

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
 Data File : VV023697.D
 Acq On : 24 Nov 2021 13:53
 Operator : SY/MD
 Sample : M4723-20DL 40X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0G67DL

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\SFAMVTR112321WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VV023697.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	16	21	37	rVB	28768	29489	4.56%	0.463%
2	1.201	44	50	58	rBV	17942	17616	2.73%	0.276%
3	1.304	74	82	95	rBV	73390	76751	11.88%	1.204%
4	1.568	156	164	172	rBV	64088	70133	10.86%	1.100%
5	1.632	175	184	197	rVB	25747	34118	5.28%	0.535%
6	1.806	230	238	248	rBV2	18151	25960	4.02%	0.407%
7	2.024	298	306	317	rBV2	19106	27254	4.22%	0.428%
8	2.108	321	332	351	rVB	118258	174509	27.01%	2.738%
9	2.481	439	448	451	rBV2	13254	19689	3.05%	0.309%
10	2.577	469	478	492	rVB2	30170	59317	9.18%	0.931%
11	2.761	522	535	550	rVB2	9771	20648	3.20%	0.324%
12	3.043	609	623	635	rBV3	8296	17668	2.73%	0.277%
13	3.275	681	695	710	rVB3	7666	17265	2.67%	0.271%
14	3.365	710	723	733	rBV6	4365	9041	1.40%	0.142%
15	3.433	733	744	756	rVV4	6282	14305	2.21%	0.224%
16	3.506	757	767	777	rVV7	3000	6933	1.07%	0.109%
17	3.580	780	790	806	rVB6	4131	9390	1.45%	0.147%
18	3.764	830	847	863	rBV3	40881	106107	16.42%	1.665%
19	3.912	882	893	916	rBV2	21298	80449	12.45%	1.262%
20	4.349	1014	1029	1052	rBV2	81734	209649	32.45%	3.289%
21	4.458	1057	1063	1079	rVB3	7332	15799	2.45%	0.248%
22	4.677	1113	1131	1148	rBV4	53294	137743	21.32%	2.161%
23	5.050	1231	1247	1254	rBV3	187402	448566	69.43%	7.038%
24	5.098	1254	1262	1282	rVB	292227	646088	100.00%	10.137%
25	5.223	1290	1301	1321	rVB4	25646	69516	10.76%	1.091%
26	5.323	1321	1332	1347	rVB8	8691	20942	3.24%	0.329%
27	5.619	1411	1424	1453	rBV	154051	331950	51.38%	5.208%
28	5.934	1508	1522	1536	rBV9	4405	11856	1.84%	0.186%
29	6.069	1551	1564	1575	rBV	103246	213956	33.12%	3.357%
30	6.130	1576	1583	1600	rVB4	45164	92261	14.28%	1.448%
31	6.362	1645	1655	1668	rBV6	4003	7239	1.12%	0.114%
32	6.548	1699	1713	1714	rBV4	7438	9519	1.47%	0.149%
33	6.805	1780	1793	1805	rVB2	10007	18287	2.83%	0.287%
34	6.989	1840	1850	1872	rBV	50680	98216	15.20%	1.541%
35	7.169	1895	1906	1921	rVB3	13842	24713	3.83%	0.388%
36	7.317	1941	1952	1968	rVV	222031	387470	59.97%	6.079%

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 Title : TRACE VOA SFAM1.0

37	7.394	1968	1976	1986	rVB	25433	43943	6.80%	0.689%
38	7.625	2039	2048	2069	rBV2	28200	54629	8.46%	0.857%
39	7.747	2078	2086	2095	rBV8	3963	7646	1.18%	0.120%
40	8.091	2182	2193	2229	rBV	150910	377700	58.46%	5.926%
41	8.854	2417	2430	2452	rBV	229959	382512	59.20%	6.001%
42	9.014	2468	2480	2502	rBV	131583	211256	32.70%	3.314%
43	9.143	2505	2520	2542	rVB	38595	70151	10.86%	1.101%
44	9.548	2636	2646	2658	rBV3	6707	11197	1.73%	0.176%
45	9.934	2754	2766	2781	rBV	15030	24993	3.87%	0.392%
46	10.217	2843	2854	2871	rBV	72811	119323	18.47%	1.872%
47	10.355	2887	2897	2912	rBV	21060	34246	5.30%	0.537%
48	10.477	2928	2935	2945	rVB	7408	10906	1.69%	0.171%
49	10.542	2945	2955	2971	rVB2	5973	9901	1.53%	0.155%
50	10.738	3006	3016	3026	rBV2	13467	20518	3.18%	0.322%
51	11.249	3163	3175	3194	rBV	239152	376144	58.22%	5.901%
52	11.336	3194	3202	3213	rVB2	8406	13001	2.01%	0.204%
53	11.529	3251	3262	3281	rBV	89027	151570	23.46%	2.378%
54	11.625	3281	3292	3310	rVV	232419	376544	58.28%	5.908%
55	11.747	3323	3330	3338	rVB5	5448	8089	1.25%	0.127%
56	12.017	3399	3414	3420	rBV	28902	46370	7.18%	0.728%
57	12.117	3434	3445	3467	rBV2	35032	75522	11.69%	1.185%
58	12.300	3495	3502	3510	rVB2	4879	7467	1.16%	0.117%
59	12.413	3521	3537	3547	rBV	21939	34813	5.39%	0.546%
60	12.464	3547	3553	3571	rVB2	10748	16809	2.60%	0.264%
61	12.725	3626	3634	3643	rVB2	14954	21107	3.27%	0.331%
62	12.882	3669	3683	3695	rBV2	76446	121595	18.82%	1.908%
63	13.046	3727	3734	3746	rVV10	4616	8508	1.32%	0.133%
64	13.217	3779	3787	3794	rVB3	5207	6873	1.06%	0.108%
65	13.268	3794	3803	3812	rBV4	10784	17010	2.63%	0.267%
66	13.709	3931	3940	3945	rBV7	4505	7615	1.18%	0.119%
67	13.914	3996	4004	4014	rVB10	3817	7794	1.21%	0.122%
68	14.098	4049	4061	4074	rBV7	8643	20585	3.19%	0.323%
69	14.429	4158	4164	4180	rBV5	5921	9800	1.52%	0.154%
70	14.635	4216	4228	4238	rBV	32090	52340	8.10%	0.821%
71	14.818	4274	4285	4299	rBV	27722	46989	7.27%	0.737%
72	15.811	4583	4594	4604	rBV	3665	7860	1.22%	0.123%

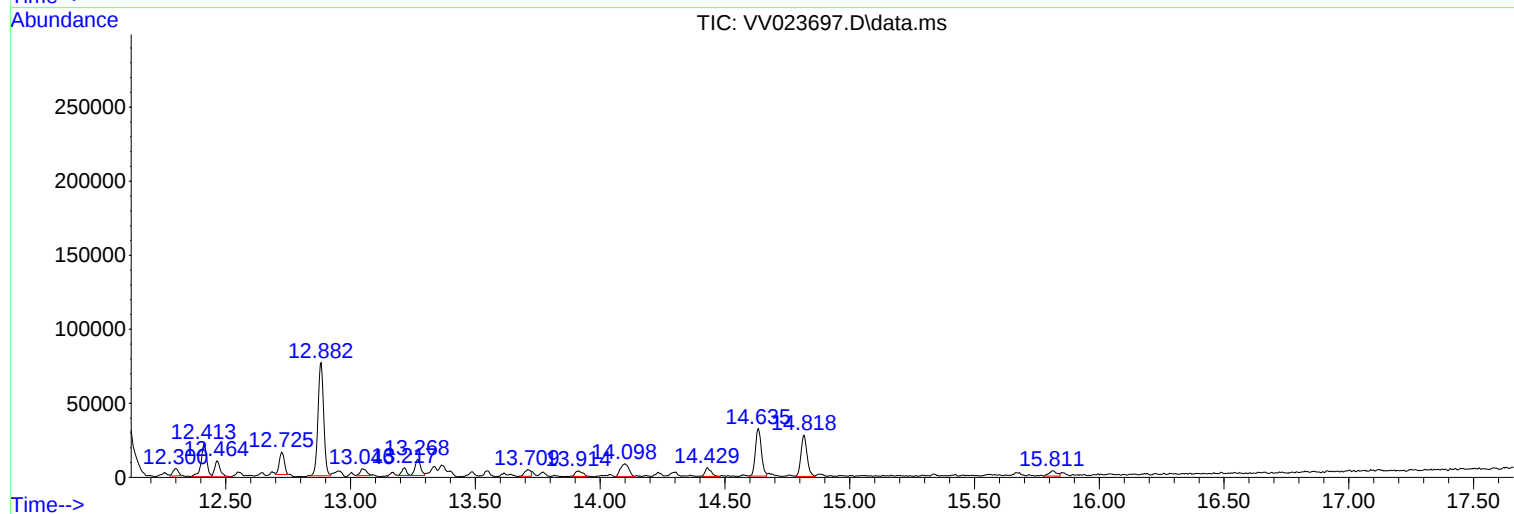
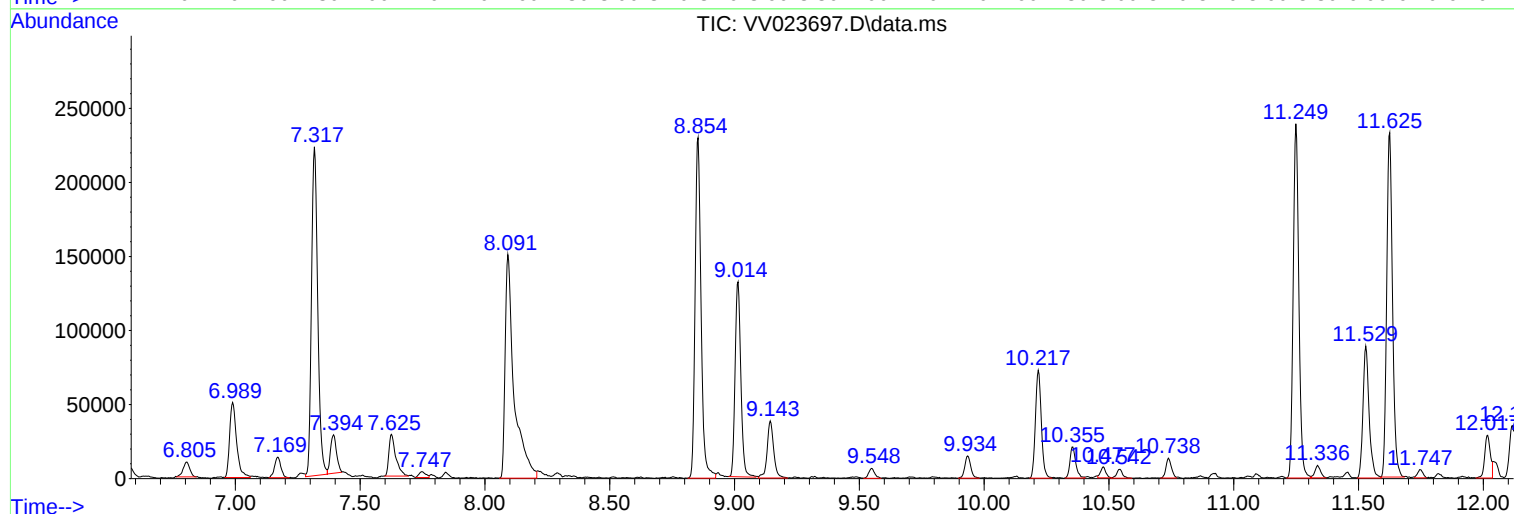
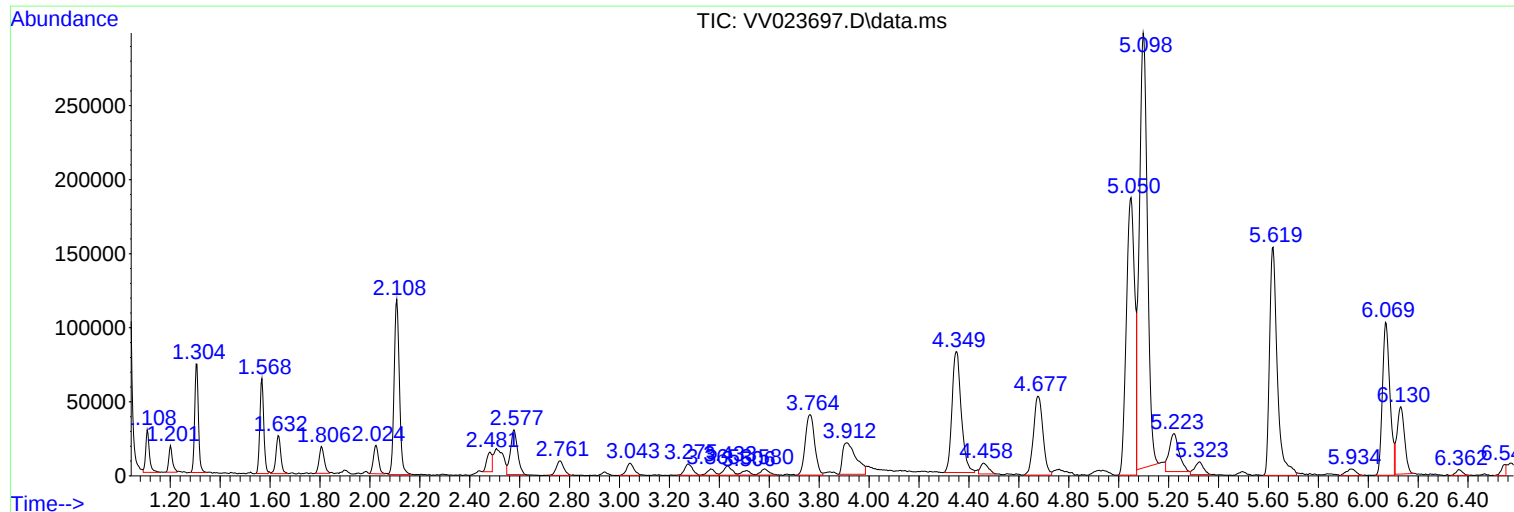
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Instrument :
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ClientSampleId :
C0G67DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.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_V/WATER
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Instrument :
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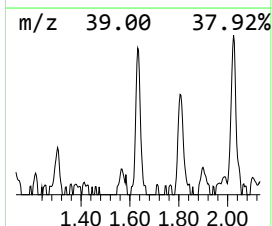
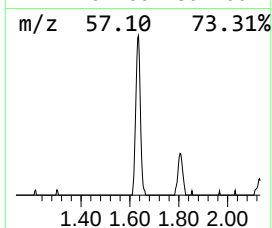
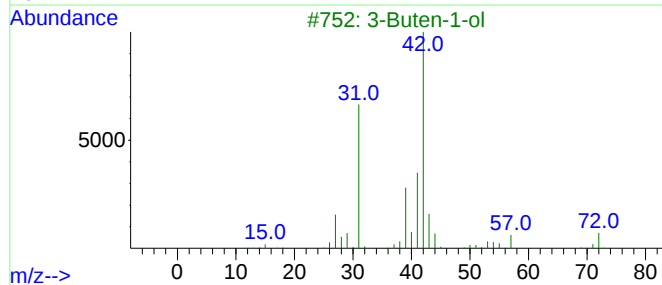
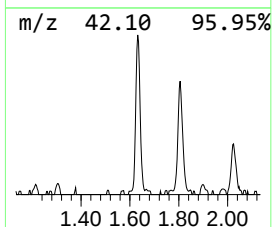
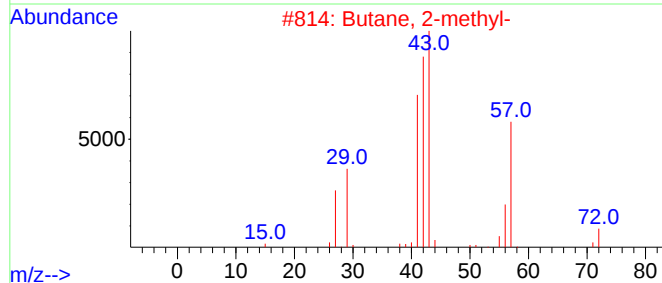
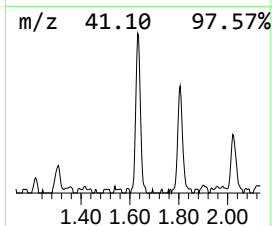
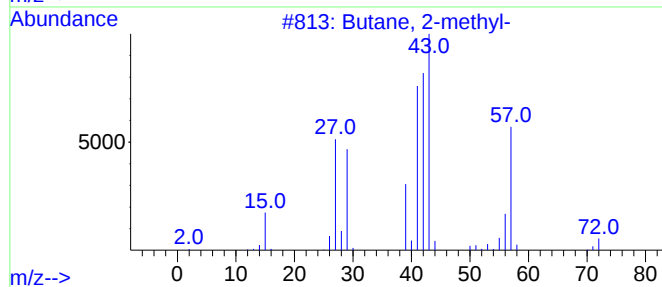
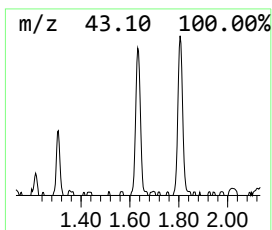
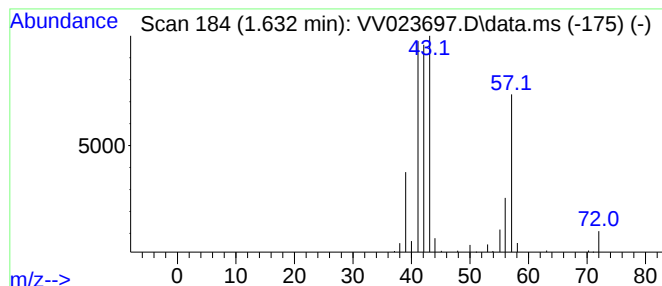
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.632	0.51 ug/L	34118	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	80
2		Butane, 2-methyl-	72	C5H12	000078-78-4	72
3		3-Buten-1-ol	72	C4H8O	000627-27-0	43
4		Pentane	72	C5H12	000109-66-0	28
5		Isobutane	58	C4H10	000075-28-5	25



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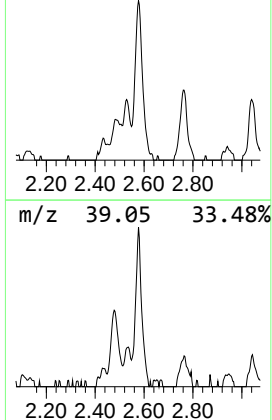
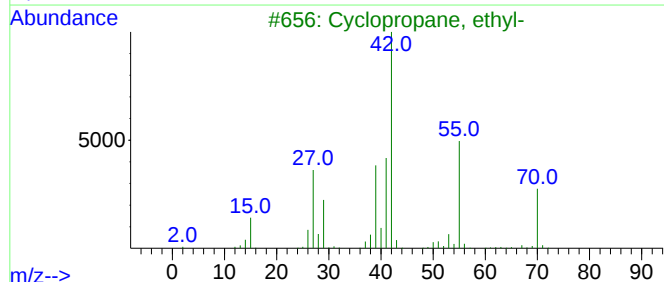
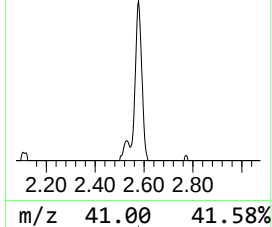
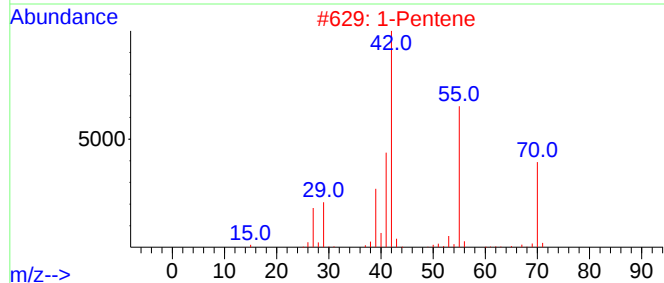
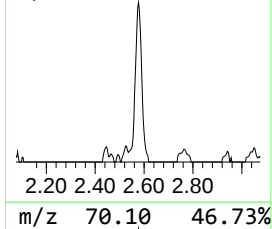
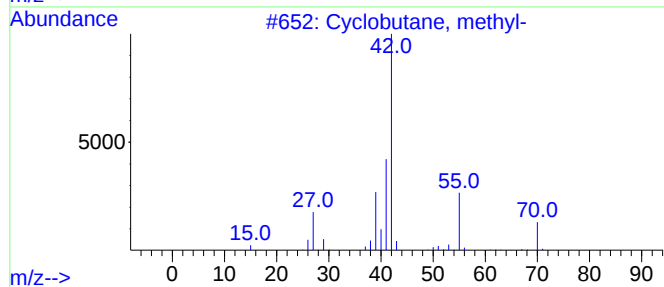
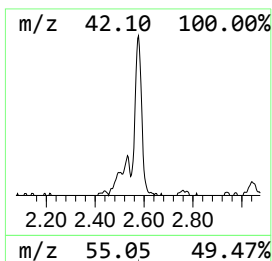
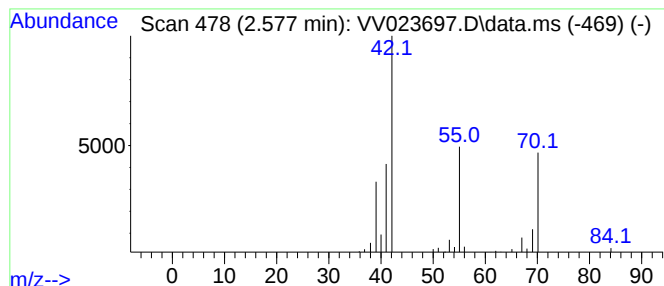
Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Cyclic2.577 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.577	0.89 ug/L	59317	1,4-Difluorobenzene	5.619	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
2	1-Pentene	70	C5H10	000109-67-1	72
3	Cyclopropane, ethyl-	70	C5H10	001191-96-4	59
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	53
5	1-Pentene	70	C5H10	000109-67-1	50



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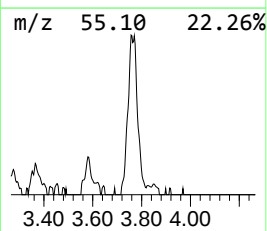
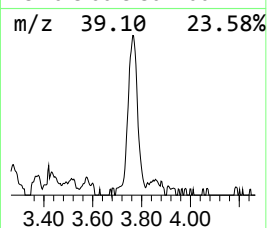
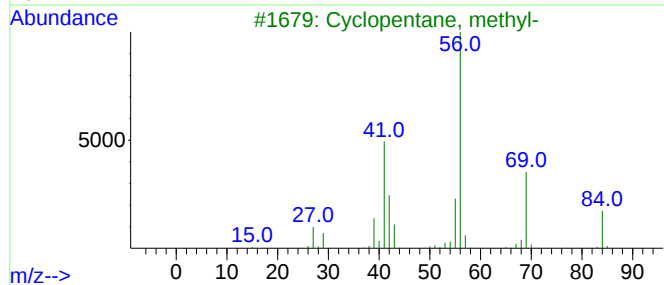
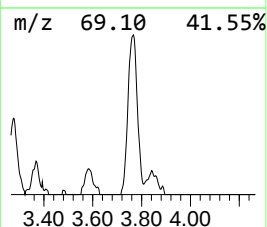
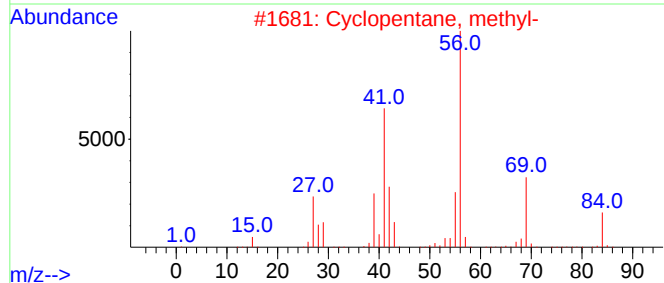
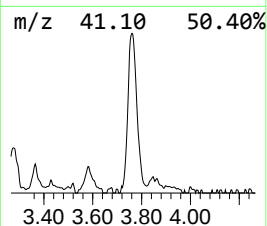
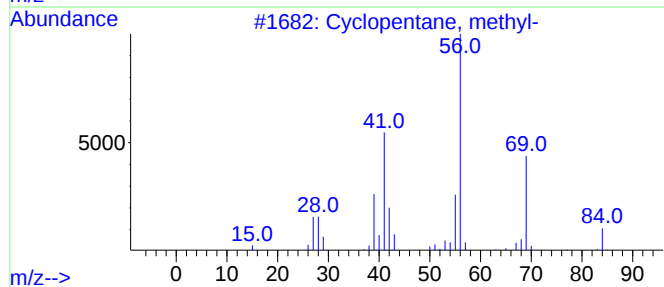
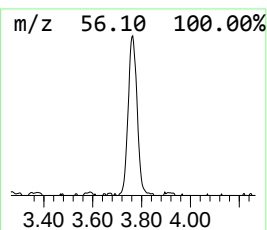
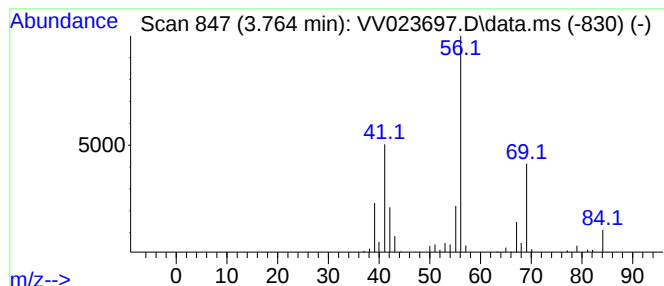
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Cyclic3.764 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.764	1.60 ug/L	106107	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2			Cyclopentane, methyl-	84	C6H12	000096-37-7	83
3			Cyclopentane, methyl-	84	C6H12	000096-37-7	72
4			Cyclopentane, methyl-	84	C6H12	000096-37-7	72
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72



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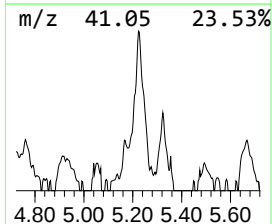
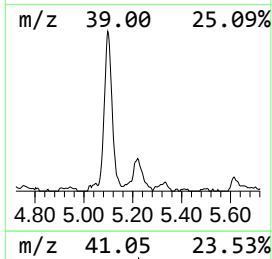
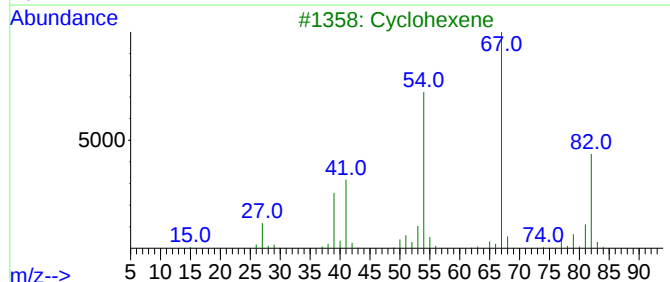
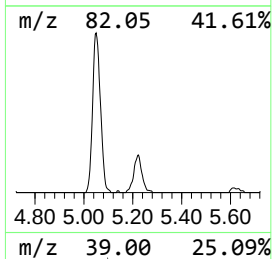
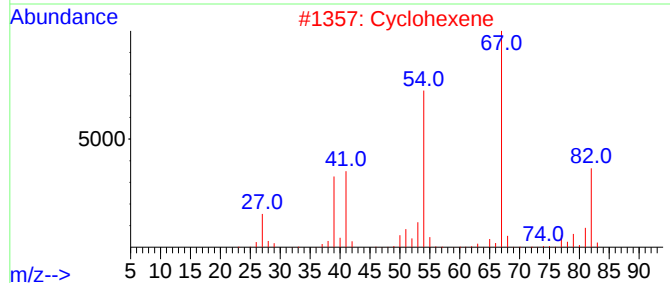
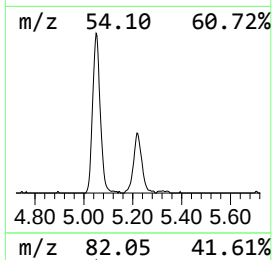
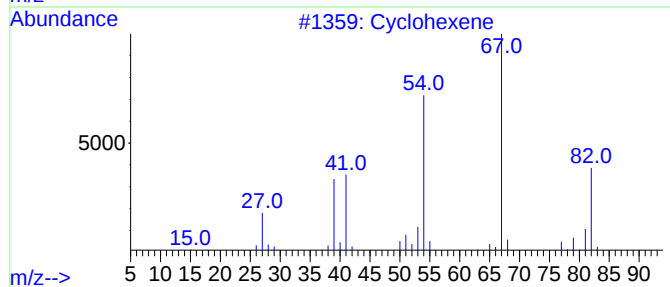
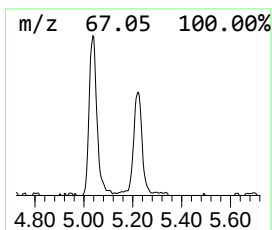
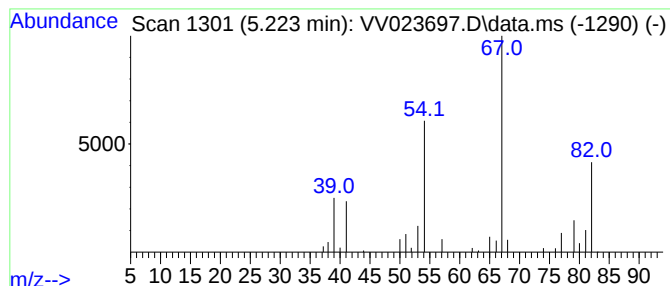
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Cyclobutane, ethenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.223	1.05 ug/L	69516	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene	82	C6H10	000110-83-8	80
2			Cyclohexene	82	C6H10	000110-83-8	80
3			Cyclohexene	82	C6H10	000110-83-8	80
4			Cyclohexene	82	C6H10	000110-83-8	80
5			Cyclobutane, ethenyl-	82	C6H10	002597-49-1	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
Data File : VV023697.D
Acq On : 24 Nov 2021 13:53
Operator : SY/MD
Sample : M4723-20DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G67DL

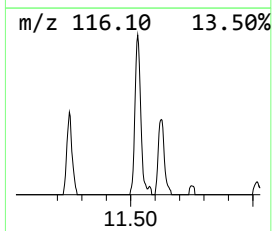
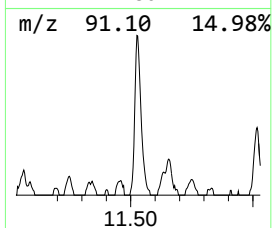
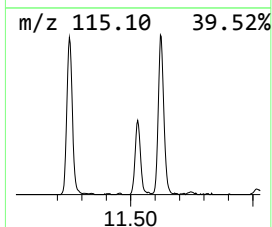
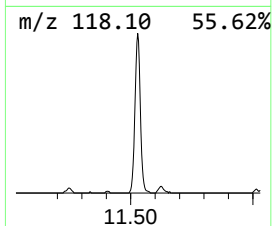
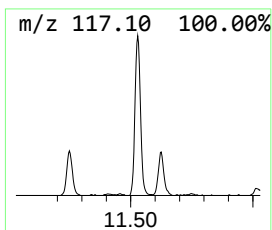
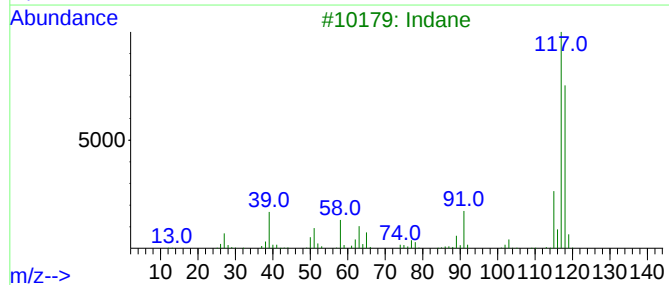
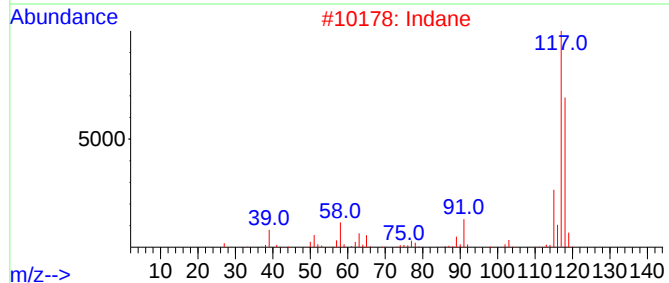
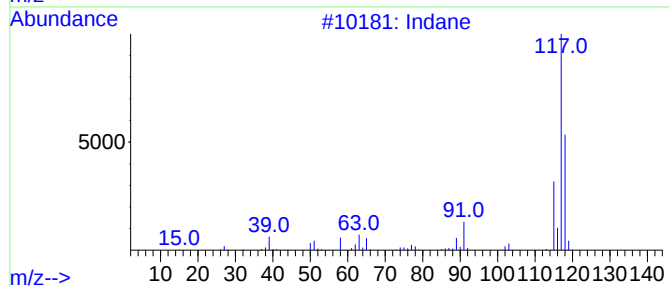
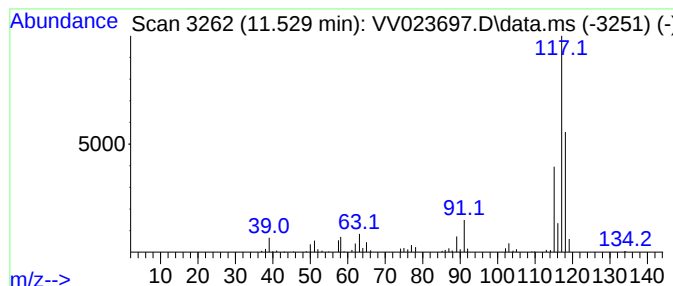
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	2.01 ug/L	151570	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	95
2			Indane	118	C9H10	000496-11-7	93
3			Indane	118	C9H10	000496-11-7	90
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
Data File : VV023697.D
Acq On : 24 Nov 2021 13:53
Operator : SY/MD
Sample : M4723-20DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G67DL

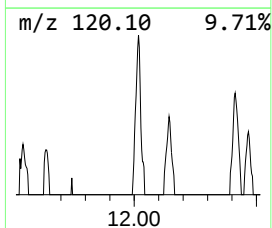
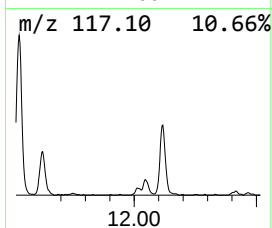
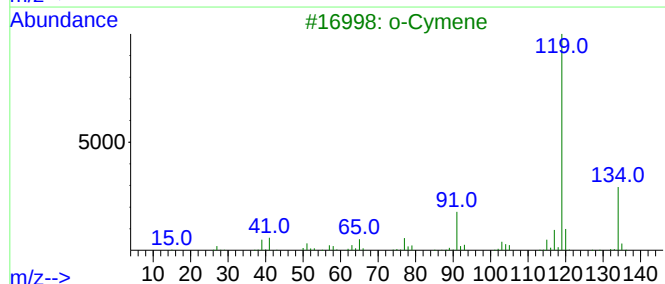
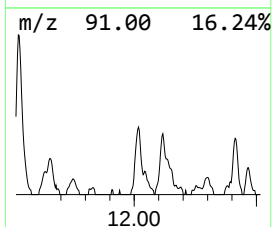
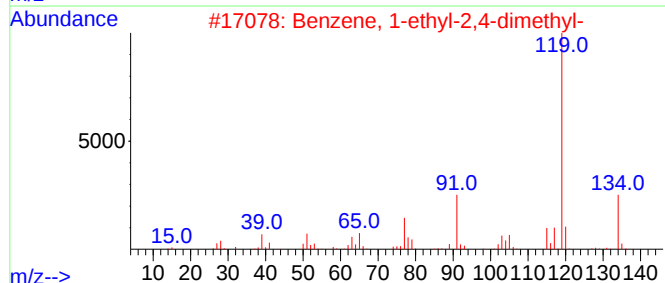
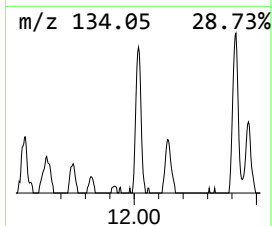
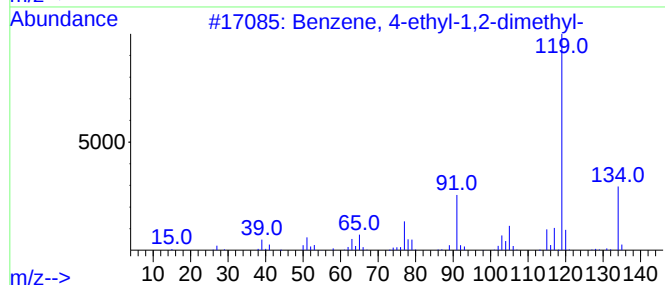
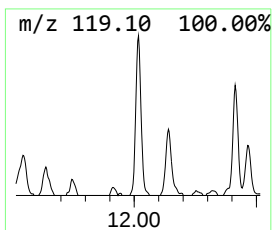
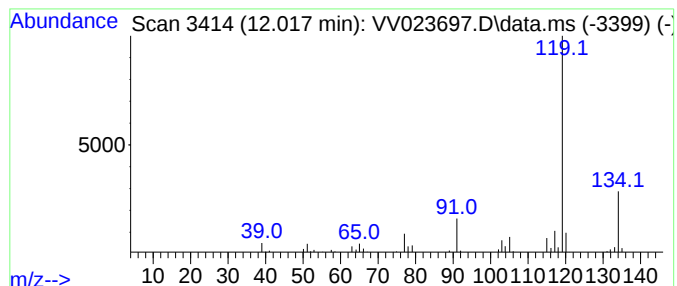
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.017	0.62 ug/L	46370	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3			o-Cymene	134	C10H14	000527-84-4	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\VV112421\
Data File : VV023697.D
Acq On : 24 Nov 2021 13:53
Operator : SY/MD
Sample : M4723-20DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G67DL

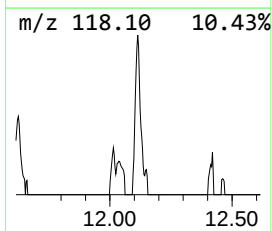
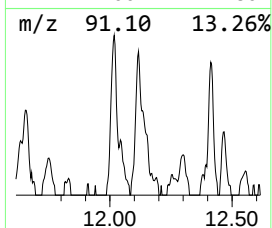
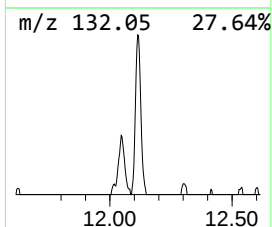
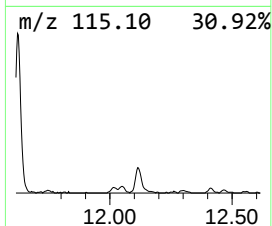
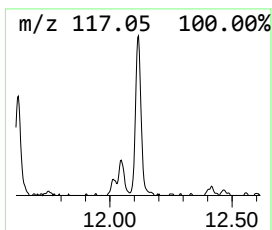
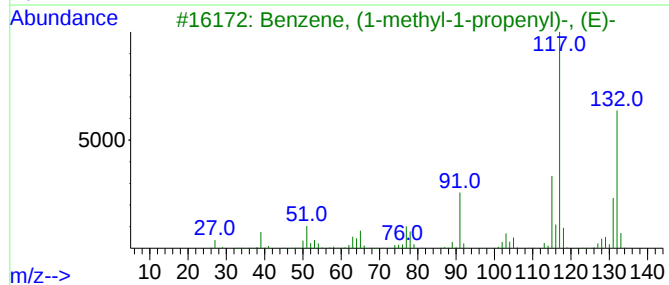
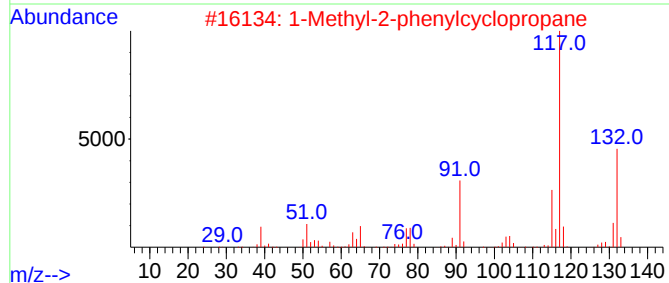
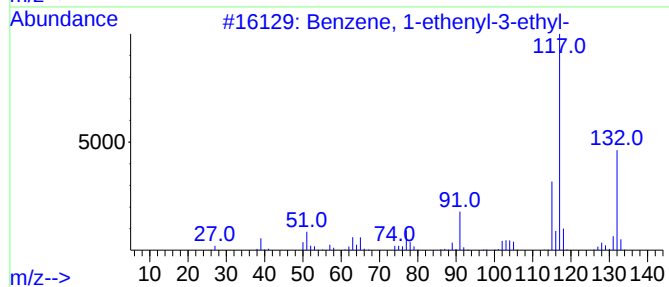
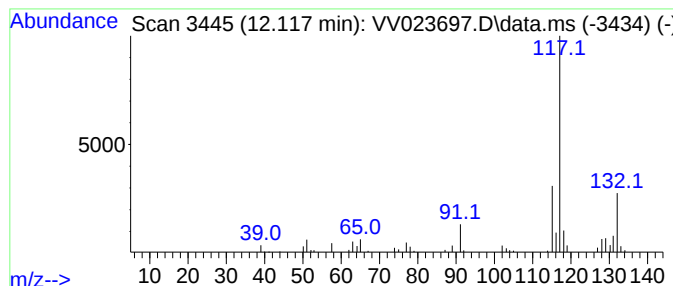
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	1.00 ug/L	75522	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
Data File : VV023697.D
Acq On : 24 Nov 2021 13:53
Operator : SY/MD
Sample : M4723-20DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G67DL

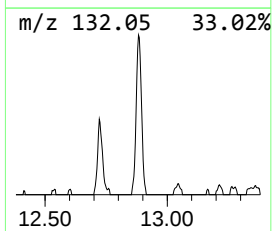
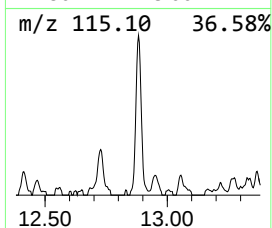
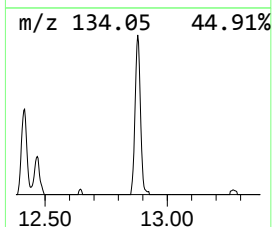
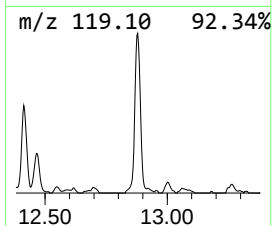
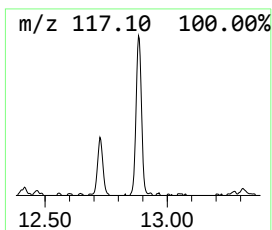
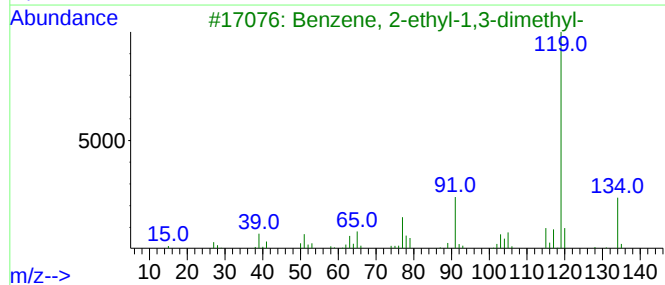
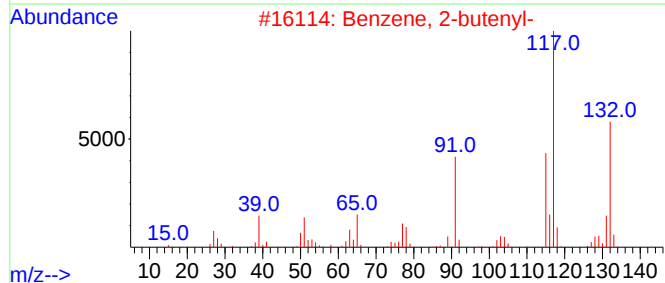
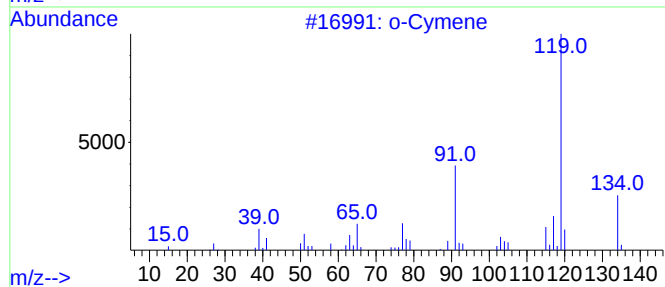
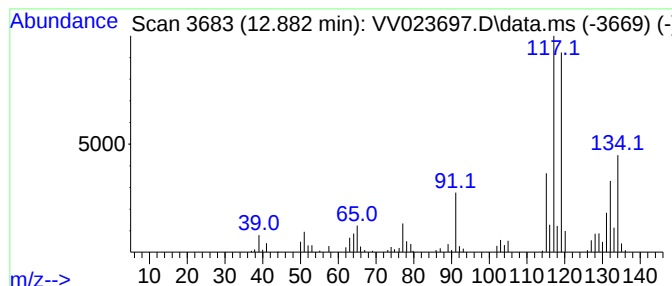
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 o-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.882	1.62 ug/L	121595	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	50
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	50
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	50
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	50
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
Data File : VV023697.D
Acq On : 24 Nov 2021 13:53
Operator : SY/MD
Sample : M4723-20DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G67DL

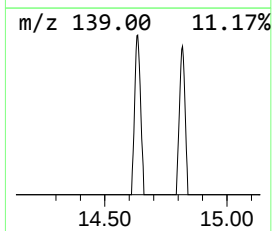
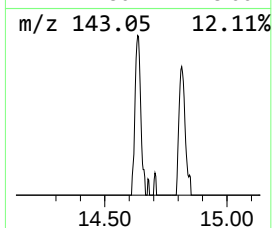
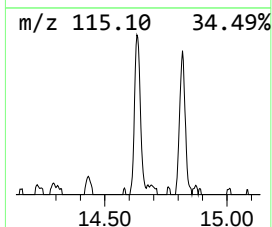
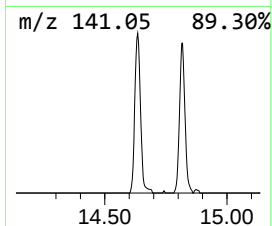
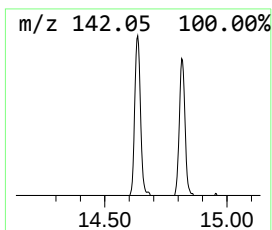
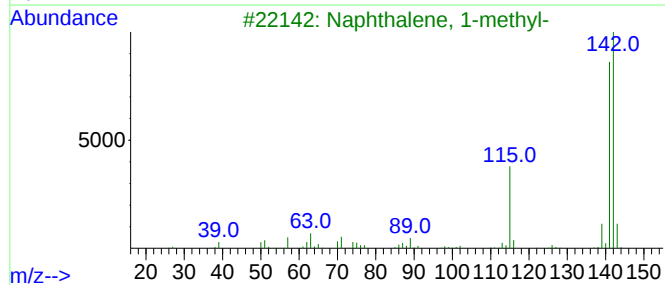
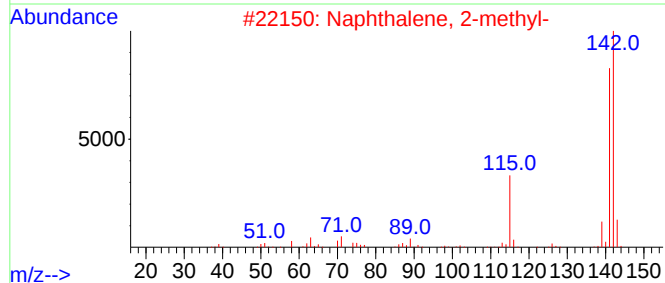
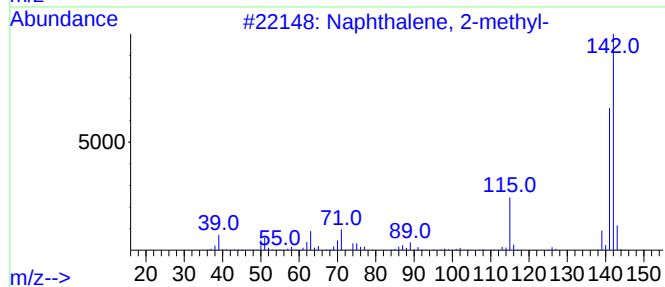
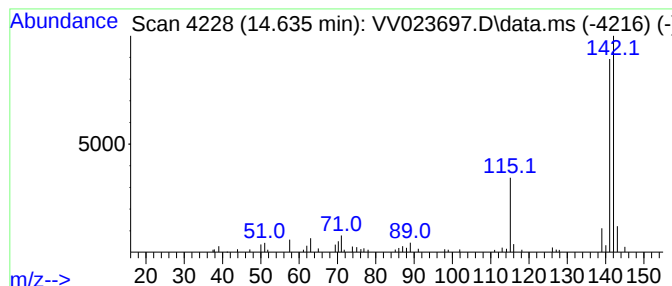
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.635	0.70 ug/L	52340	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
Data File : VV023697.D
Acq On : 24 Nov 2021 13:53
Operator : SY/MD
Sample : M4723-20DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G67DL

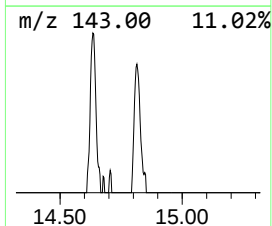
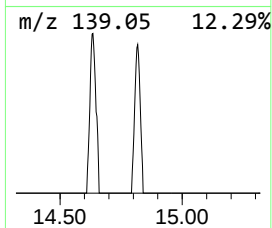
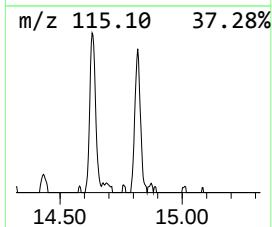
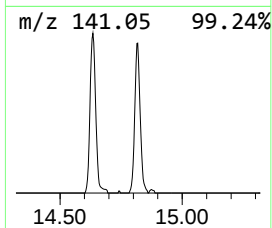
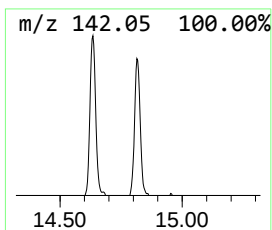
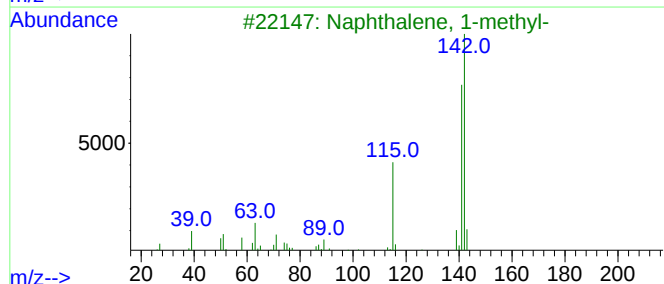
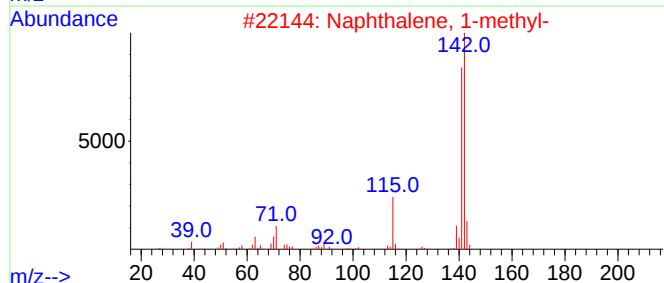
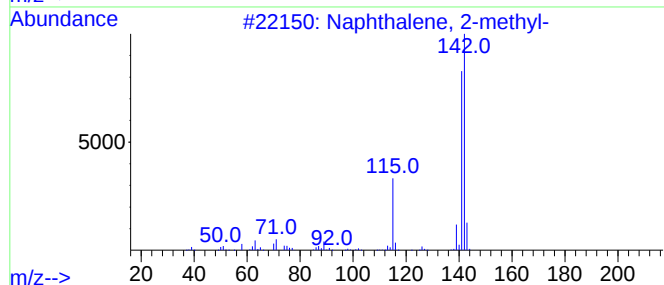
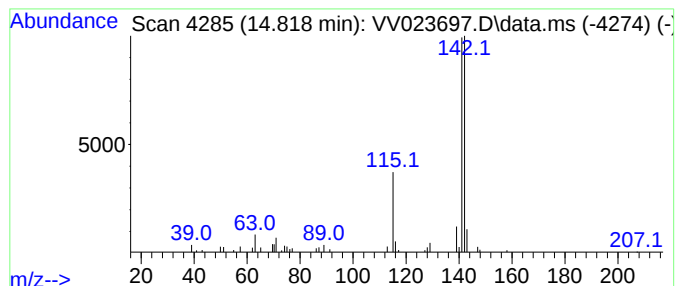
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.818	0.62 ug/L	46989	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5			Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV112421\
 Data File : VV023697.D
 Acq On : 24 Nov 2021 13:53
 Operator : SY/MD
 Sample : M4723-20DL 40X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0G67DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.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
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(DEL) Alkane: C...	2.577	0.9	ug/L	59317	1	5.619	331950	5.0
(DEL) Alkane: C...	3.764	1.6	ug/L	106107	1	5.619	331950	5.0
Cyclobutane, et...	5.223	1.1	ug/L	69516	1	5.619	331950	5.0
Indane	11.529	2.0	ug/L	151570	3	11.249	376144	5.0
Benzene, 4-ethy...	12.017	0.6	ug/L	46370	3	11.249	376144	5.0
Benzene, 1-ethe...	12.117	1.0	ug/L	75522	3	11.249	376144	5.0
o-Cymene	12.882	1.6	ug/L	121595	3	11.249	376144	5.0
Naphthalene, 2-...	14.635	0.7	ug/L	52340	3	11.249	376144	5.0
Naphthalene, 1-...	14.818	0.6	ug/L	46989	3	11.249	376144	5.0