

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW112522\
 Data File : VW029352.D
 Acq On : 25 Nov 2022 13:47
 Operator : SY/MD
 Sample : N5725-14DL 4X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BHB67DL

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_V\Method\SFAMVTR112322WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VW029352.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.105	14	20	38	rVB	3303071	3095298	100.00%	17.622%
2	1.214	49	54	72	rVB2	168335	178897	5.78%	1.018%
3	1.307	76	83	100	rVB2	257618	288423	9.32%	1.642%
4	1.565	156	163	180	rVB	105670	119661	3.87%	0.681%
5	2.105	320	331	346	rBV2	210452	309593	10.00%	1.763%
6	2.503	445	455	472	rVB2	22747	38918	1.26%	0.222%
7	3.182	651	666	691	rBV	65384	142626	4.61%	0.812%
8	3.902	868	890	942	rBV2	245295	790902	25.55%	4.503%
9	4.339	1008	1026	1047	rBV	141957	354736	11.46%	2.020%
10	4.597	1090	1106	1122	rBV3	13611	33573	1.08%	0.191%
11	5.040	1224	1244	1277	rBV2	347254	923108	29.82%	5.255%
12	5.609	1408	1421	1457	rBV	330681	689607	22.28%	3.926%
13	5.908	1501	1514	1529	rBV5	21231	46705	1.51%	0.266%
14	6.063	1543	1562	1594	rBV	195840	414814	13.40%	2.362%
15	6.979	1834	1847	1871	rBV	100213	193925	6.27%	1.104%
16	7.307	1935	1949	1966	rBV	443083	793965	25.65%	4.520%
17	7.616	2033	2045	2068	rBV	57677	107472	3.47%	0.612%
18	8.085	2175	2191	2228	rBV2	349388	742437	23.99%	4.227%
19	8.844	2415	2427	2453	rBV	466678	783232	25.30%	4.459%
20	9.005	2466	2477	2489	rBV	20727	34937	1.13%	0.199%
21	9.133	2506	2517	2533	rBV	70484	114513	3.70%	0.652%
22	9.535	2631	2642	2660	rVB	168291	263786	8.52%	1.502%
23	9.921	2751	2762	2776	rBV	57364	88117	2.85%	0.502%
24	10.207	2839	2851	2865	rBV	135747	215687	6.97%	1.228%
25	10.728	2996	3013	3030	rVB	57332	88907	2.87%	0.506%
26	11.239	3157	3172	3191	rBV	533201	828086	26.75%	4.714%
27	11.516	3243	3258	3276	rBV	661459	1029782	33.27%	5.863%
28	11.612	3277	3288	3309	rVB	458567	725027	23.42%	4.128%
29	11.735	3314	3326	3348	rBV	1216182	1900892	61.41%	10.822%
30	12.005	3400	3410	3429	rBV2	51833	93577	3.02%	0.533%
31	12.104	3429	3441	3457	rVB	247306	385285	12.45%	2.193%
32	12.349	3507	3517	3527	rBV	167119	264193	8.54%	1.504%
33	12.455	3539	3550	3560	rBV	365788	648561	20.95%	3.692%
34	12.532	3570	3574	3586	rVB	38310	54409	1.76%	0.310%
35	12.712	3620	3630	3647	rVB	90910	139609	4.51%	0.795%
36	12.870	3660	3679	3690	rBV	90029	155439	5.02%	0.885%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Title : TRACE VOA SFAM1.0

37	12.937	3691	3700	3712	rVB3	23735	39827	1.29%	0.227%
38	13.034	3719	3730	3745	rBV3	94464	157526	5.09%	0.897%
39	13.480	3859	3869	3879	rBV3	31416	65957	2.13%	0.375%
40	13.612	3895	3910	3919	rBV	95448	162828	5.26%	0.927%
41	15.760	4568	4578	4590	rBV2	37910	60455	1.95%	0.344%

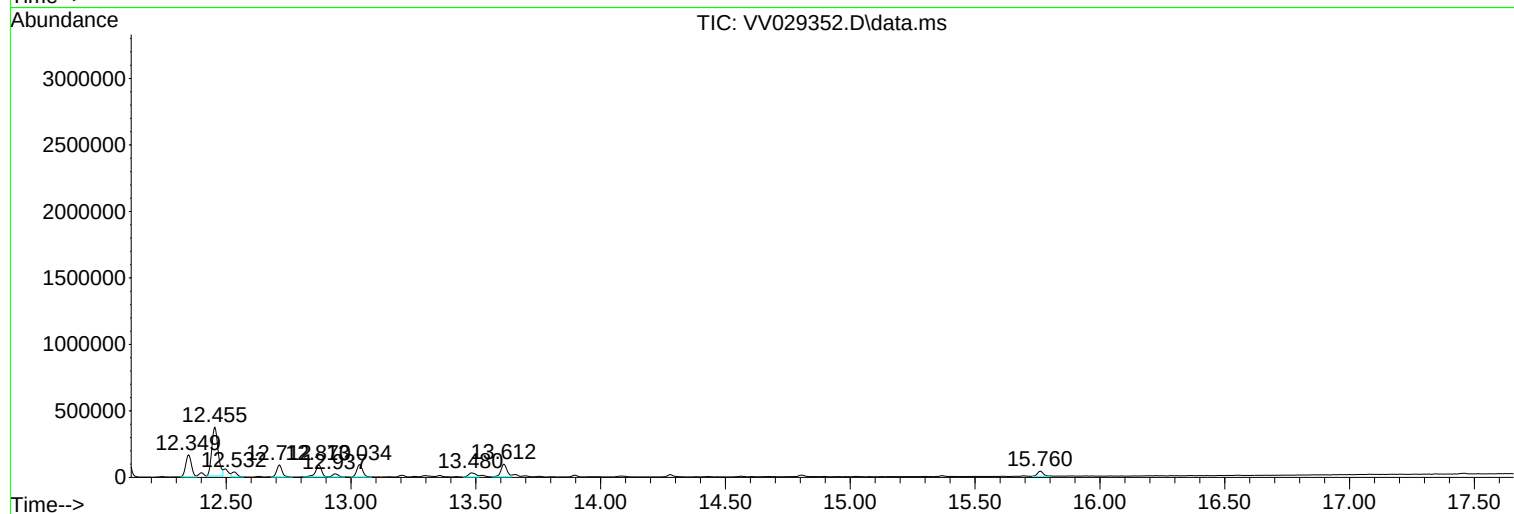
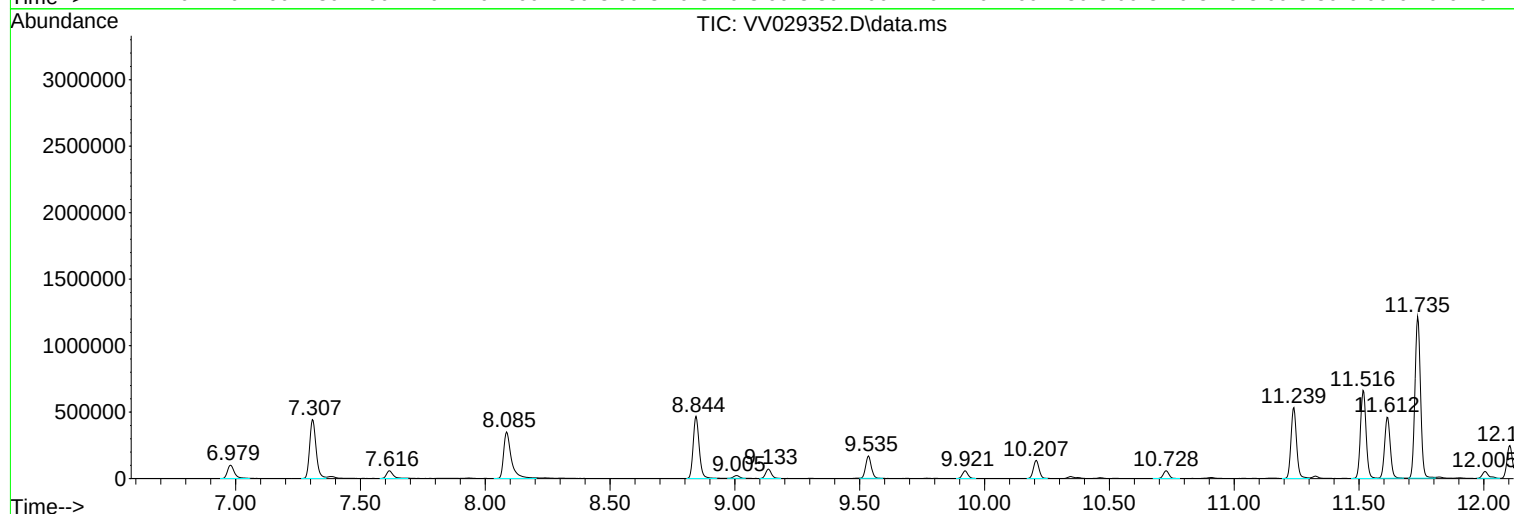
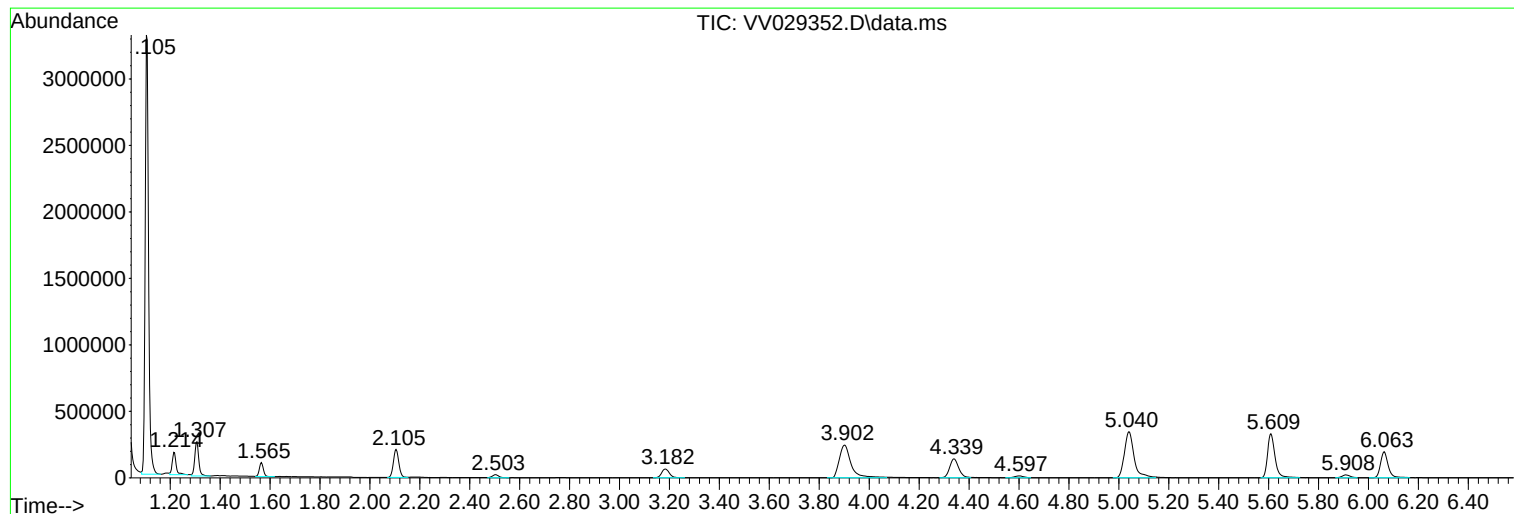
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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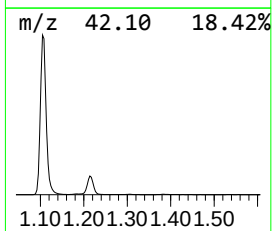
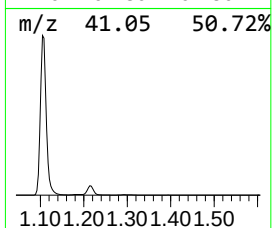
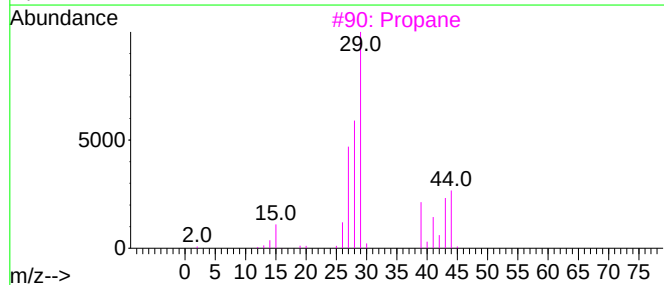
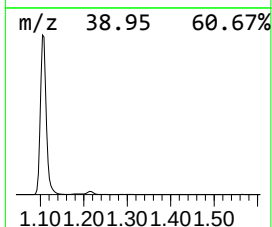
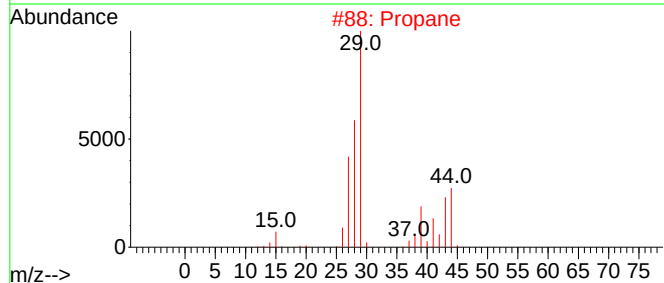
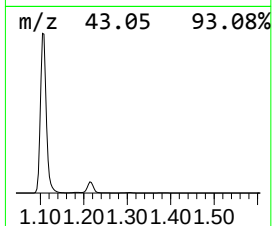
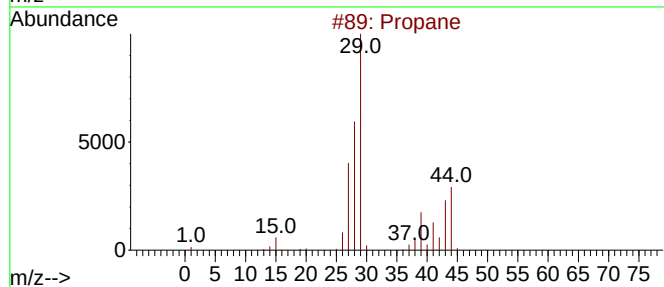
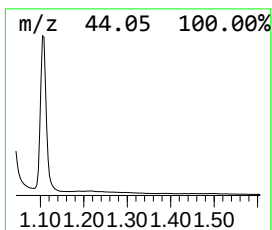
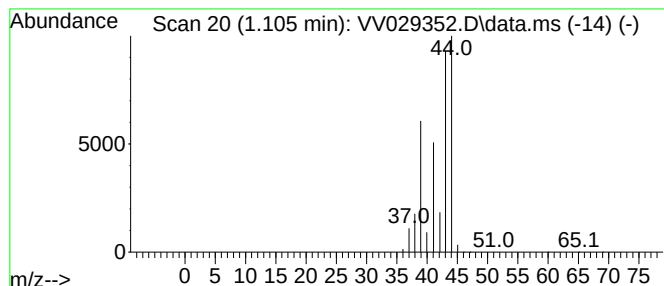
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.105	22.44 ug/L	3095300	1,4-Difluorobenzene	5.609

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane	44	C3H8	000074-98-6	83
2			Propane	44	C3H8	000074-98-6	59
3			Propane	44	C3H8	000074-98-6	9
4			Propene	42	C3H6	000115-07-1	9
5			Ethylene oxide	44	C2H4O	000075-21-8	5



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Instrument :
MSVOA_V
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BHB67DL

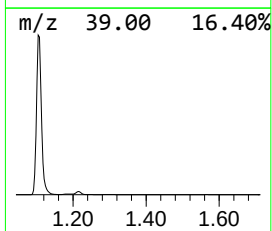
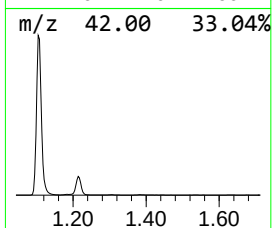
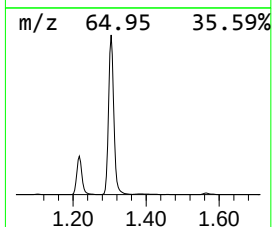
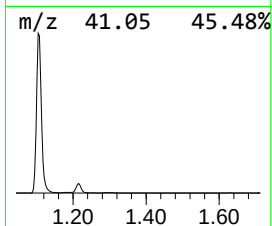
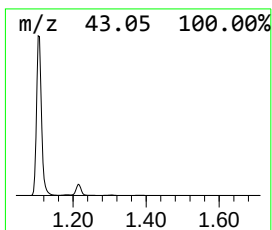
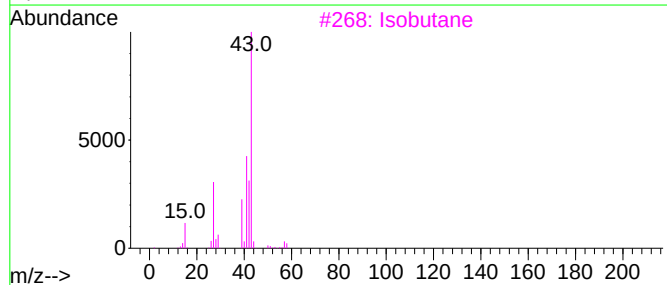
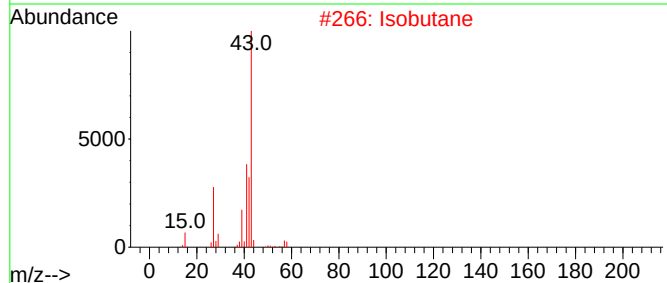
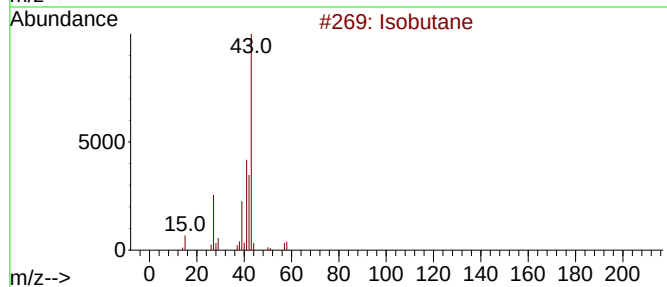
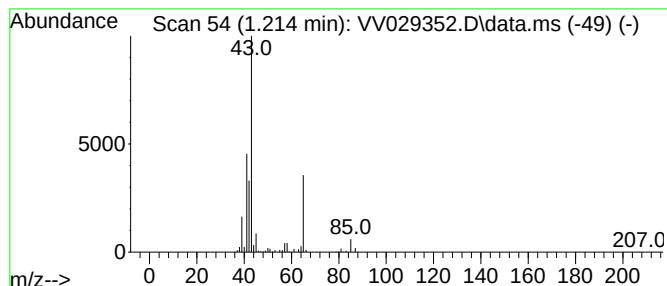
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.214	1.30 ug/L	178897	1,4-Difluorobenzene	5.609

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutane	58	C4H10	000075-28-5	46
2			Isobutane	58	C4H10	000075-28-5	46
3			Isobutane	58	C4H10	000075-28-5	43
4			Isobutane	58	C4H10	000075-28-5	43
5			Isobutane	58	C4H10	000075-28-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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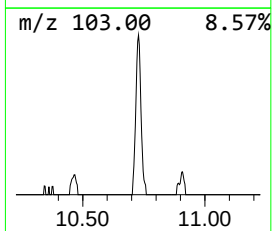
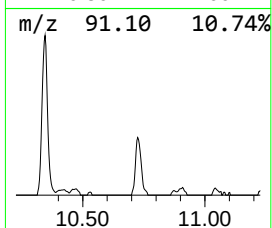
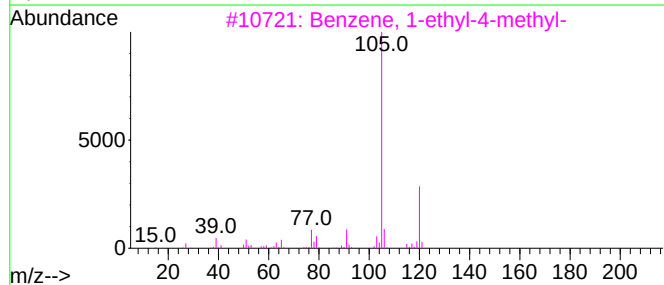
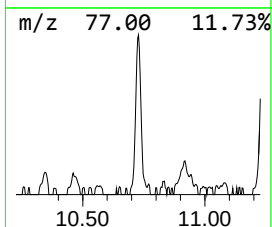
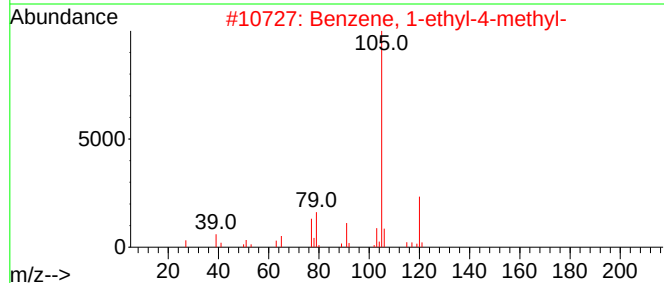
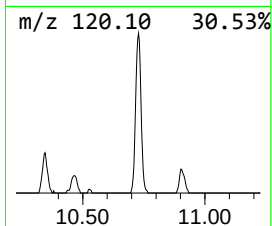
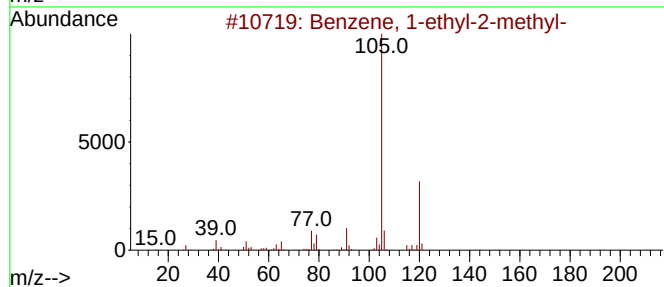
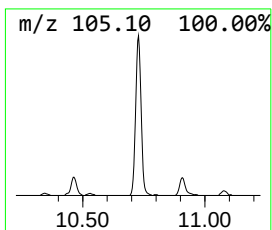
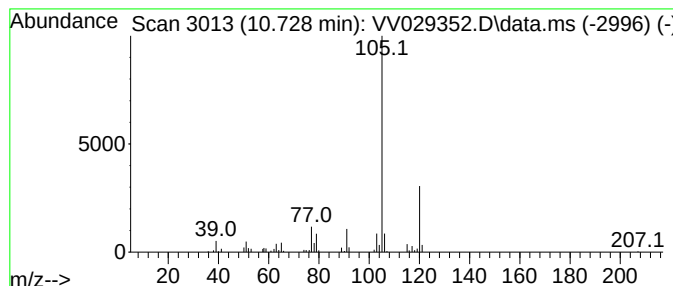
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.728	0.54 ug/L	88907	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB67DL

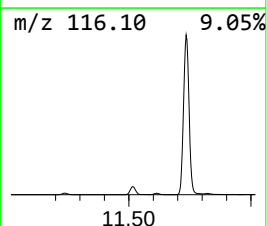
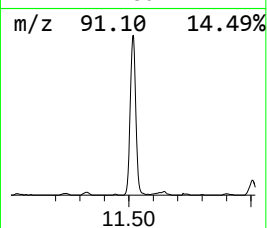
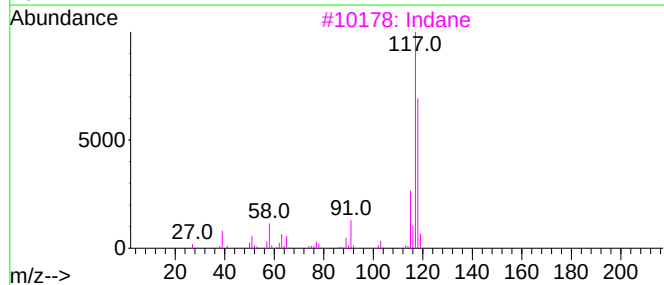
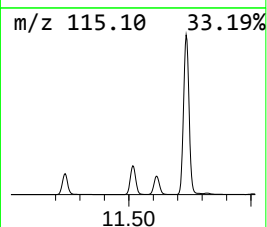
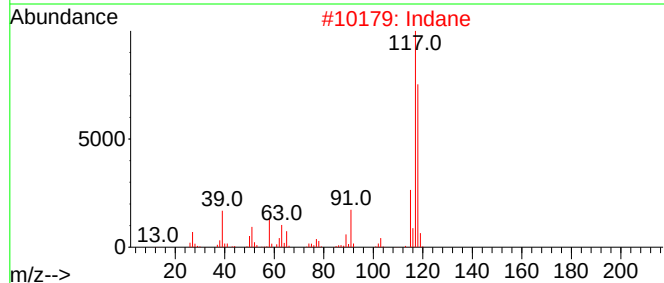
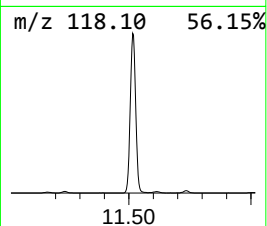
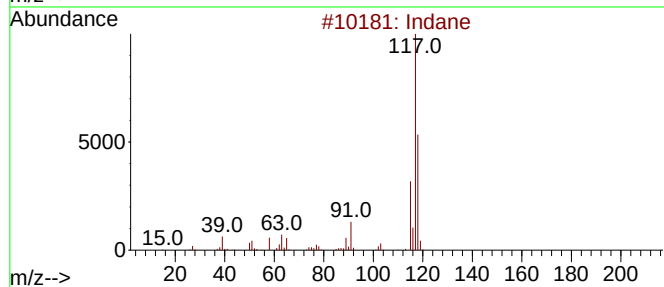
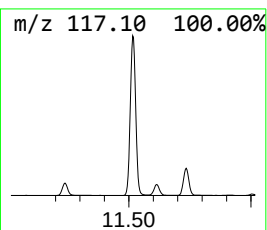
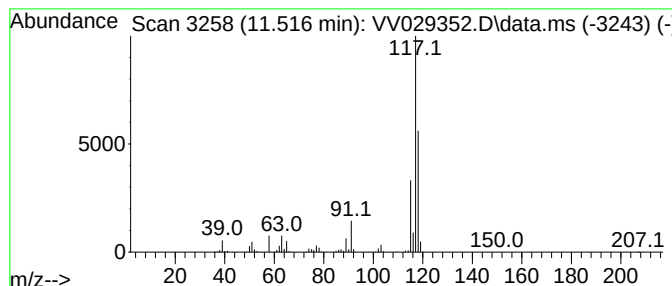
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.516	6.22 ug/L	1029780	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	95
2			Indane	118	C9H10	000496-11-7	91
3			Indane	118	C9H10	000496-11-7	87
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
5			Deltacyclene	118	C9H10	007785-10-6	80



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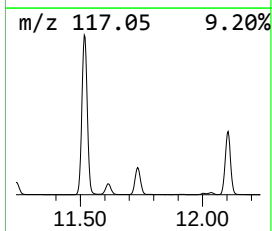
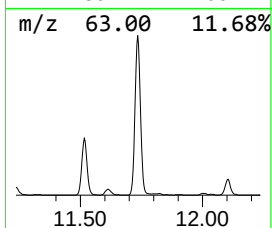
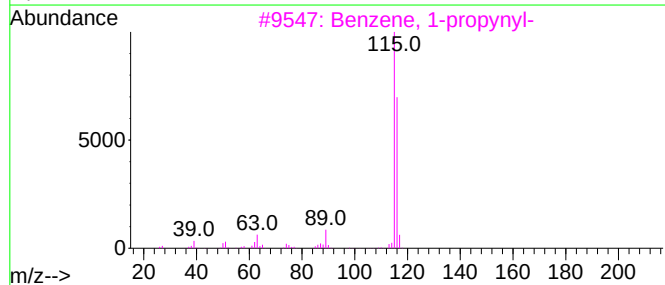
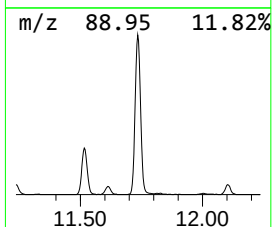
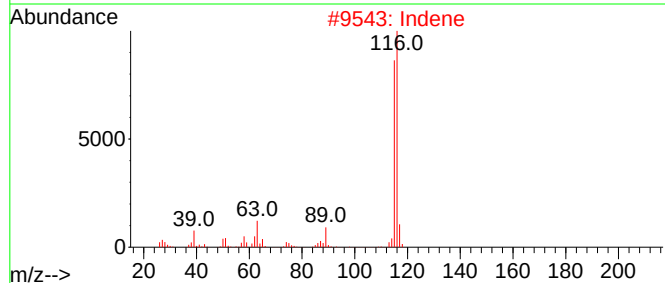
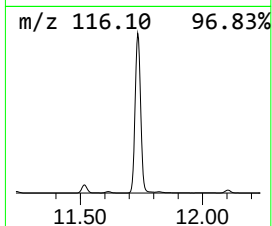
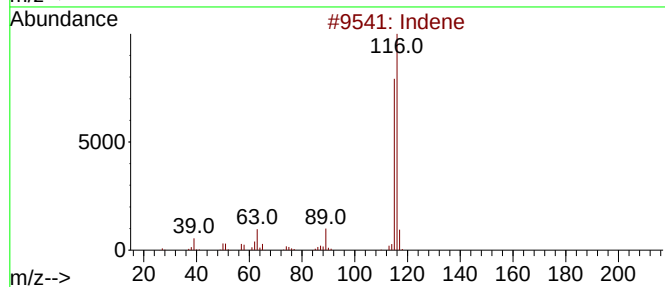
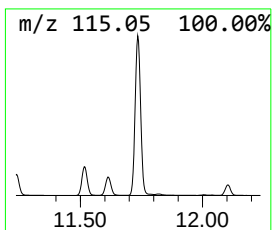
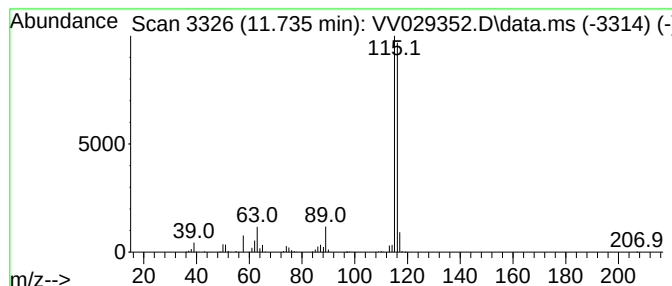
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.735	11.48 ug/L	1900890	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Indene	116	C9H8	000095-13-6	95
3			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
4			Indene	116	C9H8	000095-13-6	94
5			3-Methylphenylacetylene	116	C9H8	000766-82-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB67DL

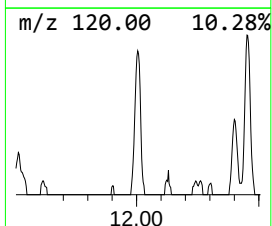
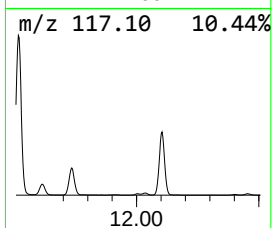
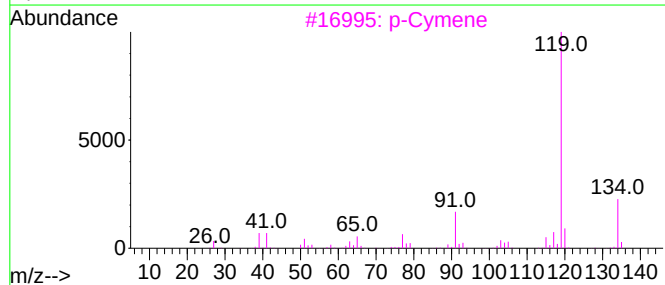
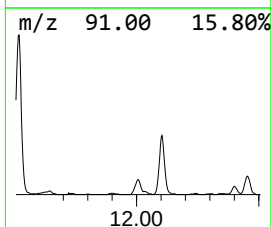
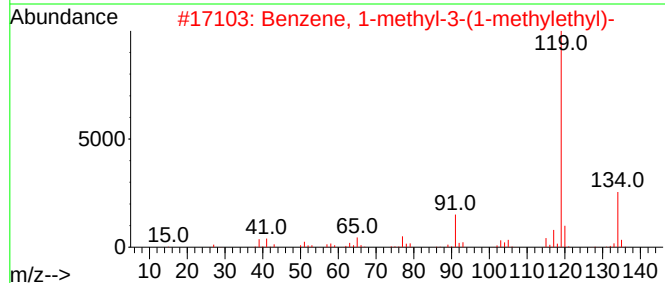
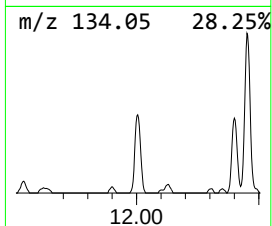
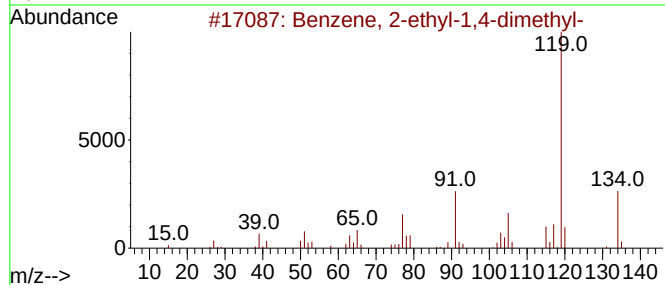
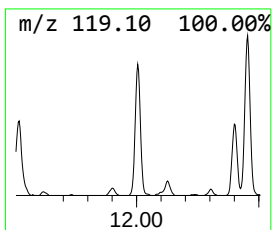
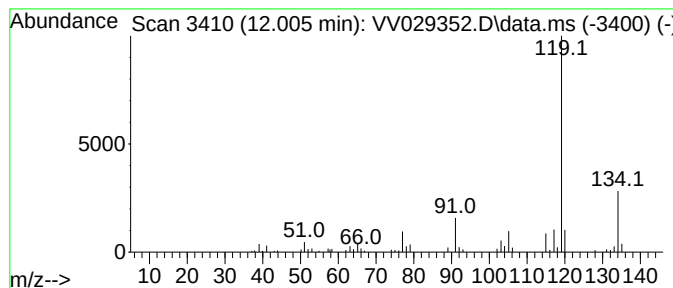
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.005	0.57 ug/L	93577	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			p-Cymene	134	C10H14	000099-87-6	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB67DL

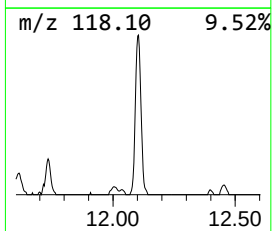
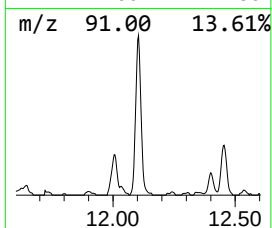
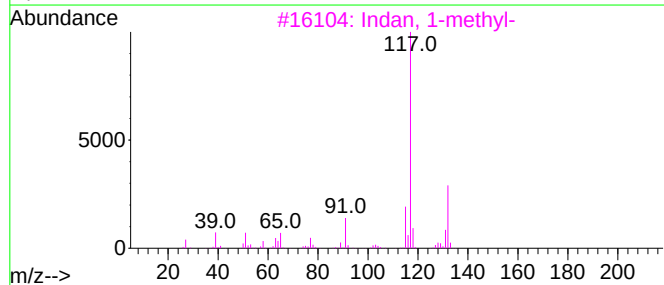
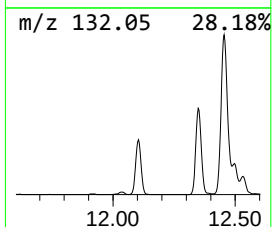
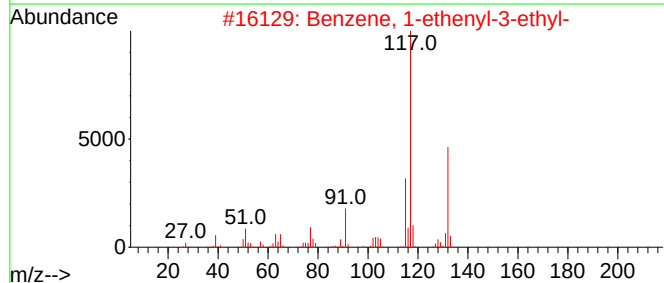
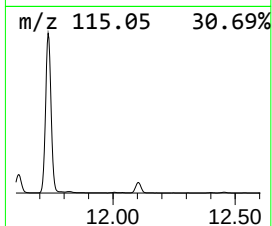
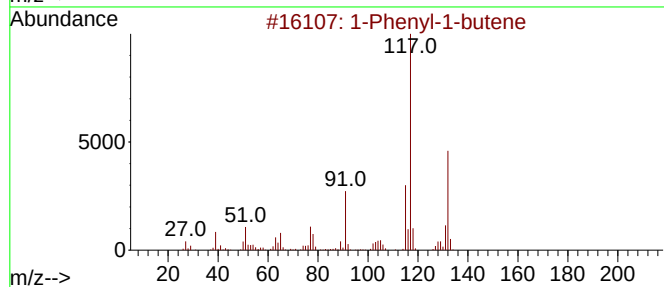
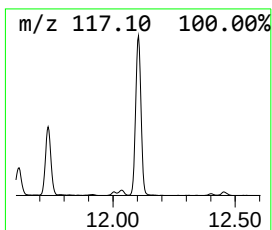
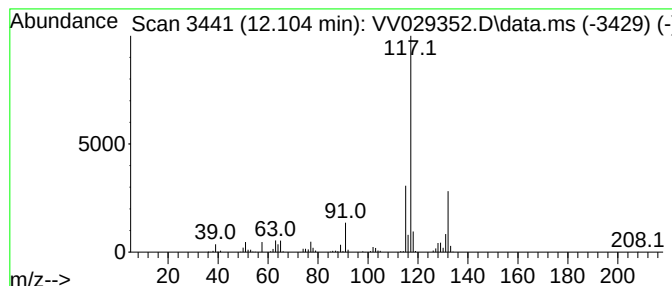
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	2.33 ug/L	385285	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	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, 2-butenyl-	132	C10H12	001560-06-1	86
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB67DL

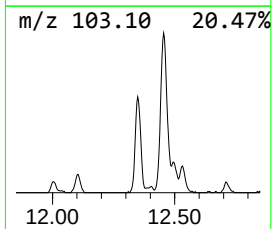
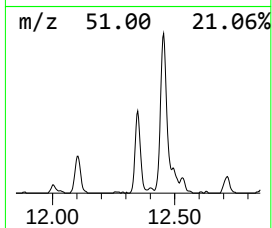
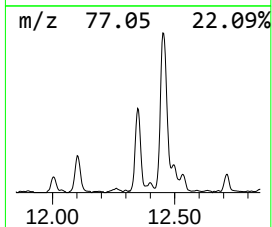
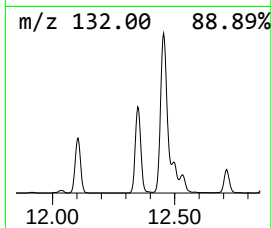
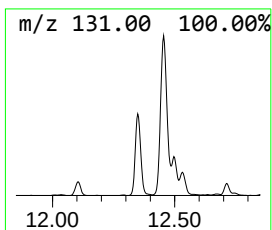
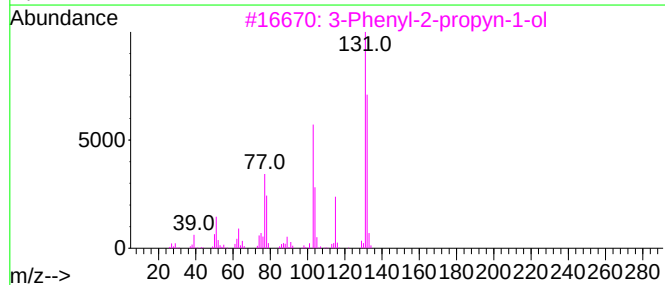
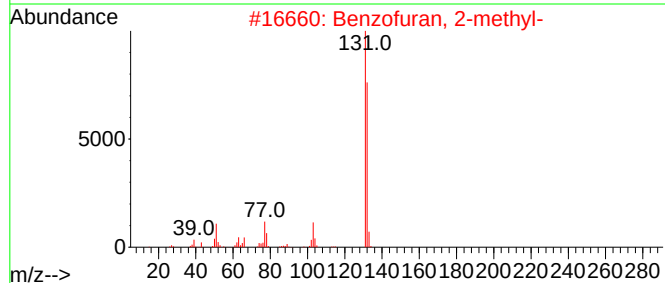
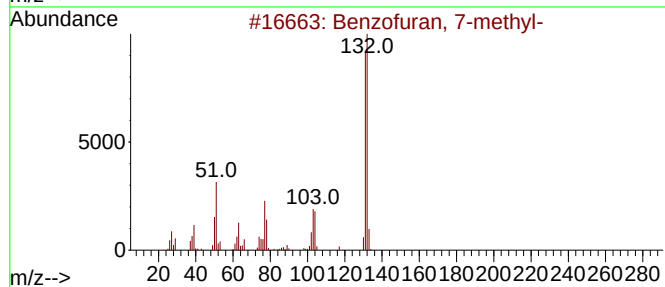
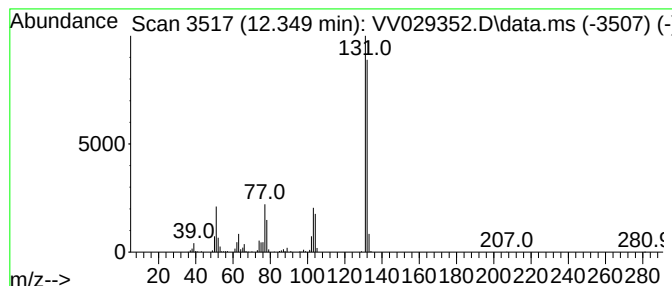
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzofuran, 7-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.349	1.60 ug/L	264193	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	96
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

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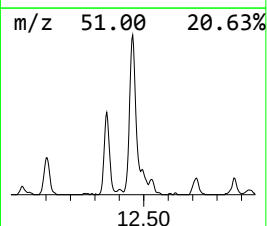
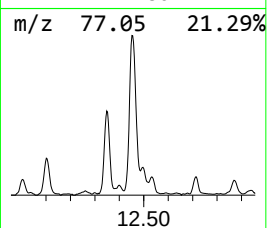
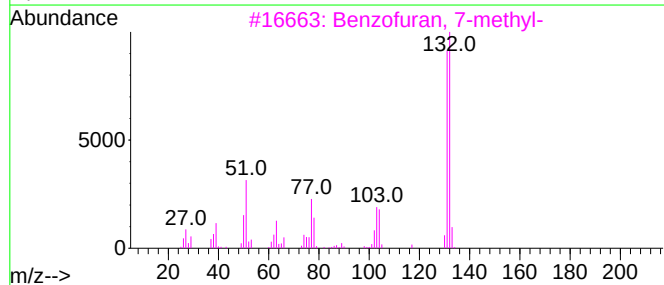
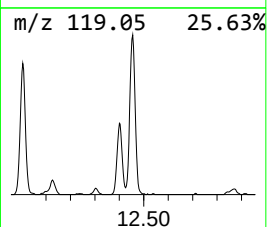
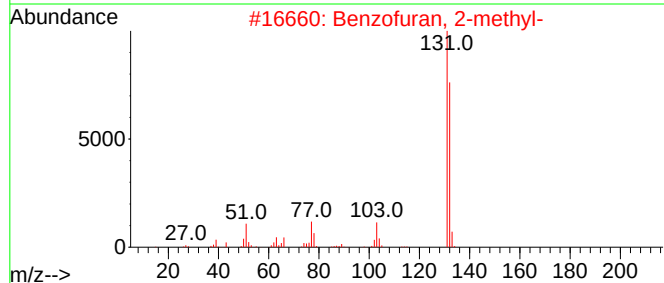
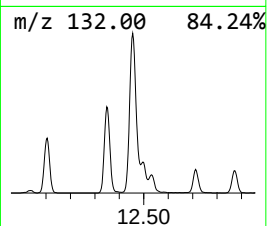
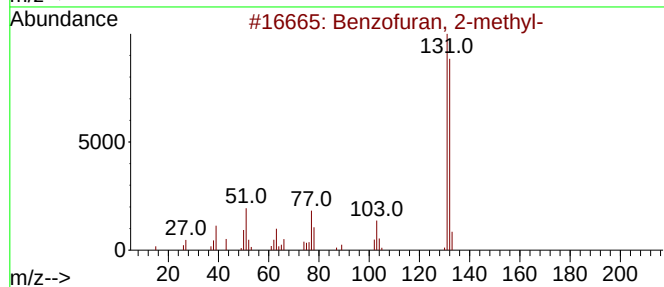
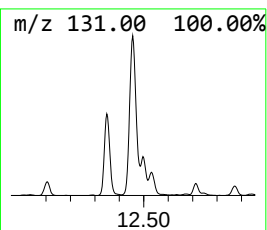
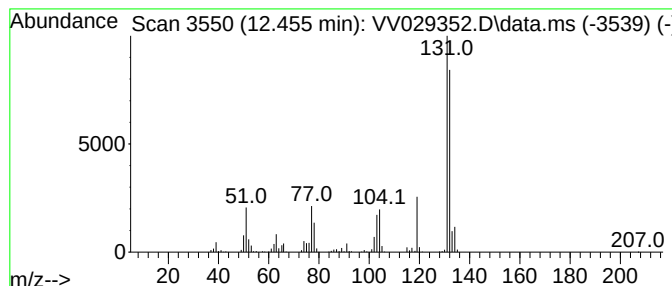
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzofuran, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.455	3.92 ug/L	648561	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
5			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

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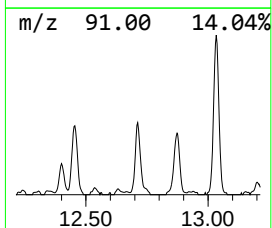
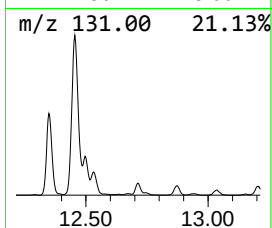
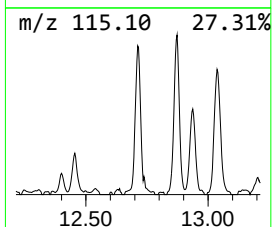
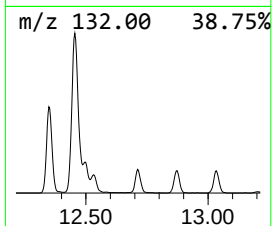
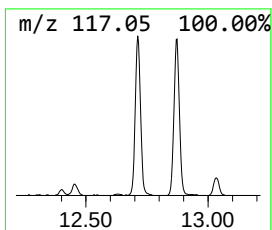
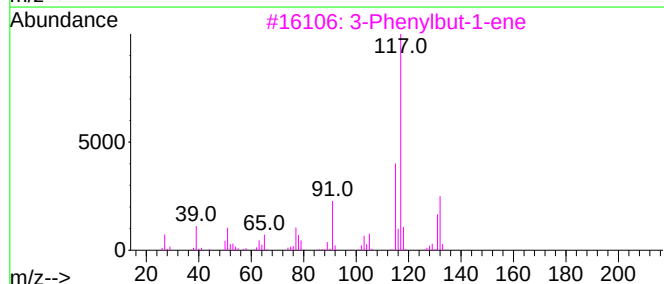
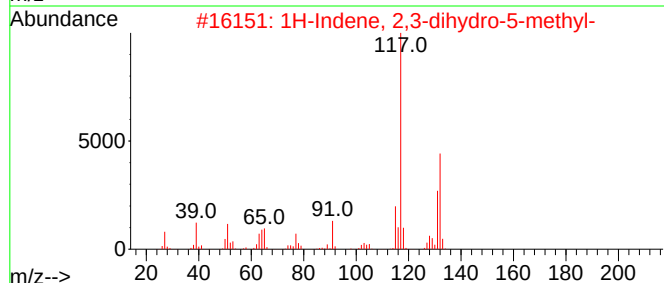
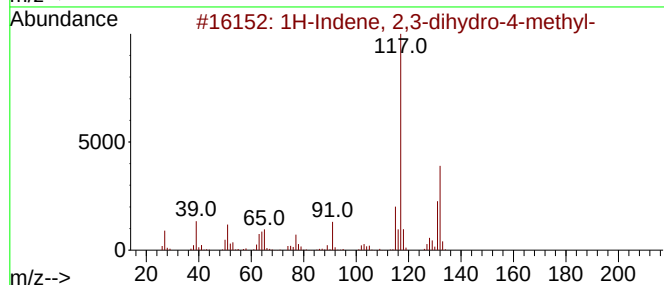
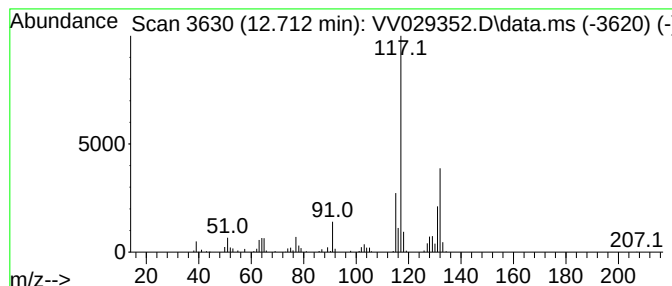
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.712	0.84 ug/L	139609	1,4-Dichlorobenzene-d4	11.239

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		Indan, 1-methyl-	132	C10H12	000767-58-8	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

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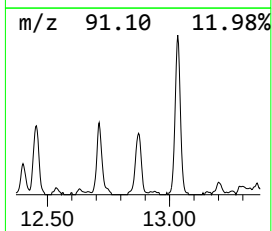
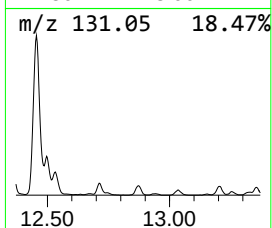
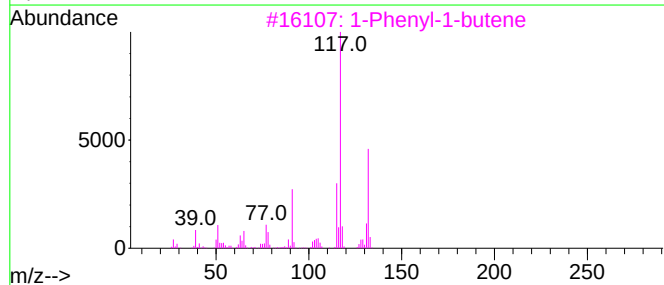
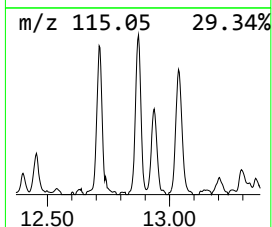
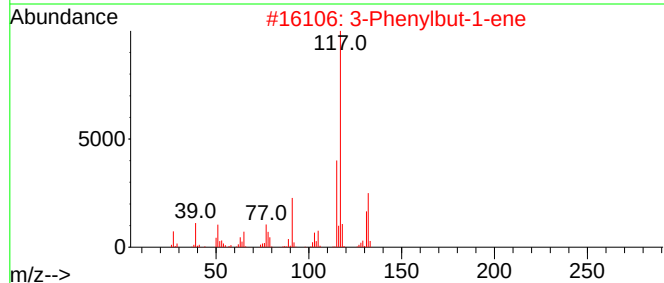
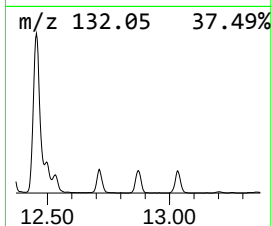
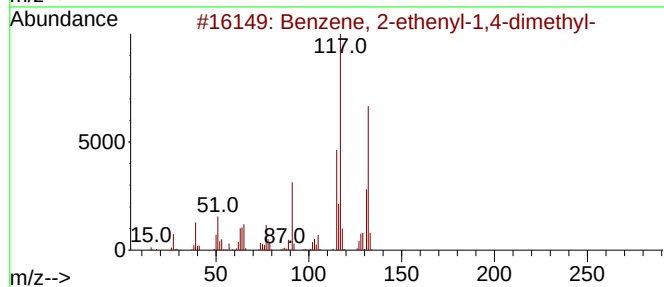
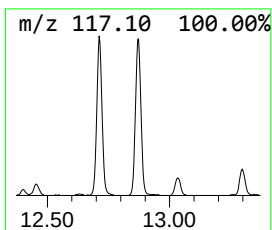
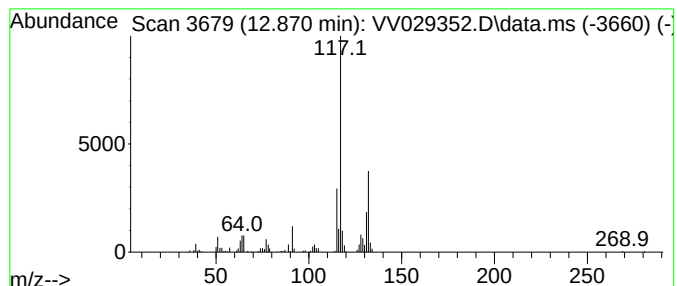
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.870	0.94 ug/L	155439	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB67DL

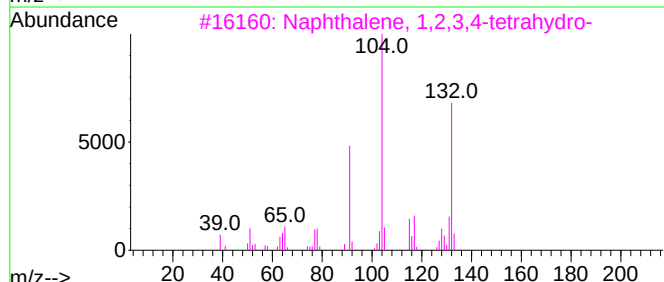
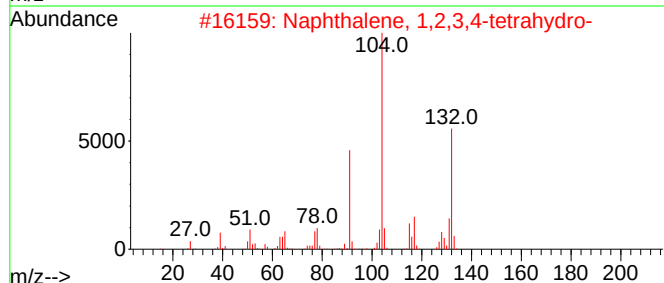
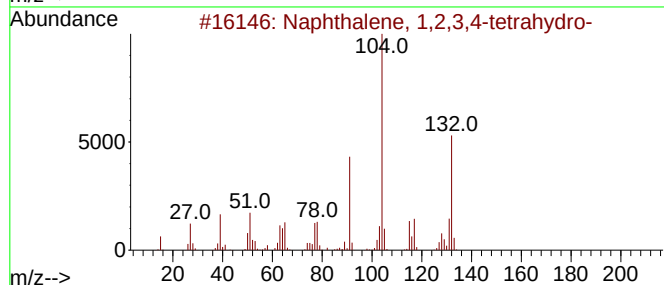
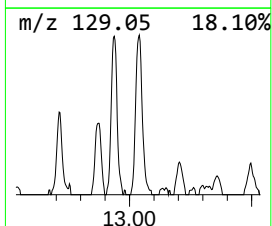
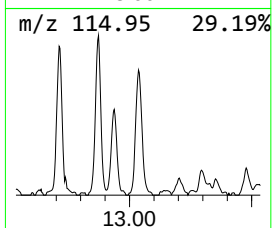
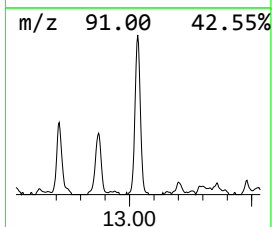
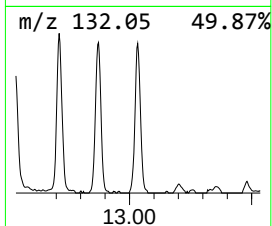
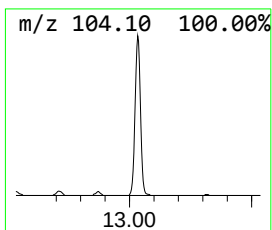
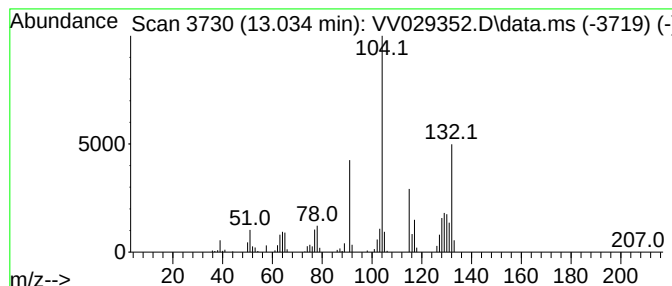
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.034	0.95 ug/L	157526	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
Data File : VV029352.D
Acq On : 25 Nov 2022 13:47
Operator : SY/MD
Sample : N5725-14DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB67DL

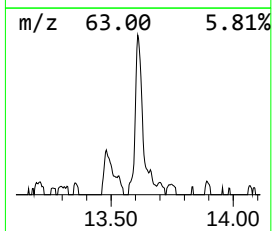
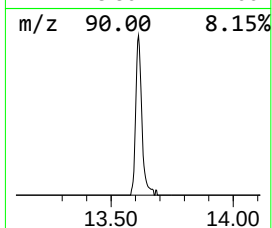
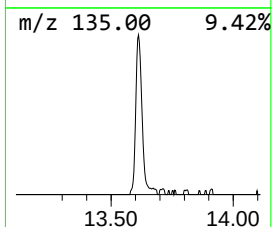
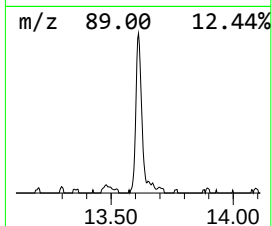
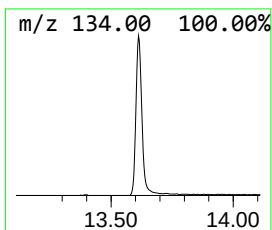
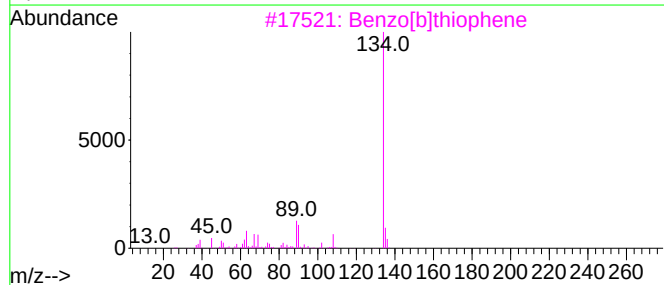
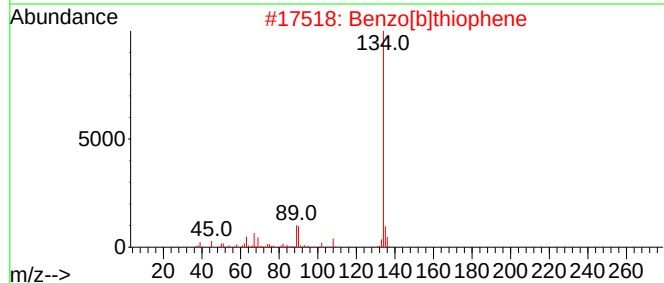
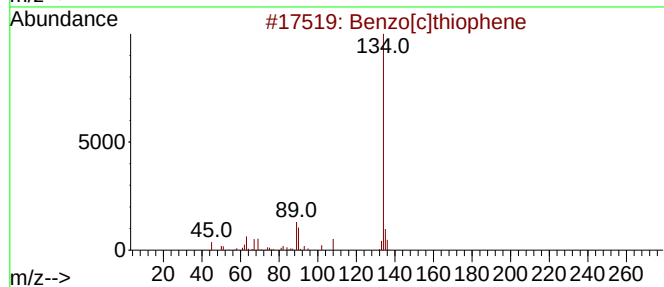
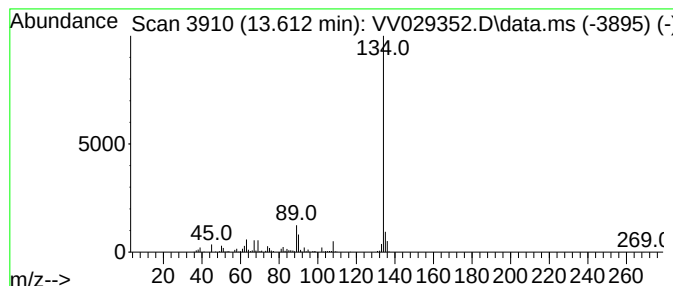
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 14 Benzo[c]thiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.612	0.98 ug/L	162828	1,4-Dichlorobenzene-d4	11.239

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	95
3		Benzo[b]thiophene	134	C8H6S	000095-15-8	94
4		Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
5		Benzo[b]thiophene	134	C8H6S	000095-15-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112522\
 Data File : VV029352.D
 Acq On : 25 Nov 2022 13:47
 Operator : SY/MD
 Sample : N5725-14DL 4X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BHB67DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.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
(DEL) Alkane: S...	1.105	22.4	ug/L	3095300	1	5.609	689607	5.0
(DEL) Alkane: S...	1.214	1.3	ug/L	178897	1	5.609	689607	5.0
Benzene, 1-ethy...	10.728	0.5	ug/L	88907	3	11.239	828086	5.0
Indane	11.516	6.2	ug/L	1029780	3	11.239	828086	5.0
Indene	11.735	11.5	ug/L	1900890	3	11.239	828086	5.0
Benzene, 2-ethy...	12.005	0.6	ug/L	93577	3	11.239	828086	5.0
1-Phenyl-1-butene	12.104	2.3	ug/L	385285	3	11.239	828086	5.0
Benzofuran, 7-m...	12.349	1.6	ug/L	264193	3	11.239	828086	5.0
Benzofuran, 2-m...	12.455	3.9	ug/L	648561	3	11.239	828086	5.0
1H-Indene, 2,3-...	12.712	0.8	ug/L	139609	3	11.239	828086	5.0
Benzene, 2-ethe...	12.870	0.9	ug/L	155439	3	11.239	828086	5.0
Naphthalene, 1,...	13.034	0.9	ug/L	157526	3	11.239	828086	5.0
Benzo[c]thiophene	13.612	1.0	ug/L	162828	3	11.239	828086	5.0