

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
 Data File : VU061098.D
 Acq On : 10 Oct 2024 18:49
 Operator : MD/SY
 Sample : P4380-14
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 E27H3

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\SFAMUTR100124WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU061098.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.354	13	20	38	rBV2	31719	37127	1.39%	0.274%
2	1.451	39	50	56	rBV	22099	44439	1.66%	0.328%
3	1.583	81	91	102	rBV2	88371	159837	5.96%	1.180%
4	1.718	125	133	153	rBV	496388	612378	22.85%	4.521%
5	1.901	184	190	203	rVB	49964	65562	2.45%	0.484%
6	2.232	286	293	310	rBV	31448	76895	2.87%	0.568%
7	2.550	382	392	406	rVB	96377	153724	5.74%	1.135%
8	3.849	784	796	817	rVB	12056	27940	1.04%	0.206%
9	4.695	1031	1059	1112	rBV3	52107	309571	11.55%	2.285%
10	5.049	1153	1169	1190	rBV	93966	206810	7.72%	1.527%
11	5.711	1356	1375	1382	rBV2	202314	467153	17.43%	3.449%
12	5.759	1383	1390	1410	rVB	255211	508396	18.97%	3.753%
13	6.010	1443	1468	1485	rBV2	42596	93124	3.47%	0.687%
14	6.238	1527	1539	1562	rVB	205893	393278	14.68%	2.903%
15	6.611	1645	1655	1665	rBV2	14697	27743	1.04%	0.205%
16	6.682	1665	1677	1691	rBV	116450	227637	8.49%	1.680%
17	7.560	1937	1950	1968	rBV	64206	116178	4.34%	0.858%
18	7.888	2037	2052	2065	rBV	232177	406759	15.18%	3.003%
19	7.959	2065	2074	2094	rVB	115274	203115	7.58%	1.499%
20	8.174	2131	2141	2155	rBV2	38291	66074	2.47%	0.488%
21	8.544	2244	2256	2272	rVB	102276	175559	6.55%	1.296%
22	8.634	2272	2284	2321	rBV	290926	573346	21.39%	4.233%
23	9.409	2513	2525	2547	rBV	289937	511328	19.08%	3.775%
24	9.563	2564	2573	2595	rVB	37393	65829	2.46%	0.486%
25	9.682	2595	2610	2627	rBV	122759	212676	7.94%	1.570%
26	10.094	2727	2738	2756	rBV	85878	142274	5.31%	1.050%
27	10.746	2931	2941	2954	rVB	93977	153810	5.74%	1.135%
28	10.991	3005	3017	3035	rBV	56384	119351	4.45%	0.881%
29	11.081	3035	3045	3062	rVB	95590	148395	5.54%	1.095%
30	11.283	3099	3108	3125	rVB	55016	87947	3.28%	0.649%
31	11.460	3152	3163	3172	rBV	148594	241006	8.99%	1.779%
32	11.630	3194	3216	3226	rBV7	14142	32530	1.21%	0.240%
33	11.733	3237	3248	3256	rBV3	60620	102950	3.84%	0.760%
34	11.804	3256	3270	3287	rVV3	325088	761290	28.41%	5.620%
35	11.884	3287	3295	3309	rVB	154855	239479	8.94%	1.768%
36	12.055	3338	3348	3351	rBV	67905	96647	3.61%	0.713%

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Min Area: 3 % of largest Peak

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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\SFAMUTR100124WMA.M
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37	12.084	3352	3357	3365	rVB2	77552	121551	4.54%	0.897%
38	12.183	3377	3388	3406	rVB	284329	481007	17.95%	3.551%
39	12.302	3413	3425	3437	rBV	167369	273186	10.19%	2.017%
40	12.367	3437	3445	3460	rVB	28966	49520	1.85%	0.366%
41	12.457	3462	3473	3478	rBV2	31285	52455	1.96%	0.387%
42	12.486	3478	3482	3494	rVB	30873	43978	1.64%	0.325%
43	12.563	3494	3506	3514	rBV	49966	75914	2.83%	0.560%
44	12.675	3531	3541	3561	rVB4	33973	85509	3.19%	0.631%
45	12.868	3590	3601	3609	rBV3	36003	60459	2.26%	0.446%
46	12.920	3609	3617	3623	rVV	18795	36583	1.37%	0.270%
47	12.965	3625	3631	3638	rVV	43910	63460	2.37%	0.468%
48	13.019	3638	3648	3661	rVB2	110554	182573	6.81%	1.348%
49	13.283	3723	3730	3743	rVB4	28427	44591	1.66%	0.329%
50	13.441	3769	3779	3793	rVB3	150043	253286	9.45%	1.870%
51	13.618	3824	3834	3851	rVB5	31498	63554	2.37%	0.469%
52	13.830	3891	3900	3915	rBV6	31069	68925	2.57%	0.509%
53	14.074	3962	3976	3997	rBV	1659885	2679829	100.00%	19.783%
54	14.199	4003	4015	4027	rBV	83722	139704	5.21%	1.031%
55	14.476	4091	4101	4108	rBV2	16539	27283	1.02%	0.201%
56	14.662	4150	4159	4171	rBV4	21109	44968	1.68%	0.332%
57	14.868	4214	4223	4235	rVB3	20193	33924	1.27%	0.250%
58	15.010	4257	4267	4278	rBV6	18794	35017	1.31%	0.259%
59	15.212	4320	4330	4343	rVV	158184	252178	9.41%	1.862%
60	15.402	4377	4389	4401	rBV	134917	227350	8.48%	1.678%
61	15.945	4552	4558	4569	rVB2	23687	36579	1.36%	0.270%
62	16.254	4638	4654	4663	rBV	28221	53830	2.01%	0.397%
63	16.402	4689	4700	4707	rBV	52024	86119	3.21%	0.636%
64	16.441	4707	4712	4722	rVB	25946	39115	1.46%	0.289%
65	16.878	4839	4848	4858	rBV3	18281	33328	1.24%	0.246%
66	17.087	4905	4913	4925	rVB	18924	32158	1.20%	0.237%

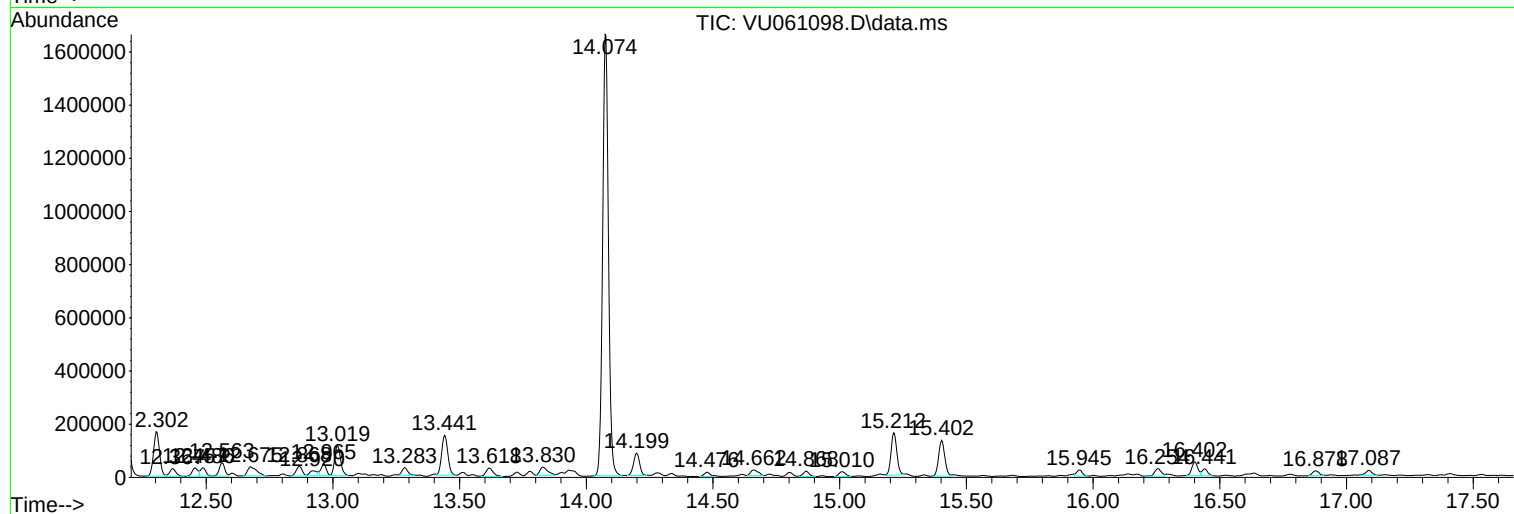
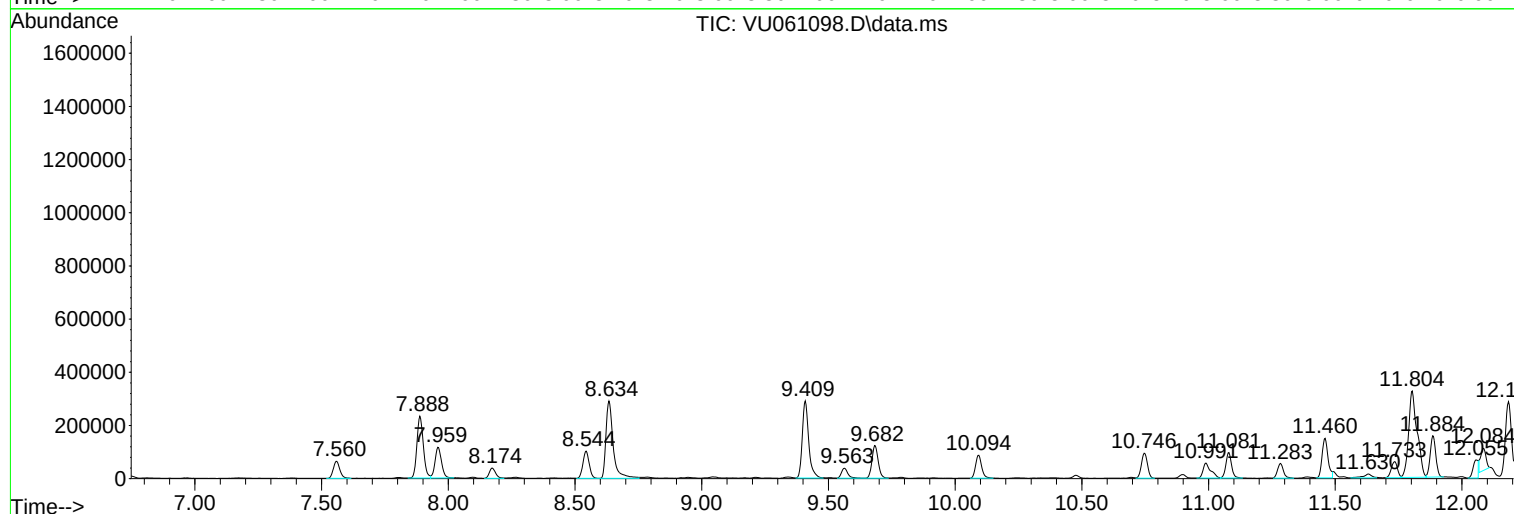
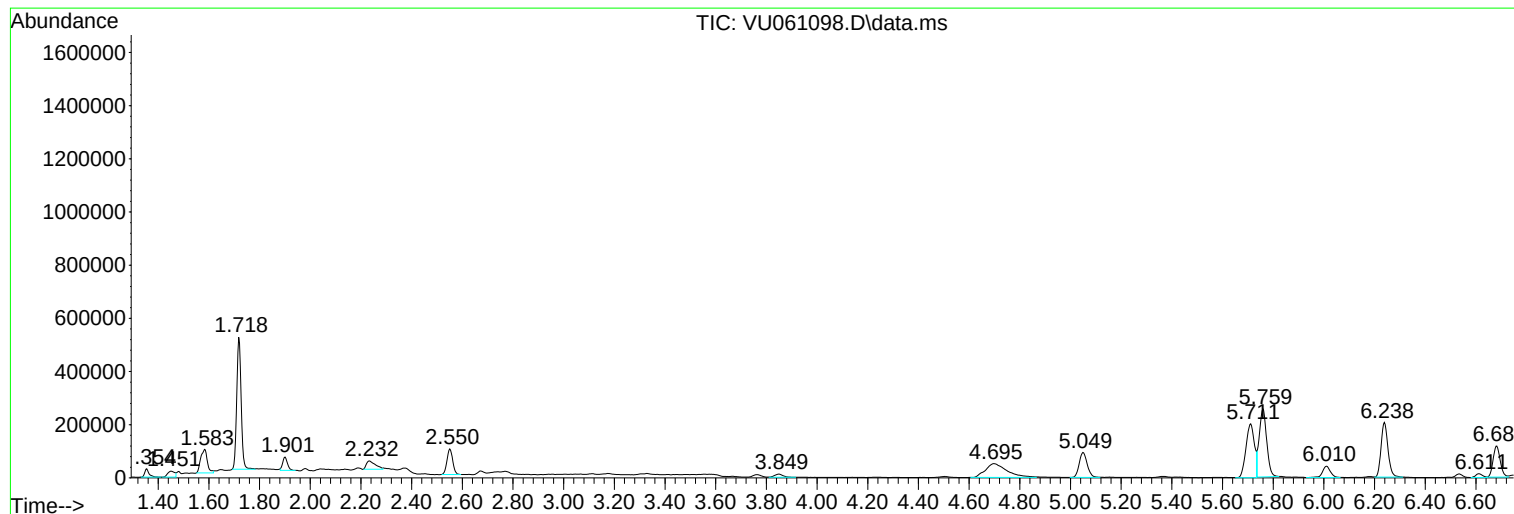
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100124WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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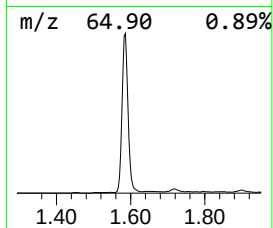
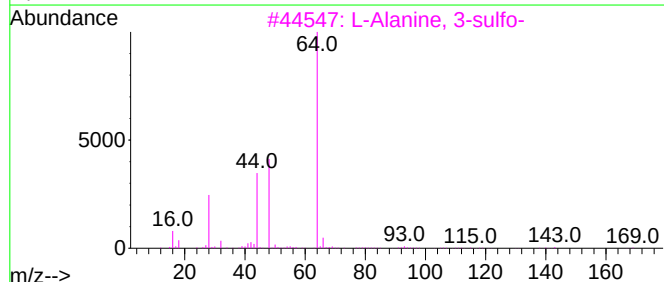
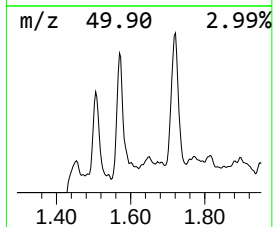
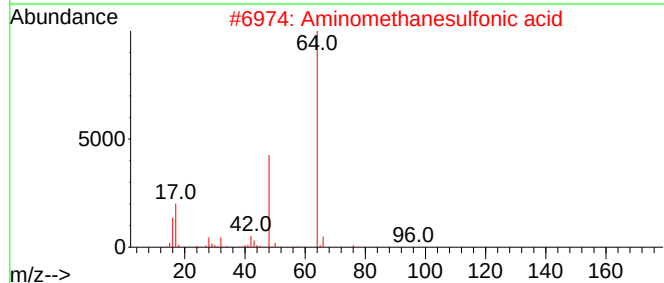
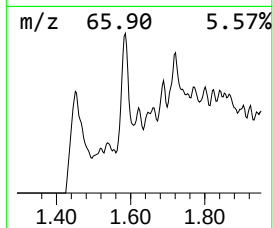
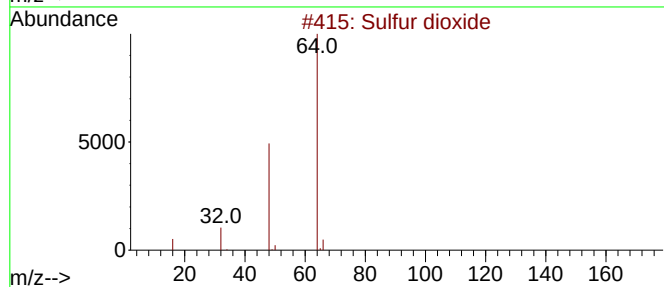
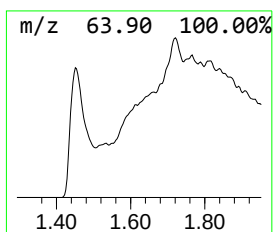
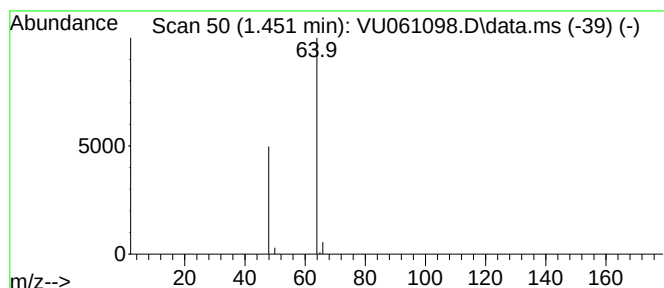
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100124WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Sulfur dioxide Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.451	0.56 ug/L	44439	1,4-Difluorobenzene	6.238

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfur dioxide	64	O2S	007446-09-5	90
2			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	64
3			L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
5			Sulfur dioxide	64	O2S	007446-09-5	4



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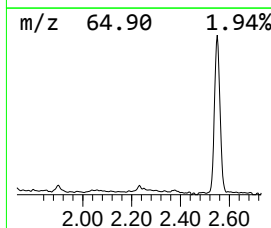
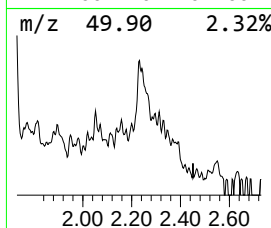
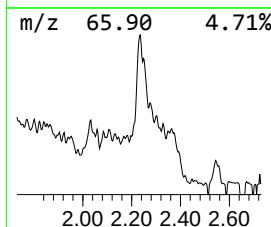
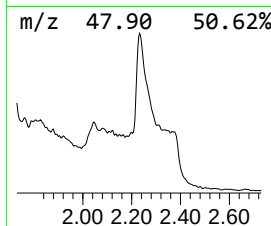
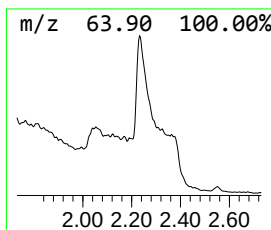
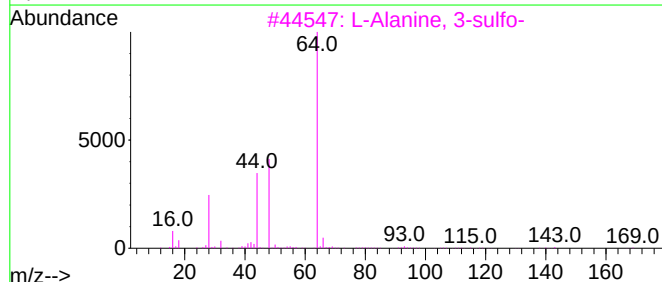
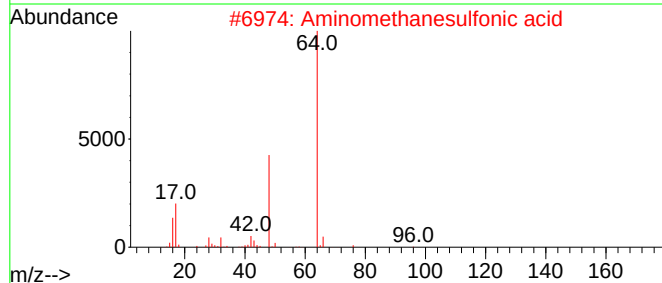
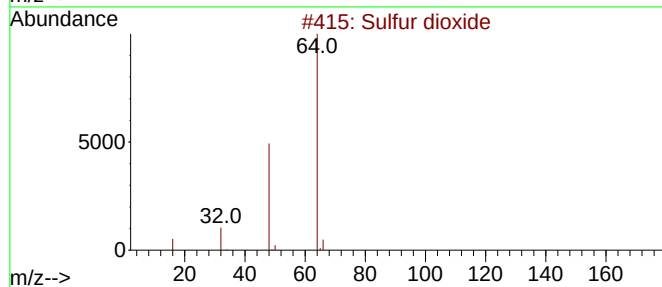
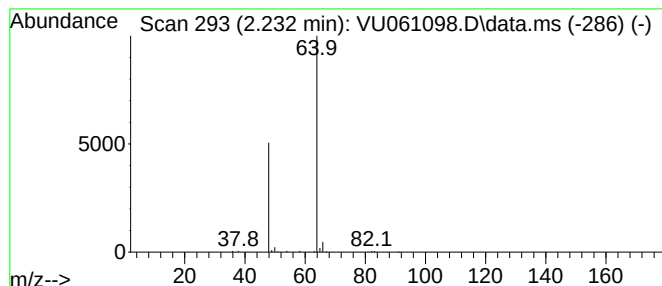
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100124WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.232	0.98 ug/L	76895	1,4-Difluorobenzene	6.238

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfur dioxide	64	O2S	007446-09-5	74
2			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
3			L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4			2-Aminoethyl hydrogen sulfate	141	C2H7NO4S	000926-39-6	9
5			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9



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ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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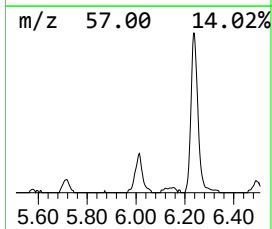
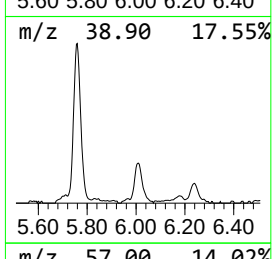
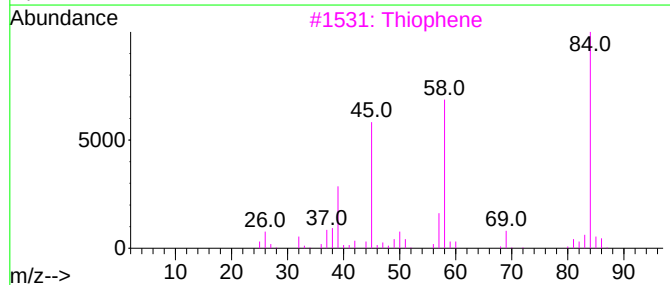
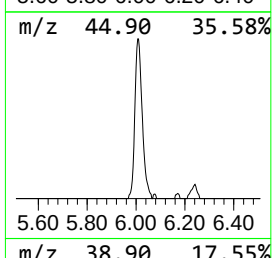
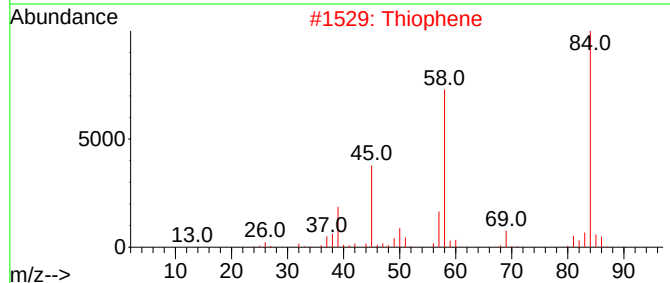
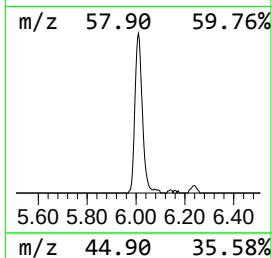
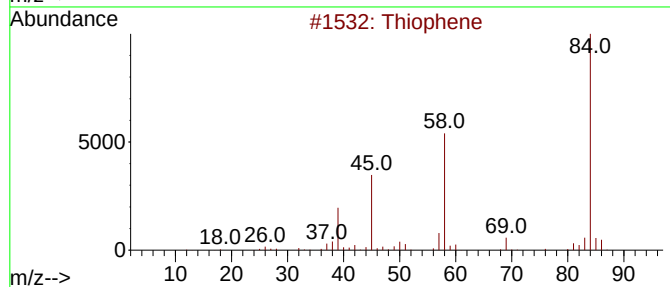
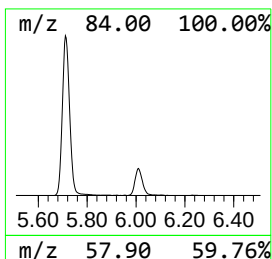
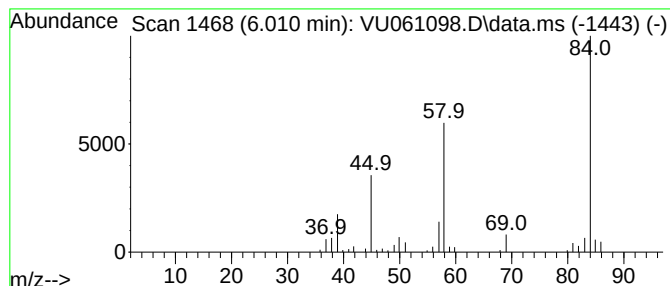
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100124WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Thiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.010	1.18 ug/L	93124	1,4-Difluorobenzene	6.238

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiophene	84	C4H4S	000110-02-1	97
2			Thiophene	84	C4H4S	000110-02-1	97
3			Thiophene	84	C4H4S	000110-02-1	91
4			Thiophene	84	C4H4S	000110-02-1	91
5			4H-1,2,4-Triazol-4-amine	84	C2H4N4	000584-13-4	25



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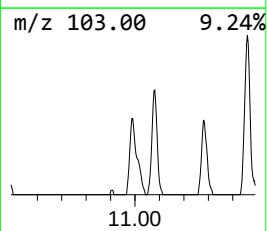
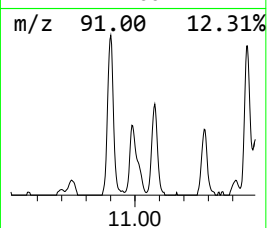
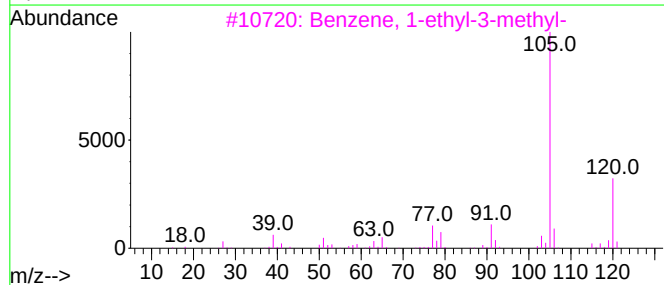
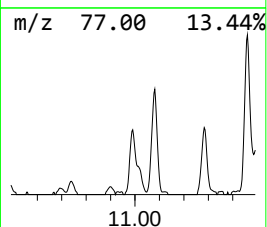
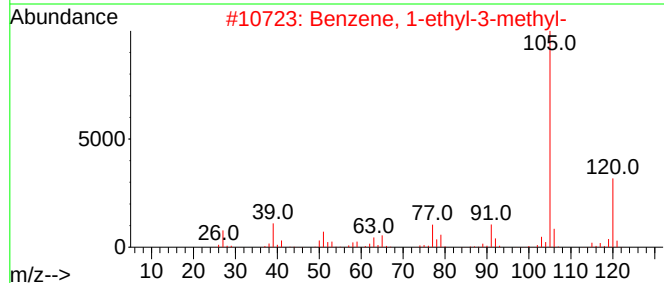
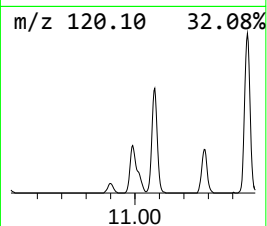
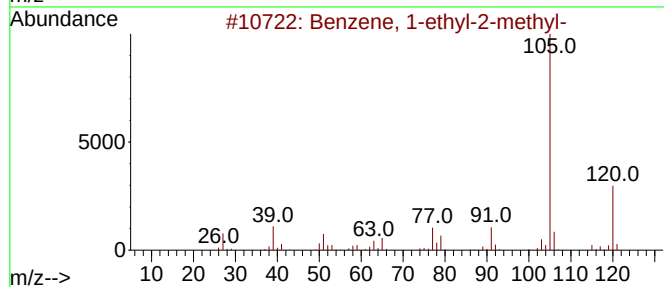
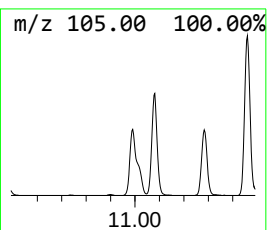
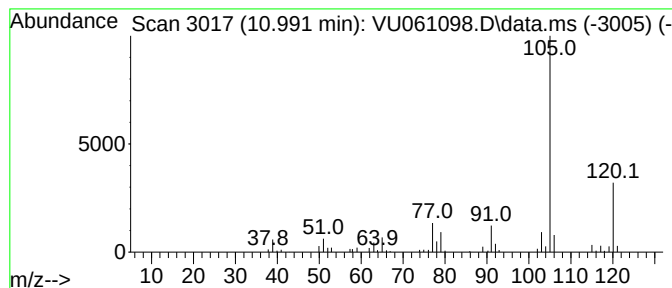
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100124WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.991	0.78 ug/L	119351	1,4-Dichlorobenzene-d4	11.804

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



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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

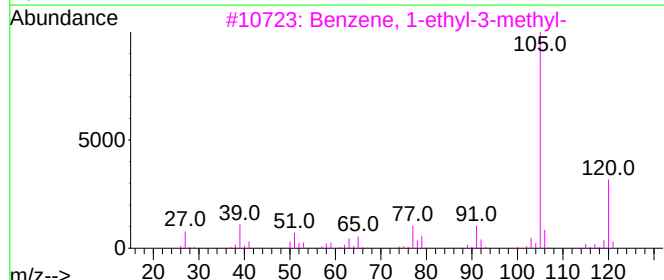
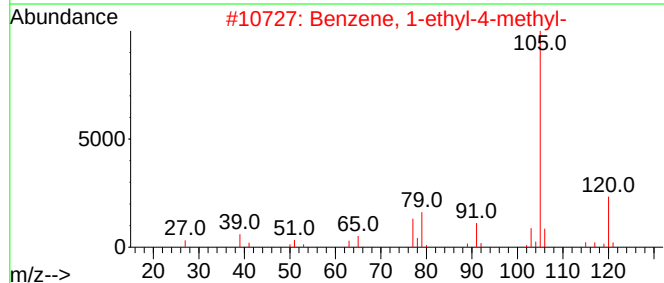
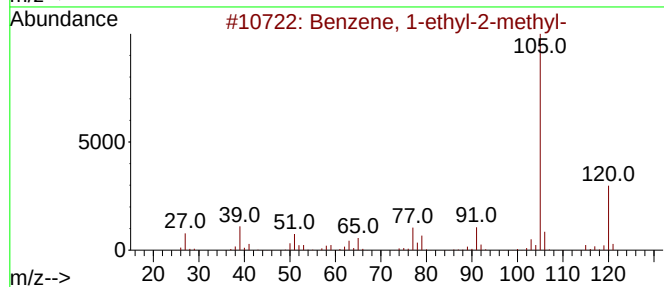
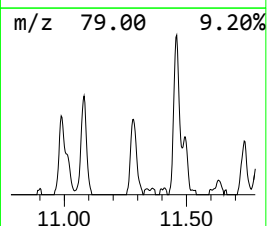
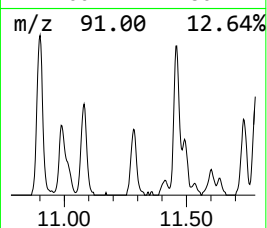
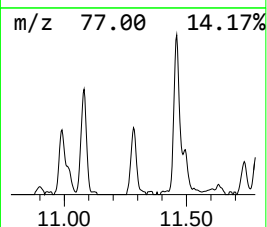
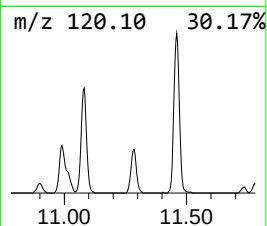
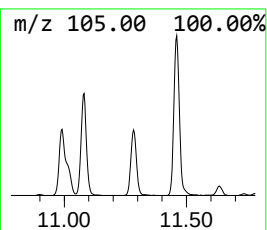
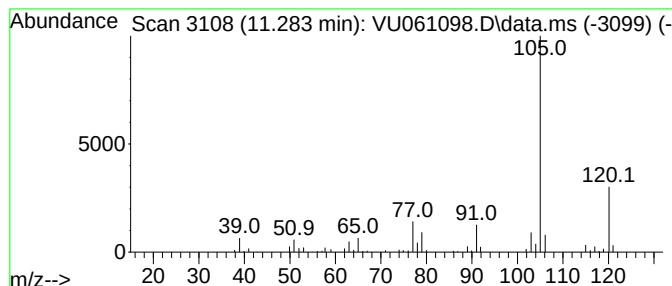
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.283	0.58 ug/L	87947	1,4-Dichlorobenzene-d4	11.804

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-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

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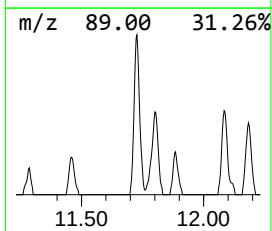
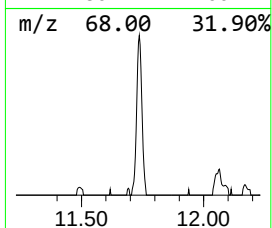
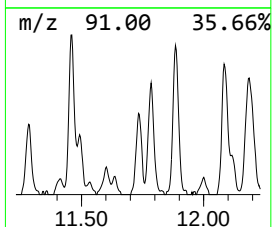
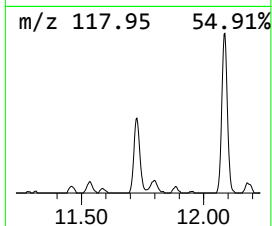
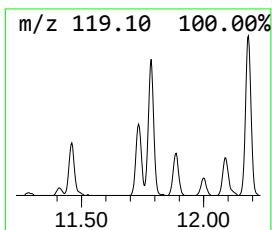
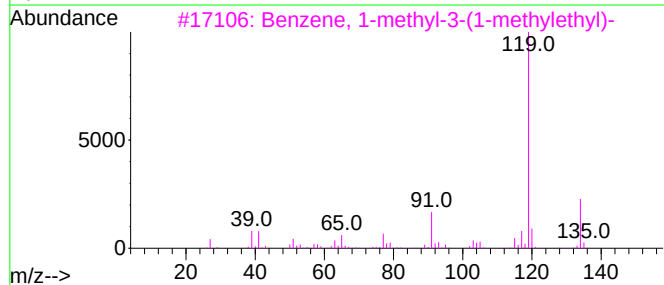
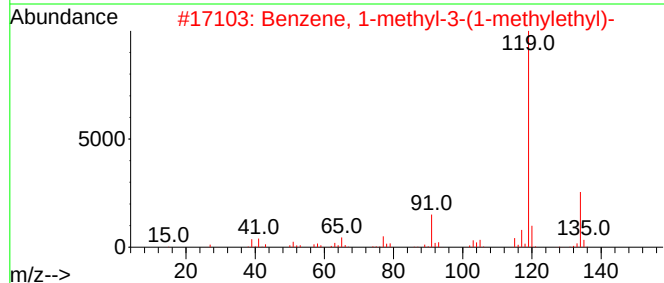
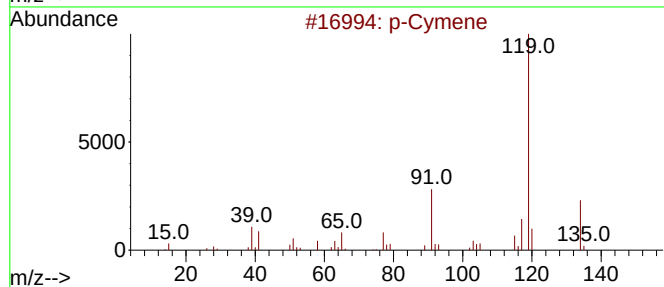
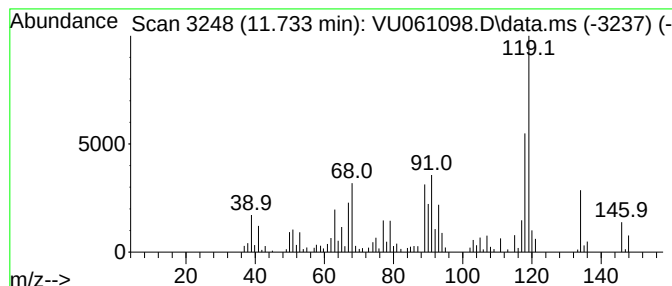
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 p-Cymene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.733	0.68 ug/L	102950	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	86
2			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	55
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	55
4			p-Cymene	134	C10H14	000099-87-6	55
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

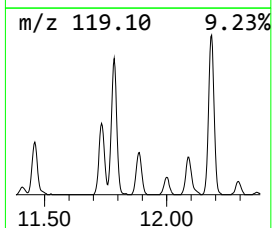
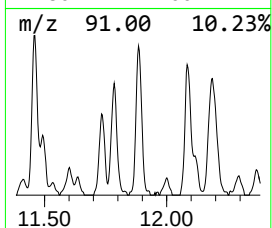
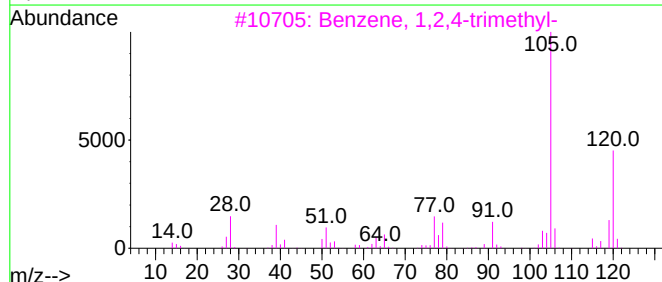
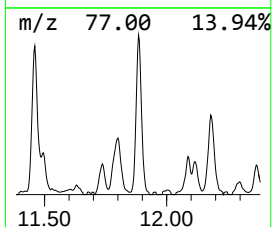
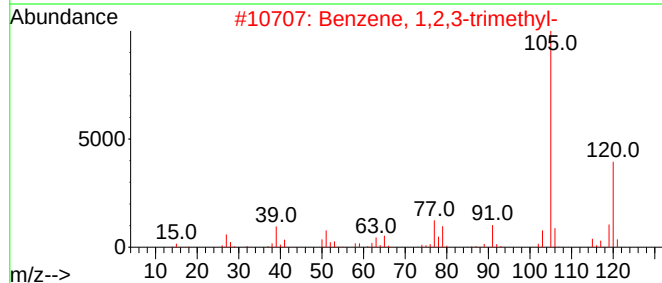
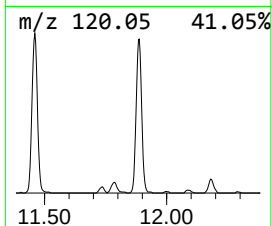
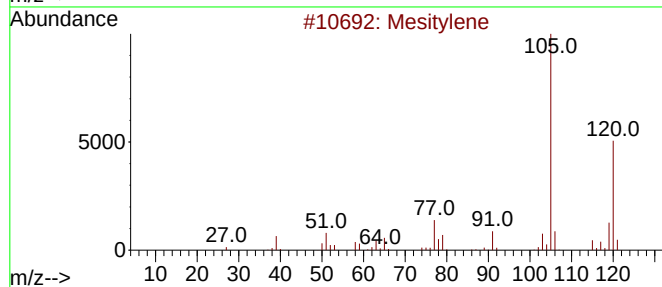
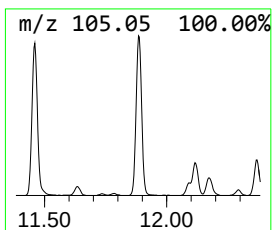
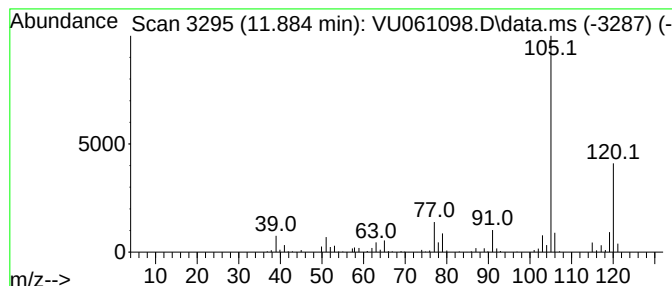
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.884	1.57 ug/L	239479	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	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

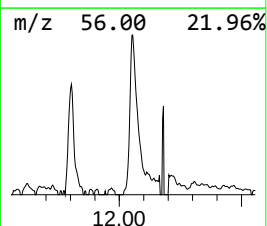
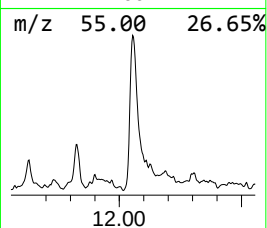
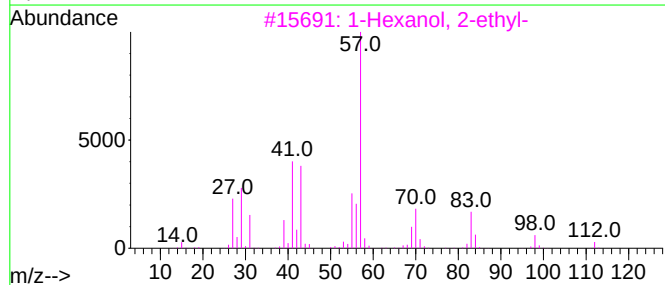
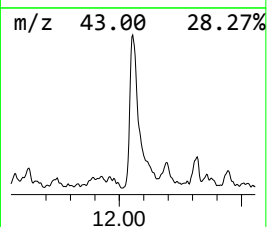
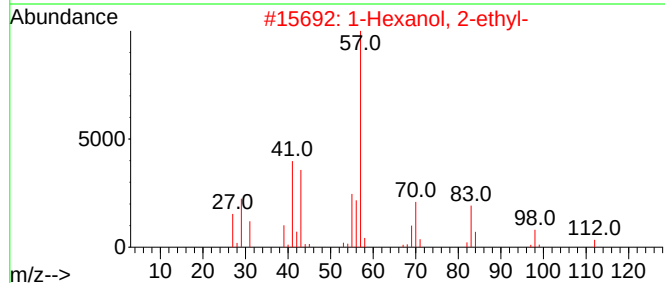
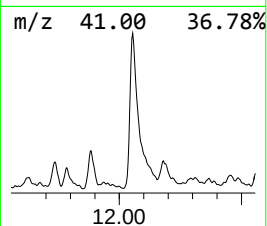
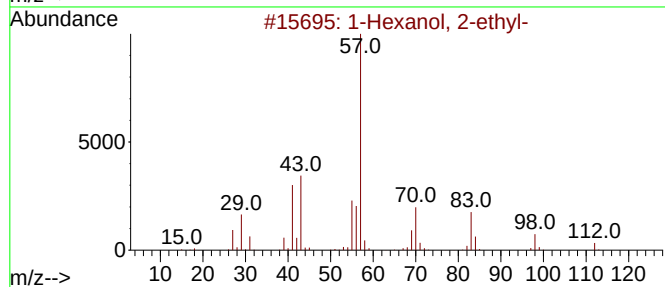
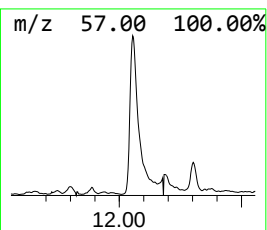
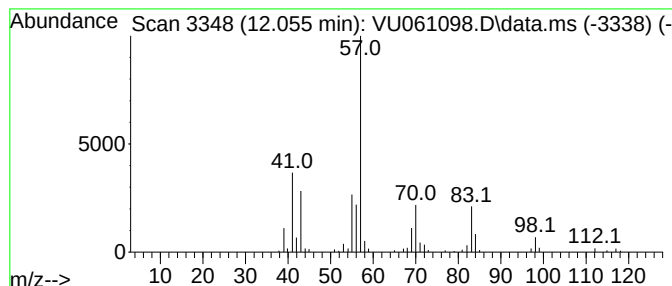
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 1-Hexanol, 2-ethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.055	0.63 ug/L	96647	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
5			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

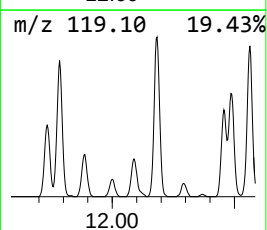
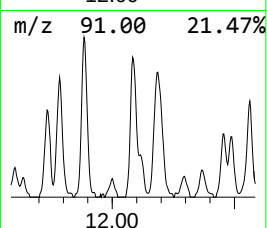
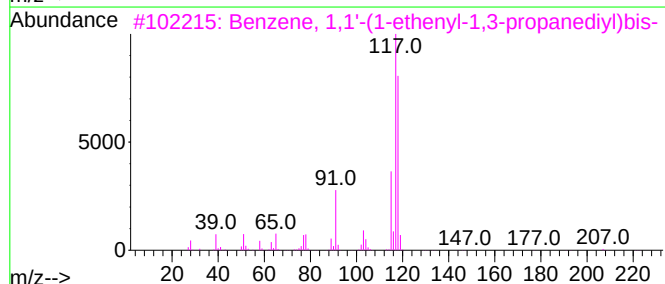
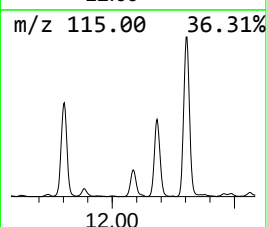
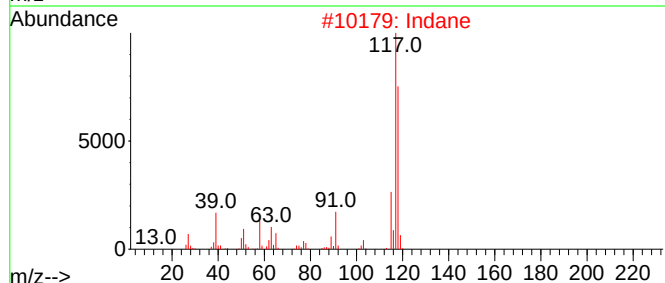
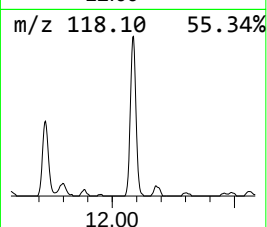
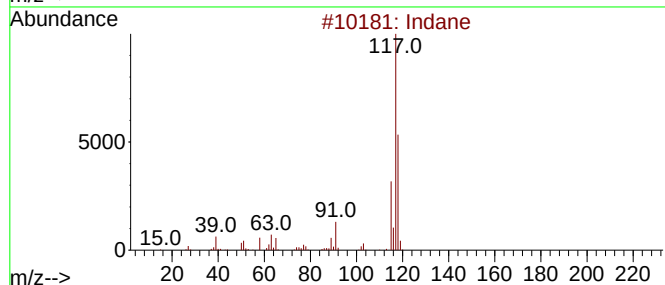
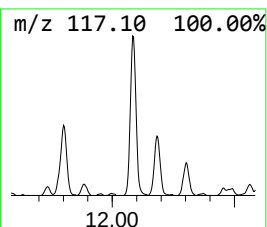
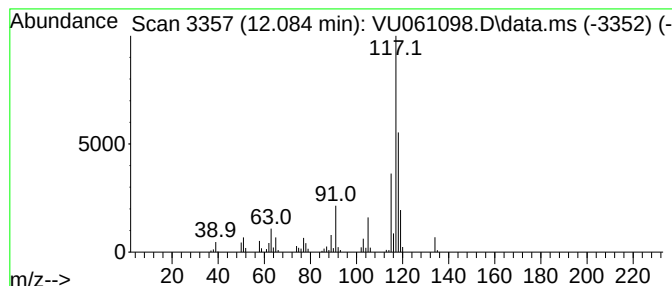
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Indane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.084	0.80 ug/L	121551	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	92
2	Indane			118	C9H10	000496-11-7	76
3	Benzene, 1,1'-(1-ethenyl-1,3-pro...			222	C17H18	061141-97-7	72
4	Indane			118	C9H10	000496-11-7	70
5	Benzene, 2-propenyl-			118	C9H10	000300-57-2	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

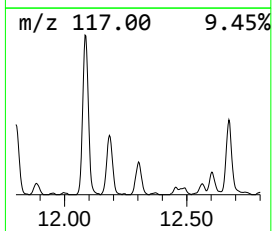
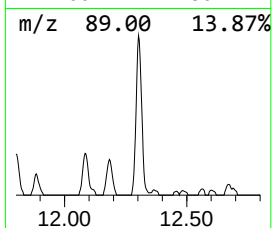
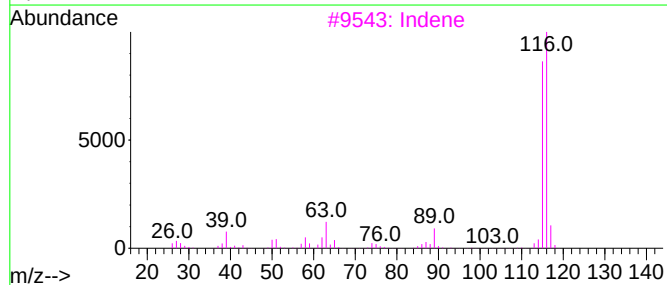
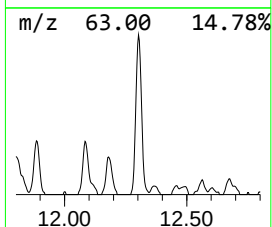
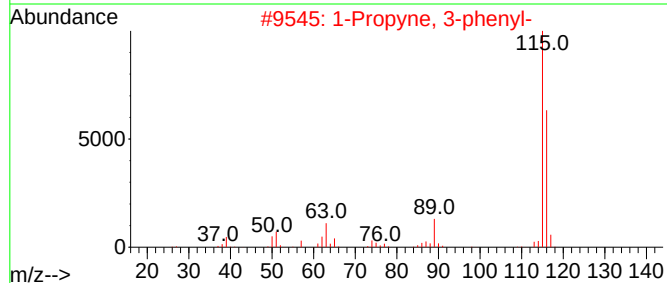
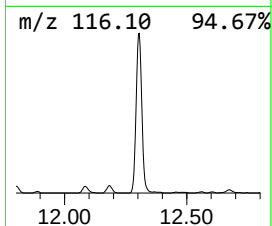
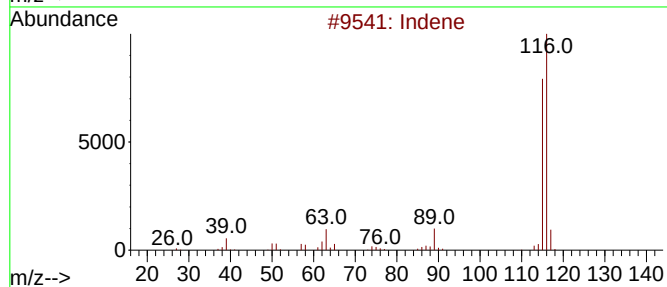
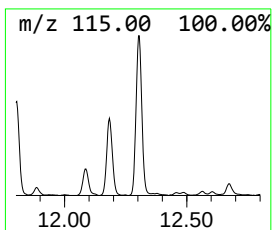
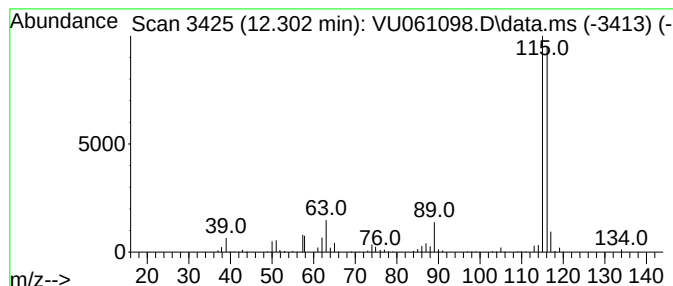
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Indene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.302	1.79 ug/L	273186	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	95
2			1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	95
3			Indene	116	C9H8	000095-13-6	94
4			Indene	116	C9H8	000095-13-6	94
5			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

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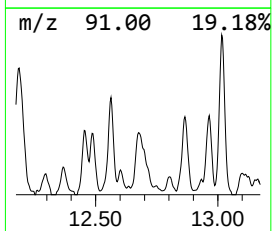
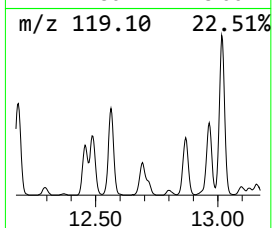
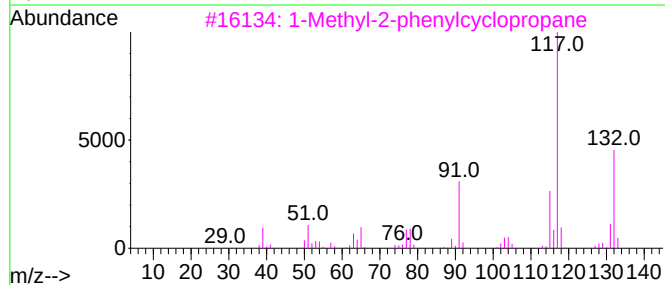
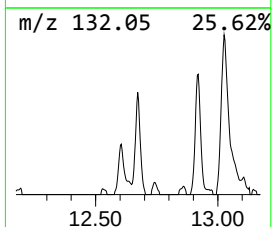
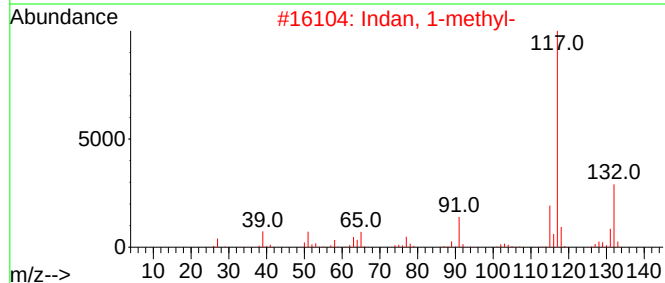
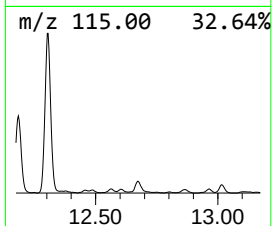
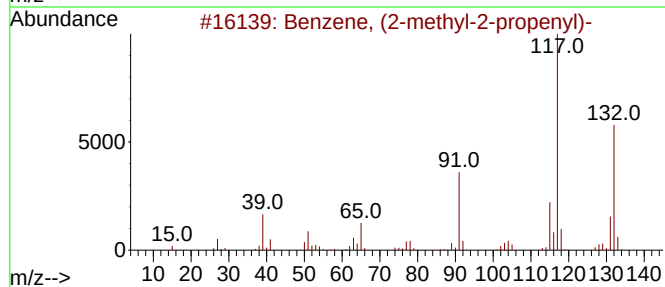
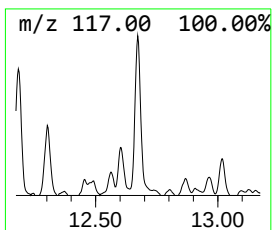
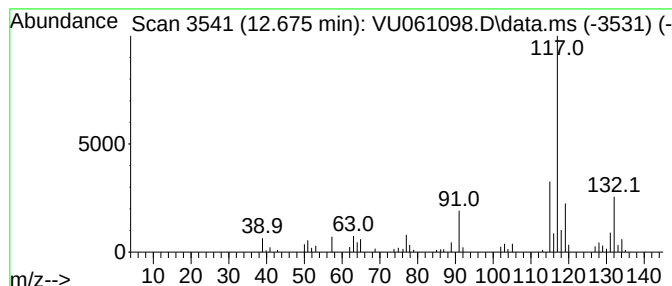
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Benzene, (2-methyl-2-propen... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.675	0.56 ug/L	85509	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
2			Indan, 1-methyl-	132	C10H12	000767-58-8	81
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81
4			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	76
5			Indan, 1-methyl-	132	C10H12	000767-58-8	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

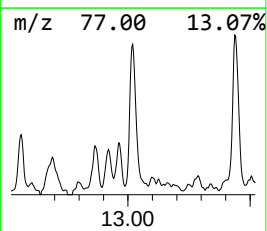
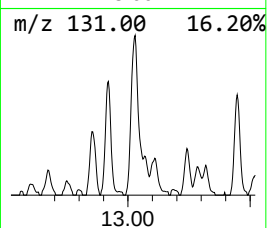
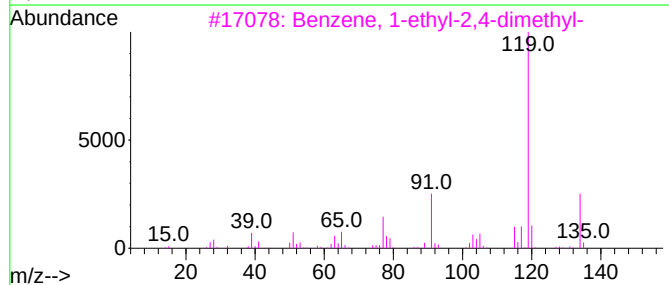
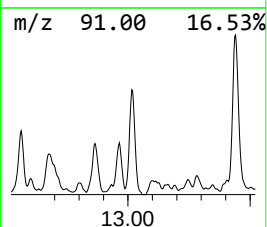
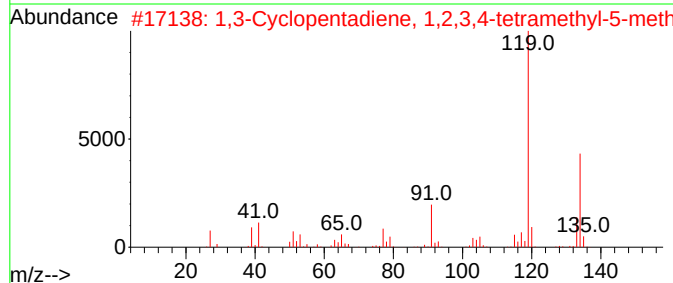
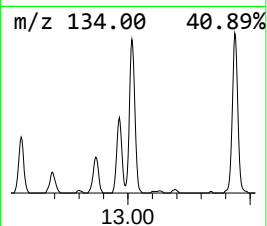
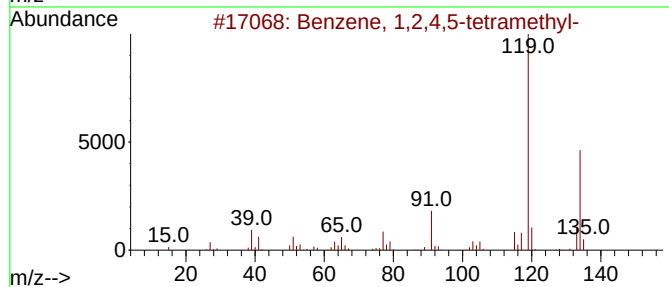
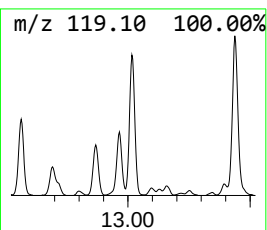
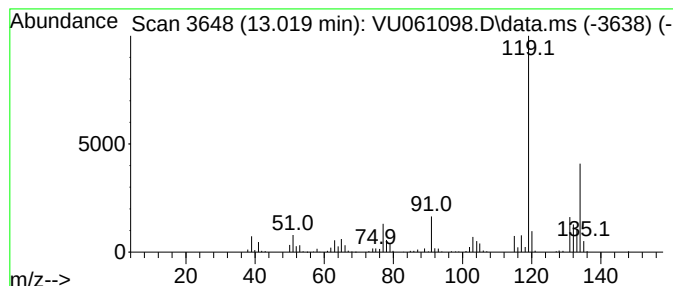
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 14 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.020	1.20 ug/L	182573	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	95
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

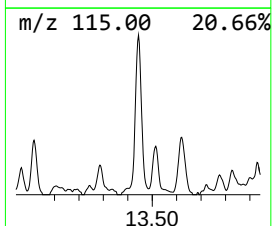
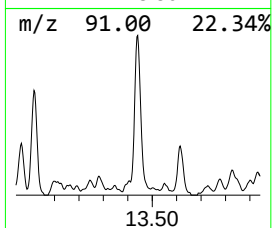
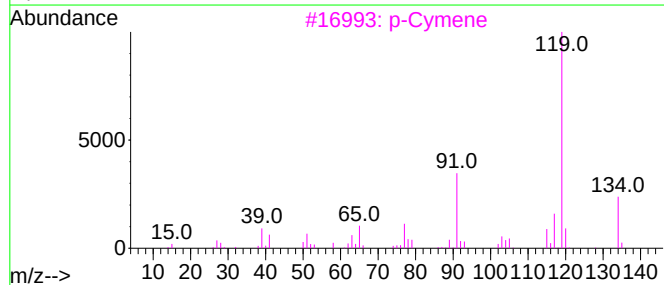
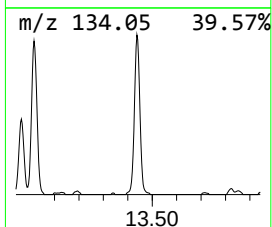
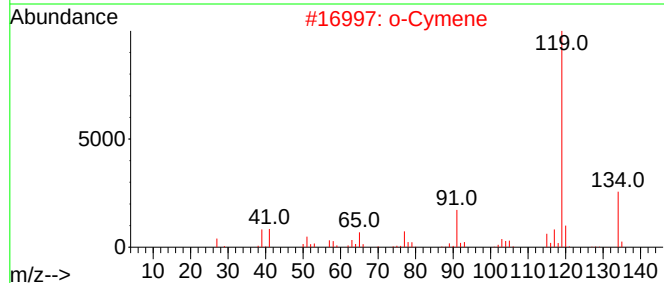
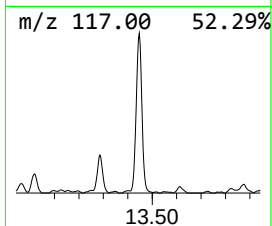
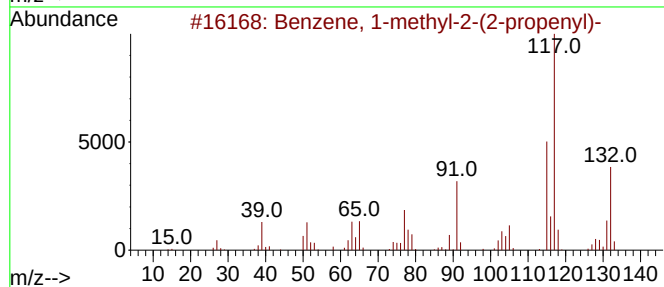
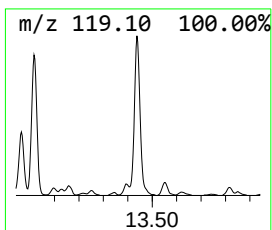
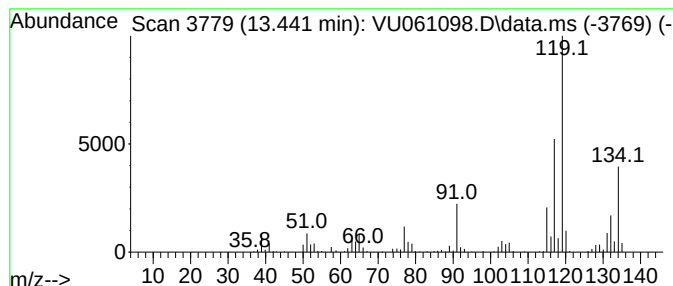
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Benzene, 1-methyl-2-(2-prop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.441	1.66 ug/L	253286	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
2			o-Cymene	134	C10H14	000527-84-4	64
3			p-Cymene	134	C10H14	000099-87-6	64
4			p-Cymene	134	C10H14	000099-87-6	64
5			o-Cymene	134	C10H14	000527-84-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

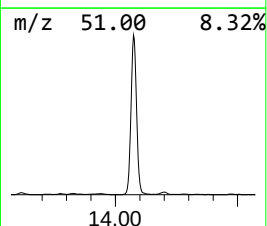
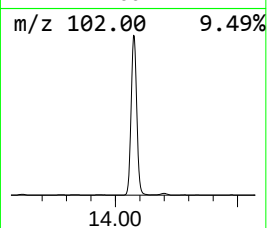
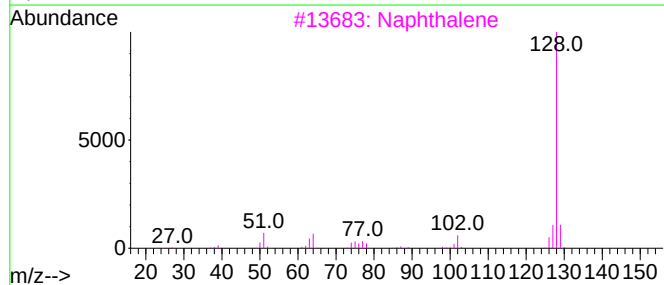
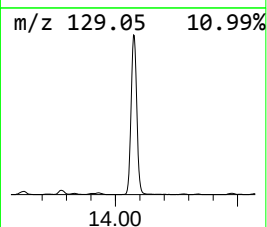
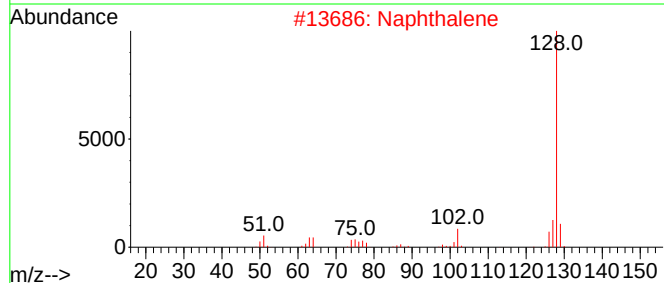
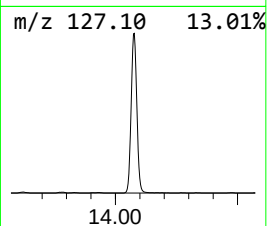
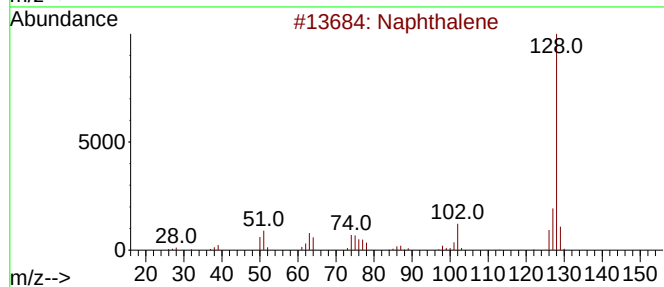
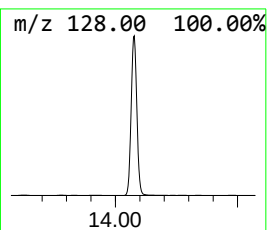
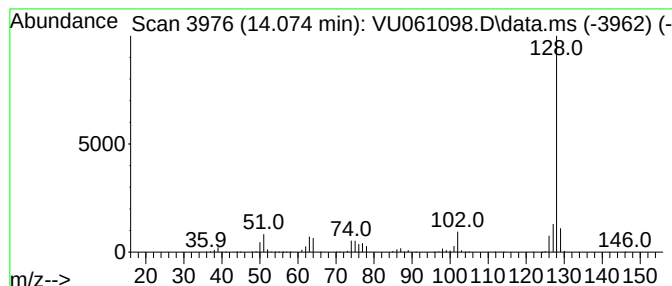
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 16 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.074	17.60 ug/L	2679830	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	97
2			Naphthalene	128	C10H8	000091-20-3	95
3			Naphthalene	128	C10H8	000091-20-3	94
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

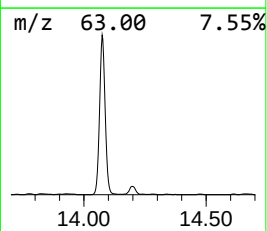
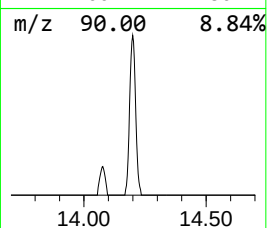
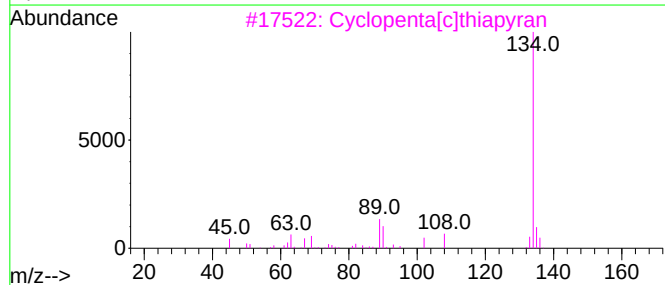
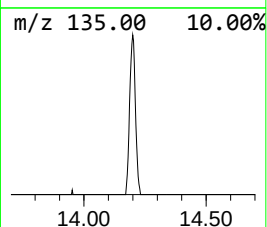
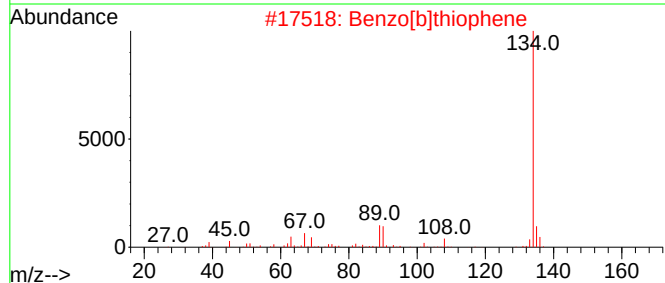
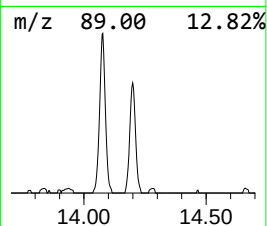
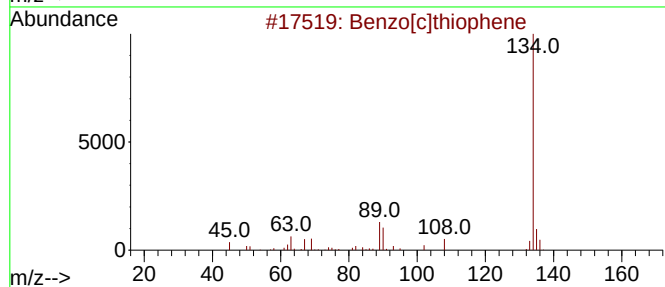
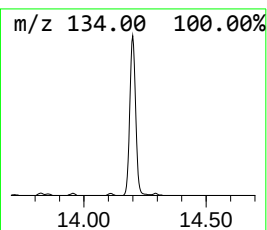
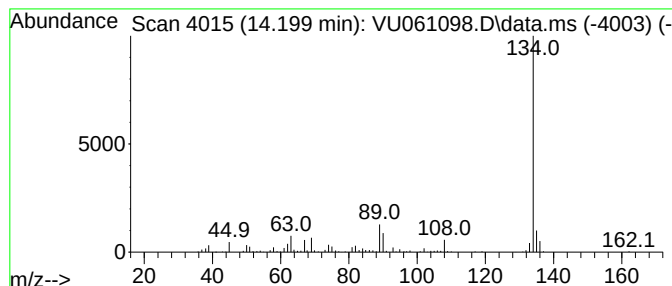
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 17 Benzo[c]thiophene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.200	0.92 ug/L	139704	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	93
5			Benzo[b]thiophene	134	C8H6S	000095-15-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

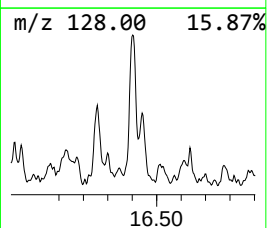
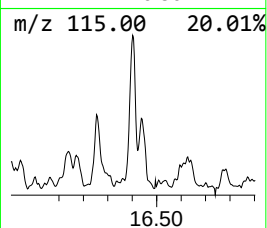
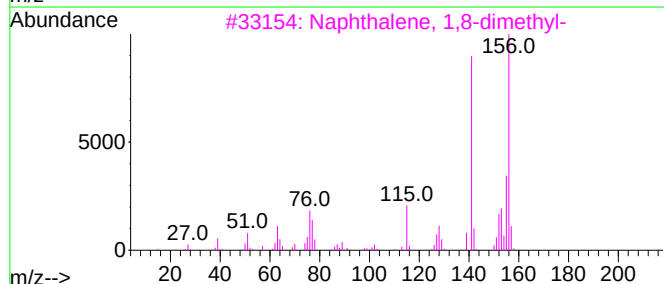
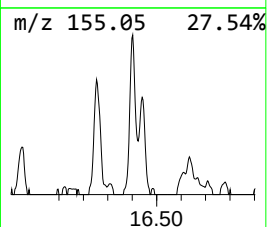
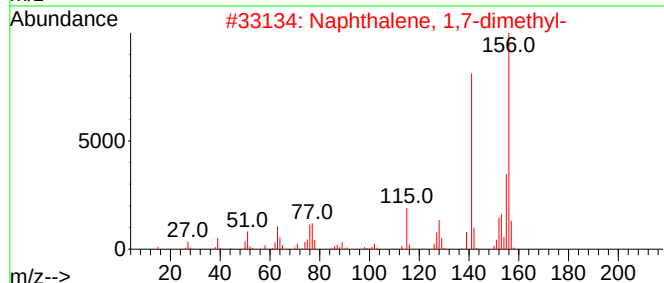
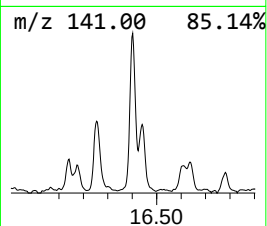
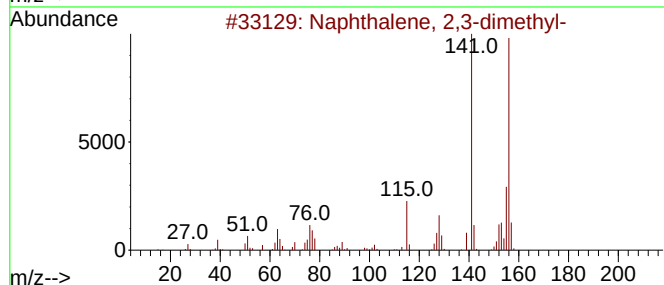
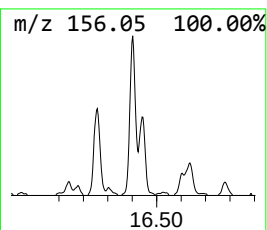
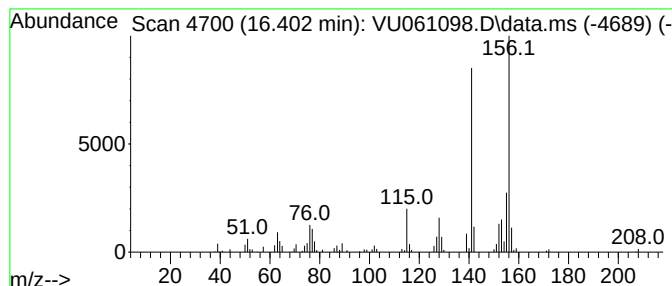
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 20 Naphthalene, 2,3-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.402	0.57 ug/L	86119	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	98
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101024\
Data File : VU061098.D
Acq On : 10 Oct 2024 18:49
Operator : MD/SY
Sample : P4380-14
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E27H3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100124WMA.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
Sulfur dioxide	1.451	0.6	ug/L	44439	1	6.238	393278	5.0
unknown-01	2.232	1.0	ug/L	76895	1	6.238	393278	5.0
Thiophene	6.010	1.2	ug/L	93124	1	6.238	393278	5.0
Benzene, 1-ethy...	10.991	0.8	ug/L	119351	3	11.804	761290	5.0
Benzene, 1-ethy...	11.283	0.6	ug/L	87947	3	11.804	761290	5.0
p-Cymene	11.733	0.7	ug/L	102950	3	11.804	761290	5.0
Benzene, 1,2,3-...	11.884	1.6	ug/L	239479	3	11.804	761290	5.0
1-Hexanol, 2-et...	12.055	0.6	ug/L	96647	3	11.804	761290	5.0
Indane	12.084	0.8	ug/L	121551	3	11.804	761290	5.0
Indene	12.302	1.8	ug/L	273186	3	11.804	761290	5.0
Benzene, (2-met...	12.675	0.6	ug/L	85509	3	11.804	761290	5.0
Benzene, 1,2,4,...	13.020	1.2	ug/L	182573	3	11.804	761290	5.0
Benzene, 1-meth...	13.441	1.7	ug/L	253286	3	11.804	761290	5.0
Azulene	14.074	17.6	ug/L	2679830	3	11.804	761290	5.0
Benzo[c]thiophene	14.200	0.9	ug/L	139704	3	11.804	761290	5.0
Naphthalene, 2,...	16.402	0.6	ug/L	86119	3	11.804	761290	5.0