

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
 Data File : VU058757.D
 Acq On : 18 Apr 2024 13:04
 Operator : MD/SY
 Sample : P2181-01
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AK9

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\SFAMUTR040424WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VU058757.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	14	20	34	rVB2	9106	13498	2.11%	0.317%
2	1.589	84	93	104	rVB	23486	27699	4.33%	0.651%
3	1.904	183	191	203	rVB	18037	22502	3.52%	0.529%
4	2.557	383	394	410	rVB	40564	66851	10.46%	1.571%
5	4.634	1028	1040	1070	rBV	18572	56816	8.89%	1.335%
6	5.052	1156	1170	1187	rBV	40853	90634	14.18%	2.130%
7	5.717	1359	1377	1389	rBV2	78733	203933	31.91%	4.792%
8	6.242	1528	1540	1564	rBV	100160	192504	30.12%	4.524%
9	6.682	1666	1677	1693	rBV	48740	94144	14.73%	2.212%
10	7.560	1941	1950	1970	rBV	21862	38882	6.08%	0.914%
11	7.894	2042	2054	2068	rBV	88496	158822	24.85%	3.732%
12	7.968	2069	2077	2093	rVB2	5095	9759	1.53%	0.229%
13	8.177	2132	2142	2157	rBV3	11431	20663	3.23%	0.486%
14	8.631	2273	2283	2315	rBV2	72455	146327	22.90%	3.438%
15	9.412	2514	2526	2542	rBV	140573	234574	36.70%	5.512%
16	9.495	2542	2552	2565	rVV	143035	222648	34.84%	5.232%
17	9.566	2565	2574	2587	rVB	17563	27964	4.38%	0.657%
18	9.692	2599	2613	2632	rBB2	19699	33318	5.21%	0.783%
19	10.097	2730	2739	2756	rVB2	13805	22640	3.54%	0.532%
20	10.479	2847	2858	2875	rBV	75613	119582	18.71%	2.810%
21	10.749	2931	2942	2958	rBB	30442	48659	7.61%	1.143%
22	10.901	2979	2989	3002	rBV	9402	14517	2.27%	0.341%
23	11.020	3012	3026	3037	rBV2	6558	11630	1.82%	0.273%
24	11.290	3101	3110	3115	rBV	5374	7636	1.19%	0.179%
25	11.463	3155	3164	3175	rVB	7402	11456	1.79%	0.269%
26	11.637	3210	3218	3227	rVB	14489	20909	3.27%	0.491%
27	11.807	3258	3271	3286	rBV	144016	236174	36.95%	5.550%
28	11.888	3289	3296	3307	rVB2	5353	7377	1.15%	0.173%
29	12.087	3347	3358	3373	rBV	96367	151758	23.75%	3.566%
30	12.187	3378	3389	3404	rBV	96709	157869	24.70%	3.710%
31	12.299	3414	3424	3437	rBV4	24831	43785	6.85%	1.029%
32	12.373	3439	3447	3461	rVB2	6683	10319	1.61%	0.242%
33	12.675	3527	3541	3559	rBV	43259	69384	10.86%	1.630%
34	12.923	3609	3618	3624	rBV2	10669	17548	2.75%	0.412%
35	12.968	3624	3632	3639	rVV	16882	28213	4.41%	0.663%
36	13.019	3639	3648	3664	rVV4	22808	55386	8.67%	1.301%

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR040424WMA.M
 Title : TRACE VOA SFAM1.0

37	13.103	3668	3674	3690	rVB6	8252	16969	2.66%	0.399%
38	13.248	3711	3719	3725	rVV4	6035	8594	1.34%	0.202%
39	13.290	3725	3732	3738	rVV2	11865	18647	2.92%	0.438%
40	13.318	3738	3741	3755	rVB2	5306	7878	1.23%	0.185%
41	13.447	3771	3781	3793	rVB3	20734	33827	5.29%	0.795%
42	13.515	3793	3802	3813	rBV4	11990	18354	2.87%	0.431%
43	13.621	3822	3835	3848	rVB3	24444	41734	6.53%	0.981%
44	13.781	3874	3885	3892	rBV4	15951	25751	4.03%	0.605%
45	13.830	3892	3900	3911	rVV6	6632	11410	1.79%	0.268%
46	13.904	3916	3923	3928	rVV2	8421	12972	2.03%	0.305%
47	13.952	3928	3938	3951	rVB7	19781	39205	6.13%	0.921%
48	14.080	3958	3978	3998	rBV	370237	639098	100.00%	15.018%
49	14.199	3998	4015	4022	rVV3	26523	59282	9.28%	1.393%
50	14.238	4022	4027	4034	rVV5	16131	27555	4.31%	0.648%
51	14.280	4035	4040	4050	rVV5	12672	19491	3.05%	0.458%
52	14.335	4050	4057	4066	rVB8	5979	8753	1.37%	0.206%
53	14.669	4151	4161	4170	rBV8	7143	16026	2.51%	0.377%
54	14.801	4190	4202	4211	rBV2	22026	34866	5.46%	0.819%
55	14.875	4211	4225	4237	rVB5	10958	24894	3.90%	0.585%
56	15.016	4258	4269	4278	rBV5	4155	6916	1.08%	0.163%
57	15.135	4293	4306	4307	rBV3	8557	10193	1.59%	0.240%
58	15.164	4308	4315	4326	rVB	16829	28630	4.48%	0.673%
59	15.254	4338	4343	4354	rVB8	8480	14023	2.19%	0.330%
60	15.405	4378	4390	4416	rBV7	44463	104833	16.40%	2.463%
61	15.904	4537	4545	4557	rBV7	3695	7055	1.10%	0.166%
62	16.006	4569	4577	4588	rVB5	5191	6922	1.08%	0.163%
63	16.177	4622	4630	4643	rVB3	14818	22085	3.46%	0.519%
64	16.605	4754	4763	4769	rBV2	13241	21637	3.39%	0.508%
65	16.640	4769	4774	4788	rVB2	10345	15214	2.38%	0.358%
66	17.090	4902	4914	4936	rBV	144445	246648	38.59%	5.796%
67	17.270	4960	4970	4979	rVB3	6270	9710	1.52%	0.228%

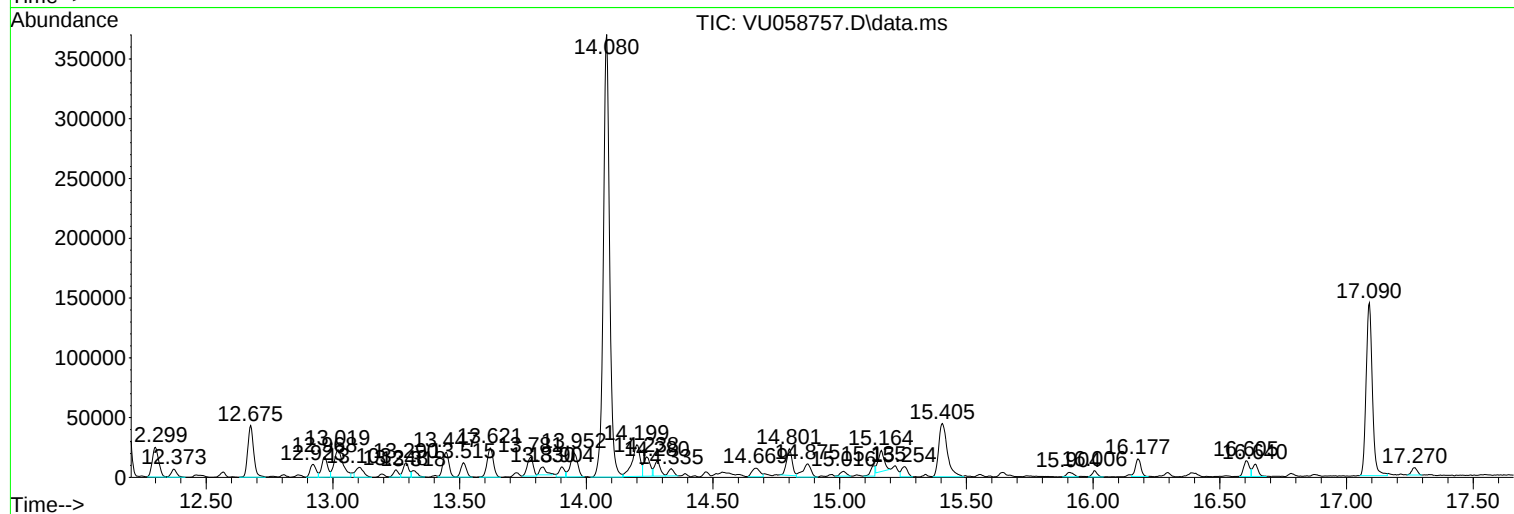
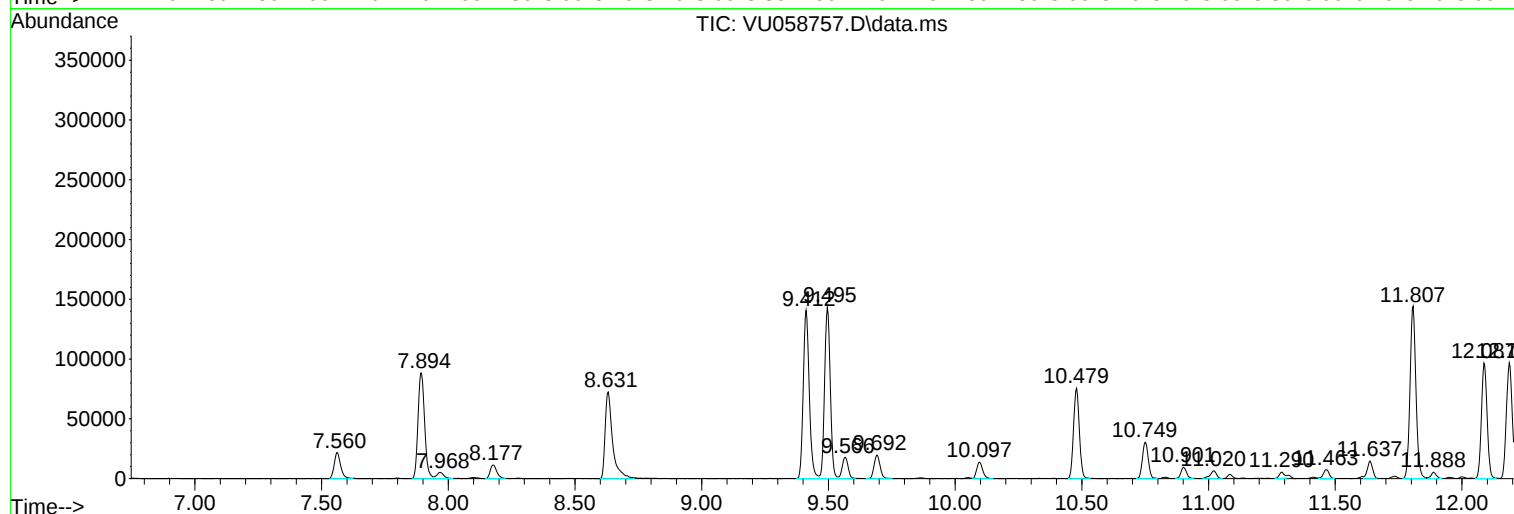
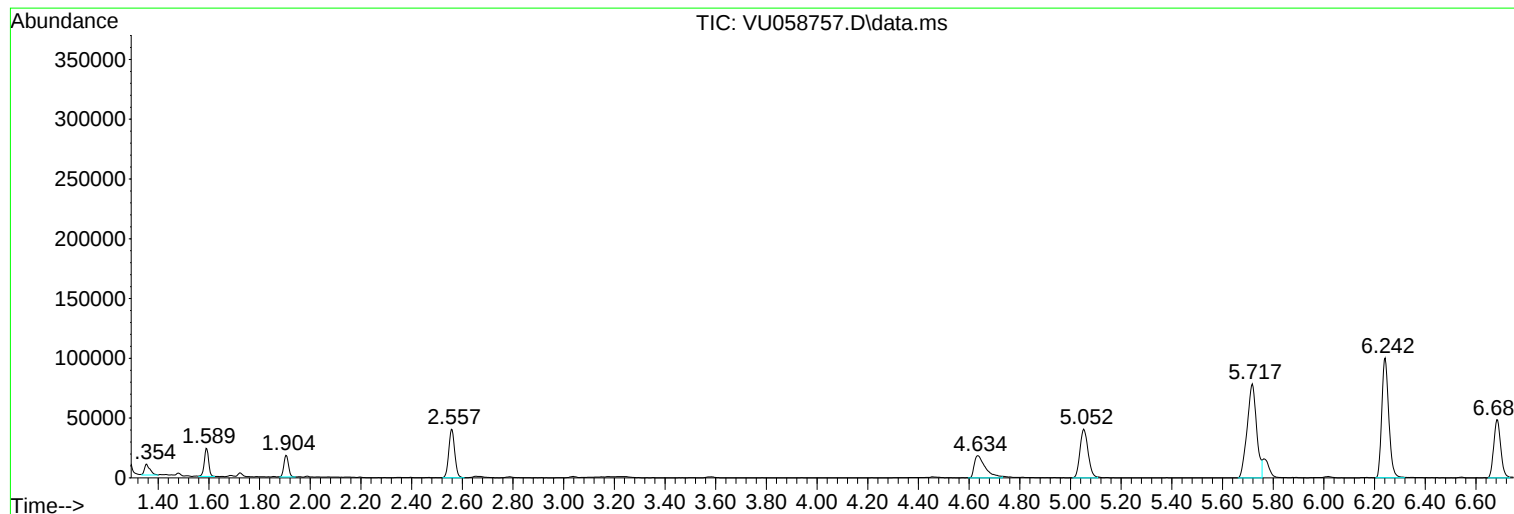
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Instrument :
MSVOA_U
ClientSampleId :
C0AK9

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR040424WMA.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 :
C0AK9

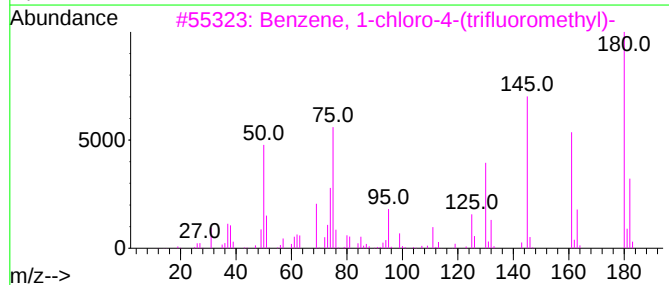
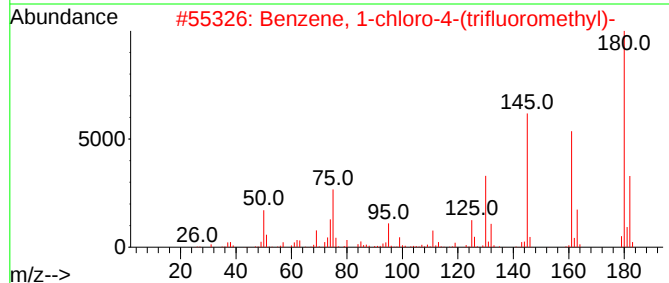
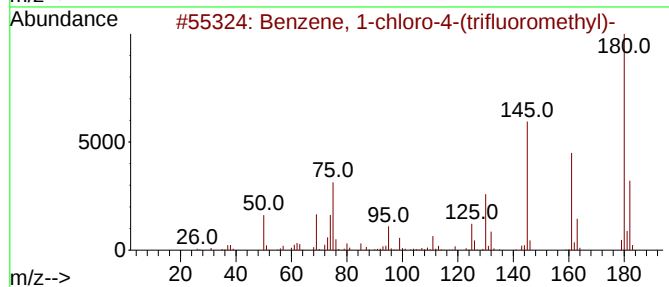
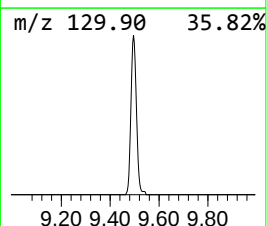
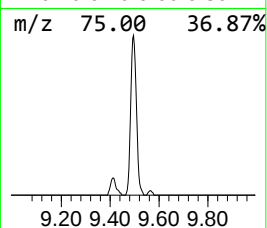
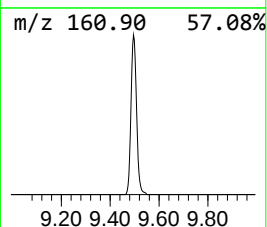
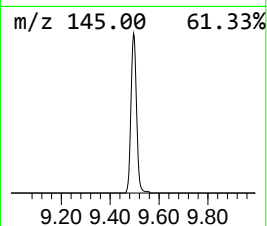
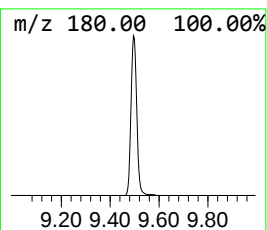
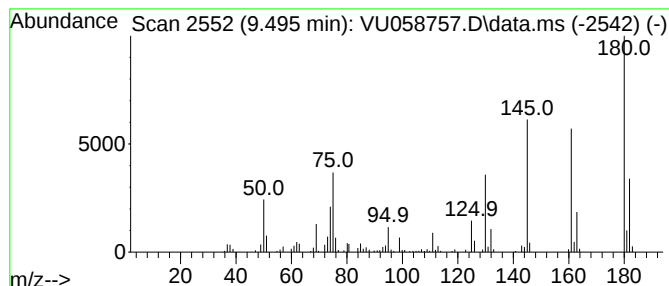
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR040424WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Benzene, 1-chloro-4-(triflu... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.495	4.75 ug/L	222648	Chlorobenzene-d5	9.412

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-4-(trifluorome...	180	C7H4ClF3	000098-56-6	98
2			Benzene, 1-chloro-4-(trifluorome...	180	C7H4ClF3	000098-56-6	98
3			Benzene, 1-chloro-4-(trifluorome...	180	C7H4ClF3	000098-56-6	95
4			Benzene, 1-chloro-4-(trifluorome...	180	C7H4ClF3	000098-56-6	93
5			Benzene, 1-chloro-3-(trifluorome...	180	C7H4ClF3	000098-15-7	93



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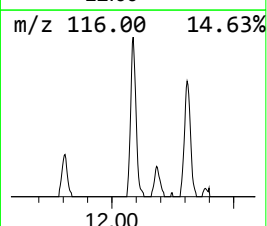
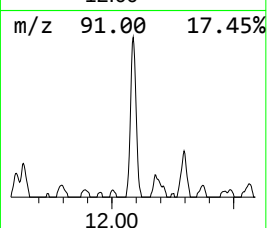
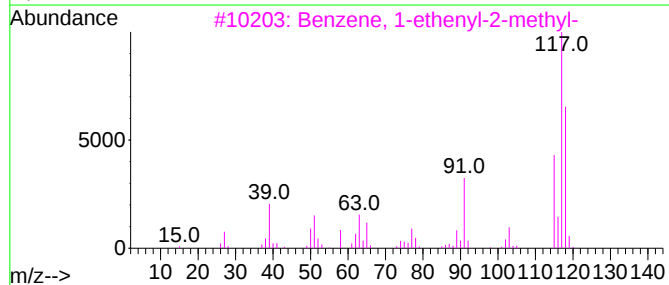
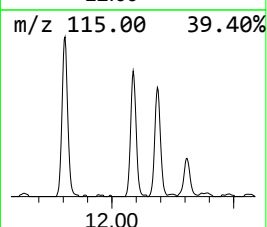
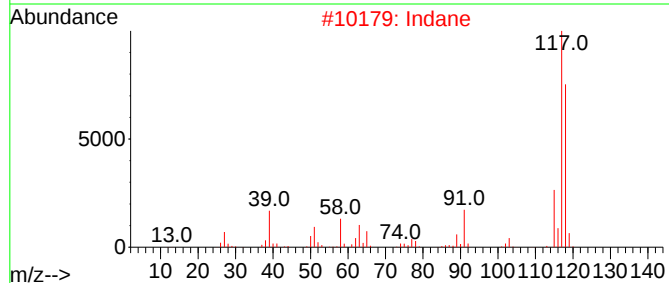
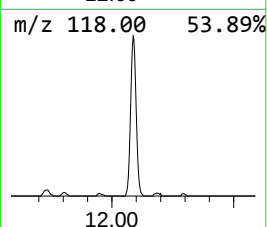
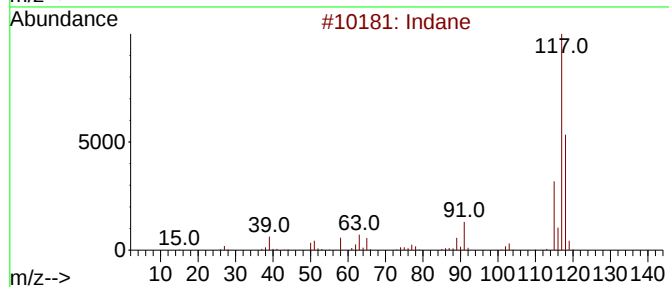
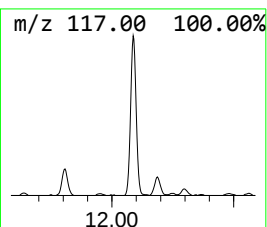
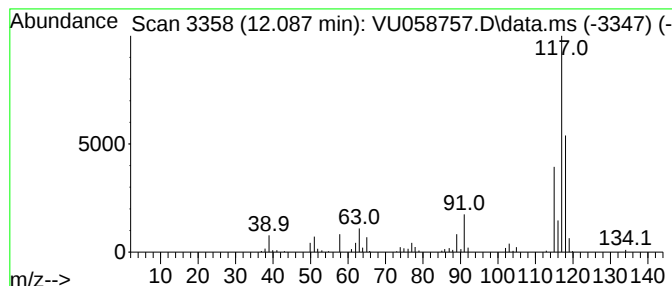
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR040424WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.087	3.21 ug/L	151758	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	76
3	Benzene, 1-ethenyl-2-methyl-			118	C9H10	000611-15-4	70
4	Indane			118	C9H10	000496-11-7	70
5	Benzene, 1-ethenyl-4-methyl-			118	C9H10	000622-97-9	68



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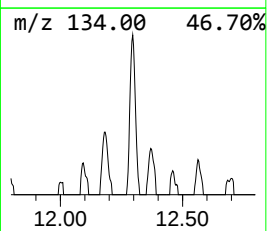
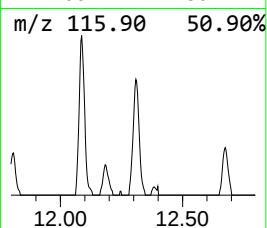
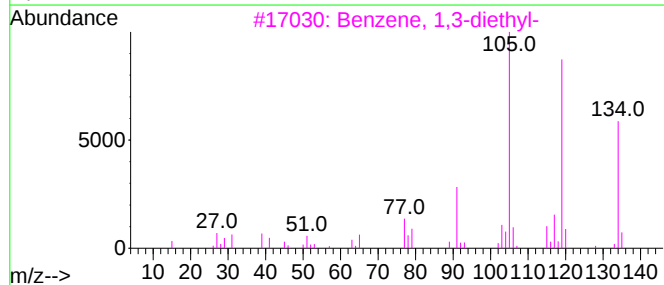
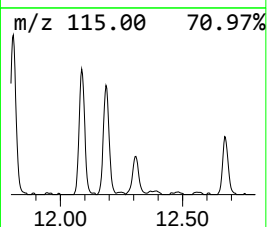
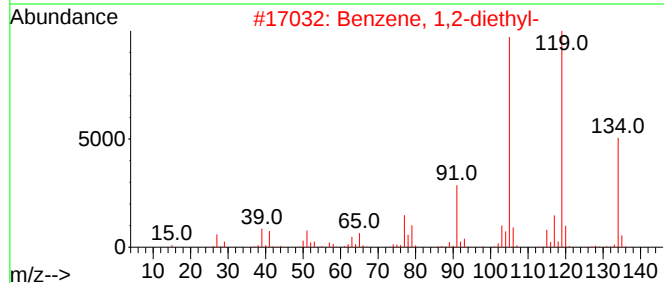
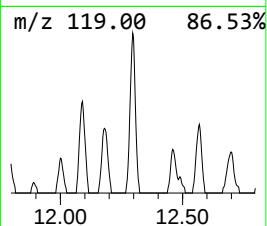
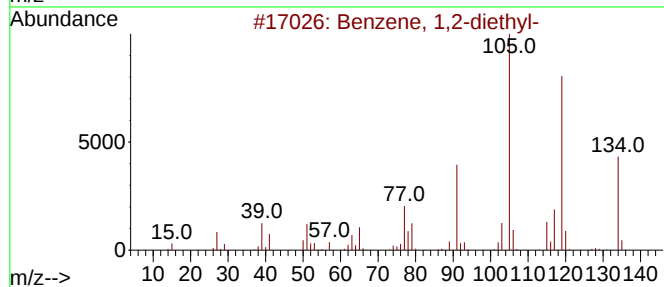
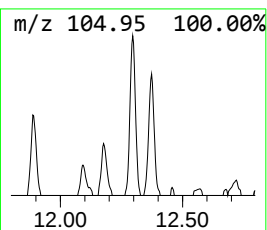
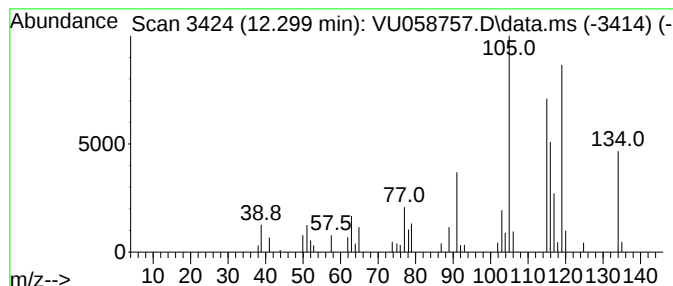
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.299	0.93 ug/L	43785	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	64
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	60
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	58
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	53
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	53



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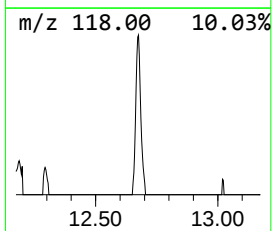
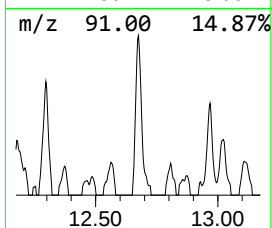
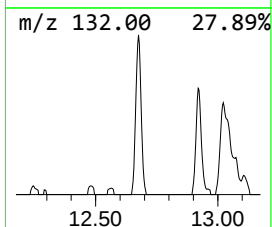
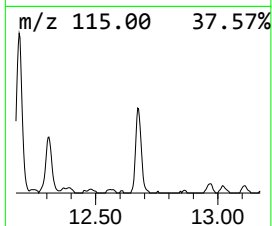
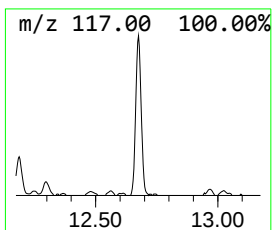
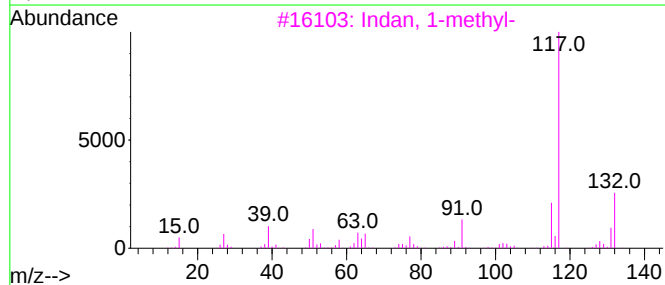
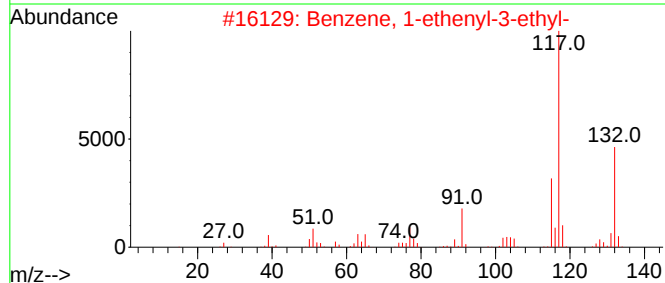
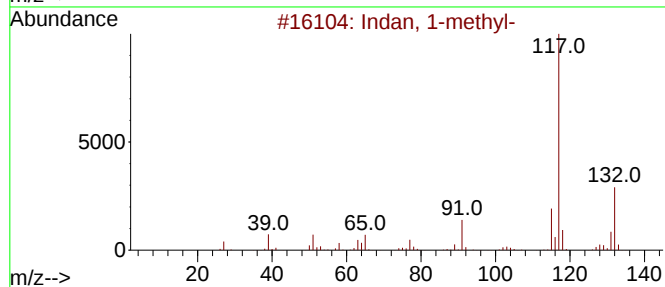
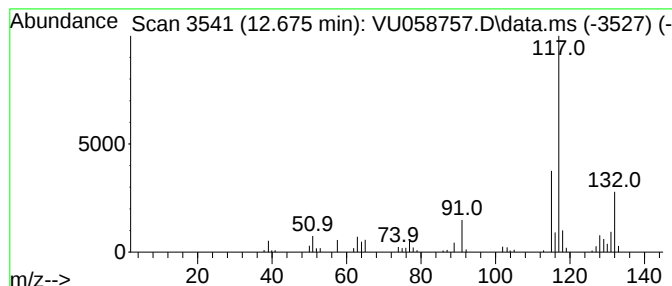
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR040424WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.675	1.47 ug/L	69384	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
5			Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

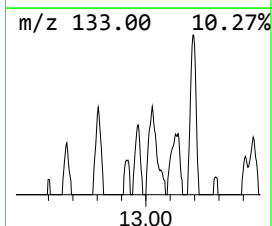
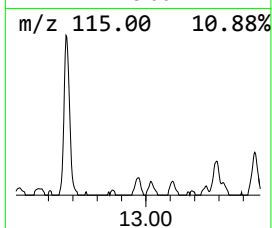
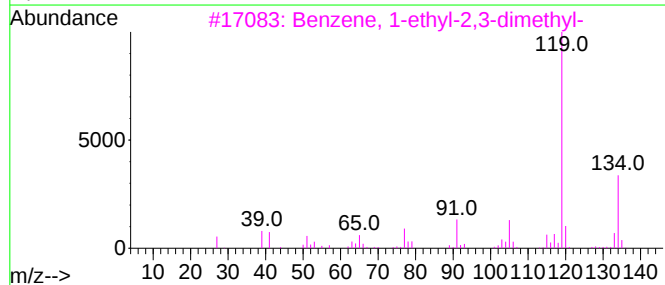
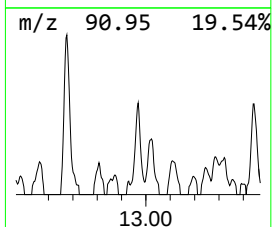
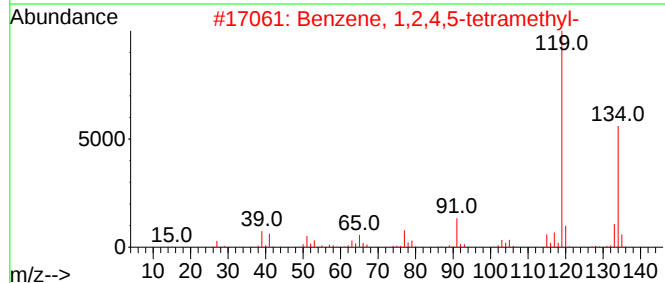
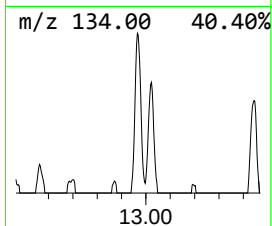
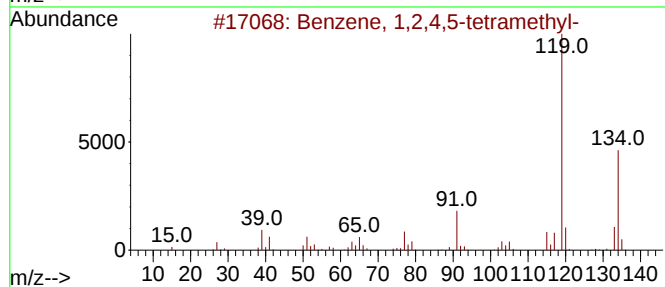
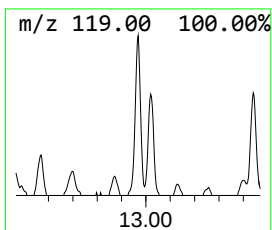
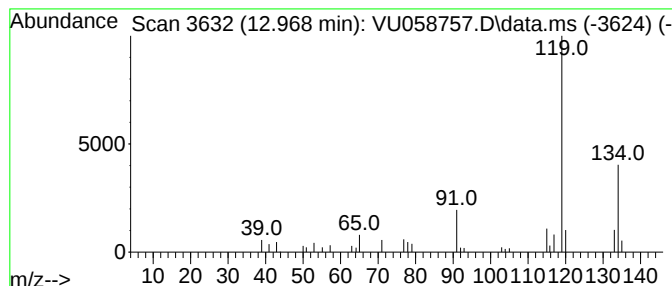
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.968	0.60 ug/L	28213	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

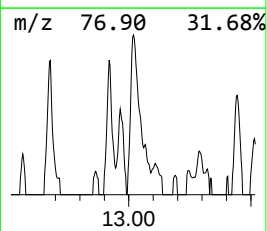
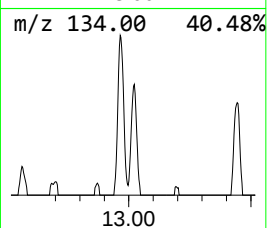
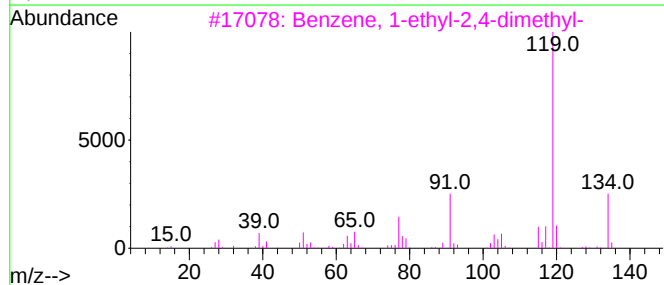
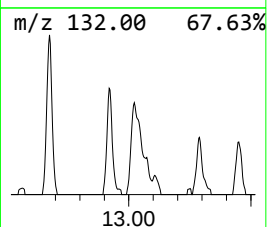
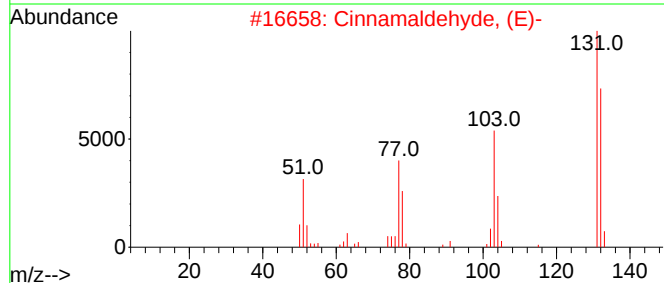
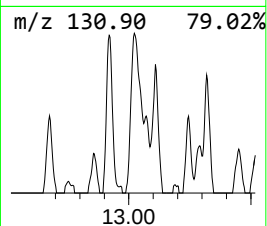
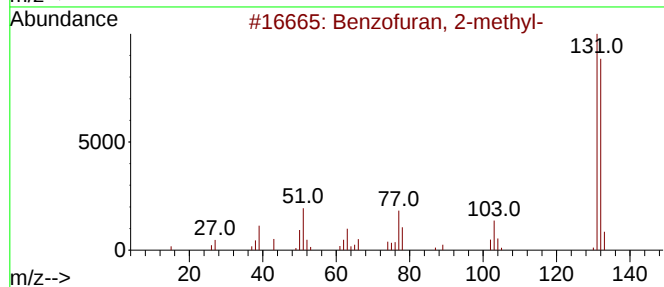
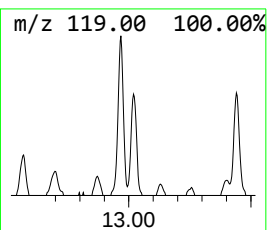
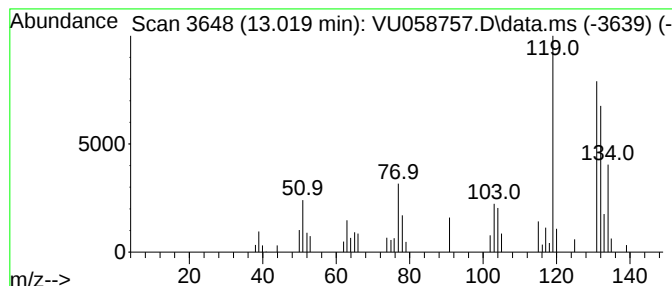
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzofuran, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.019	1.17 ug/L	55386	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	86
2			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	64
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	50
4			2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	50
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

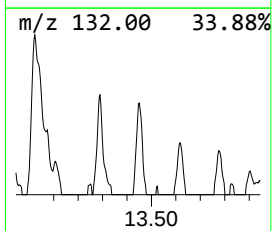
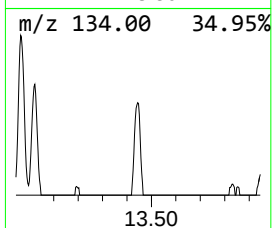
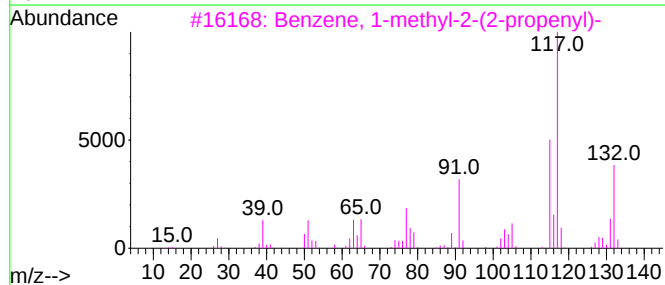
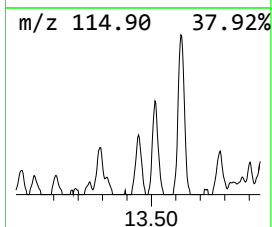
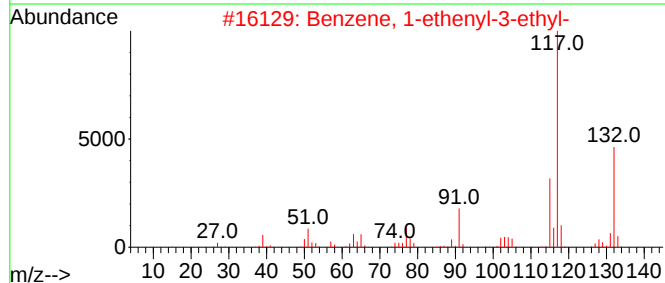
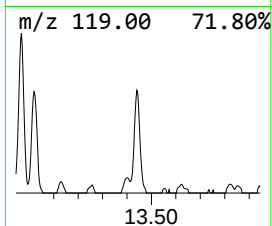
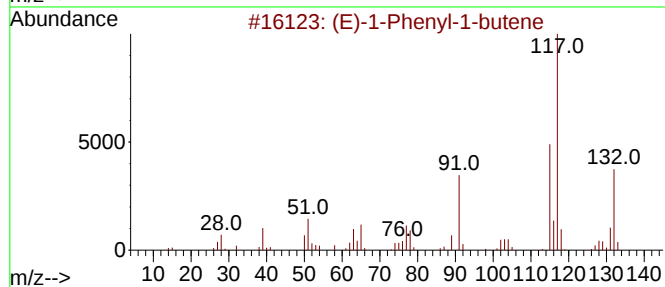
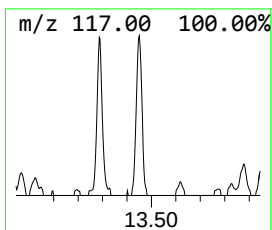
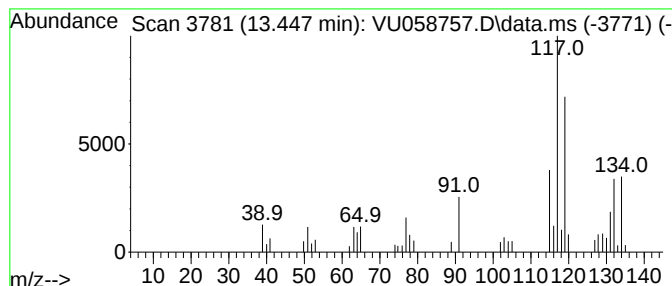
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 (E)-1-Phenyl-1-butene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.447	0.72 ug/L	33827	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	93		
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83		
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83		
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70		
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

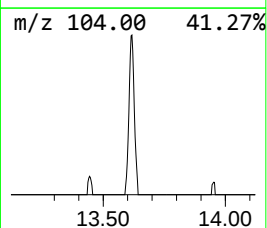
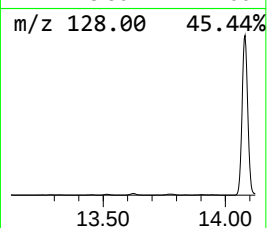
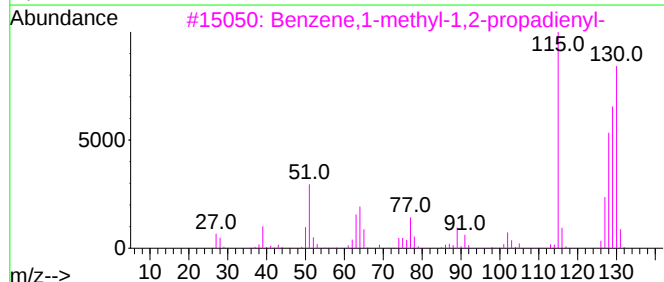
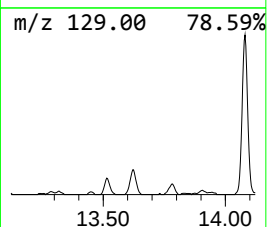
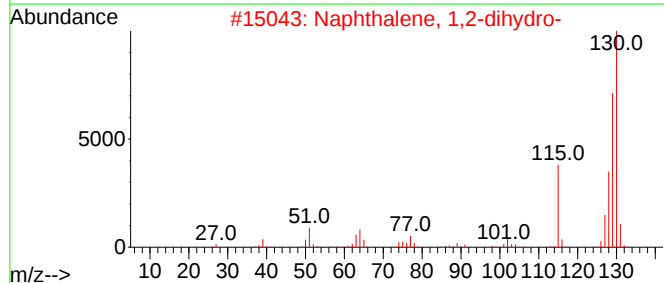
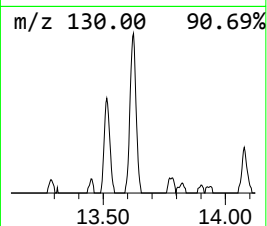
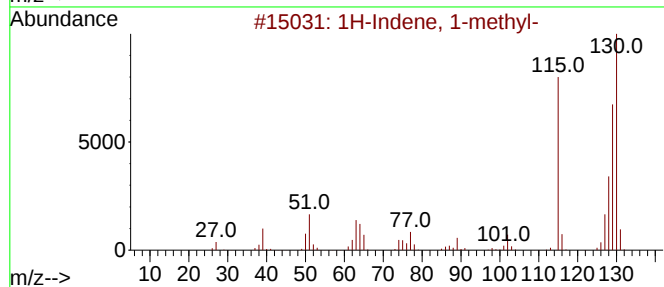
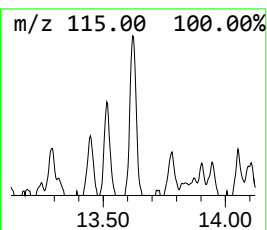
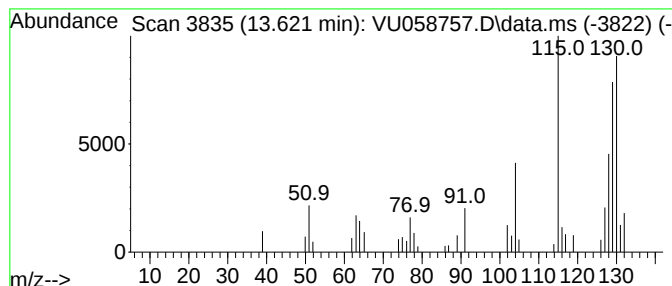
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 1H-Indene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.621	0.88 ug/L	41734	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2			Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	95
3			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
4			Benzene, 1-butyryl-	130	C10H10	000622-76-4	94
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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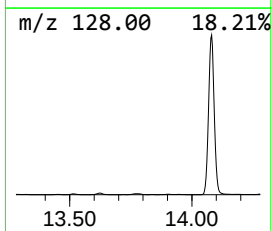
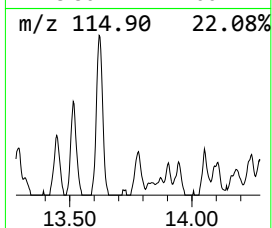
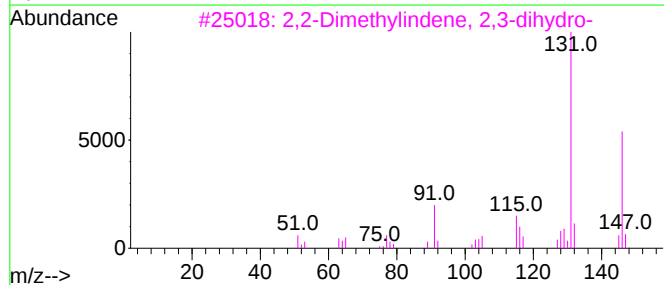
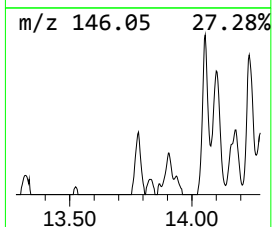
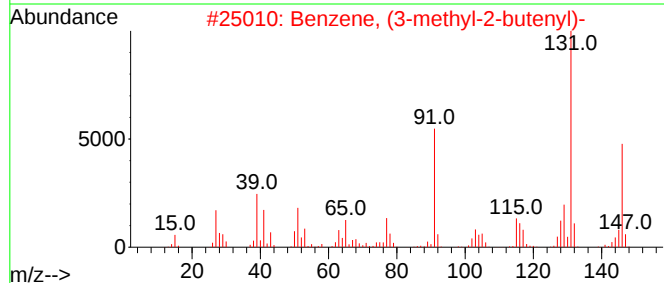
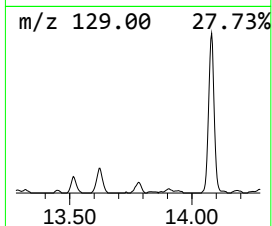
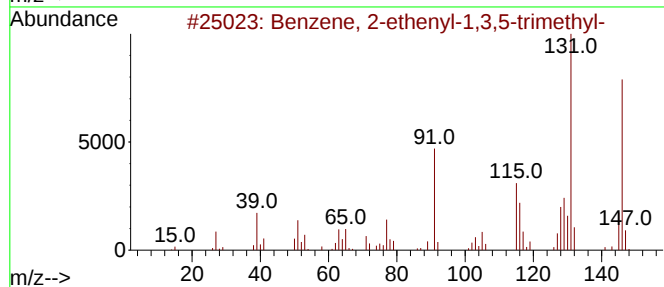
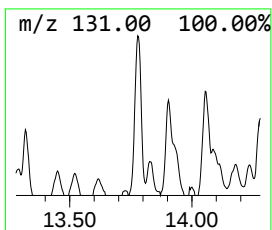
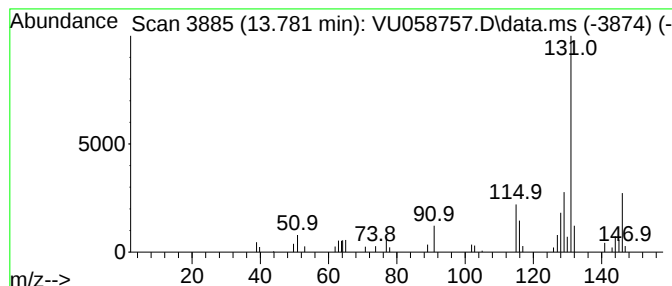
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.781	0.55 ug/L	25751	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	90
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	87
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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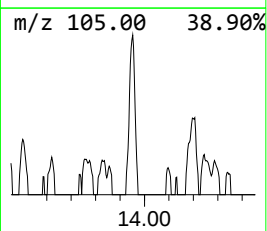
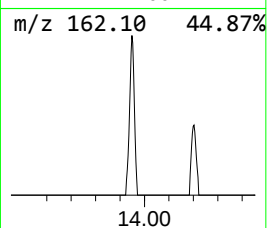
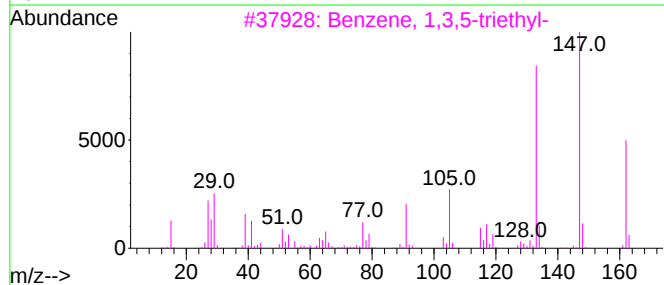
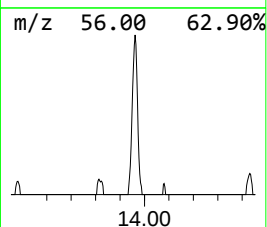
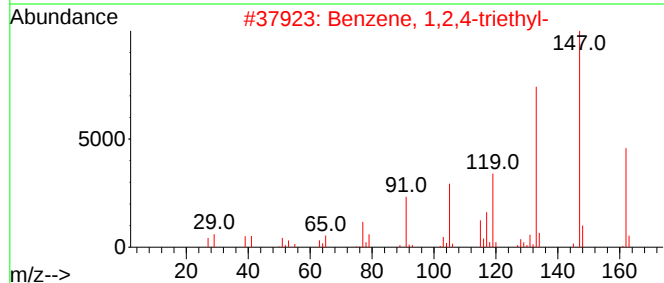
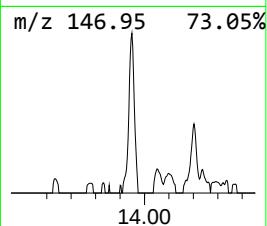
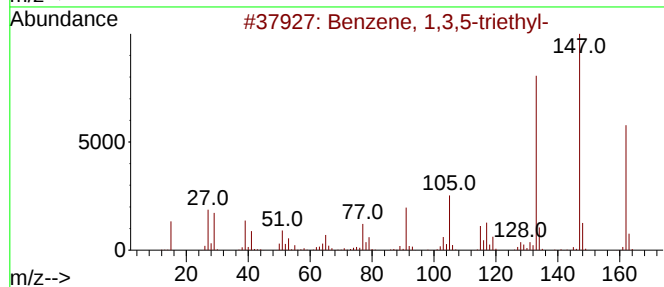
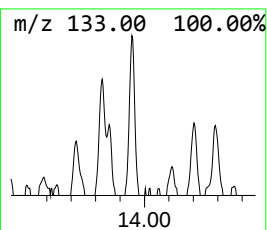
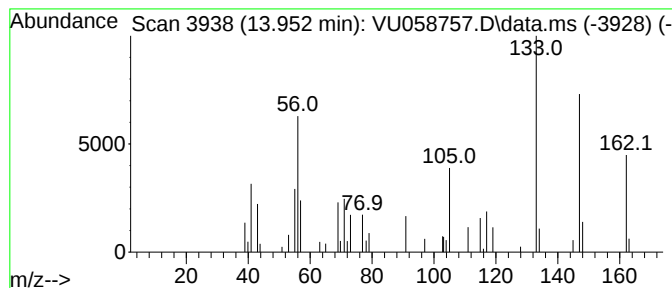
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.952	0.83 ug/L	39205	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-triethyl-	162	C12H18	000102-25-0	49
2			Benzene, 1,2,4-triethyl-	162	C12H18	000877-44-1	49
3			Benzene, 1,3,5-triethyl-	162	C12H18	000102-25-0	49
4			Benzene, 1,2,4-triethyl-	162	C12H18	000877-44-1	47
5			Benzene, 2,4-dimethyl-1-(1-methy...	162	C12H18	001483-60-9	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

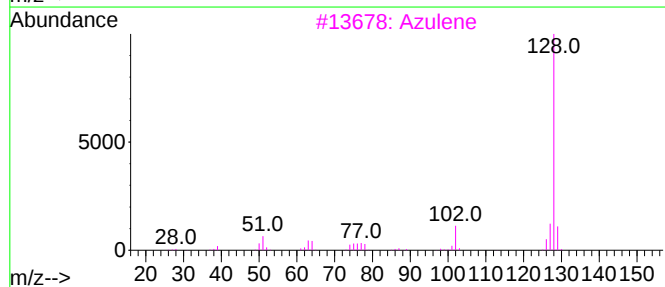
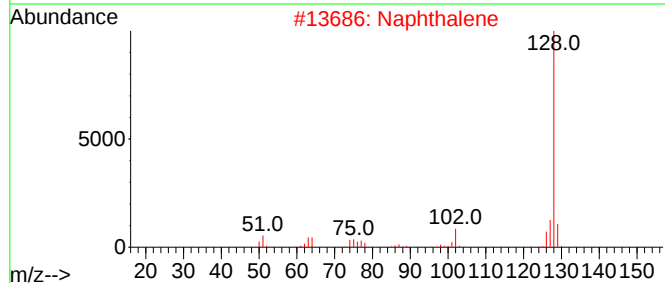
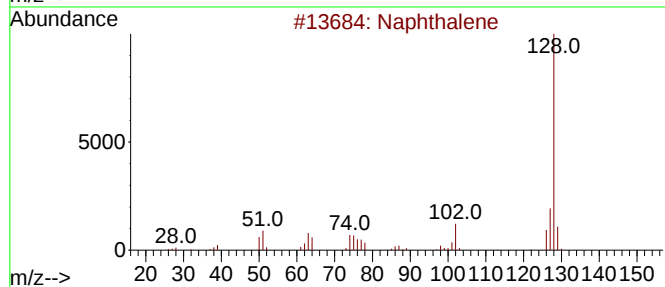
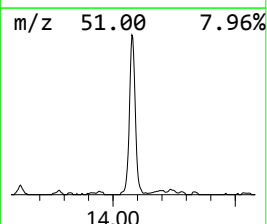
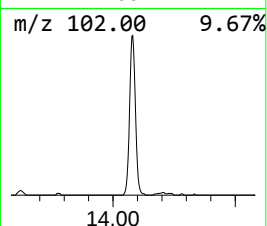
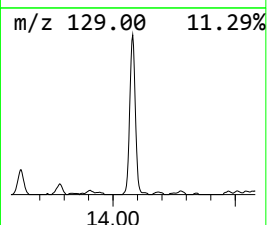
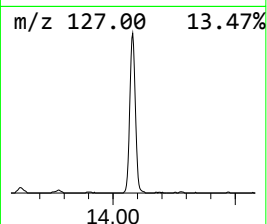
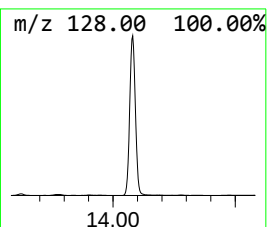
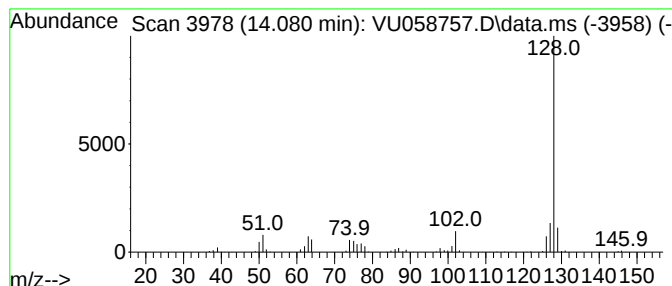
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	13.53 ug/L	639098	1,4-Dichlorobenzene-d4	11.807

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			Azulene	128	C10H8	000275-51-4	91
4			Naphthalene	128	C10H8	000091-20-3	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

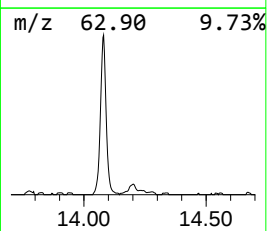
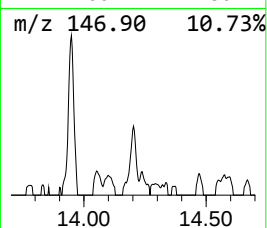
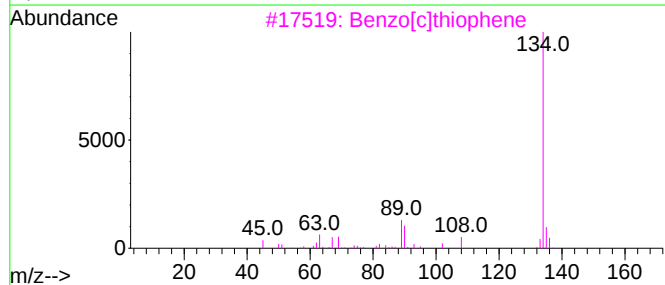
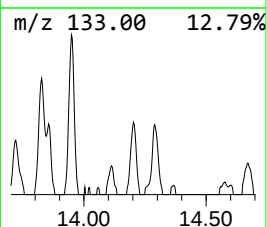
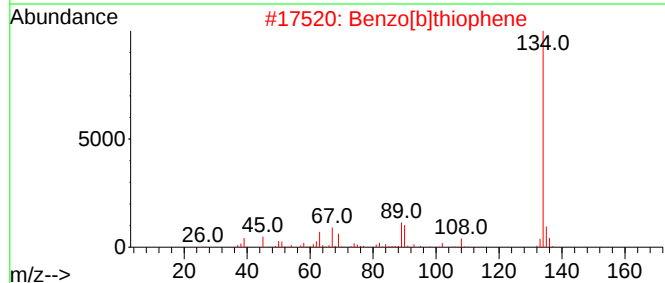
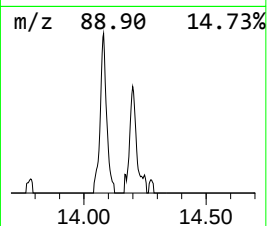
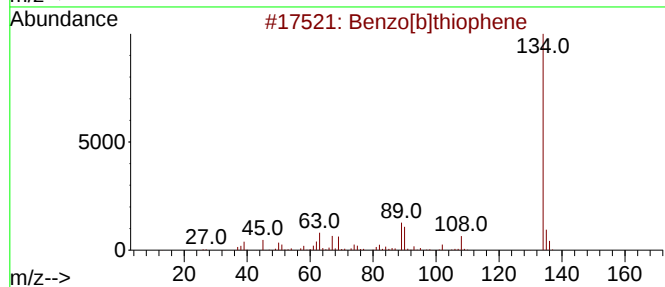
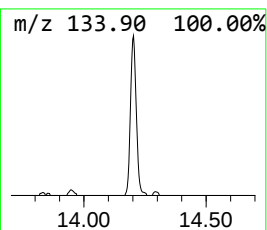
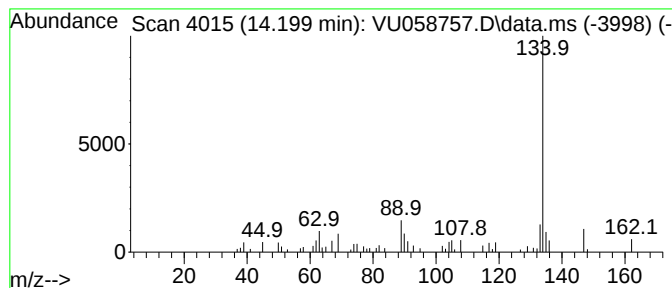
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Benzo[b]thiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.200	1.26 ug/L	59282	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
2			Benzo[b]thiophene	134	C8H6S	000095-15-8	81
3			Benzo[c]thiophene	134	C8H6S	000270-82-6	81
4			Benzo[b]thiophene	134	C8H6S	000095-15-8	74
5			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

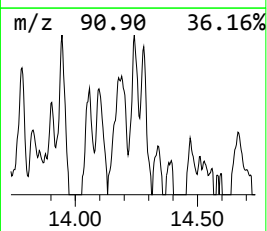
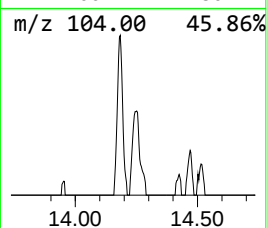
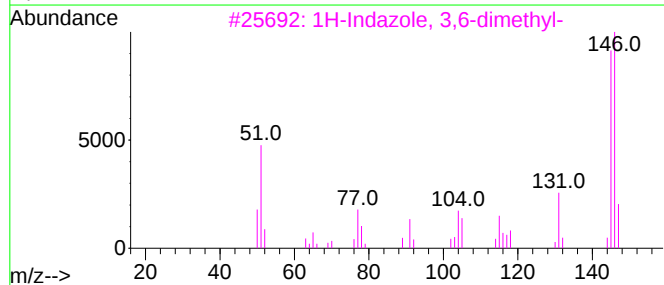
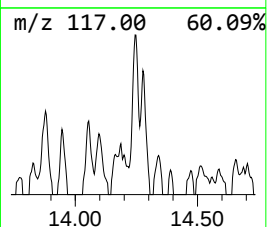
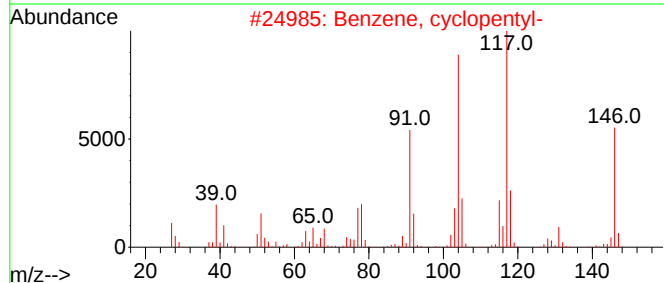
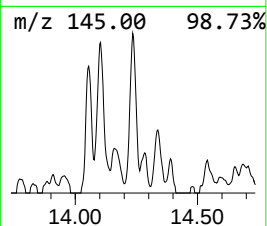
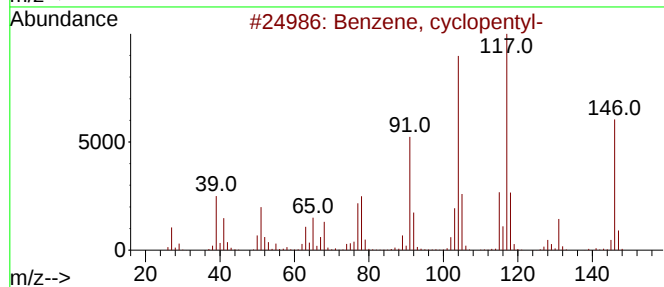
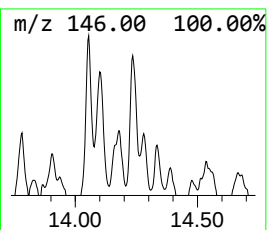
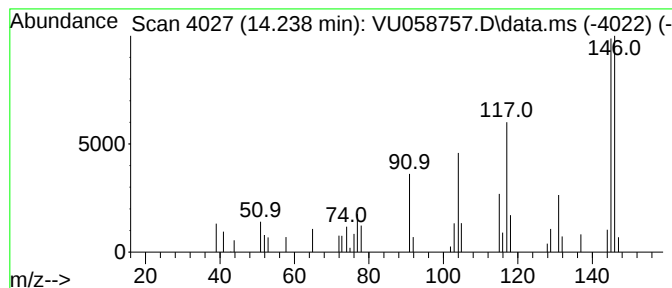
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 14 Benzene, cyclopentyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.238	0.58 ug/L	27555	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, cyclopentyl-	146	C11H14	000700-88-9	64
2			Benzene, cyclopentyl-	146	C11H14	000700-88-9	60
3			1H-Indazole, 3,6-dimethyl-	146	C9H10N2	007746-28-3	53
4			1H-1,3-Benzodiazole-5-carbaldehyde	146	C8H6N2O	058442-17-4	50
5			1H-Indole-5-methanamine	146	C9H10N2	1000362-58-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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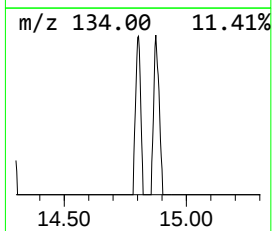
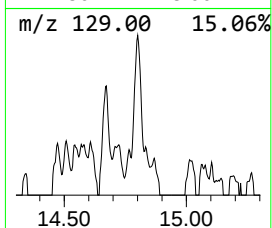
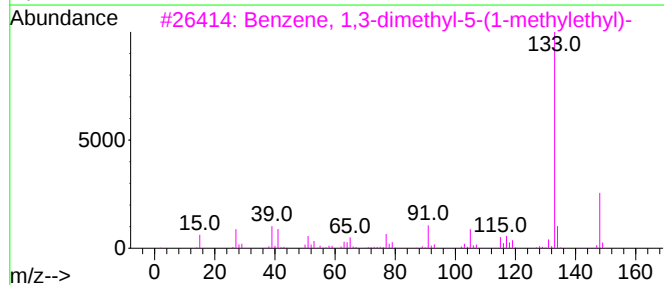
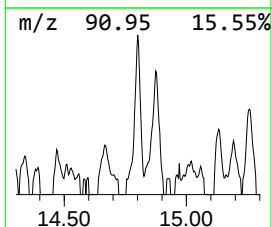
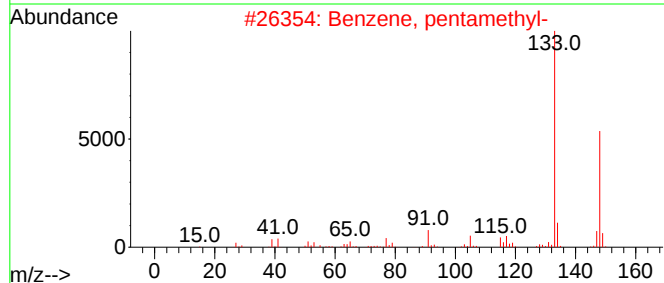
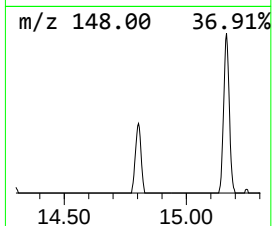
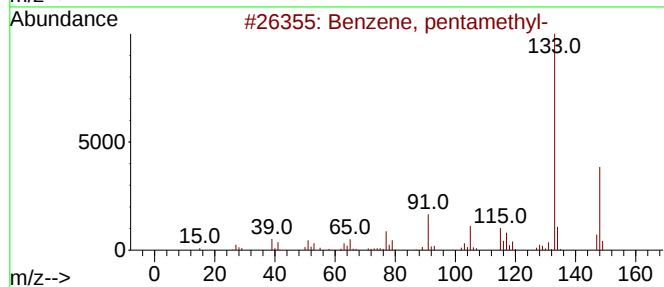
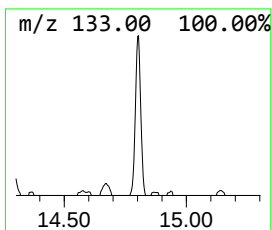
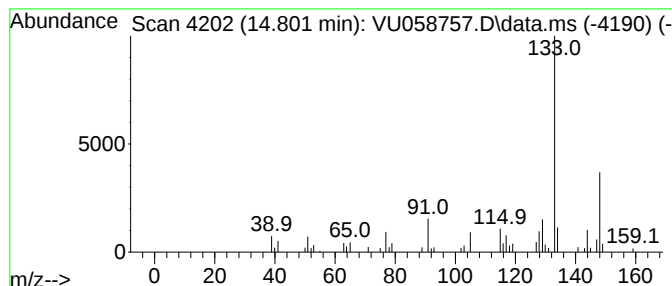
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Benzene, pentamethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.801	0.74 ug/L	34866	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	89
2			Benzene, pentamethyl-	148	C11H16	000700-12-9	76
3			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	70
4			3,4-Dimethylcumene	148	C11H16	1000370-34-1	70
5			2-tert-Butyltoluene	148	C11H16	001074-92-6	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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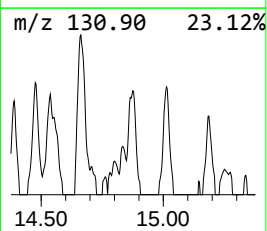
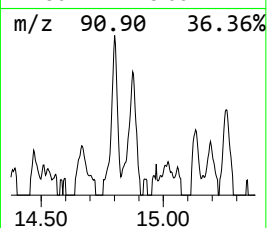
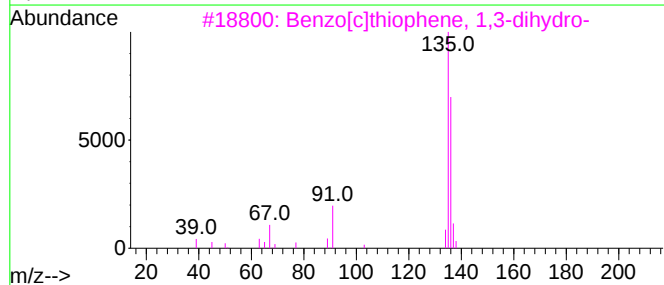
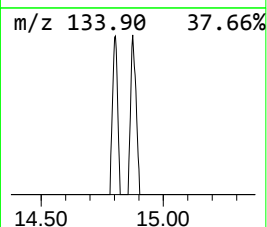
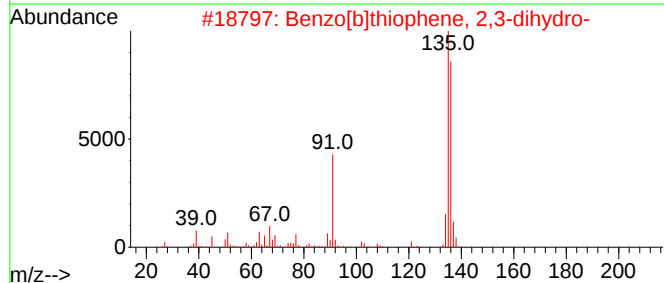
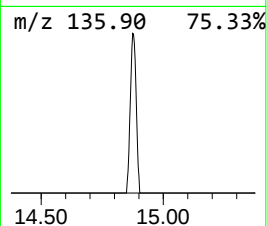
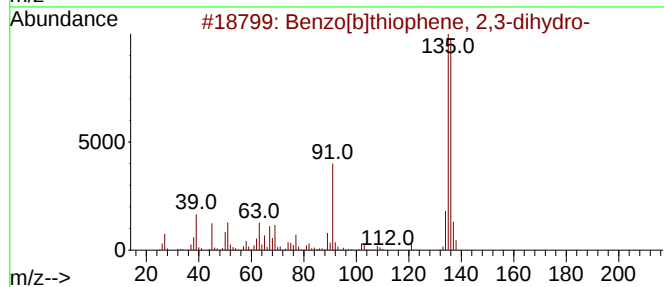
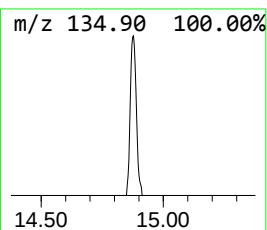
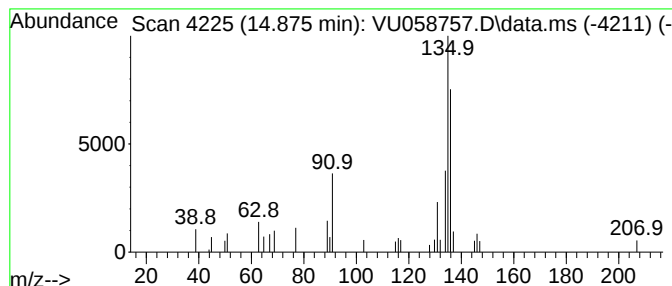
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 16 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.875	0.53 ug/L	24894	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	64
2			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	59
3			Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	59
4			4-Hydroxy-2-methylbenzaldehyde	136	C8H8O2	041438-18-0	58
5			N-Methyl-4-pyridinecarboxamide	136	C7H8N2O	006843-37-4	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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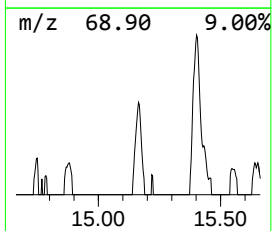
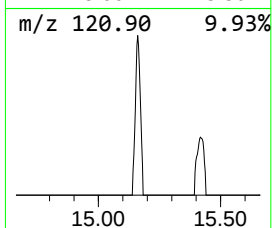
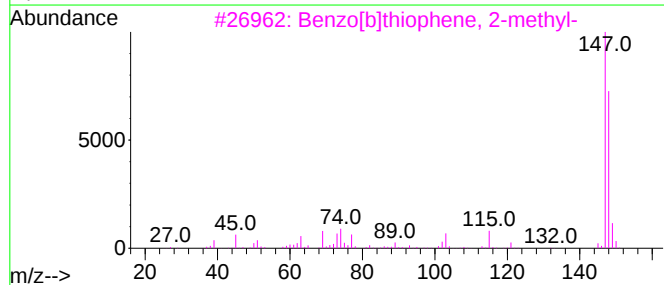
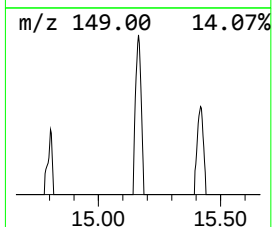
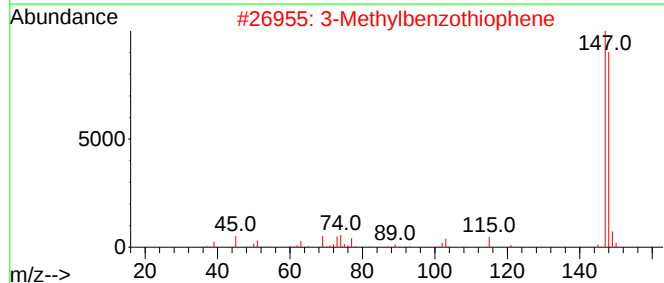
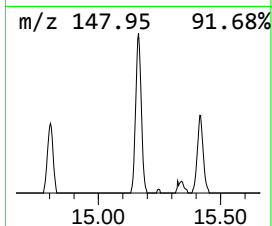
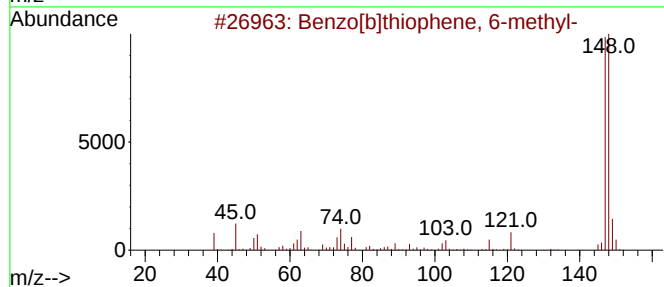
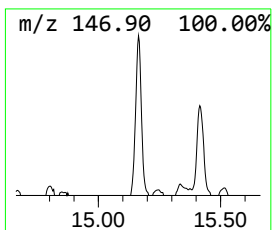
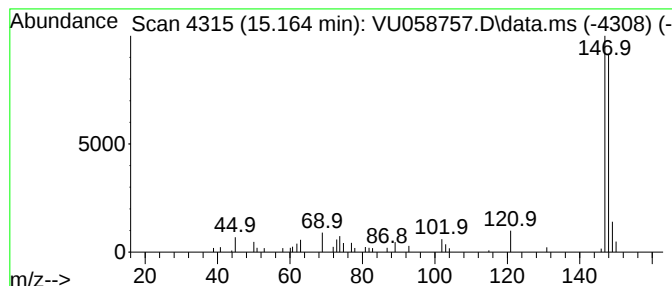
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 17 Benzo[b]thiophene, 6-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.164	0.61 ug/L	28630	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	91
2			3-Methylbenzothiophene	148	C9H8S	001455-18-1	87
3			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	80
4			5-Methylthiophene-2,3-dicarbonit...	148	C7H4N2S	1000305-40-8	80
5			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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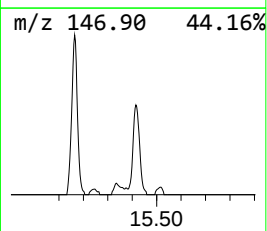
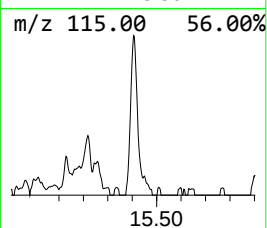
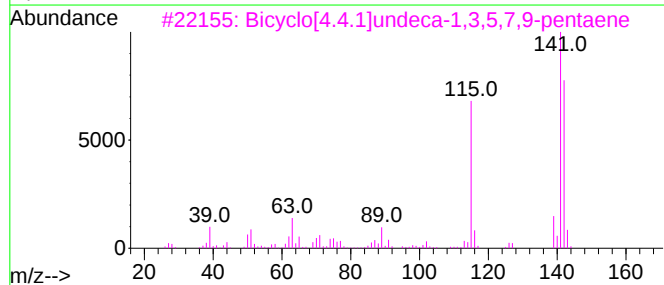
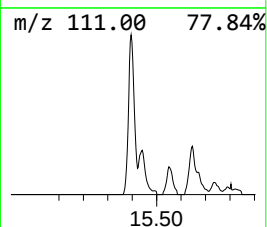
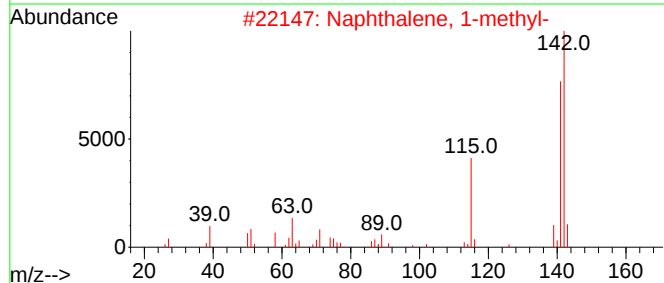
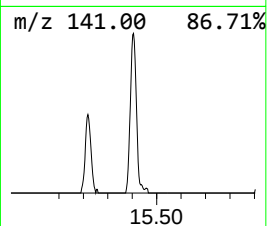
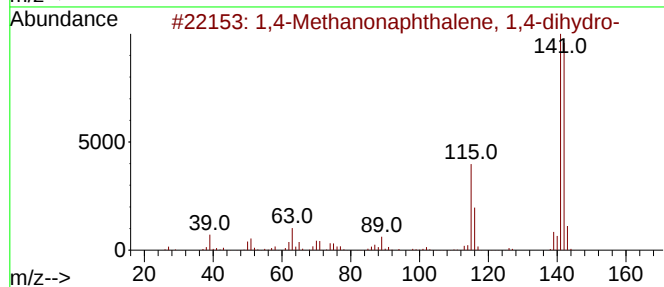
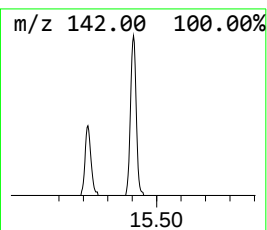
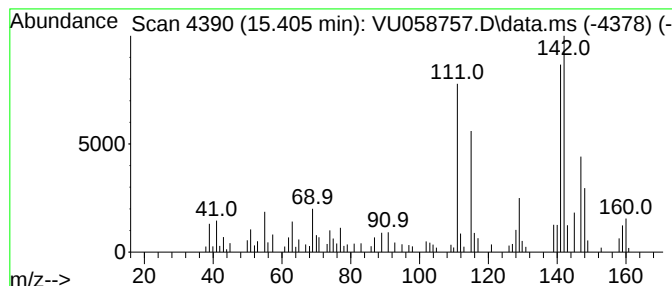
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 18 1,4-Methanonaphthalene, 1,4... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.405	2.22 ug/L	104833	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	50
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	50
3			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	50
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	50
5			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
Data File : VU058757.D
Acq On : 18 Apr 2024 13:04
Operator : MD/SY
Sample : P2181-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AK9

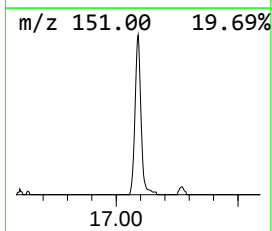
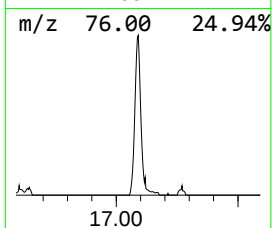
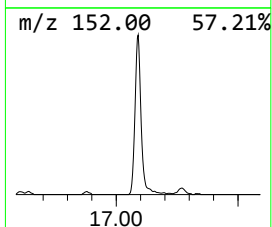
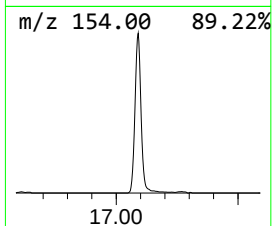
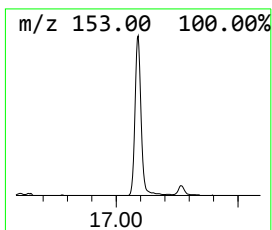
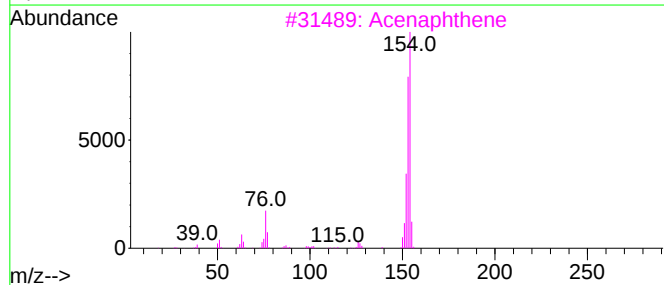
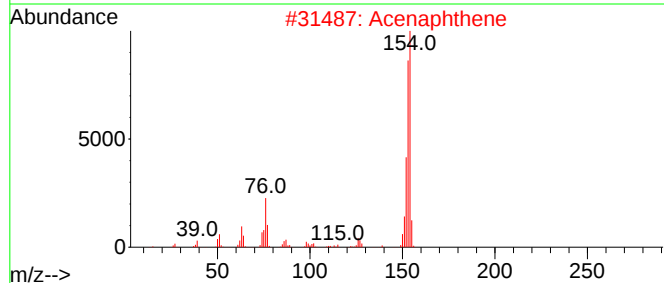
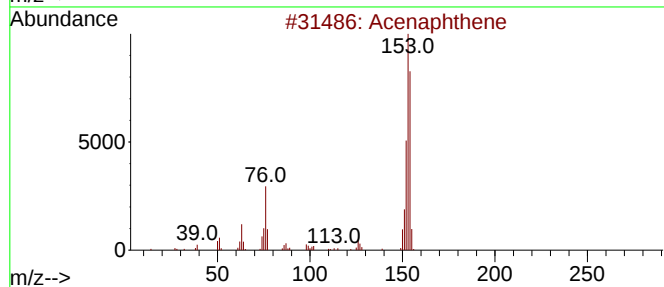
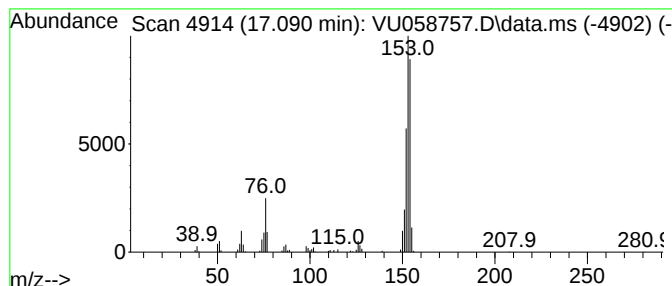
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 19 1,4-Ethenonaphthalene, 1,4-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.090	5.22 ug/L	246648	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	96
2			Acenaphthene	154	C12H10	000083-32-9	90
3			Acenaphthene	154	C12H10	000083-32-9	76
4			Acenaphthene	154	C12H10	000083-32-9	64
5			1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041824\
 Data File : VU058757.D
 Acq On : 18 Apr 2024 13:04
 Operator : MD/SY
 Sample : P2181-01
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AK9

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR040424WMA.M
 Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-chlo...	9.495	4.8	ug/L	222648	2	9.412	234574	5.0
Indane	12.087	3.2	ug/L	151758	3	11.807	236174	5.0
Benzene, 1,2-di...	12.299	0.9	ug/L	43785	3	11.807	236174	5.0
Indan, 1-methyl-	12.675	1.5	ug/L	69384	3	11.807	236174	5.0
Benzene, 1,2,4,...	12.968	0.6	ug/L	28213	3	11.807	236174	5.0
Benzofuran, 2-m...	13.019	1.2	ug/L	55386	3	11.807	236174	5.0
(E)-1-Phenyl-1-...	13.447	0.7	ug/L	33827	3	11.807	236174	5.0
1H-Indene, 1-me...	13.621	0.9	ug/L	41734	3	11.807	236174	5.0
Benzene, 2-ethe...	13.781	0.6	ug/L	25751	3	11.807	236174	5.0
unknown-01	13.952	0.8	ug/L	39205	3	11.807	236174	5.0
Azulene	14.081	13.5	ug/L	639098	3	11.807	236174	5.0
Benzo[b]thiophene	14.200	1.3	ug/L	59282	3	11.807	236174	5.0
Benzene, cyclop...	14.238	0.6	ug/L	27555	3	11.807	236174	5.0
Benzene, pentam...	14.801	0.7	ug/L	34866	3	11.807	236174	5.0
Benzo[b]thiophe...	14.875	0.5	ug/L	24894	3	11.807	236174	5.0
Benzo[b]thiophe...	15.164	0.6	ug/L	28630	3	11.807	236174	5.0
1,4-Methanonaph...	15.405	2.2	ug/L	104833	3	11.807	236174	5.0
1,4-Ethenonaph...	17.090	5.2	ug/L	246648	3	11.807	236174	5.0