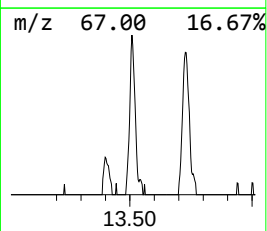
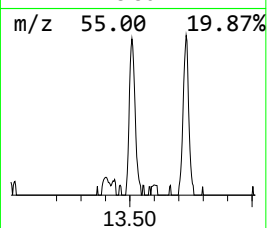
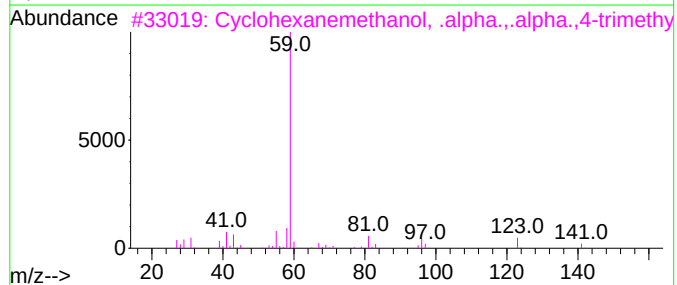
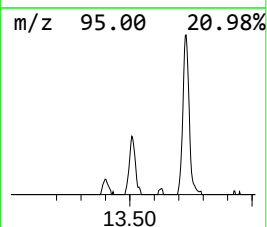
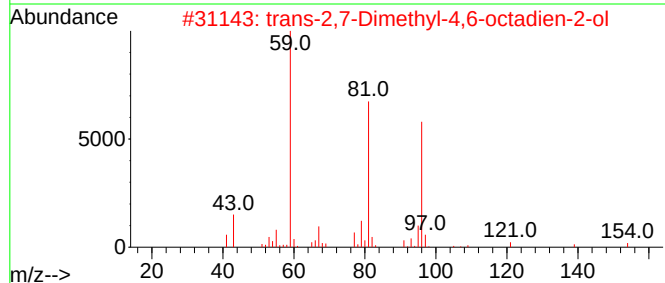
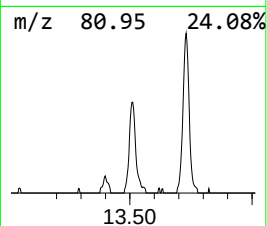
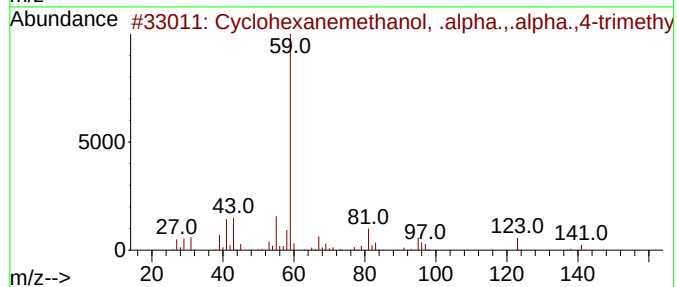
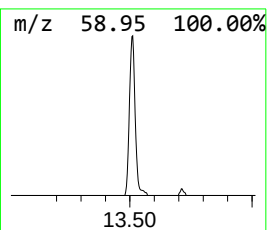
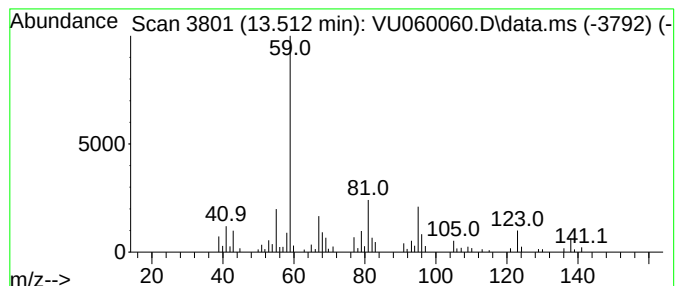
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Cyclohexanemethanol, .alpha... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.512	0.51 ug/L	49790	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	50
2			trans-2,7-Dimethyl-4,6-octadien-...	154	C10H18O	1010281-69-5	40
3			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	40
4			2-(3,4-Dibromo-4-methylcyclohexy...	312	C10H18Br2O	110202-13-6	40
5			7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060060.D
Acq On : 19 Jul 2024 15:08
Operator : MD/SY
Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

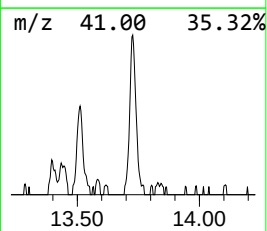
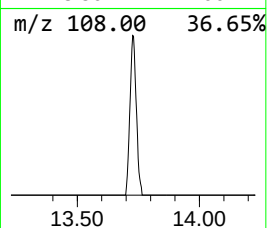
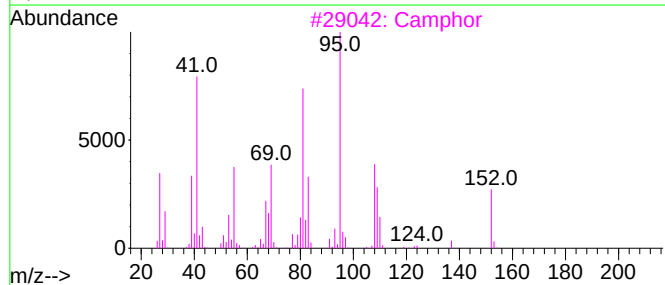
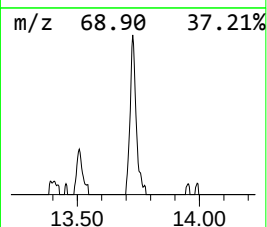
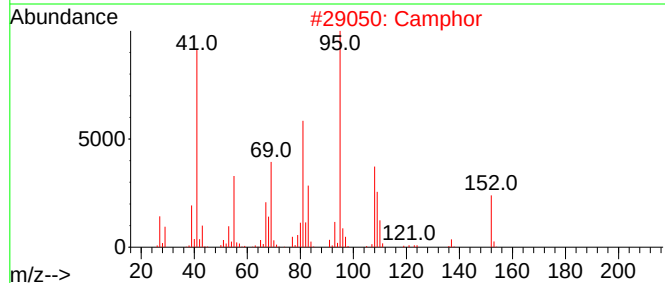
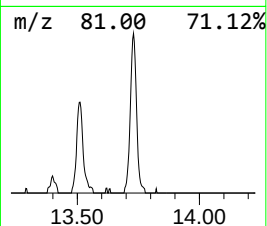
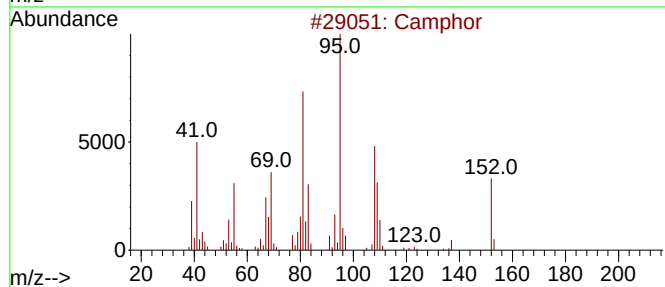
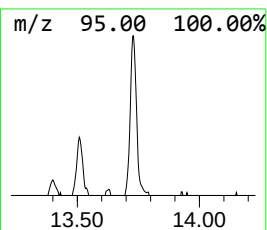
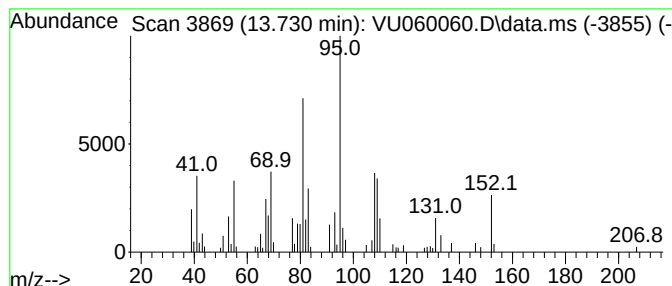
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Camphor Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.730	0.74 ug/L	72456	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphor	152	C10H16O	000076-22-2	97
2			Camphor	152	C10H16O	000076-22-2	97
3			Camphor	152	C10H16O	000076-22-2	97
4			(+)-2-Bornanone	152	C10H16O	000464-49-3	97
5			(+)-2-Bornanone	152	C10H16O	000464-49-3	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060060.D
Acq On : 19 Jul 2024 15:08
Operator : MD/SY
Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

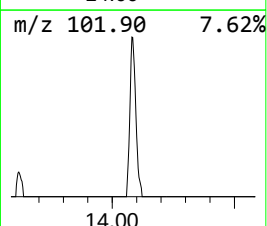
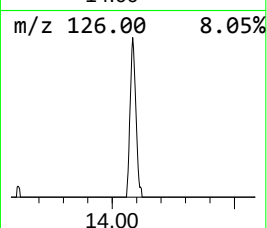
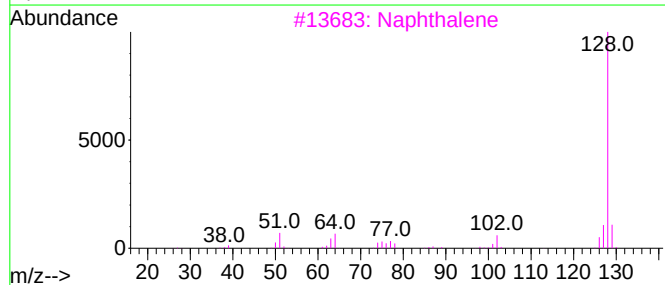
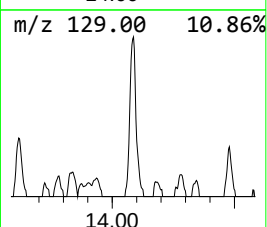
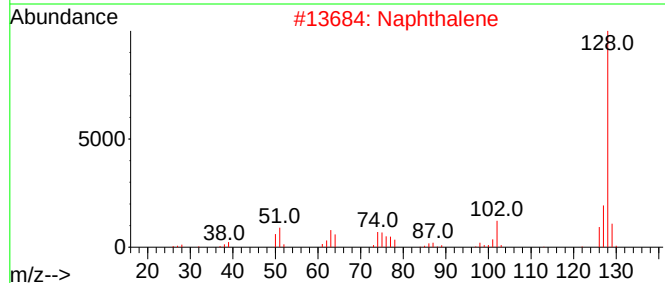
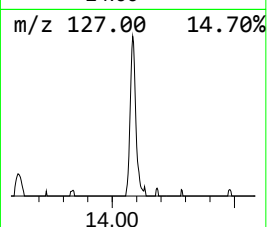
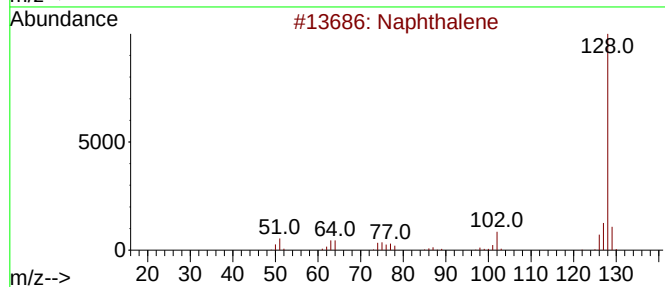
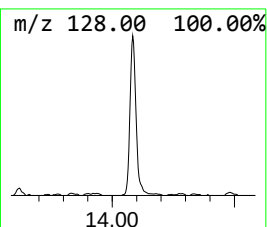
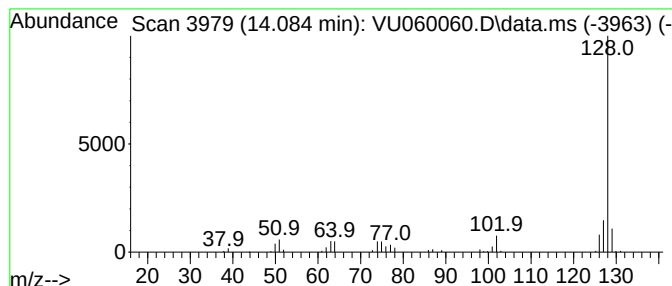
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Azulene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.084	0.87 ug/L	84636	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			Naphthalene	128	C10H8	000091-20-3	91
4			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060060.D
Acq On : 19 Jul 2024 15:08
Operator : MD/SY
Sample : P3244-03 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZD4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.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
Benzene, 1-ethy...	10.991	2.5	ug/L	238338	3	11.807	487359	5.0
Benzene, 1-ethy...	11.287	1.3	ug/L	127691	3	11.807	487359	5.0
7-Oxabicyclo[2...	11.631	0.7	ug/L	64010	3	11.807	487359	5.0
Benzene, 1,2,3-...	11.888	3.0	ug/L	296272	3	11.807	487359	5.0
Indane	12.090	0.6	ug/L	58823	3	11.807	487359	5.0
Fenchone	12.939	0.9	ug/L	85193	3	11.807	487359	5.0
Benzene, 1,2,3,...	13.447	0.9	ug/L	84753	3	11.807	487359	5.0
Cyclohexanemeth...	13.512	0.5	ug/L	49790	3	11.807	487359	5.0
Camphor	13.730	0.7	ug/L	72456	3	11.807	487359	5.0
Azulene	14.084	0.9	ug/L	84636	3	11.807	487359	5.0