

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
 Data File : VV025776.D
 Acq On : 02 May 2022 19:15
 Operator : SY/MD
 Sample : N2625-09
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EW6Y1

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\SFAMVTR040822WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV025776.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.111	15	22	38	rBV	1272381	1181140	95.57%	13.780%
2	1.182	39	44	49	rVV	21056	18846	1.52%	0.220%
3	1.217	49	55	69	rVV	156040	140626	11.38%	1.641%
4	1.301	73	81	95	rVV2	866922	1235936	100.00%	14.419%
5	1.359	95	99	104	rVV	25990	26345	2.13%	0.307%
6	1.391	104	109	112	rVV	22389	26068	2.11%	0.304%
7	1.420	112	118	129	rVB	124871	140682	11.38%	1.641%
8	1.568	152	164	176	rVB2	118305	163438	13.22%	1.907%
9	1.635	177	185	198	rBV	160978	203377	16.46%	2.373%
10	1.770	217	227	233	rBV	233726	290704	23.52%	3.391%
11	1.812	233	240	259	rVV3	208799	388757	31.45%	4.535%
12	1.902	261	268	278	rVB	21622	28539	2.31%	0.333%
13	1.982	283	293	301	rBV	83177	120641	9.76%	1.407%
14	2.027	301	307	316	rVB2	18242	24681	2.00%	0.288%
15	2.111	322	333	344	rBV	137968	204851	16.57%	2.390%
16	2.352	400	408	420	rVB2	11078	18381	1.49%	0.214%
17	2.449	420	438	446	rBV4	23379	55985	4.53%	0.653%
18	2.532	456	464	472	rVV2	14547	24395	1.97%	0.285%
19	2.577	473	478	498	rVB5	10867	22611	1.83%	0.264%
20	2.767	524	537	554	rVB3	21558	43783	3.54%	0.511%
21	2.953	581	595	611	rVB5	49347	109365	8.85%	1.276%
22	3.040	613	622	642	rVB4	22289	47843	3.87%	0.558%
23	3.153	645	657	666	rBV9	8334	18373	1.49%	0.214%
24	3.458	738	752	765	rBV7	12320	30438	2.46%	0.355%
25	3.902	877	890	932	rBV2	47354	189690	15.35%	2.213%
26	4.352	1013	1030	1052	rBV2	99090	257251	20.81%	3.001%
27	4.912	1192	1204	1221	rVB6	7796	18226	1.47%	0.213%
28	5.050	1229	1247	1280	rBV3	206153	545719	44.15%	6.367%
29	5.326	1322	1333	1341	rBV6	6039	13008	1.05%	0.152%
30	5.381	1342	1350	1373	rVB6	11969	31243	2.53%	0.364%
31	5.493	1377	1385	1397	rBV8	6426	13795	1.12%	0.161%
32	5.619	1410	1424	1454	rBV	160708	351773	28.46%	4.104%
33	6.069	1550	1564	1594	rBV	124103	273561	22.13%	3.191%
34	6.992	1835	1851	1871	rBV2	35557	73021	5.91%	0.852%
35	7.317	1941	1952	1968	rBV	216262	384057	31.07%	4.481%
36	7.394	1968	1976	1989	rVB	20994	39043	3.16%	0.455%

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37	7.628	2041	2049	2067	rBV2	17992	38000	3.07%	0.443%
38	8.088	2183	2192	2222	rBV	227621	454558	36.78%	5.303%
39	8.853	2418	2430	2449	rBV	256572	421144	34.07%	4.913%
40	9.017	2473	2481	2491	rBV2	9240	15494	1.25%	0.181%
41	9.143	2512	2520	2533	rBV2	11057	19023	1.54%	0.222%
42	10.217	2841	2854	2869	rBV	85676	139192	11.26%	1.624%
43	11.249	3165	3175	3189	rBV	217660	342078	27.68%	3.991%
44	11.529	3252	3262	3274	rBV2	25102	44822	3.63%	0.523%
45	11.622	3281	3291	3311	rVB	207764	324169	26.23%	3.782%
46	12.345	3508	3516	3532	rVB7	10656	16984	1.37%	0.198%

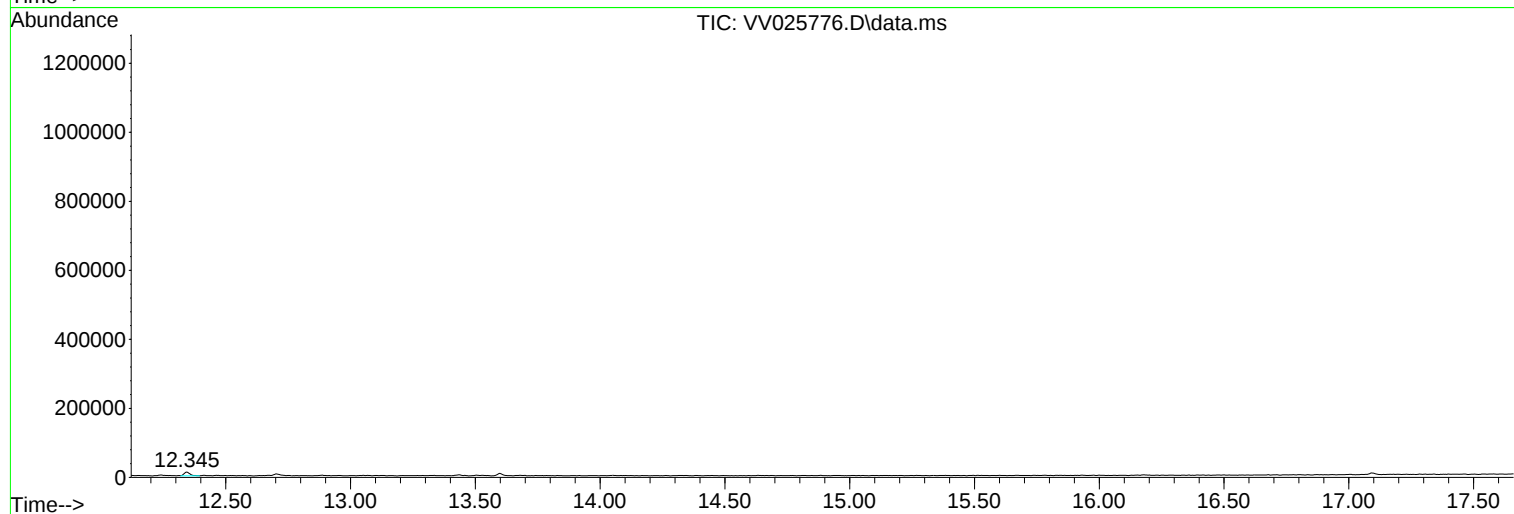
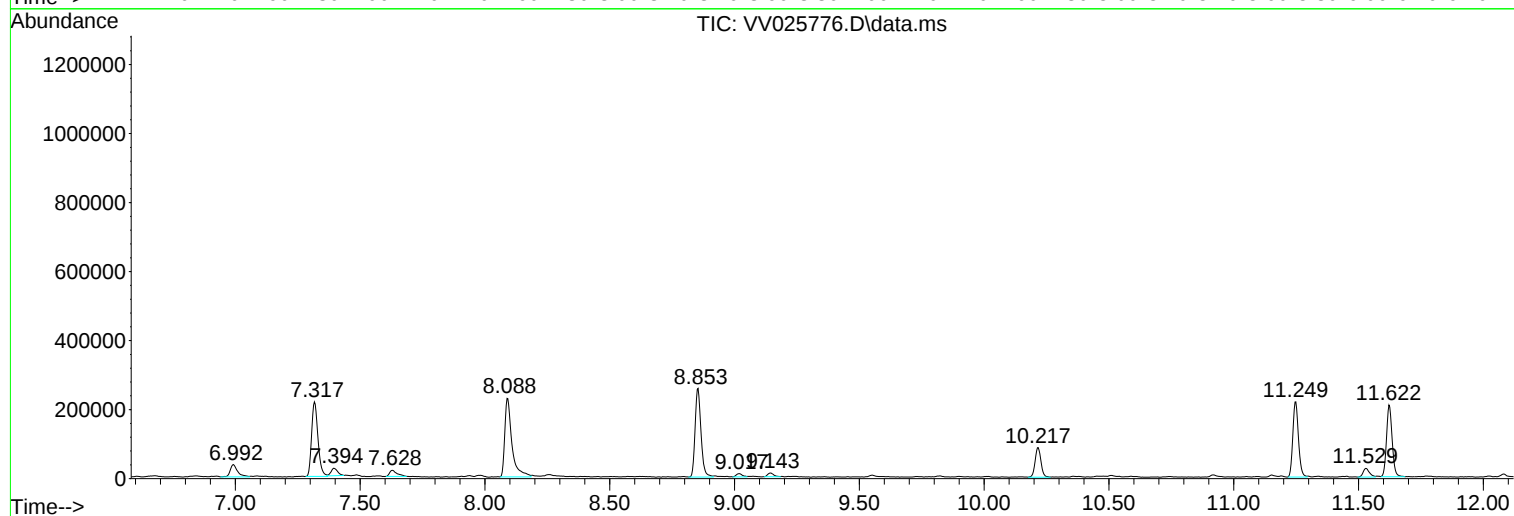
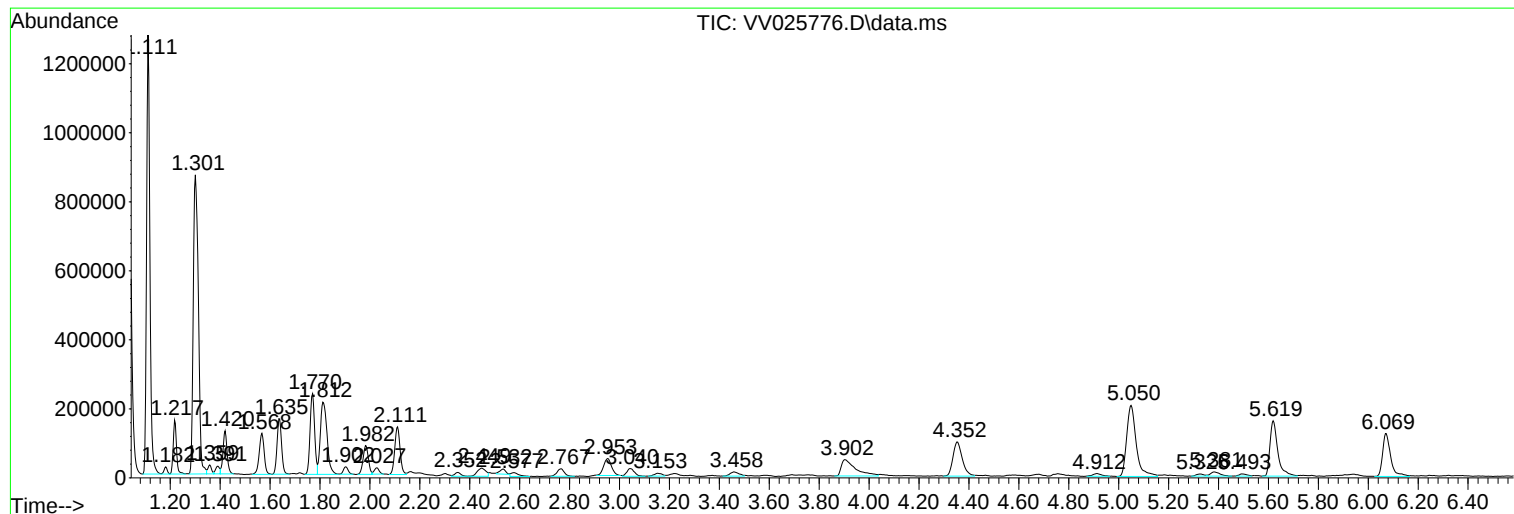
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Quant Title : TRACE VOA SFAM1.0

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



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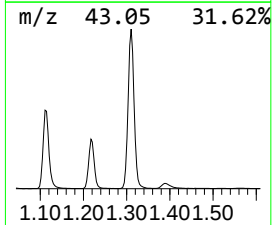
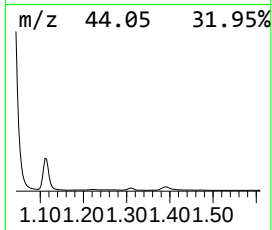
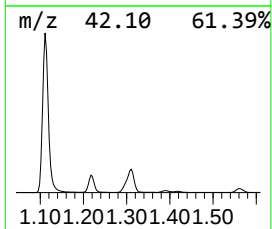
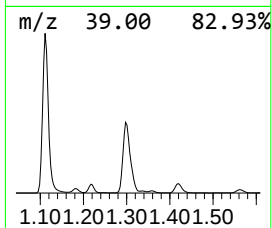
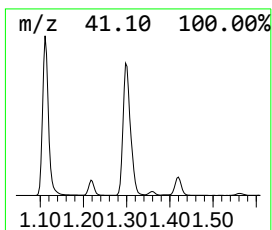
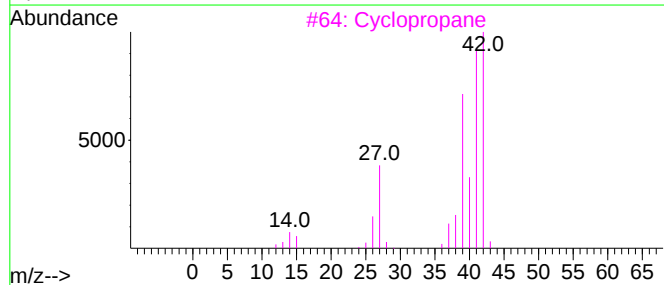
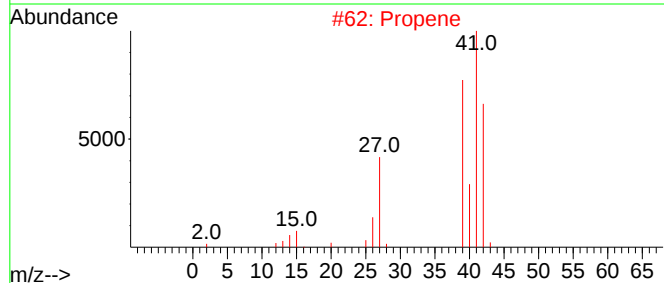
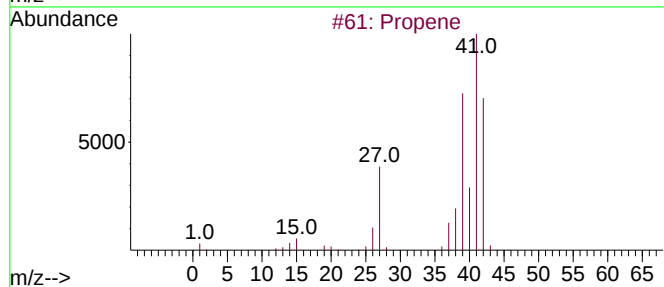
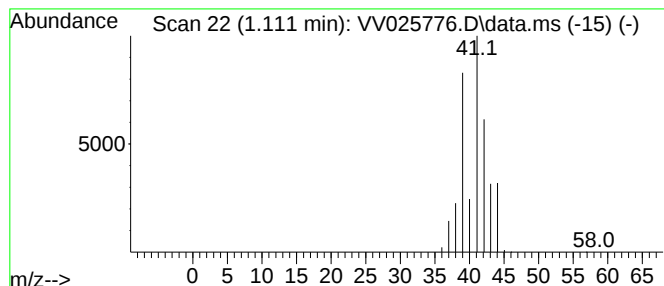
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Quant Title : TRACE VOA SFAM1.0

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.
1.111	16.79 ug/L	1181140	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	90
2			Propene	42	C3H6	000115-07-1	42
3			Cyclopropane	42	C3H6	000075-19-4	9
4			Cyclopropene	40	C3H4	002781-85-3	9
5			Cyclopropane	42	C3H6	000075-19-4	7



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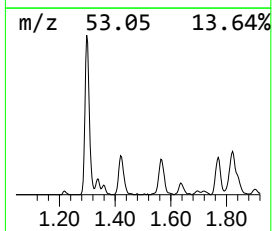
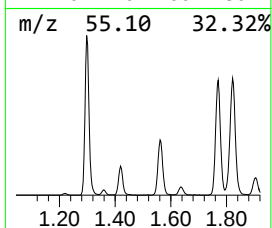
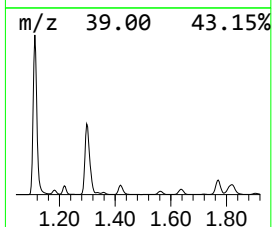
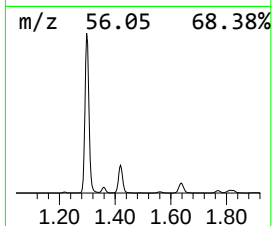
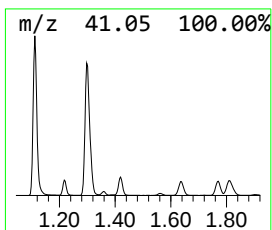
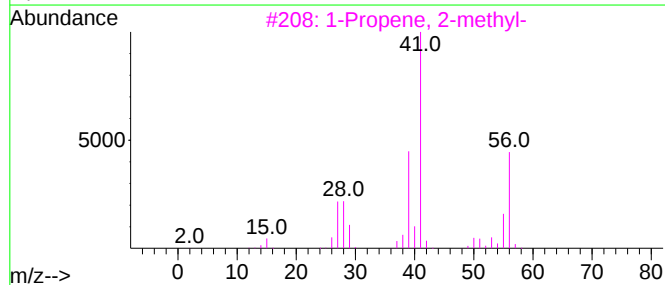
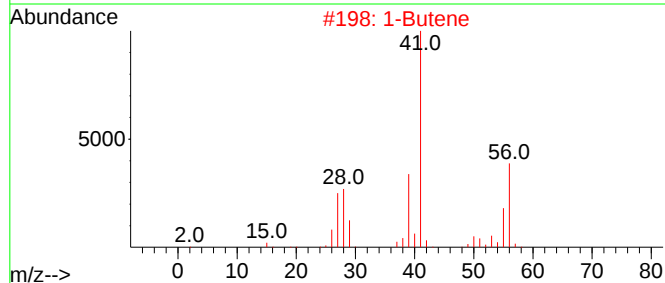
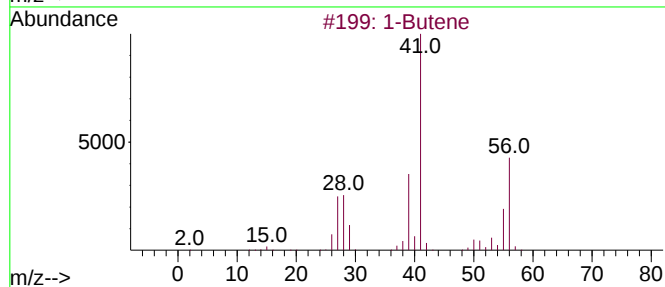
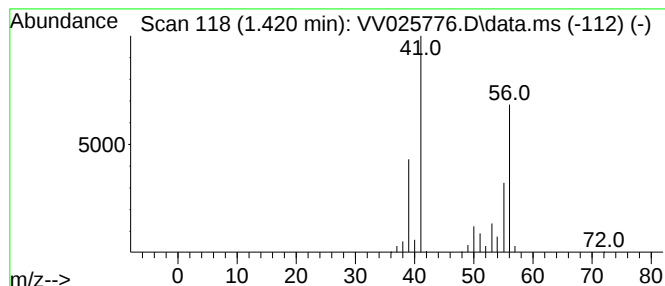
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.420	2.00 ug/L	140682	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Butene	56	C4H8	000106-98-9	83
2		1-Butene	56	C4H8	000106-98-9	80
3		1-Propene, 2-methyl-	56	C4H8	000115-11-7	72
4		1-Butene	56	C4H8	000106-98-9	72
5		2-Butene, (Z)-	56	C4H8	000590-18-1	68



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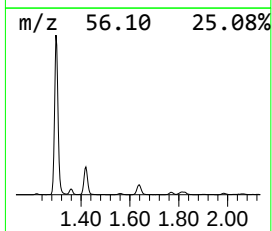
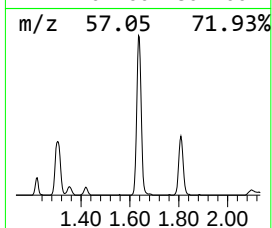
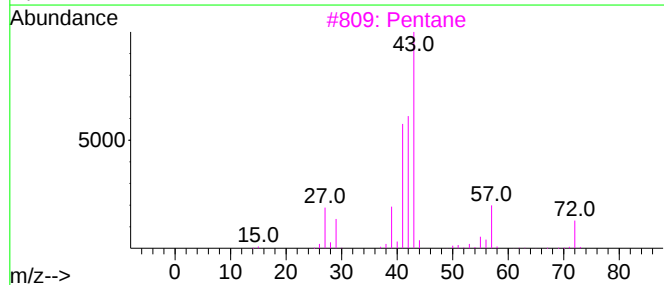
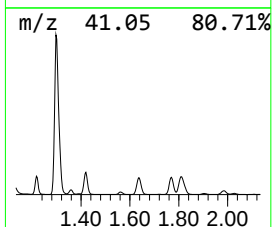
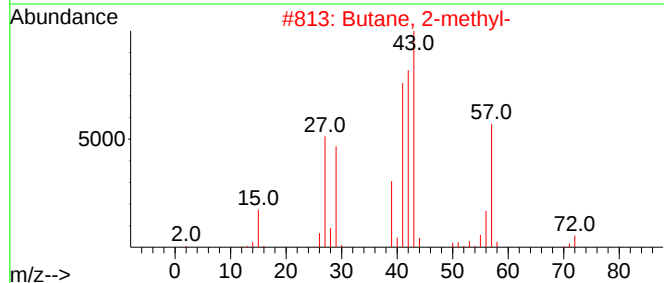
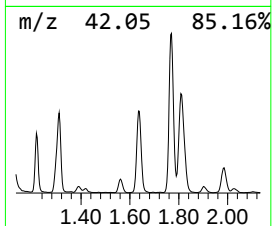
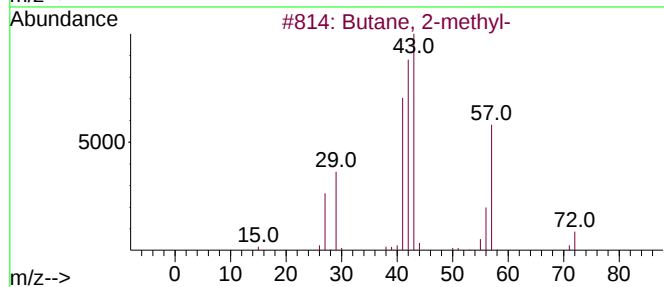
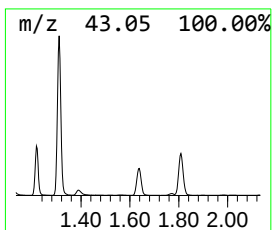
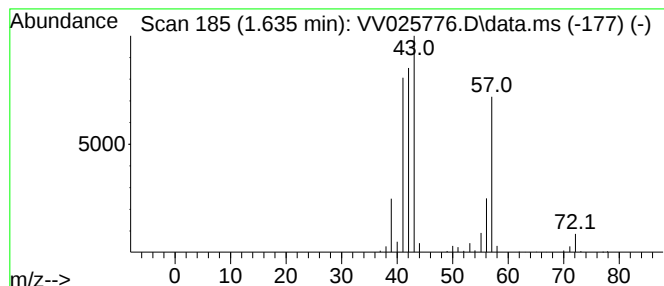
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.635	2.89 ug/L	203377	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-			72	C5H12	000078-78-4	91
2	Butane, 2-methyl-			72	C5H12	000078-78-4	91
3	Pentane			72	C5H12	000109-66-0	37
4	3-Buten-1-ol			72	C4H8O	000627-27-0	28
5	1-Pentene			70	C5H10	000109-67-1	25



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Instrument :
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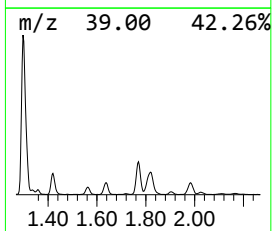
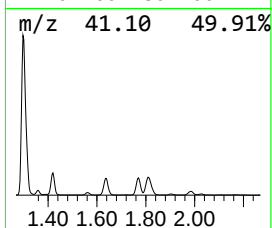
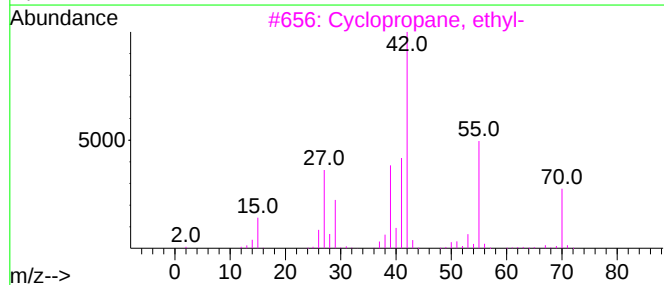
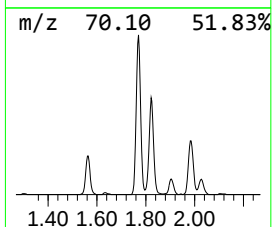
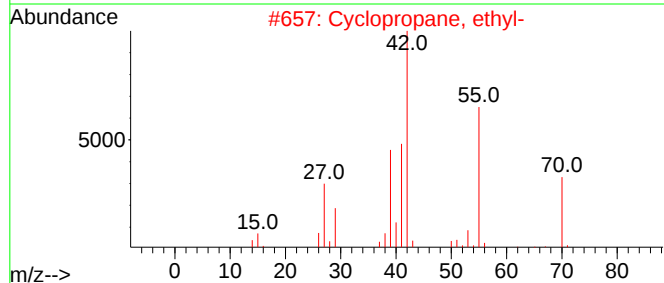
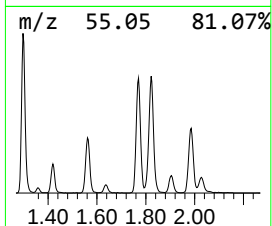
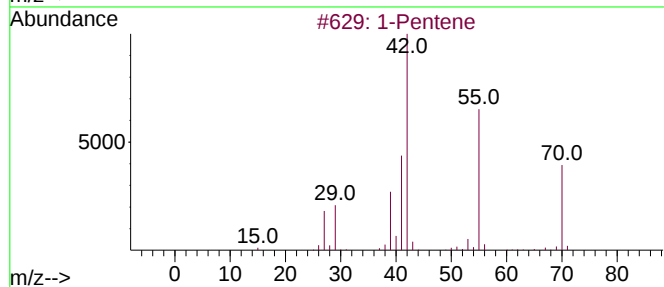
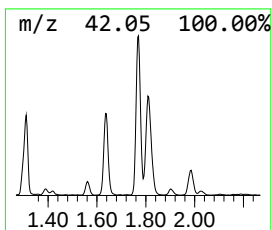
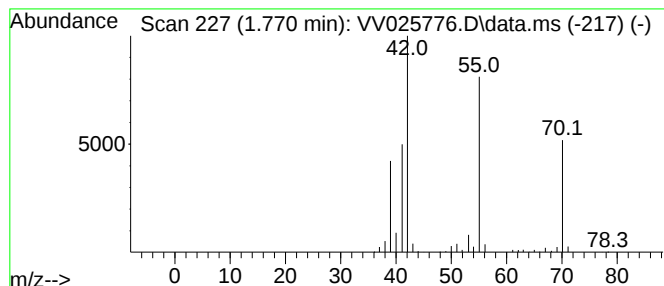
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.770	4.13 ug/L	290704	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene	70	C5H10	000109-67-1	81
2		Cyclopropane, ethyl-	70	C5H10	001191-96-4	78
3		Cyclopropane, ethyl-	70	C5H10	001191-96-4	78
4		2-Pentene, (Z)-	70	C5H10	000627-20-3	74
5		1-Pentene	70	C5H10	000109-67-1	68



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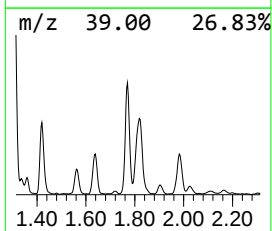
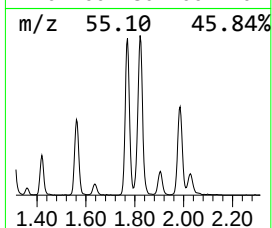
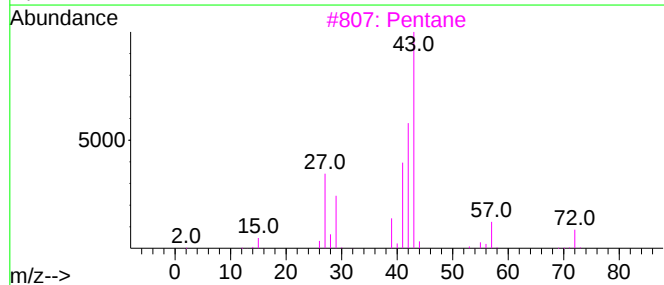
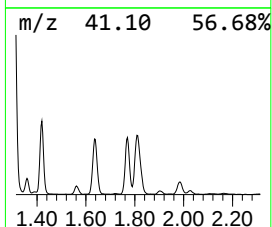
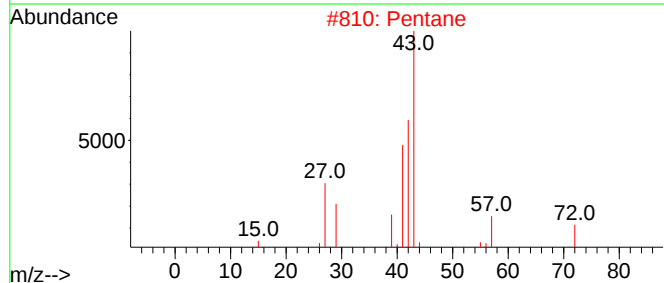
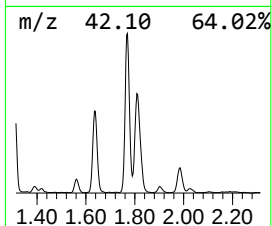
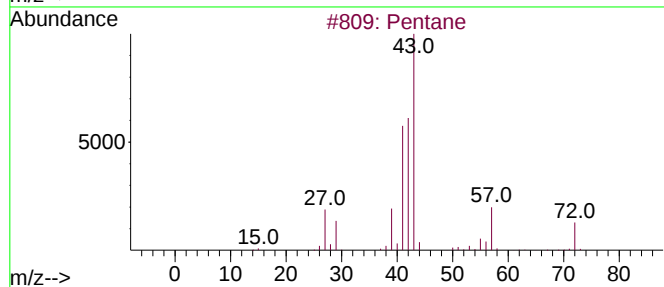
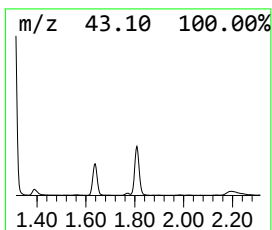
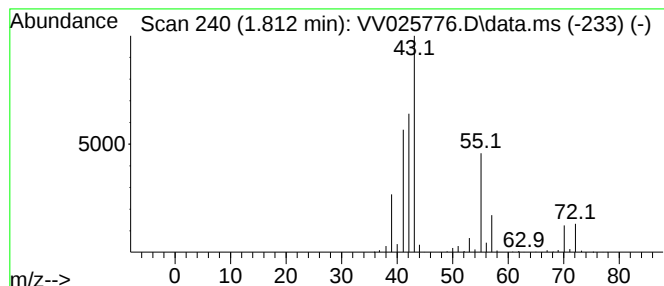
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.812	5.53 ug/L	388757	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane	72	C5H12	000109-66-0	72
2		Pentane	72	C5H12	000109-66-0	72
3		Pentane	72	C5H12	000109-66-0	72
4		Pentane	72	C5H12	000109-66-0	72
5		Pentane	72	C5H12	000109-66-0	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

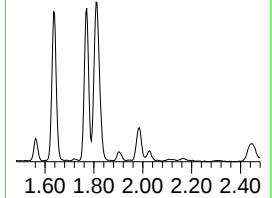
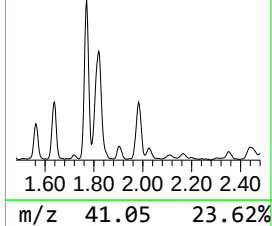
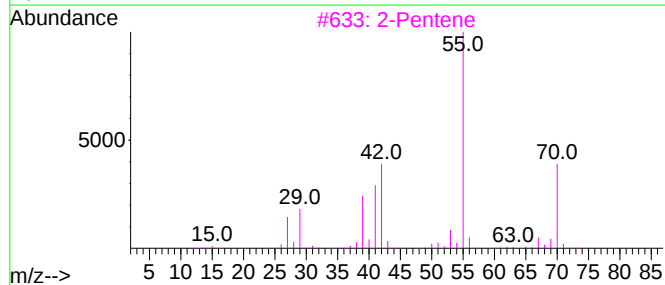
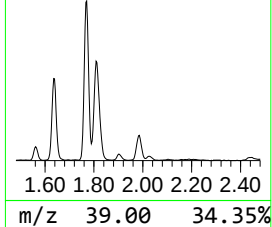
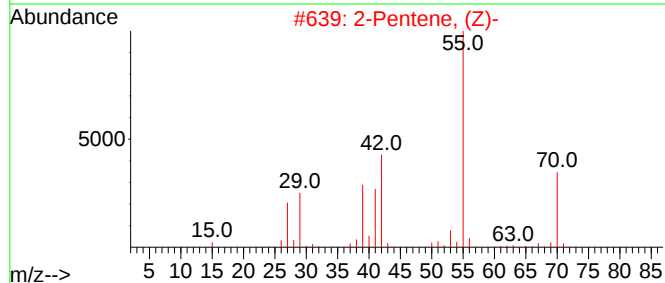
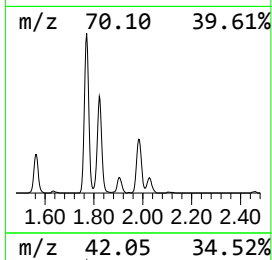
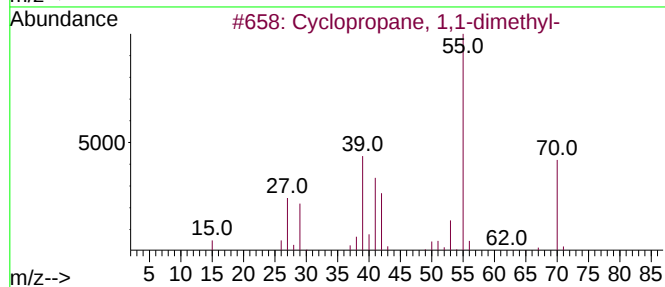
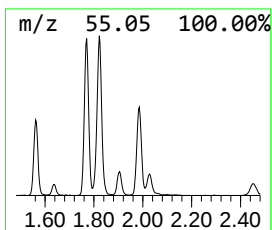
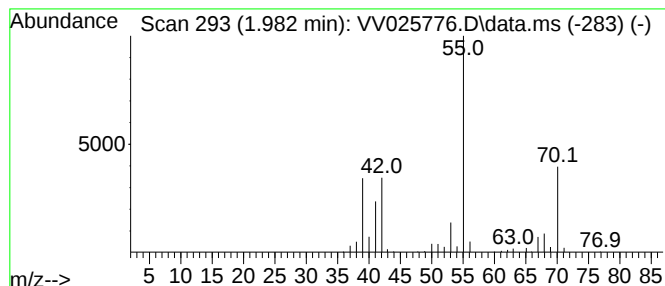
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Cyclic1.982 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.982	1.71 ug/L	120641	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	87
2			2-Pentene, (Z)-	70	C5H10	000627-20-3	87
3			2-Pentene	70	C5H10	000109-68-2	86
4			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	86
5			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

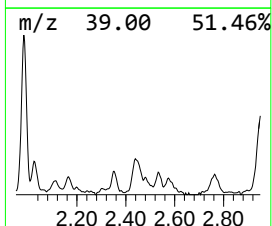
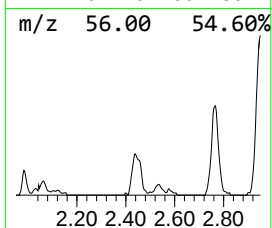
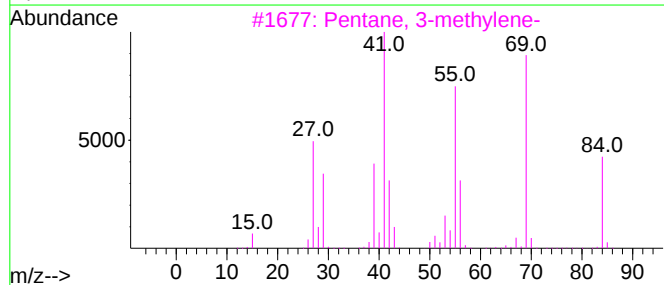
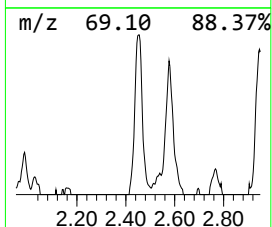
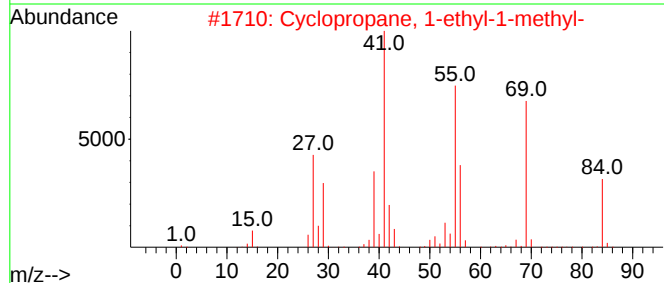
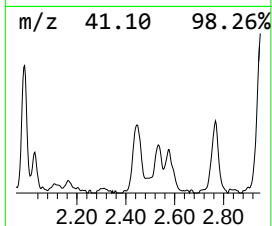
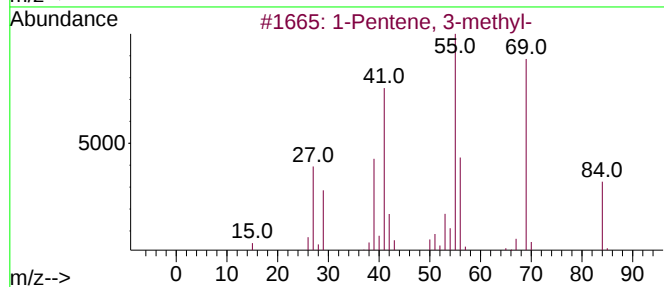
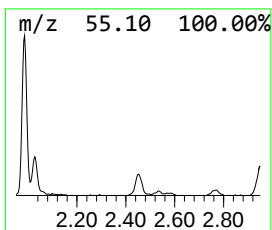
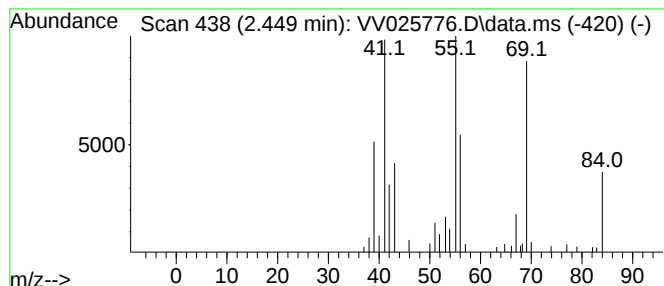
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.449	0.80 ug/L	55985	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	90
2		Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	76
3		Pentane, 3-methylene-	84	C6H12	000760-21-4	72
4		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	68
5		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

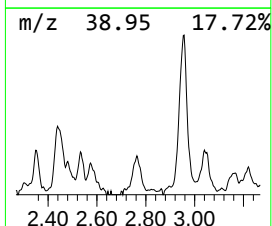
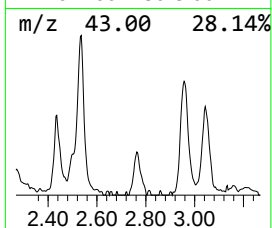
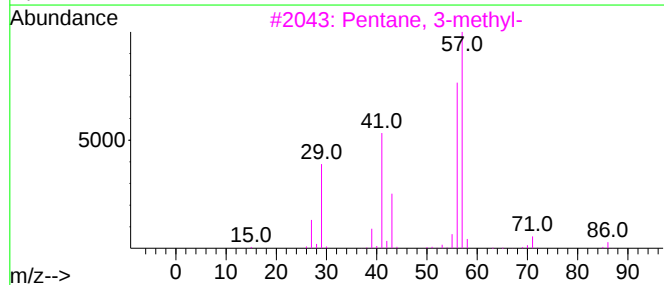
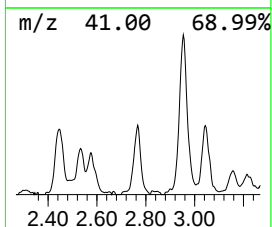
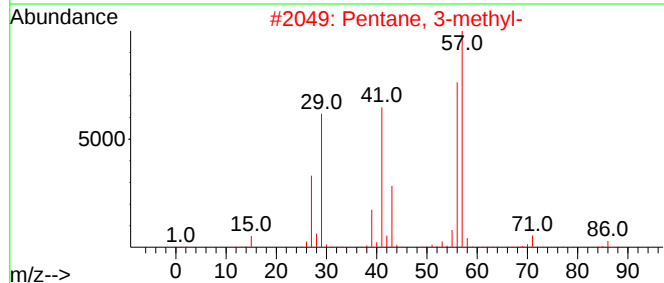
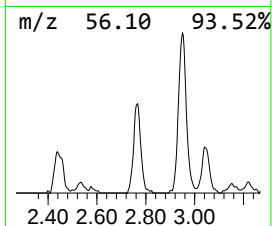
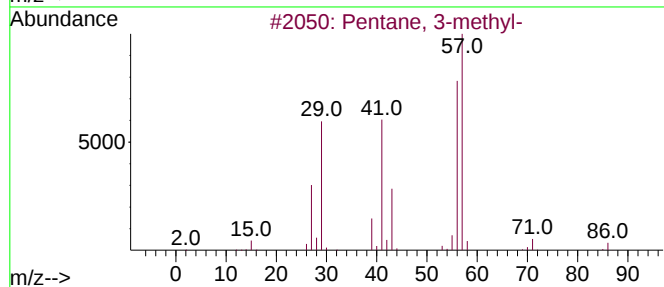
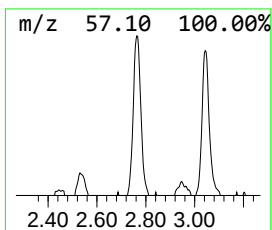
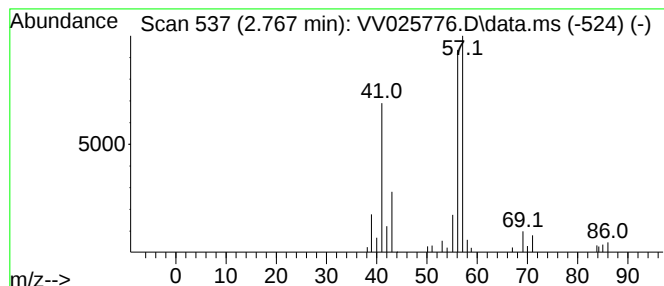
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.767	0.62 ug/L	43783	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
2			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
3			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
5			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

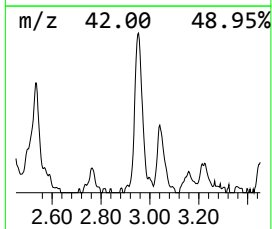
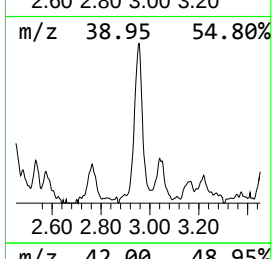
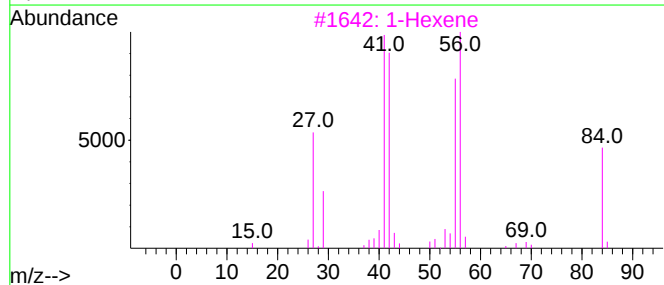
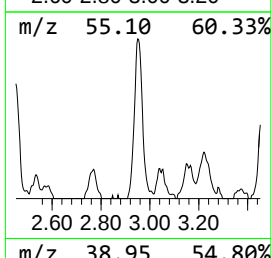
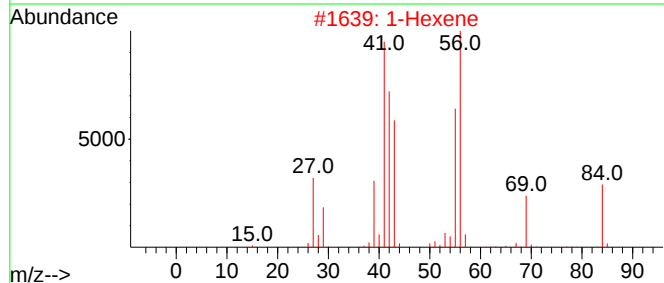
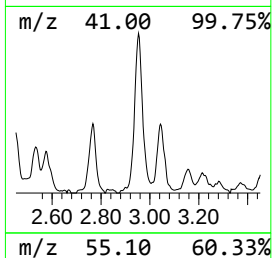
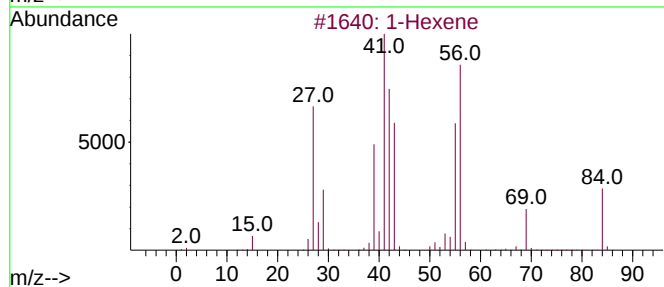
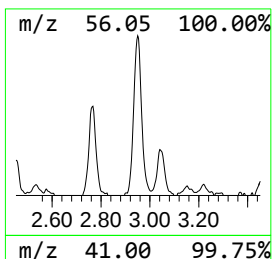
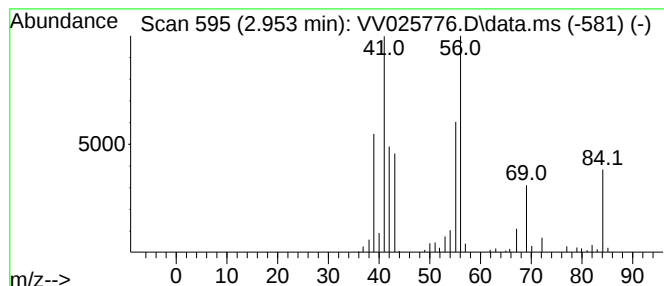
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.953	1.55 ug/L	109365	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene		84	C6H12	000592-41-6	76
2	1-Hexene		84	C6H12	000592-41-6	58
3	1-Hexene		84	C6H12	000592-41-6	58
4	1-Pentene, 2-methyl-		84	C6H12	000763-29-1	58
5	2-Hexene, (Z)-		84	C6H12	007688-21-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

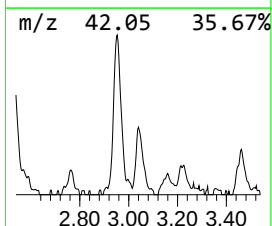
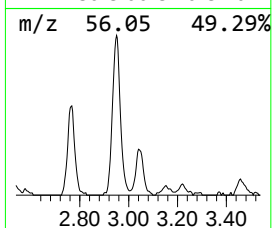
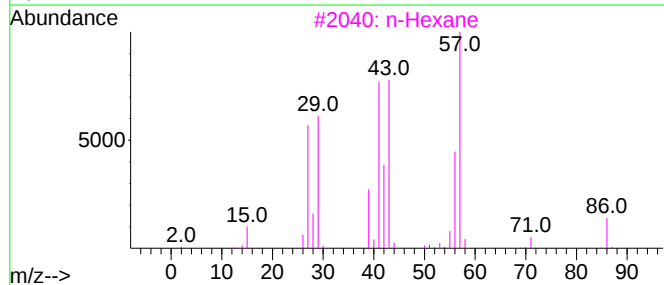
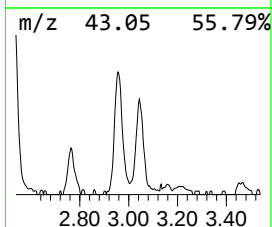
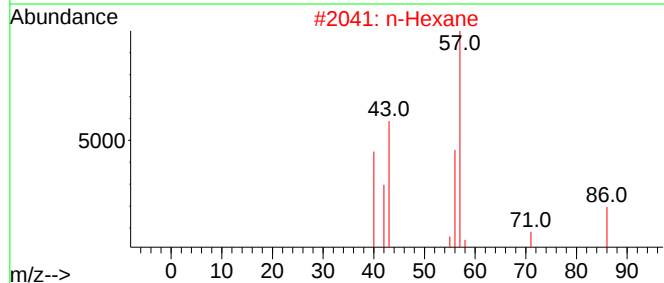
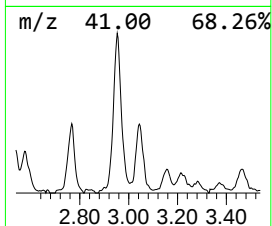
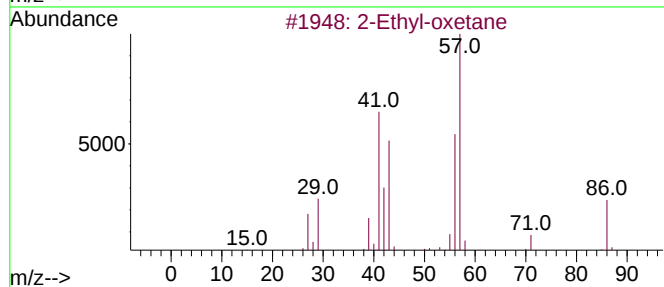
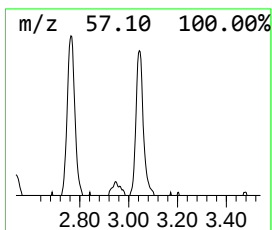
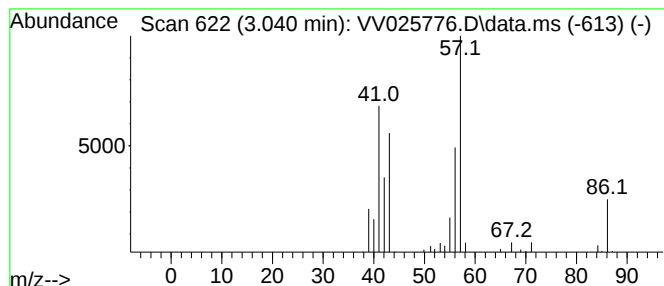
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 2-Ethyl-oxetane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	0.68 ug/L	47843	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Ethyl-oxetane	86	C5H10O	1010386-40-2	87
2		n-Hexane	86	C6H14	000110-54-3	80
3		n-Hexane	86	C6H14	000110-54-3	64
4		n-Hexane	86	C6H14	000110-54-3	50
5		n-Hexane	86	C6H14	000110-54-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

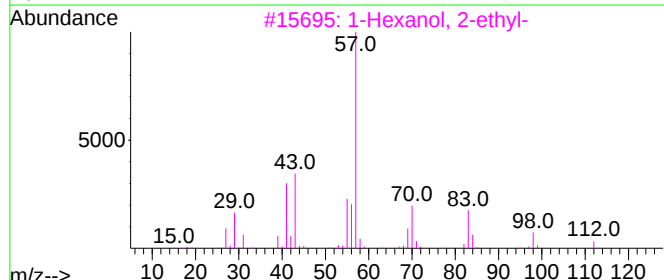
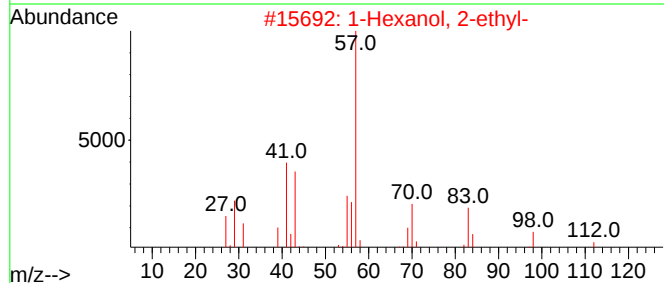
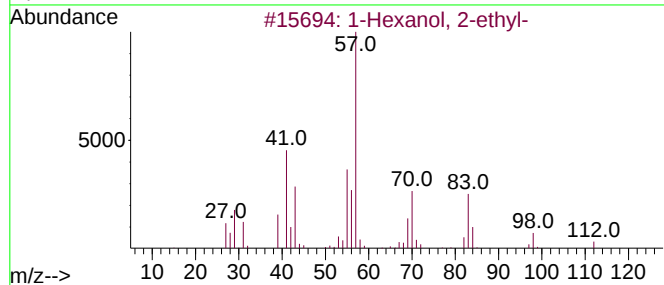
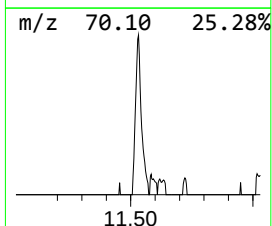
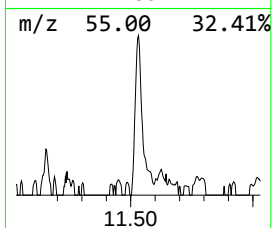
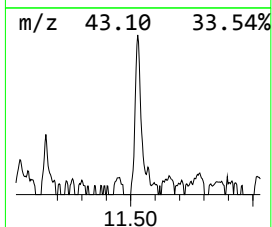
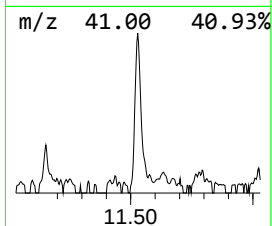
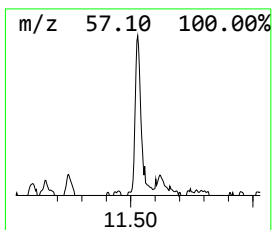
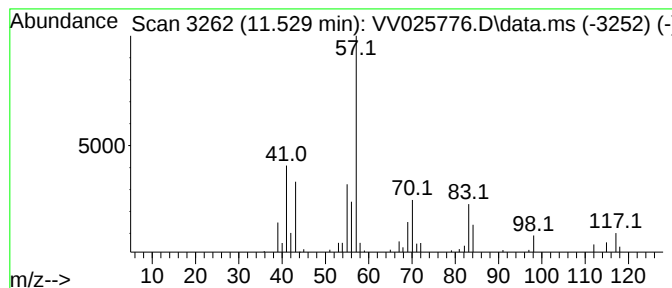
Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 1-Hexanol, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.529	0.66 ug/L	44822	1,4-Dichlorobenzene-d4			11.249
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	80
2	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	64
3	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	64
4	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025776.D
Acq On : 02 May 2022 19:15
Operator : SY/MD
Sample : N2625-09
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

					--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	1.111	16.8	ug/L	1181140	1	5.619	351773	5.0
(DEL) Alkane: S...	1.420	2.0	ug/L	140682	1	5.619	351773	5.0
(DEL) Alkane: S...	1.635	2.9	ug/L	203377	1	5.619	351773	5.0
(DEL) Alkane: S...	1.770	4.1	ug/L	290704	1	5.619	351773	5.0
(DEL) Alkane: S...	1.812	5.5	ug/L	388757	1	5.619	351773	5.0
(DEL) Alkane: C...	1.982	1.7	ug/L	120641	1	5.619	351773	5.0
(DEL) Alkane: S...	2.449	0.8	ug/L	55985	1	5.619	351773	5.0
(DEL) Alkane: S...	2.767	0.6	ug/L	43783	1	5.619	351773	5.0
(DEL) Alkane: S...	2.953	1.6	ug/L	109365	1	5.619	351773	5.0
2-Ethyl-oxetane	3.040	0.7	ug/L	47843	1	5.619	351773	5.0
1-Hexanol, 2-et...	11.529	0.7	ug/L	44822	3	11.249	342078	5.0