

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
 Data File : VV029328.D
 Acq On : 24 Nov 2022 05:38
 Operator : SY/MD
 Sample : N5694-06DL 5X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BHB42DL

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: VV029328.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.108	15	21	35	rVB	2020502	1830904	100.00%	13.813%
2	1.217	49	55	67	rVB	138890	133413	7.29%	1.007%
3	1.304	76	82	100	rVB2	170490	191261	10.45%	1.443%
4	1.565	156	163	175	rVB	108672	125086	6.83%	0.944%
5	2.105	320	331	349	rBV2	228405	351117	19.18%	2.649%
6	2.503	445	455	469	rBV	16442	28460	1.55%	0.215%
7	3.185	648	667	693	rBV	79843	169611	9.26%	1.280%
8	3.902	874	890	943	rBV2	322522	989379	54.04%	7.464%
9	4.339	1009	1026	1050	rBV	142146	362199	19.78%	2.733%
10	4.603	1091	1108	1124	rBV3	27386	70102	3.83%	0.529%
11	5.040	1226	1244	1256	rBV2	356157	882732	48.21%	6.660%
12	5.609	1407	1421	1447	rBV	264355	545127	29.77%	4.113%
13	5.908	1501	1514	1535	rBV2	109016	223945	12.23%	1.690%
14	6.063	1548	1562	1596	rBV	191426	404635	22.10%	3.053%
15	6.979	1832	1847	1877	rBV2	92589	179542	9.81%	1.355%
16	7.307	1937	1949	1967	rBV	451423	805215	43.98%	6.075%
17	7.616	2035	2045	2065	rBV	54208	100870	5.51%	0.761%
18	8.088	2179	2192	2223	rBV2	363254	736232	40.21%	5.555%
19	8.844	2410	2427	2454	rBV	378071	625352	34.16%	4.718%
20	9.133	2506	2517	2526	rBV	11841	20707	1.13%	0.156%
21	9.535	2633	2642	2659	rVB	40208	63321	3.46%	0.478%
22	9.921	2749	2762	2782	rBV	51436	81587	4.46%	0.616%
23	10.207	2839	2851	2871	rVB	130737	207372	11.33%	1.565%
24	11.239	3160	3172	3190	rBV	415707	654069	35.72%	4.935%
25	11.516	3247	3258	3277	rBV	111211	190703	10.42%	1.439%
26	11.612	3277	3288	3307	rVV	449441	705543	38.54%	5.323%
27	11.735	3315	3326	3340	rBV	141713	218660	11.94%	1.650%
28	11.821	3342	3353	3365	rVB	16518	28759	1.57%	0.217%
29	12.104	3427	3441	3455	rBV	88706	136249	7.44%	1.028%
30	12.349	3501	3517	3538	rBV	458065	703873	38.44%	5.310%
31	12.455	3538	3550	3558	rVV	100342	190098	10.38%	1.434%
32	12.497	3558	3563	3569	rVV	56360	90127	4.92%	0.680%
33	12.532	3569	3574	3584	rVB2	47461	72723	3.97%	0.549%
34	12.712	3621	3630	3636	rBV	13809	21093	1.15%	0.159%
35	13.490	3845	3872	3891	rBV	131822	294190	16.07%	2.220%
36	13.612	3891	3910	3934	rVB	423338	723848	39.54%	5.461%

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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	14.564	4196	4206	4216	rBV	23332	38780	2.12%	0.293%
38	14.802	4269	4280	4295	rBV2	32799	57589	3.15%	0.434%

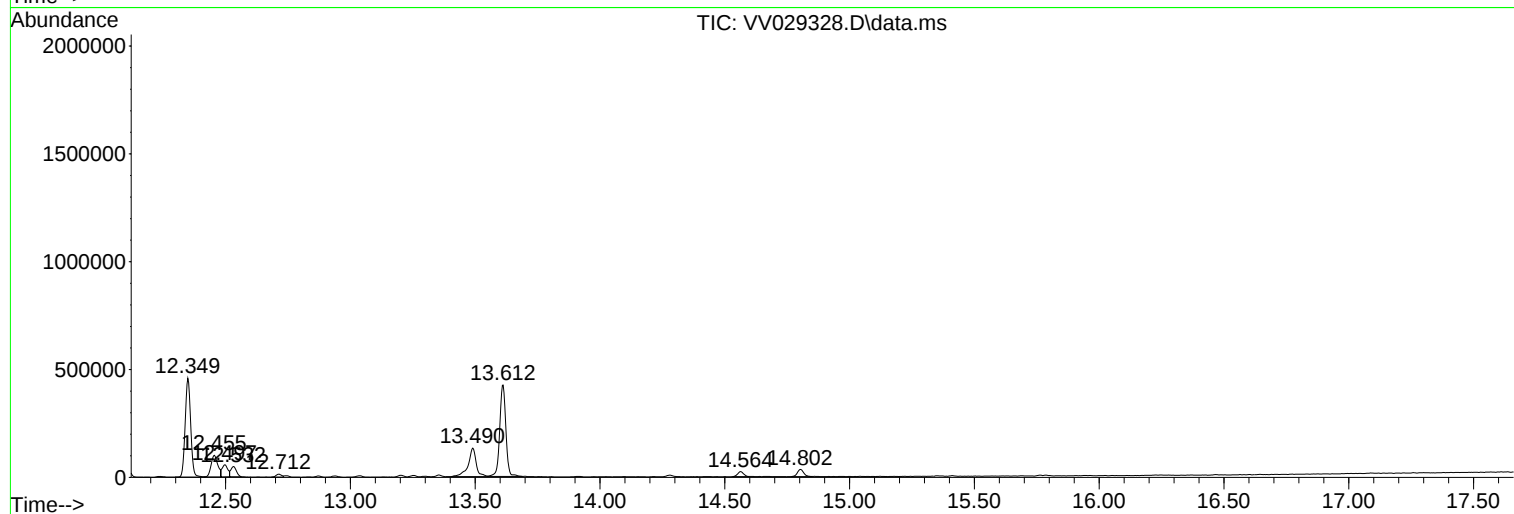
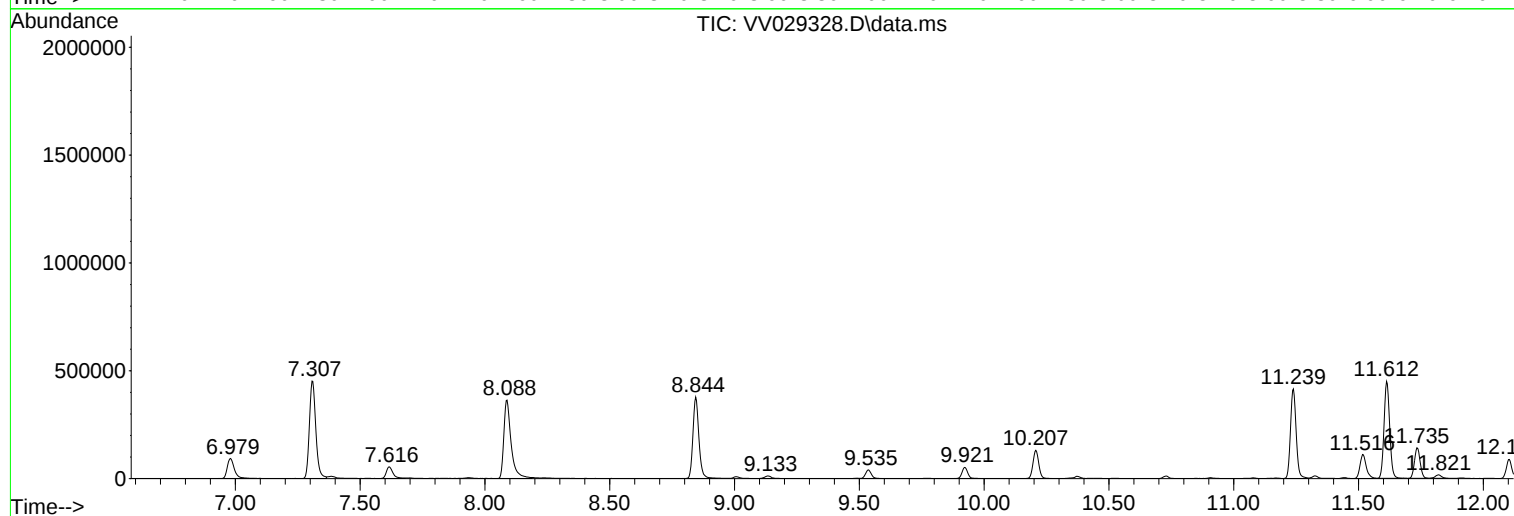
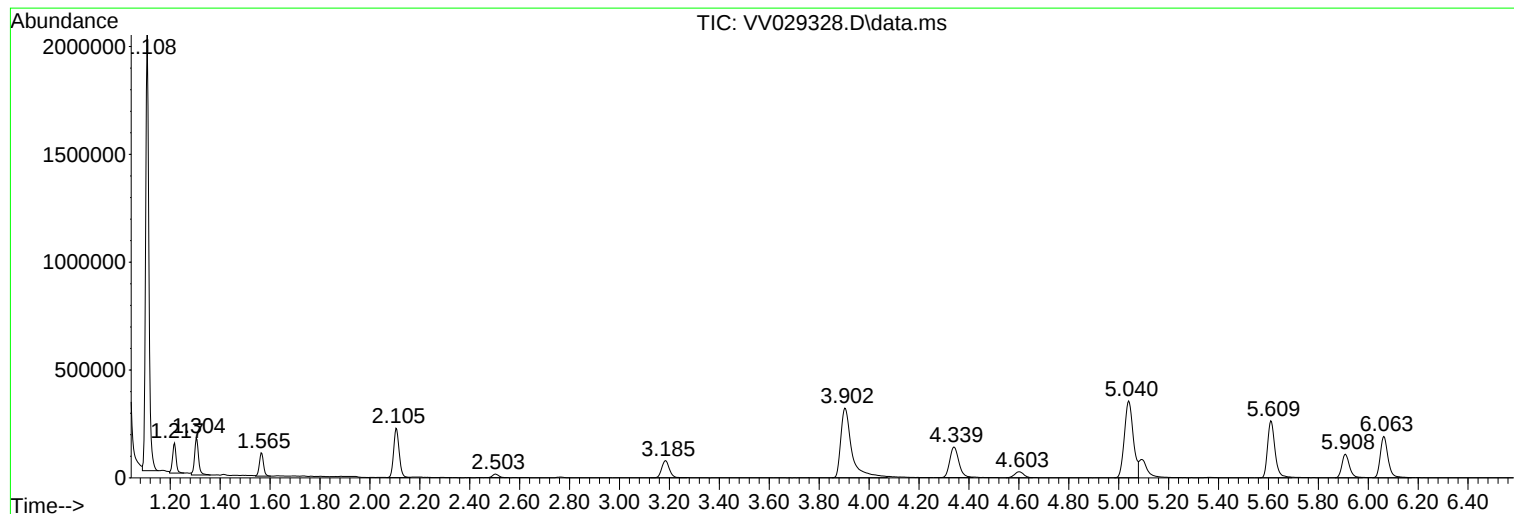
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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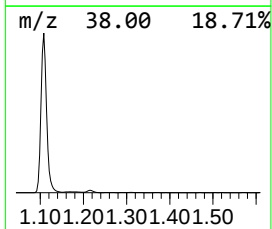
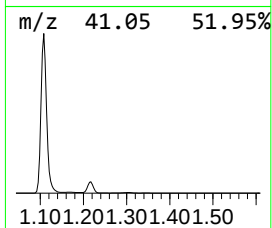
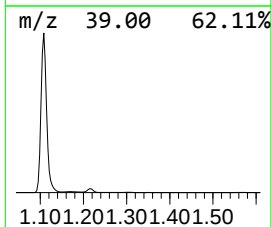
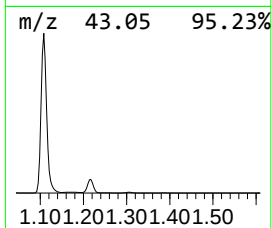
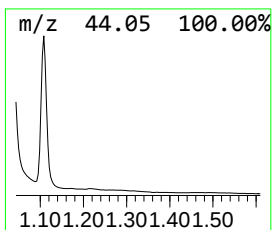
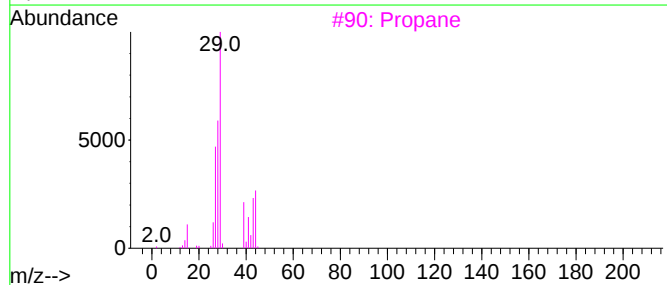
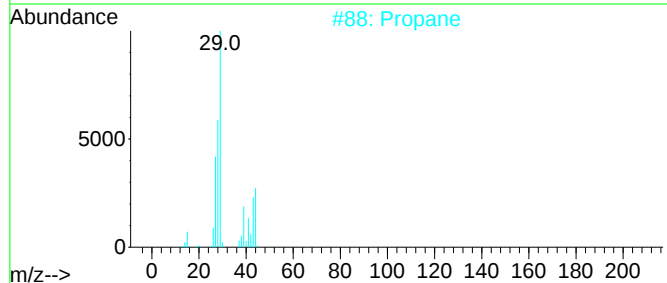
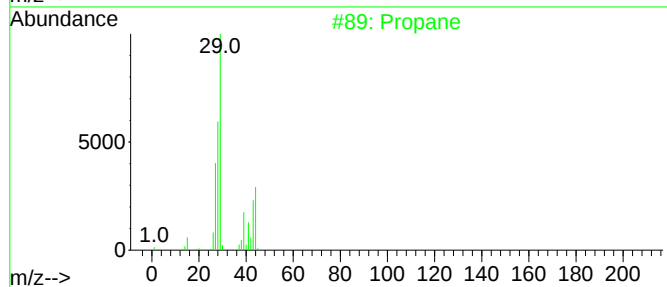
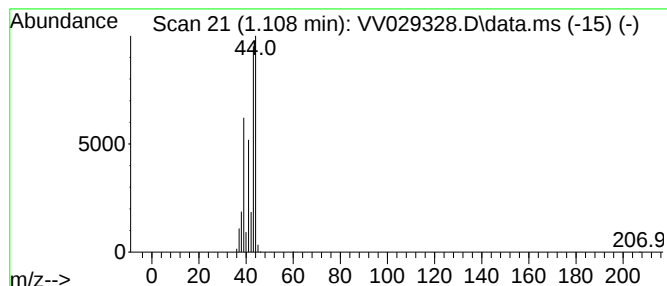
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.108	16.79 ug/L	1830900	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	78
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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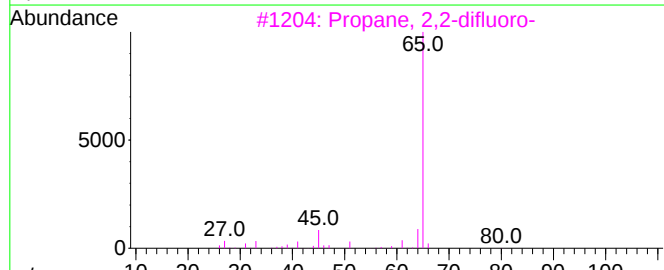
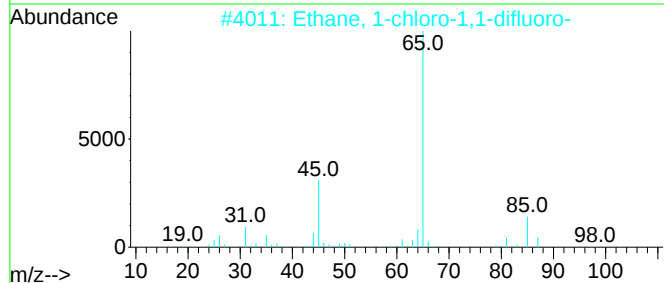
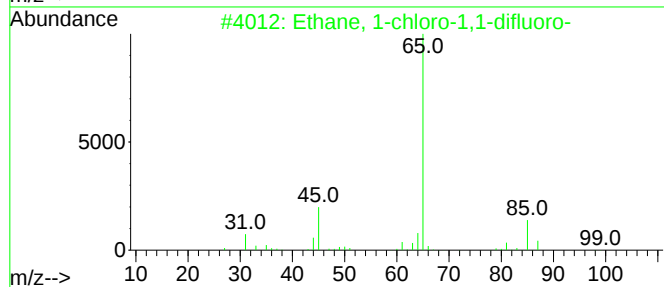
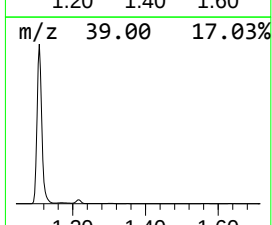
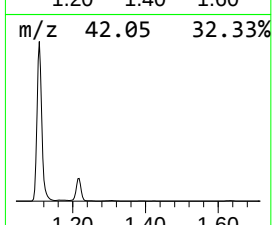
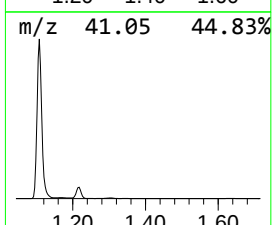
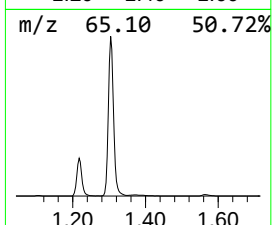
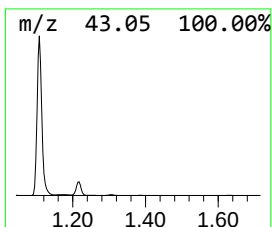
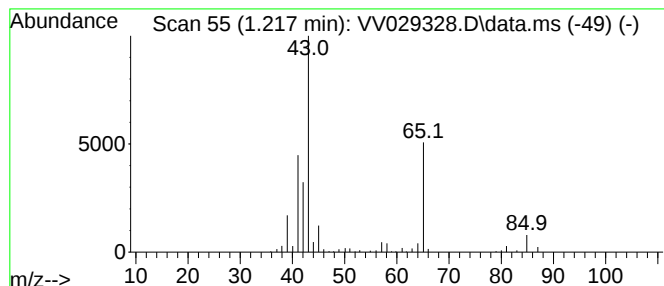
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Ethane, 1-chloro-1,1-difluoro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.217	1.22 ug/L	133413	1,4-Difluorobenzene	5.609

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1-chloro-1,1-difluoro-	100	C2H3ClF2	000075-68-3	64
2			Ethane, 1-chloro-1,1-difluoro-	100	C2H3ClF2	000075-68-3	45
3			Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	9
4			Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	9
5			Thioacetic acid	76	C2H4OS	000507-09-5	9



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ClientSampleId :
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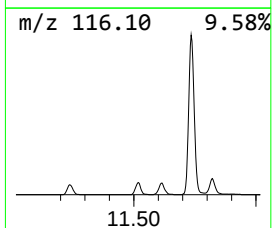
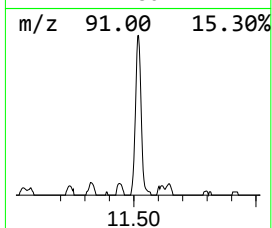
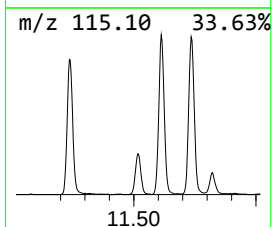
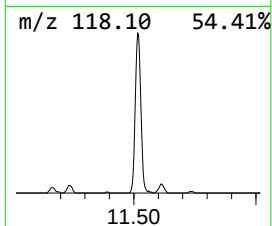
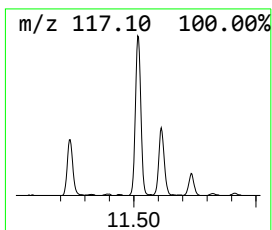
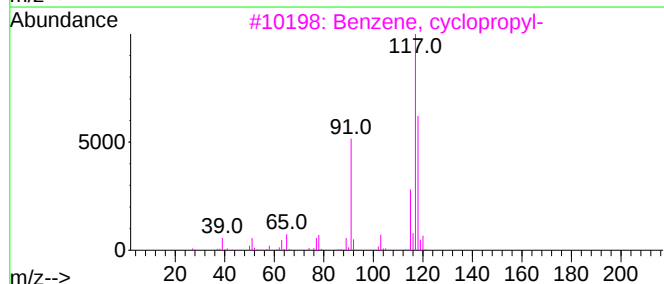
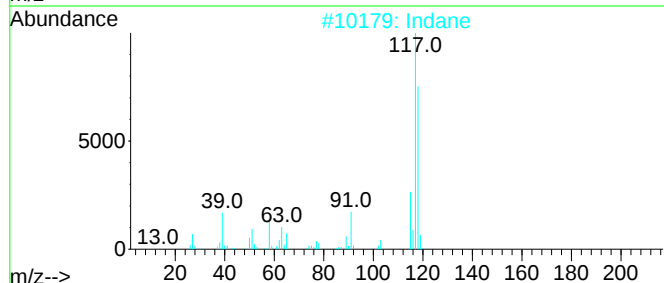
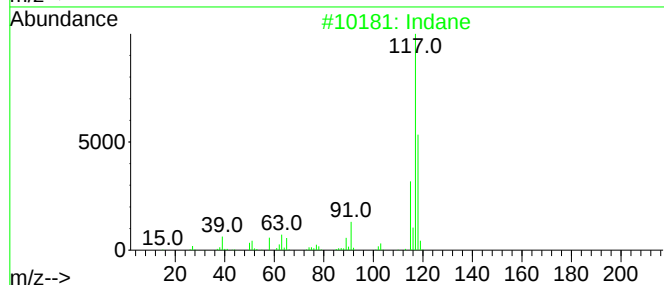
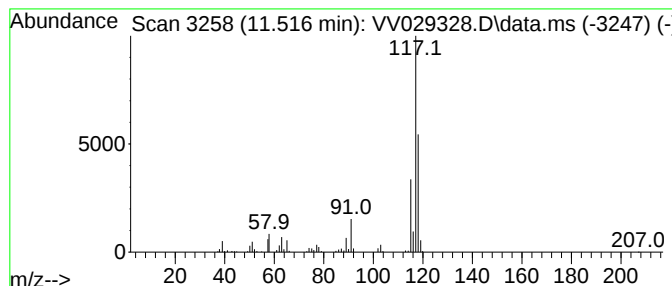
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.516	1.46 ug/L	190703	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			Benzene, cyclopropyl-	118	C9H10	000873-49-4	87
4			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	74
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74



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ClientSampleId :
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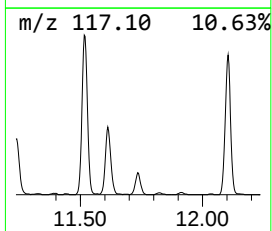
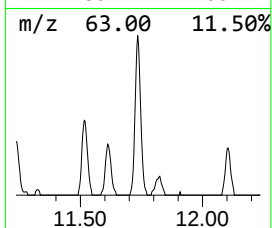
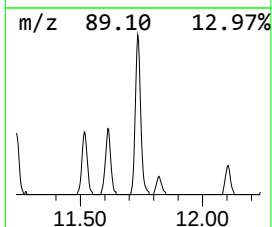
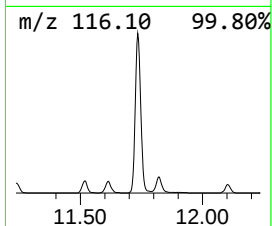
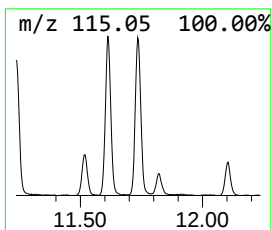
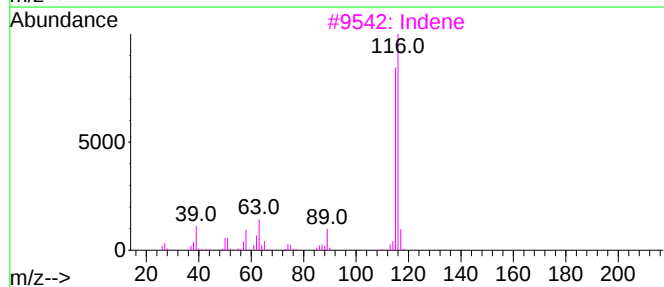
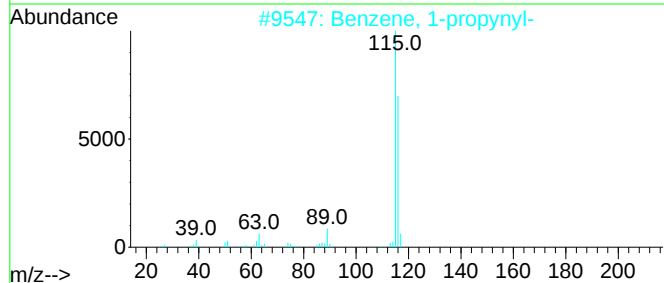
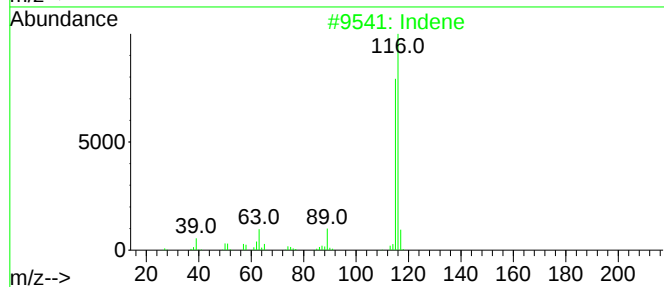
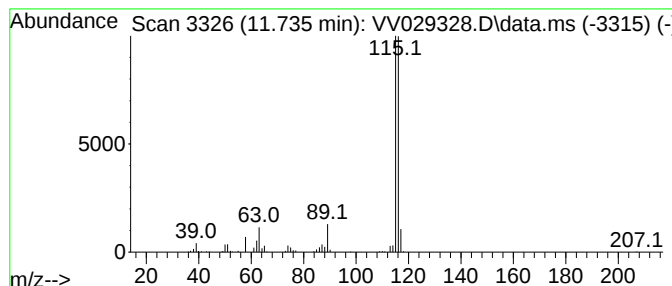
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Indene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.735	1.67 ug/L	218660	1,4-Dichlorobenzene-d4	11.239

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



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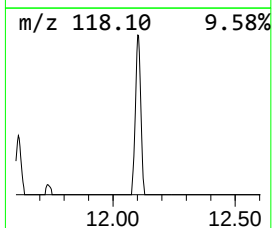
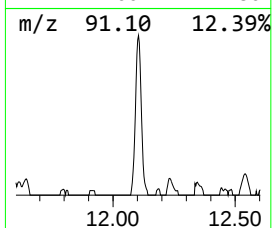
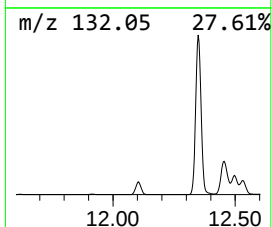
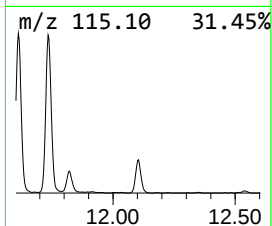
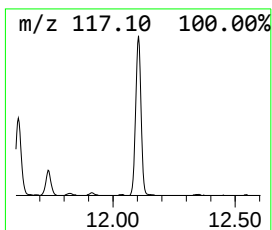
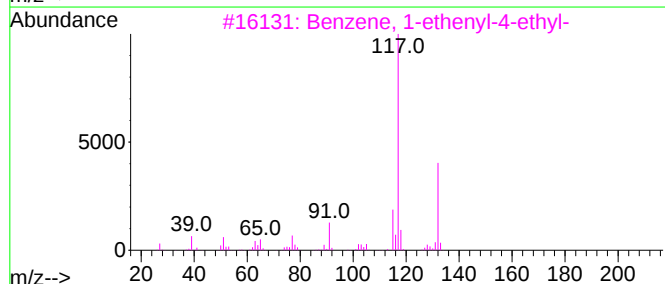
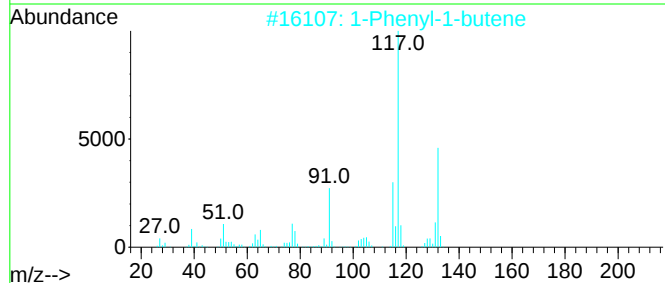
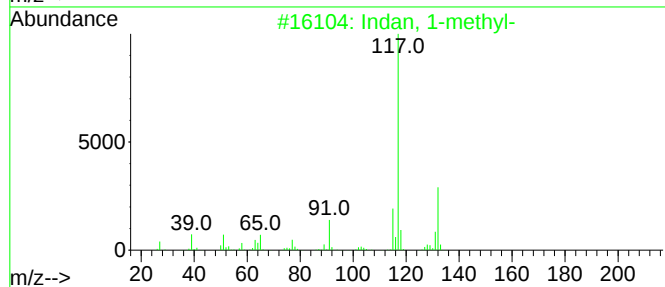
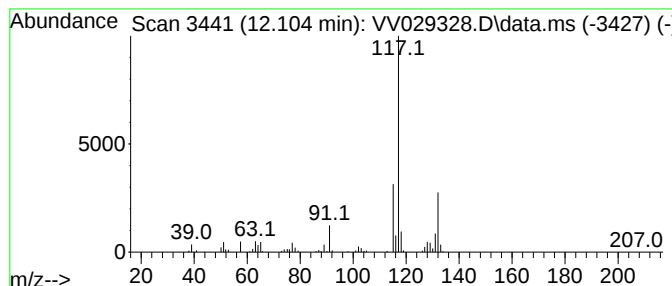
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	1.04 ug/L	136249	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	90
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
Data File : VV029328.D
Acq On : 24 Nov 2022 05:38
Operator : SY/MD
Sample : N5694-06DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB42DL

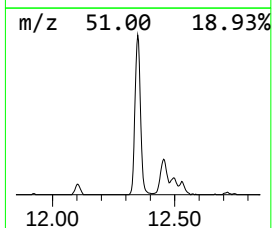
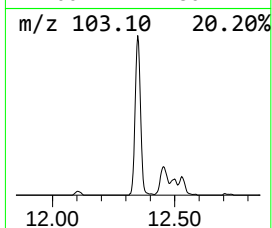
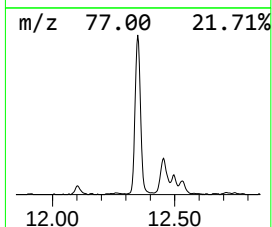
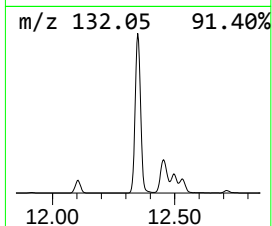
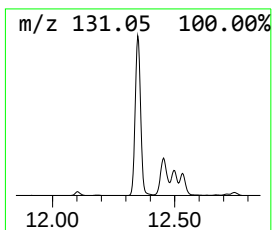
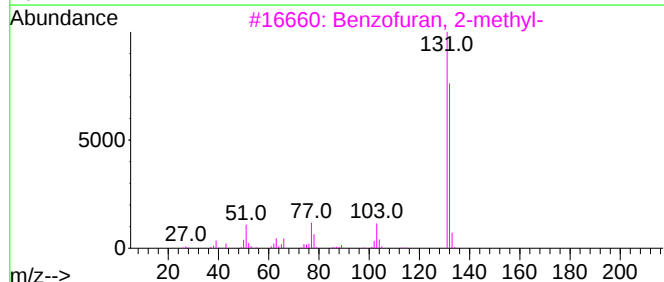
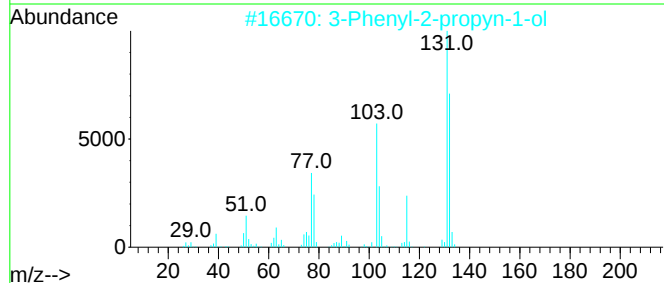
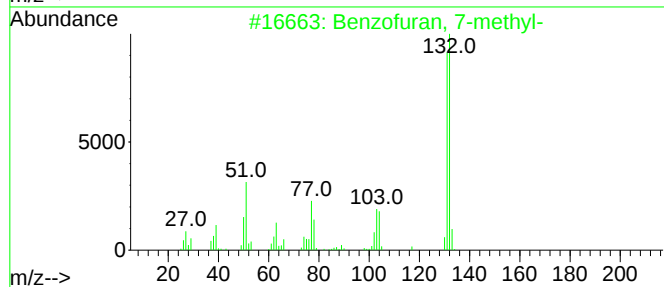
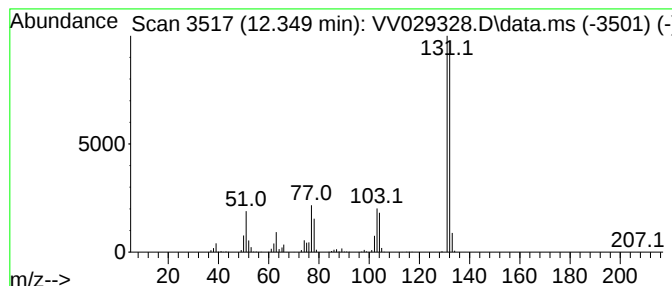
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzo[1,2-b:4,5-b']difuran, 7-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.349	5.38 ug/L	703873	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[1,2-b:4,5-b']difuran, 7-methyl-	132	C9H8O	017059-52-8	95
2			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
3			Benzo[1,2-b:4,5-b']difuran, 2-methyl-	132	C9H8O	004265-25-2	91
4			Benzo[1,2-b:4,5-b']difuran, 2-methyl-	132	C9H8O	004265-25-2	91
5			Benzo[1,2-b:4,5-b']difuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
Data File : VV029328.D
Acq On : 24 Nov 2022 05:38
Operator : SY/MD
Sample : N5694-06DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB42DL

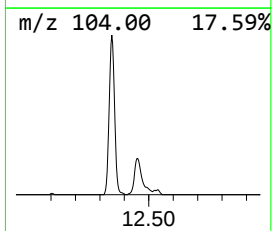
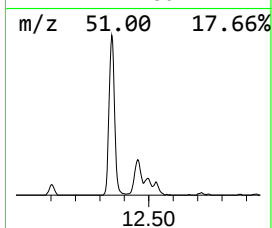
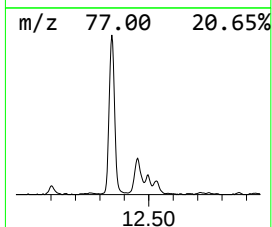
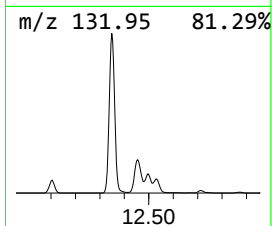
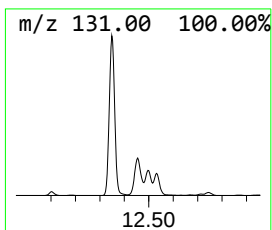
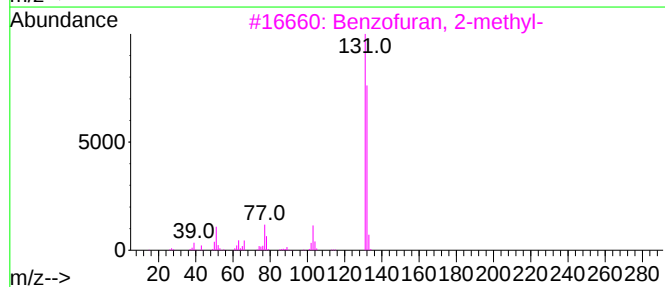
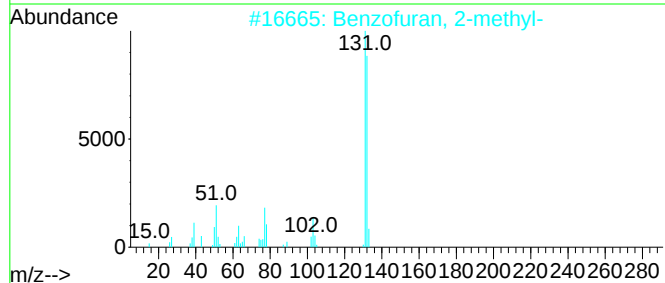
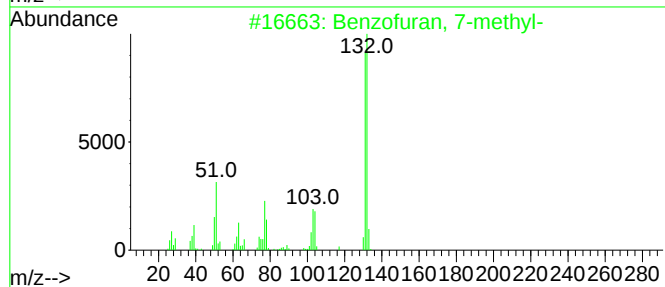
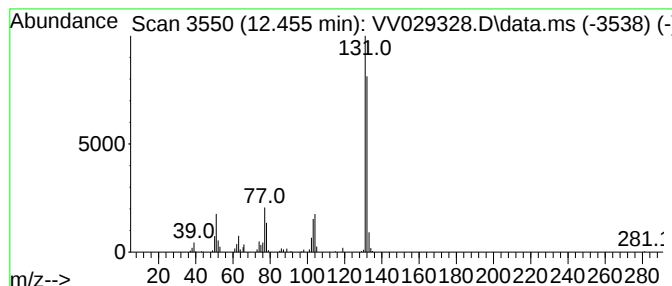
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.455	1.45 ug/L	190098	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
Data File : VV029328.D
Acq On : 24 Nov 2022 05:38
Operator : SY/MD
Sample : N5694-06DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB42DL

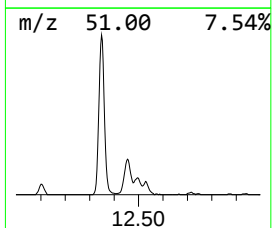
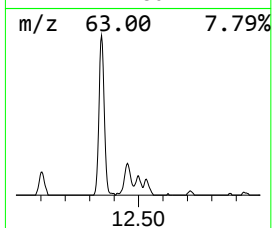
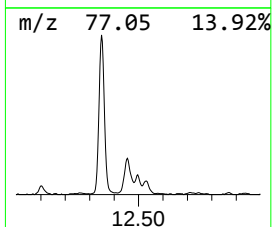
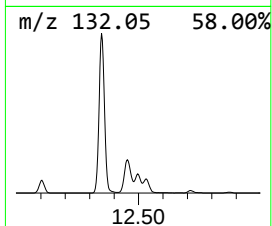
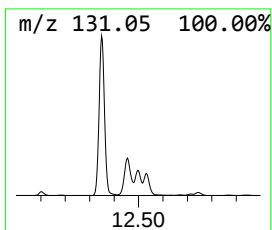
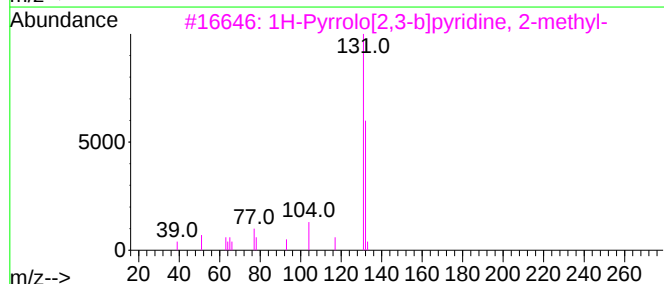
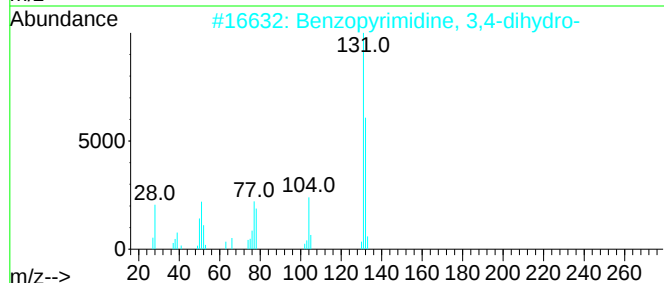
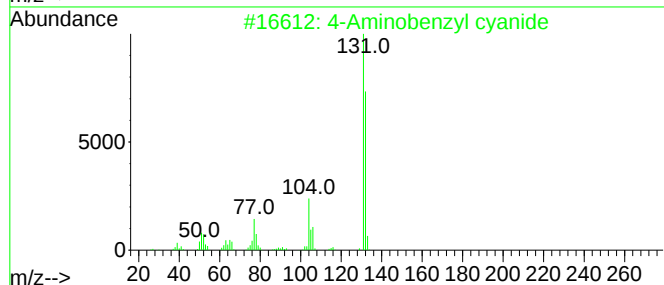
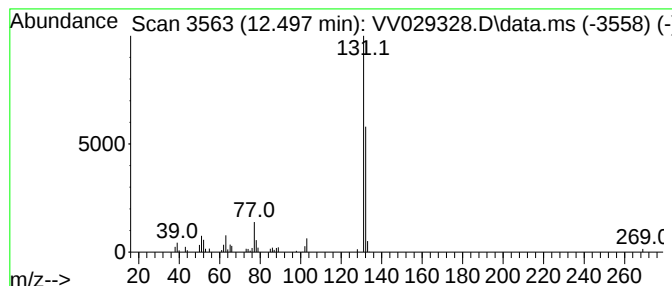
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 4-Aminobenzyl cyanide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.497	0.69 ug/L	90127	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Aminobenzyl cyanide	132	C8H8N2	003544-25-0	86
2			Benzopyrimidine, 3,4-dihydro-	132	C8H8N2	001904-64-9	72
3			1H-Pyrrolo[2,3-b]pyridine, 2-met...	132	C8H8N2	023612-48-8	72
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	58
5			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
Data File : VV029328.D
Acq On : 24 Nov 2022 05:38
Operator : SY/MD
Sample : N5694-06DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB42DL

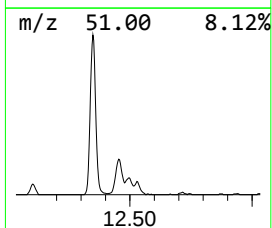
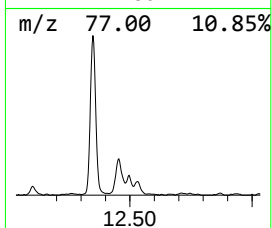
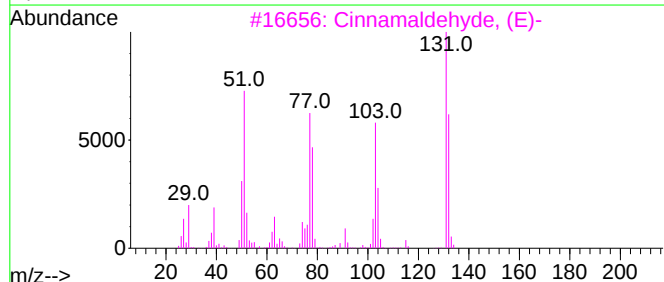
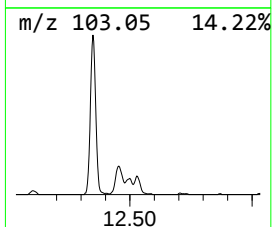
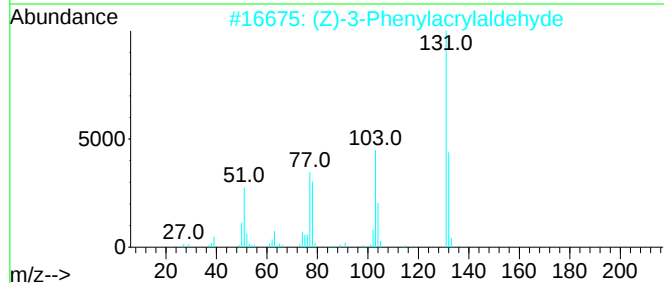
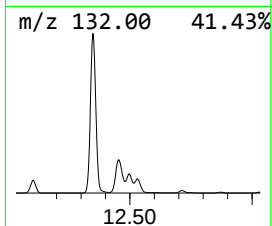
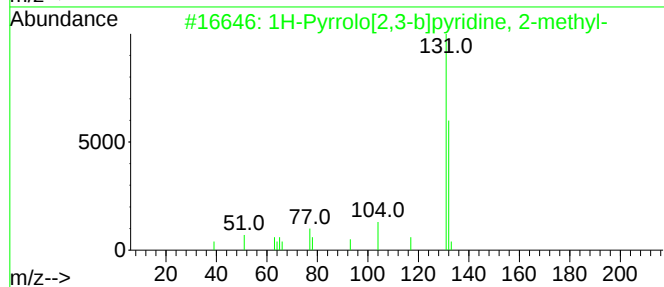
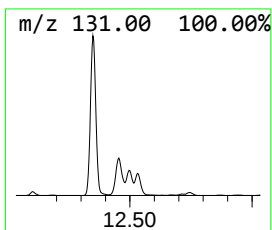
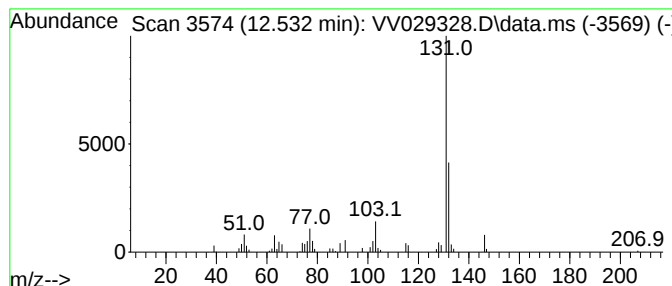
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 1H-Pyrrolo[2,3-b]pyridine, ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.532	0.56 ug/L	72723	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrrolo[2,3-b]pyridine, 2-met...	132	C8H8N2	023612-48-8	74
2			(Z)-3-Phenylacrylaldehyde	132	C9H8O	057194-69-1	64
3			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	50
4			Cinnamoyl chloride	166	C9H7ClO	000102-92-1	50
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
Data File : VV029328.D
Acq On : 24 Nov 2022 05:38
Operator : SY/MD
Sample : N5694-06DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB42DL

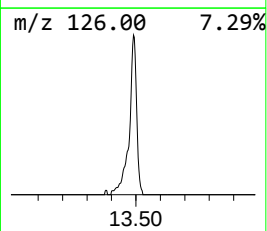
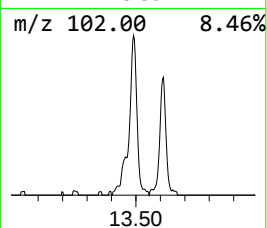
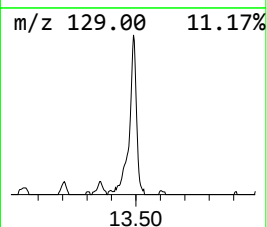
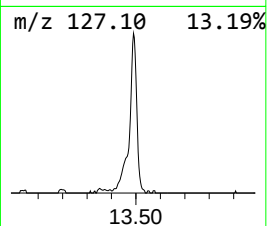
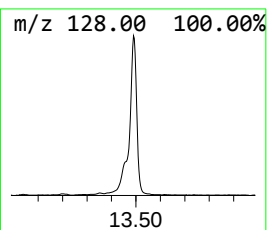
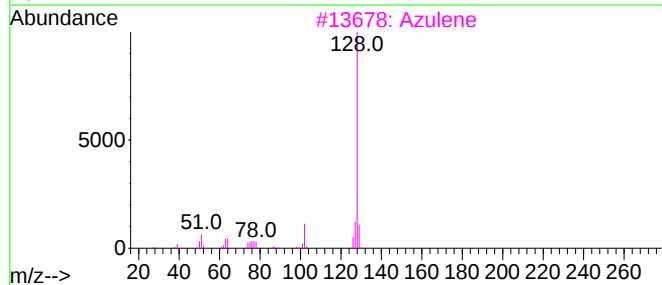
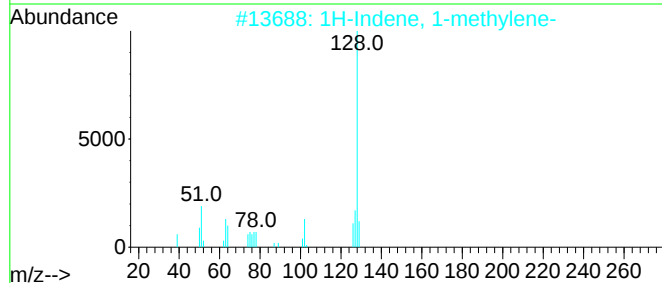
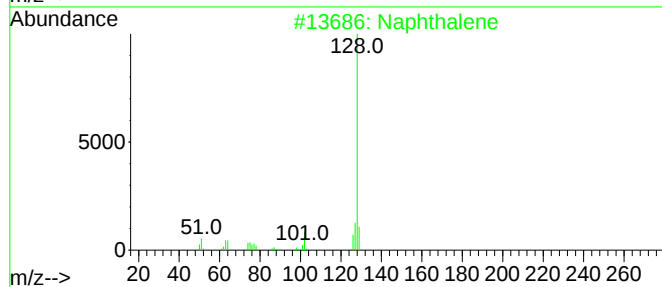
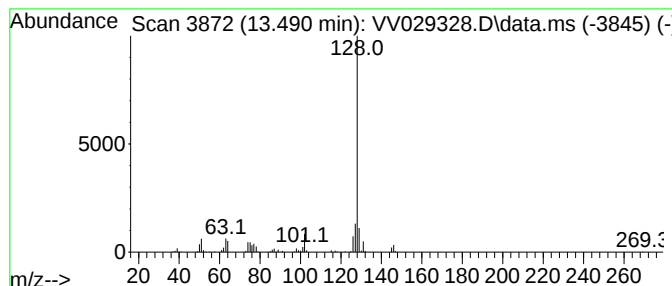
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 1H-Indene, 1-methylene- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.490	2.25 ug/L	294190	1,4-Dichlorobenzene-d4	11.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	94
2			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	93
3			Azulene	128	C10H8	000275-51-4	93
4			Naphthalene	128	C10H8	000091-20-3	93
5			Azulene	128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112322\
Data File : VV029328.D
Acq On : 24 Nov 2022 05:38
Operator : SY/MD
Sample : N5694-06DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHB42DL

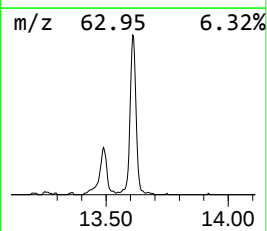
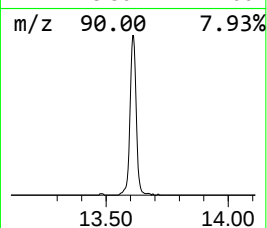
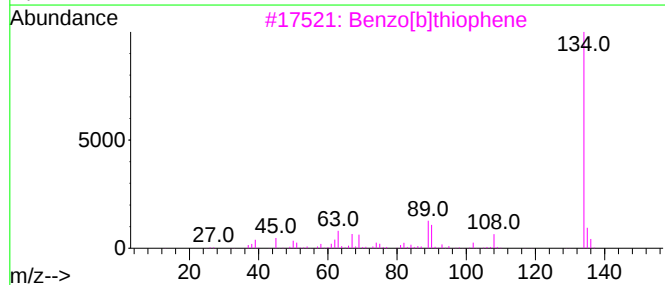
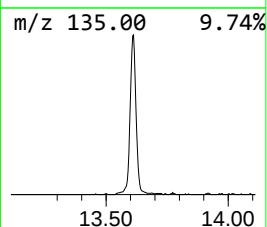
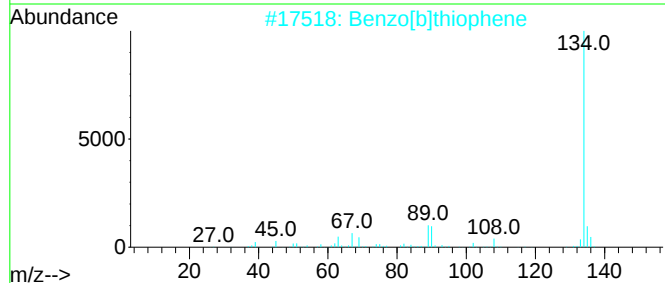
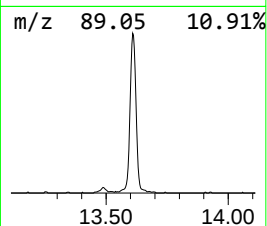
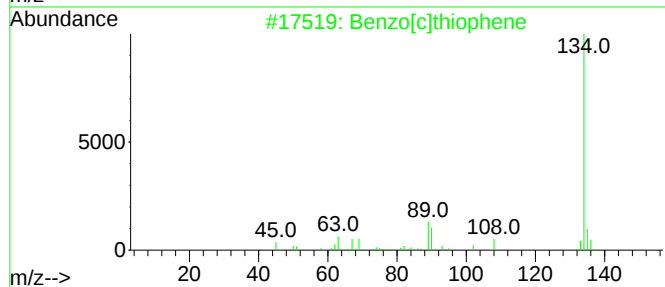
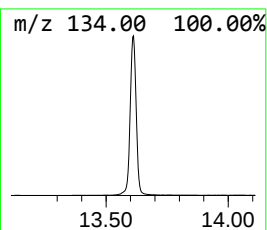
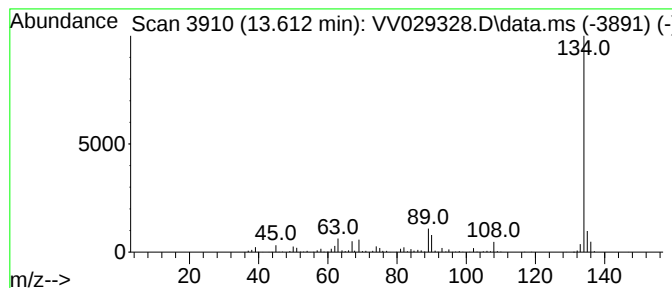
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Benzo[c]thiophene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.612	5.53 ug/L	723848	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\VW112322\
 Data File : VV029328.D
 Acq On : 24 Nov 2022 05:38
 Operator : SY/MD
 Sample : N5694-06DL 5X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BHB42DL

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...	1.108	16.8	ug/L	1830900	1	5.609	545127	5.0
Ethane, 1-chlor...	1.217	1.2	ug/L	133413	1	5.609	545127	5.0
Indane	11.516	1.5	ug/L	190703	3	11.239	654069	5.0
Indene	11.735	1.7	ug/L	218660	3	11.239	654069	5.0
Indan, 1-methyl-	12.104	1.0	ug/L	136249	3	11.239	654069	5.0
Benzofuran, 7-m...	12.349	5.4	ug/L	703873	3	11.239	654069	5.0
Benzofuran, 2-m...	12.455	1.5	ug/L	190098	3	11.239	654069	5.0
4-Aminobenzyl c...	12.497	0.7	ug/L	90127	3	11.239	654069	5.0
1H-Pyrrolo[2,3-...	12.532	0.6	ug/L	72723	3	11.239	654069	5.0
1H-Indene, 1-me...	13.490	2.3	ug/L	294190	3	11.239	654069	5.0
Benzo[c]thiophene	13.612	5.5	ug/L	723848	3	11.239	654069	5.0