

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
 Data File : VV025777.D
 Acq On : 02 May 2022 19:39
 Operator : SY/MD
 Sample : N2625-10
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EW6Y2

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: VV025777.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	rVV	1242448	1134867	93.00%	13.278%
2	1.182	39	44	49	rVB	15185	13950	1.14%	0.163%
3	1.217	49	55	70	rBV	166178	149030	12.21%	1.744%
4	1.301	73	81	95	rVV2	858331	1220228	100.00%	14.277%
5	1.359	95	99	103	rVV2	25916	25665	2.10%	0.300%
6	1.391	103	109	112	rVV	26799	32919	2.70%	0.385%
7	1.420	112	118	135	rVB	123346	143163	11.73%	1.675%
8	1.568	152	164	176	rVB2	119767	168407	13.80%	1.970%
9	1.635	177	185	197	rVB	163554	202875	16.63%	2.374%
10	1.770	218	227	233	rBV	216900	272062	22.30%	3.183%
11	1.815	233	241	259	rVB3	206258	379929	31.14%	4.445%
12	1.905	261	269	278	rVB	21754	28339	2.32%	0.332%
13	1.982	282	293	301	rBV2	82474	121406	9.95%	1.421%
14	2.027	301	307	314	rVB2	17125	22415	1.84%	0.262%
15	2.111	322	333	343	rBV	135833	201443	16.51%	2.357%
16	2.352	400	408	419	rVB2	15884	25910	2.12%	0.303%
17	2.445	424	437	450	rBV6	23456	63812	5.23%	0.747%
18	2.532	459	464	472	rVB2	13273	18253	1.50%	0.214%
19	2.577	474	478	490	rVB3	10099	17420	1.43%	0.204%
20	2.764	521	536	549	rBV3	21138	44955	3.68%	0.526%
21	2.953	583	595	611	rVB5	46388	102933	8.44%	1.204%
22	3.043	612	623	642	rVB3	19929	44609	3.66%	0.522%
23	3.156	645	658	666	rBV5	8973	19279	1.58%	0.226%
24	3.224	671	679	689	rVB6	6018	13012	1.07%	0.152%
25	3.458	740	752	767	rBV7	12032	29675	2.43%	0.347%
26	3.905	879	891	917	rBV2	42928	166817	13.67%	1.952%
27	4.352	1013	1030	1053	rBV2	102470	263717	21.61%	3.086%
28	4.757	1143	1156	1177	rBV2	6609	20048	1.64%	0.235%
29	4.915	1193	1205	1216	rBV5	6630	13502	1.11%	0.158%
30	5.050	1231	1247	1278	rBV2	203431	538116	44.10%	6.296%
31	5.320	1321	1331	1340	rBV3	7190	14983	1.23%	0.175%
32	5.500	1375	1387	1397	rBV7	5882	13382	1.10%	0.157%
33	5.619	1412	1424	1450	rBV	170608	369122	30.25%	4.319%
34	6.072	1549	1565	1581	rBV	129111	271521	22.25%	3.177%
35	6.992	1838	1851	1872	rBV2	34520	69697	5.71%	0.815%
36	7.317	1941	1952	1968	rBV	216581	376785	30.88%	4.409%

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 Title : TRACE VOA SFAM1.0

37	7.394	1969	1976	1990	rVB	18765	35466	2.91%	0.415%
38	7.628	2040	2049	2058	rBV2	17637	32961	2.70%	0.386%
39	8.091	2182	2193	2229	rBV	228391	467610	38.32%	5.471%
40	8.850	2418	2429	2447	rBV	266259	441921	36.22%	5.171%
41	9.017	2473	2481	2490	rBV2	10214	15159	1.24%	0.177%
42	9.143	2511	2520	2532	rBV3	10090	20509	1.68%	0.240%
43	10.217	2843	2854	2866	rVB	88331	141774	11.62%	1.659%
44	11.249	3163	3175	3197	rBV	237445	370867	30.39%	4.339%
45	11.529	3252	3262	3281	rBV3	29462	55958	4.59%	0.655%
46	11.622	3281	3291	3314	rVB	201353	320139	26.24%	3.746%
47	12.345	3505	3516	3526	rBV5	9524	15805	1.30%	0.185%
48	17.094	4983	4993	5001	rBV2	8400	14245	1.17%	0.167%

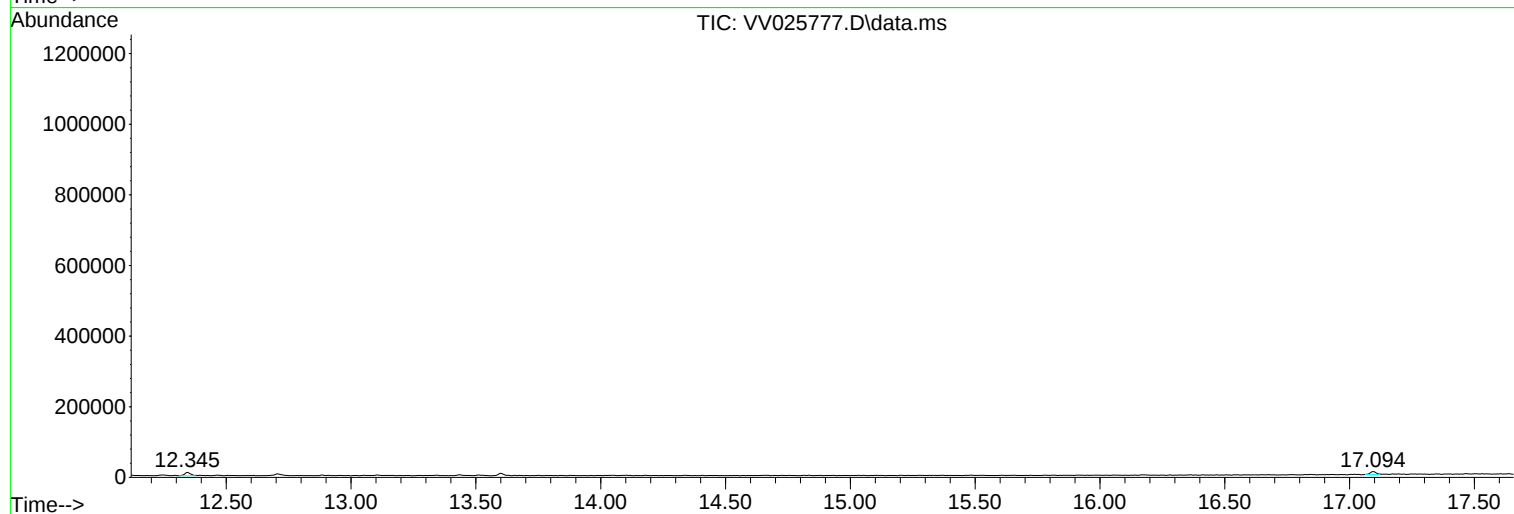
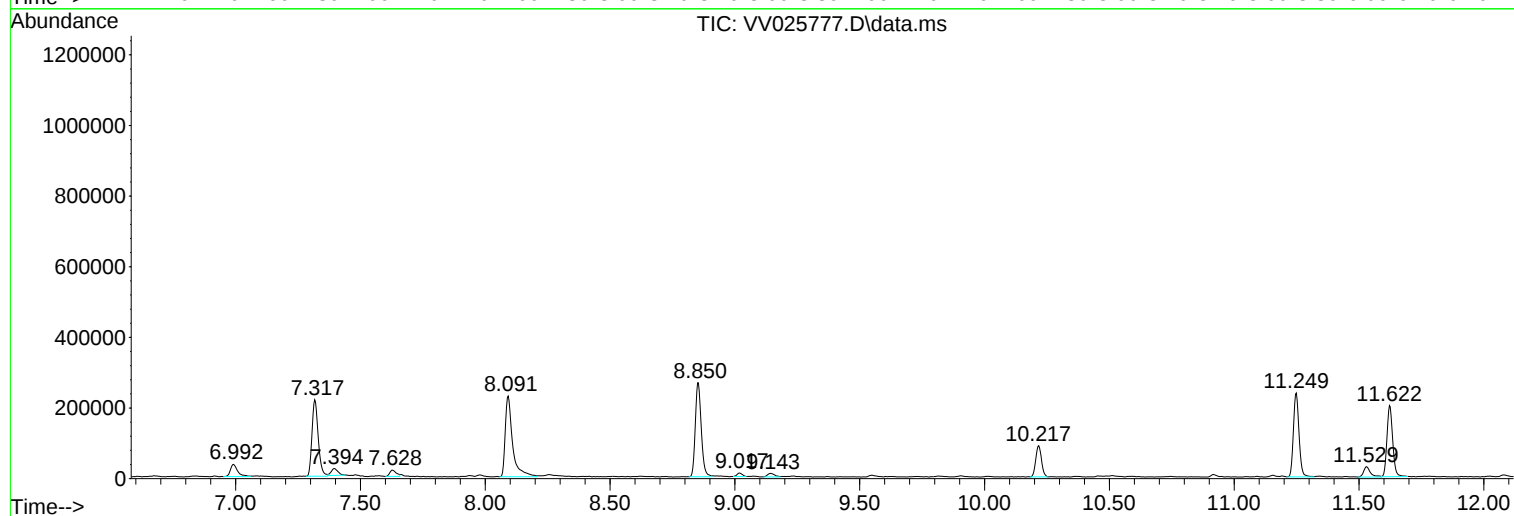
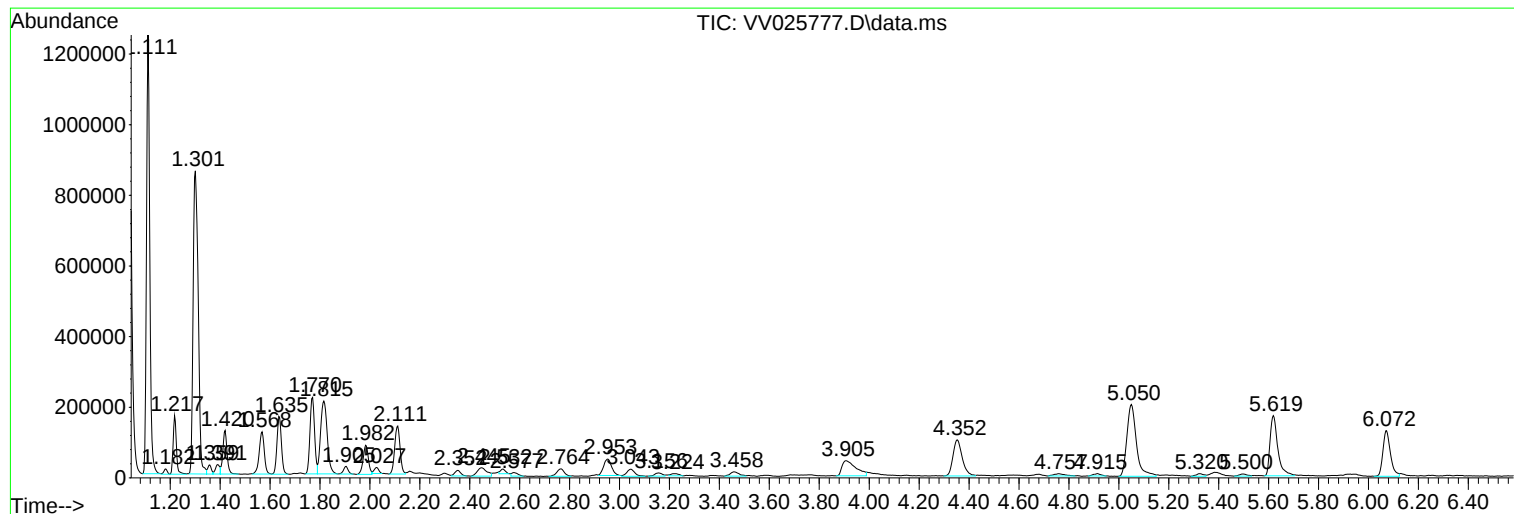
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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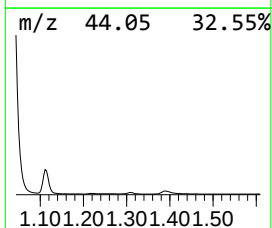
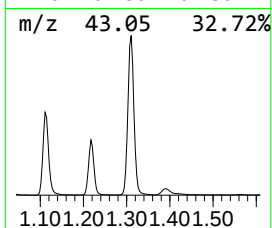
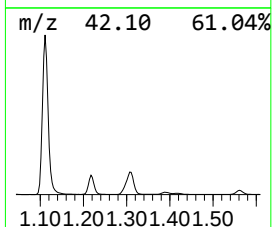
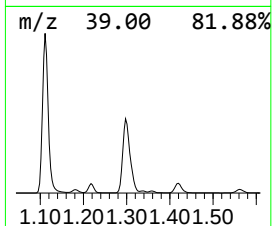
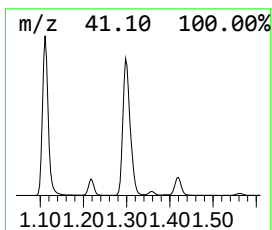
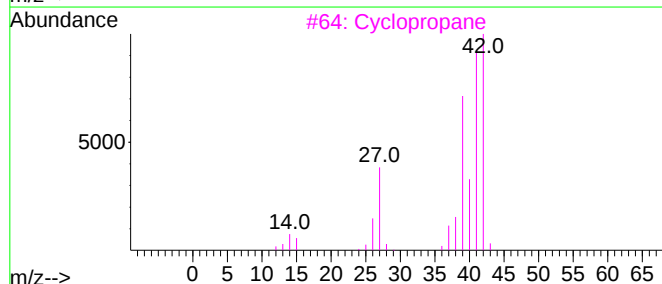
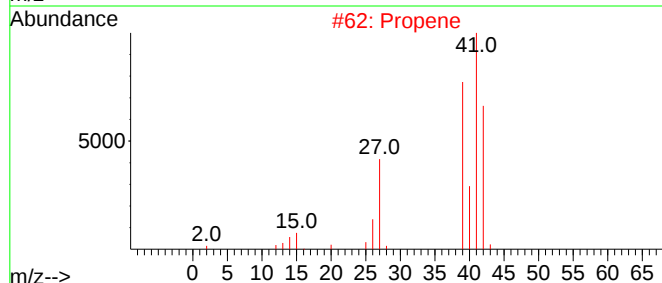
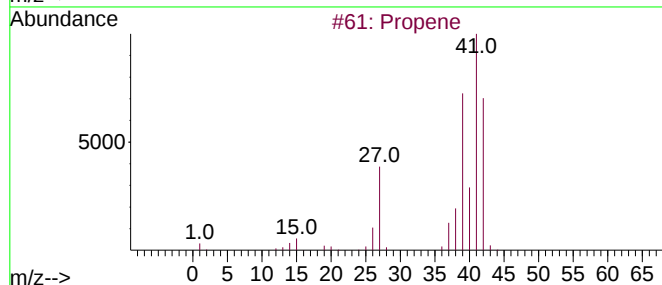
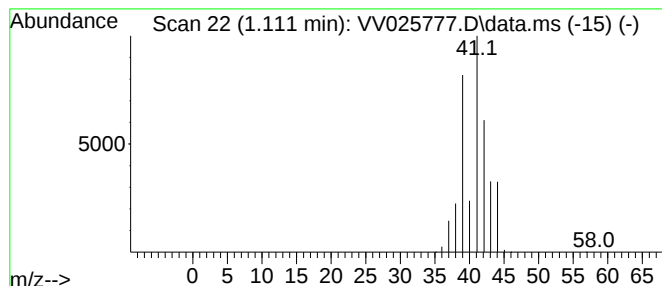
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Quant Title : TRACE VOA SFAM1.0

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	15.37 ug/L	1134870	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			Cyclopropane	42	C3H6	000075-19-4	7
5			Cyclopropene	40	C3H4	002781-85-3	5



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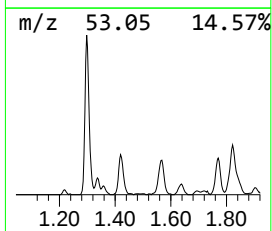
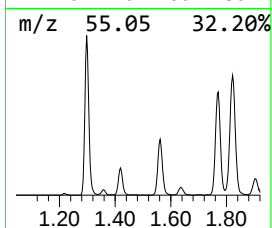
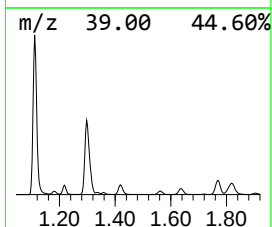
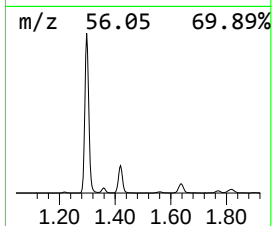
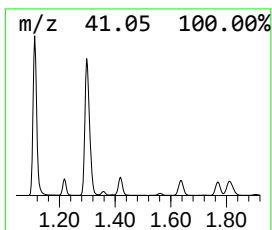
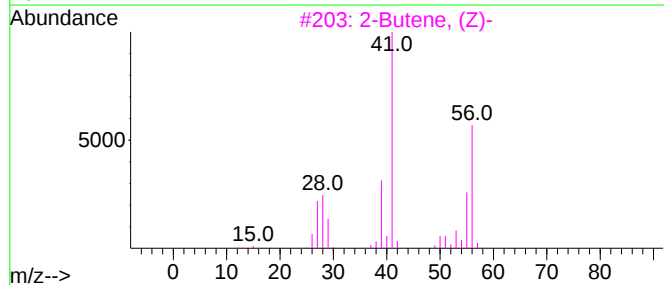
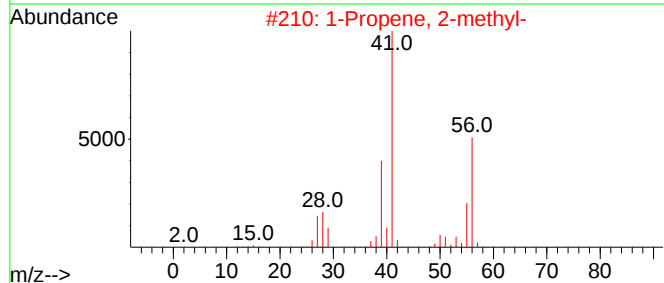
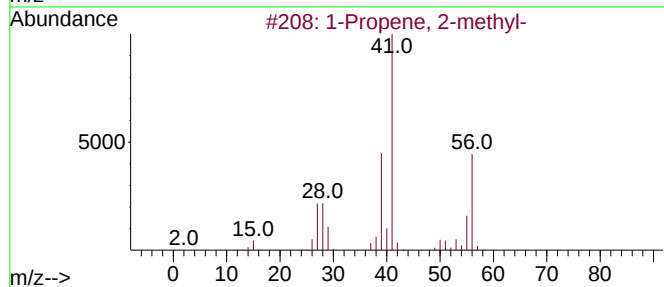
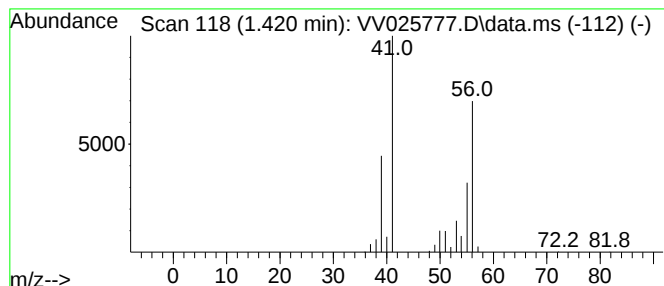
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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	1.94 ug/L	143163	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Propene, 2-methyl-	56	C4H8	000115-11-7	72
2		1-Propene, 2-methyl-	56	C4H8	000115-11-7	72
3		2-Butene, (Z)-	56	C4H8	000590-18-1	72
4		2-Butene, (Z)-	56	C4H8	000590-18-1	68
5		2-Butene	56	C4H8	000107-01-7	64



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ALS Vial : 21 Sample Multiplier: 1

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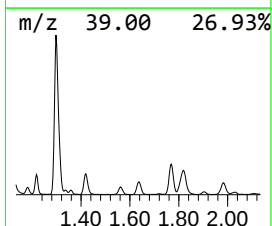
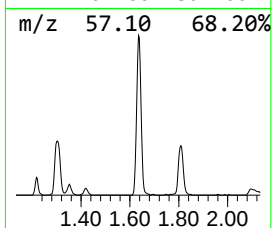
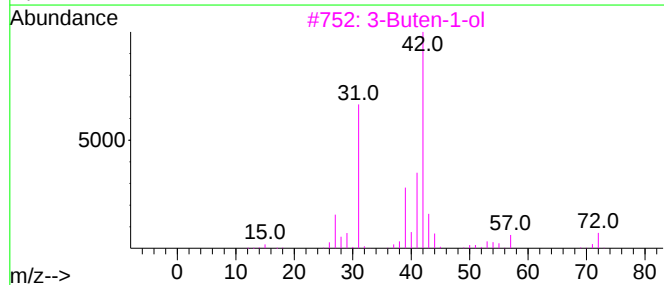
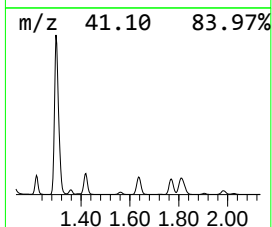
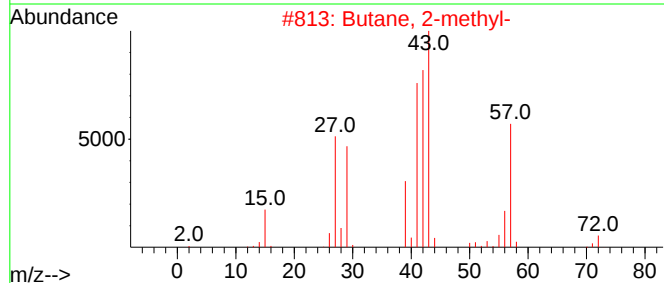
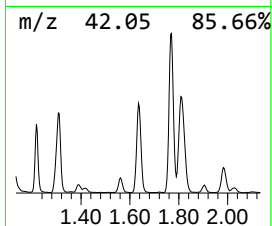
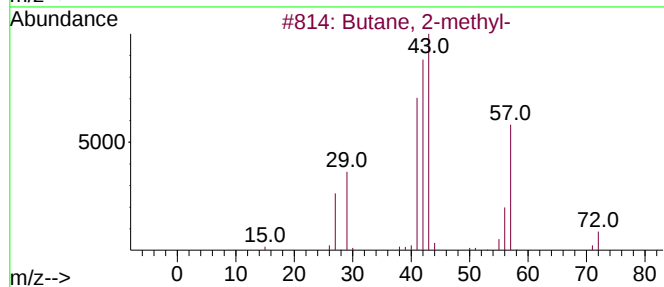
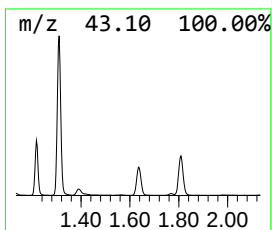
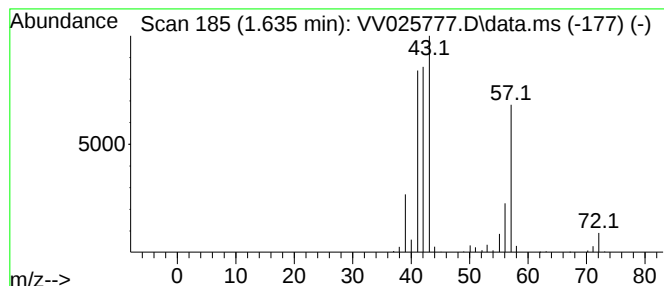
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.635	2.75 ug/L	202875	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	90
3	3-Buten-1-ol			72	C4H8O	000627-27-0	47
4	Pentane			72	C5H12	000109-66-0	37
5	1-Butanesulfonyl chloride			156	C4H9ClO2S	002386-60-9	10



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Instrument :
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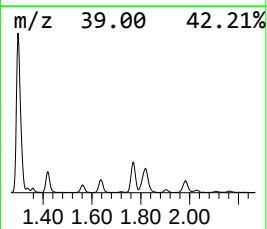
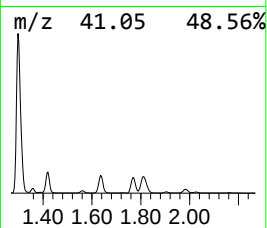
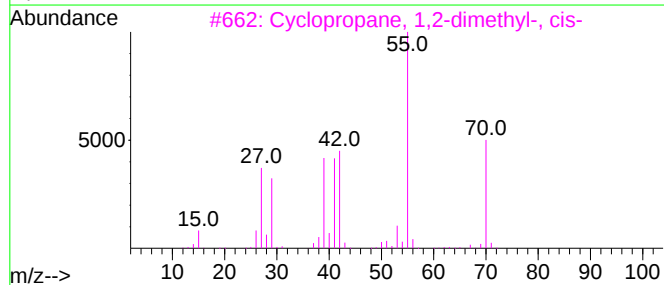
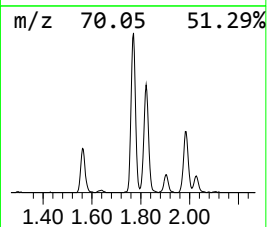
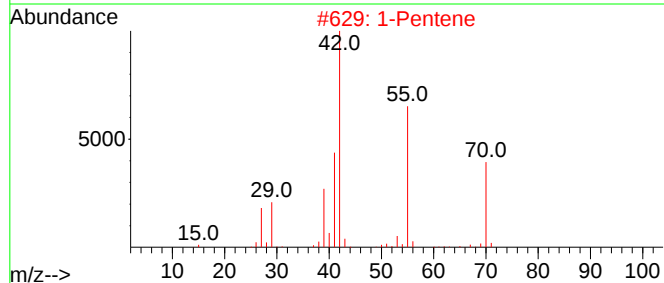
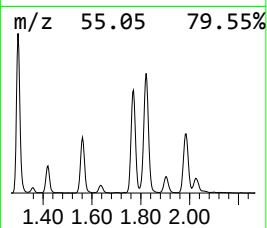
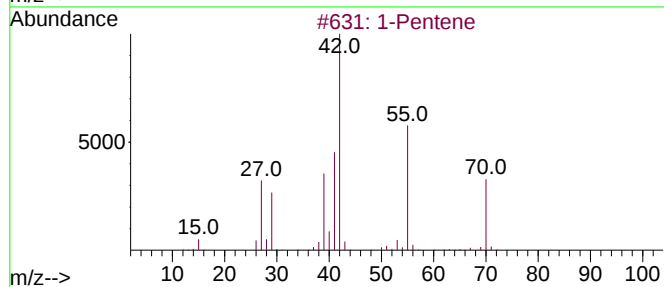
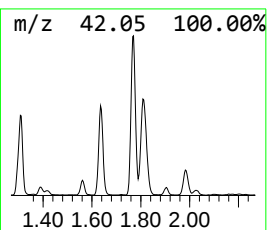
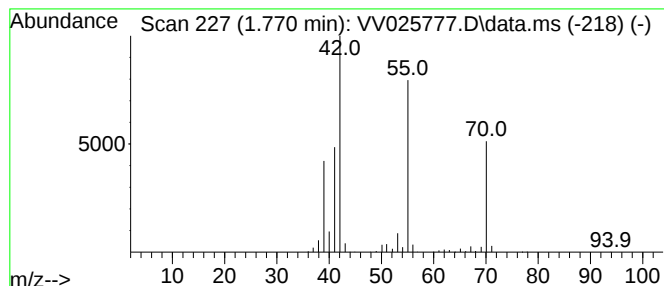
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.770	3.69 ug/L	272062	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene	70	C5H10	000109-67-1	70
2		1-Pentene	70	C5H10	000109-67-1	68
3		Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	64
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
5		Cyclobutane, methyl-	70	C5H10	000598-61-8	62



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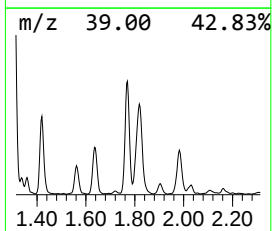
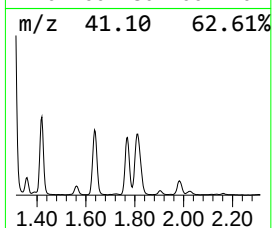
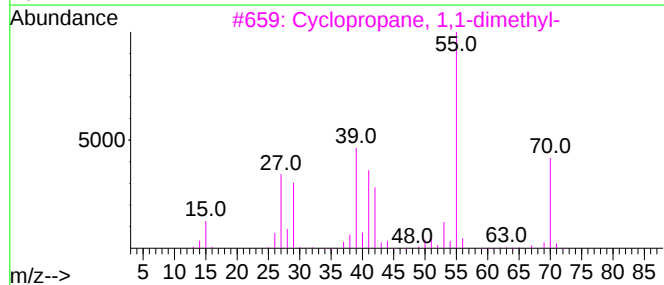
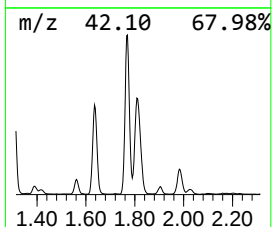
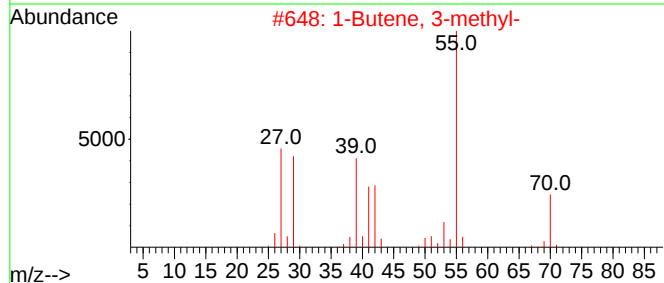
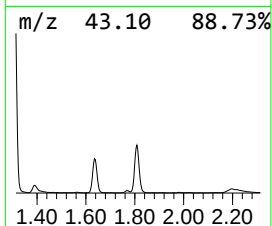
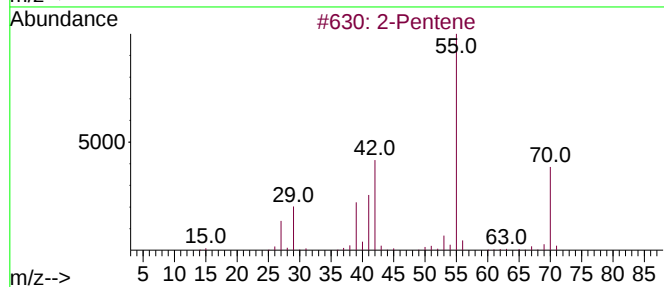
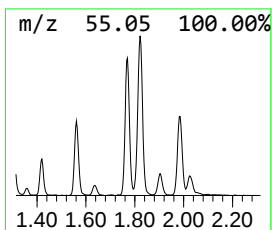
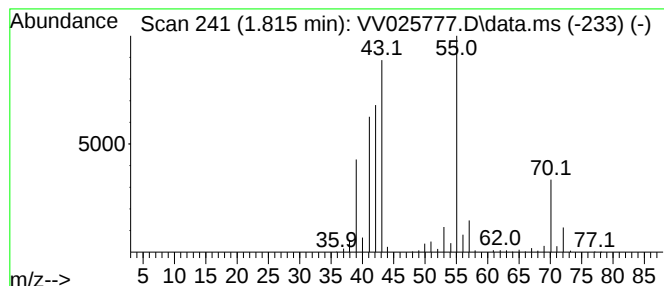
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Quant Title : TRACE VOA SFAM1.0

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.
1.815	5.15 ug/L	379929	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentene	70	C5H10	000109-68-2	64
2			1-Butene, 3-methyl-	70	C5H10	000563-45-1	58
3			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	50
4			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	49
5			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y2

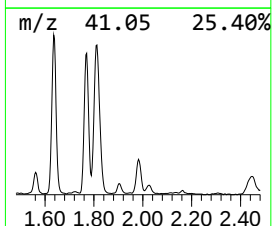
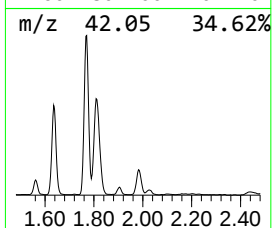
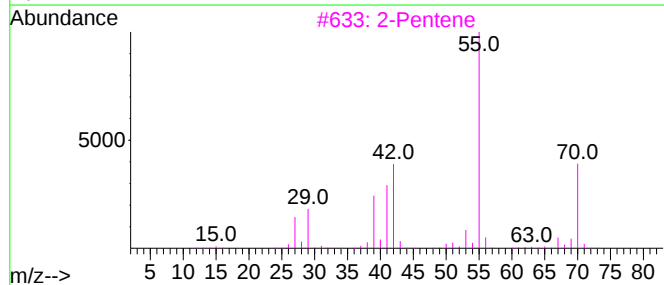
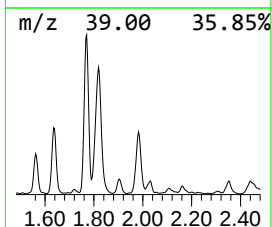
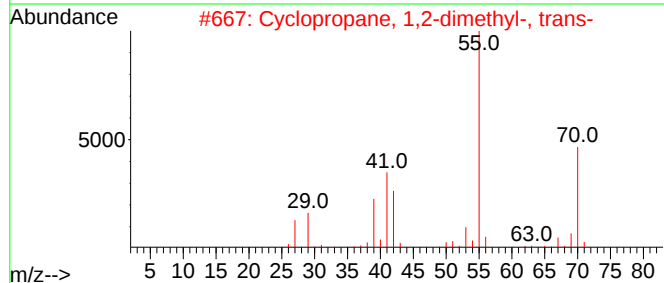
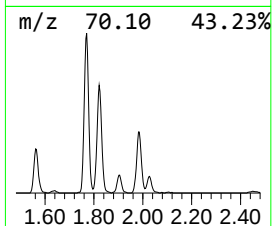
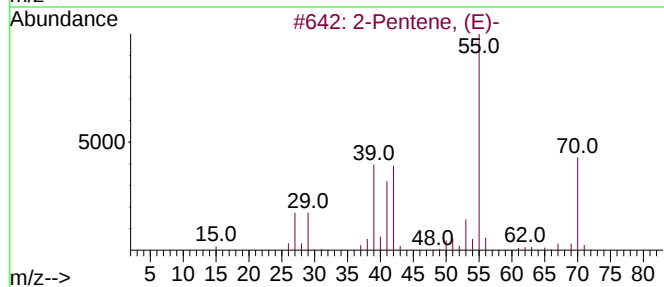
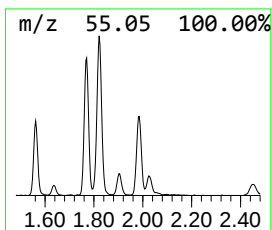
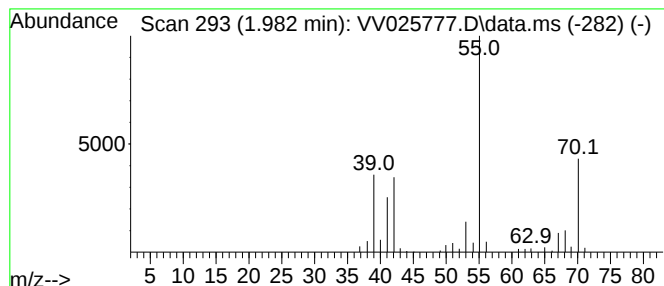
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.982	1.64 ug/L	121406	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (E)-		70	C5H10	000646-04-8	93
2	Cyclopropane, 1,2-dimethyl-, trans-		70	C5H10	002402-06-4	87
3	2-Pentene		70	C5H10	000109-68-2	86
4	Cyclopropane, 1,2-dimethyl-, cis-		70	C5H10	000930-18-7	83
5	2-Methyl-1-butene		70	C5H10	000563-46-2	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y2

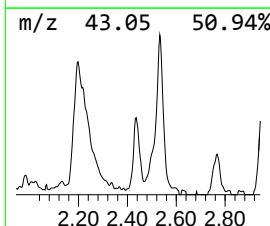
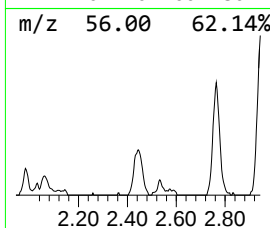
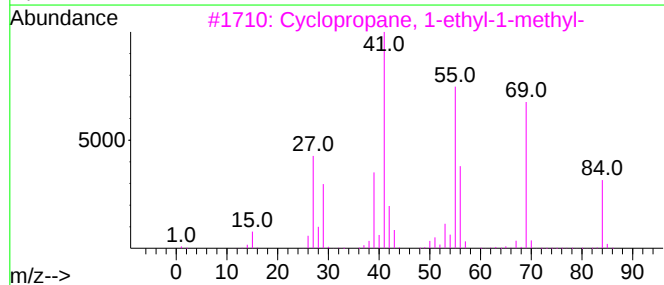
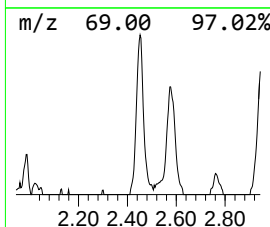
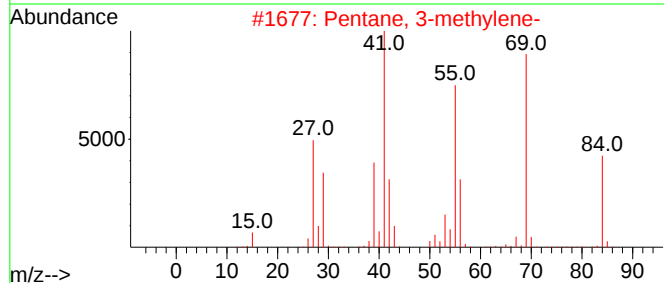
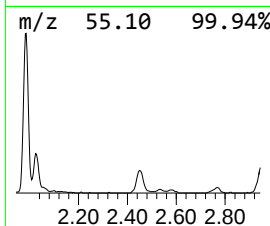
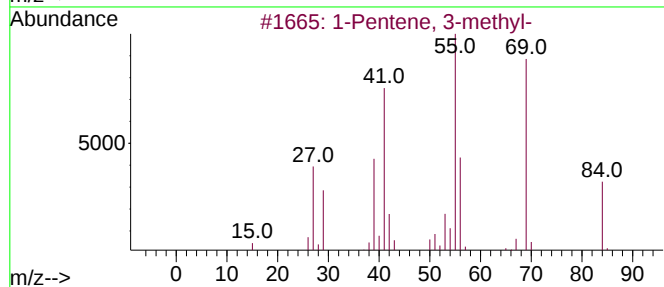
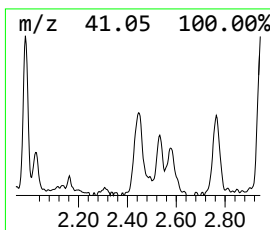
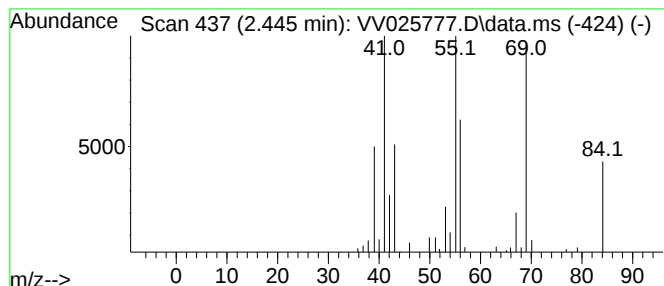
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Quant Title : TRACE VOA SFAM1.0

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.
2.445	0.86 ug/L	63812	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	87
2		Pentane, 3-methylene-	84	C6H12	000760-21-4	81
3		Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	76
4		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	62
5		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

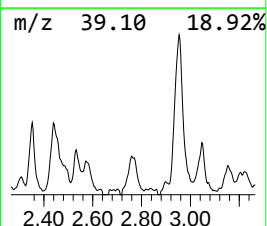
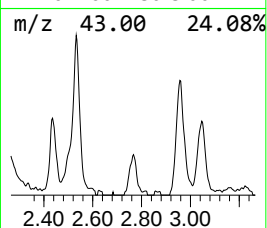
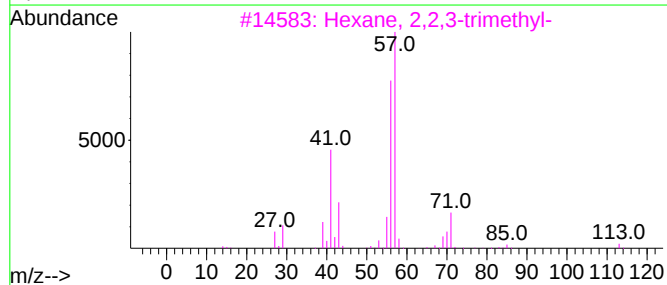
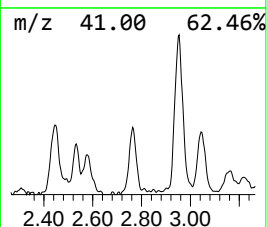
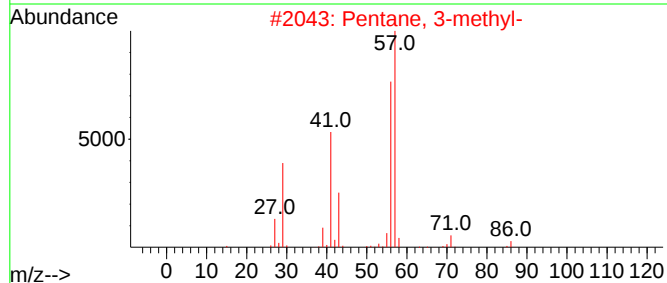
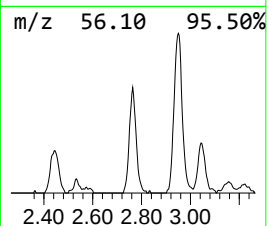
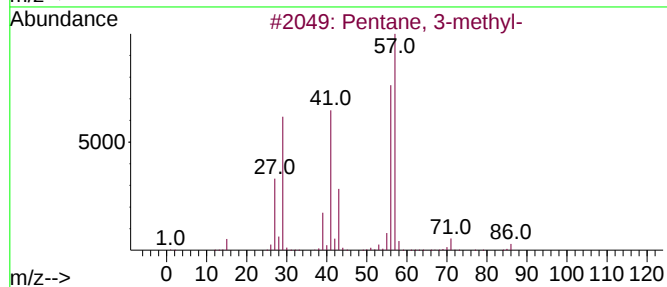
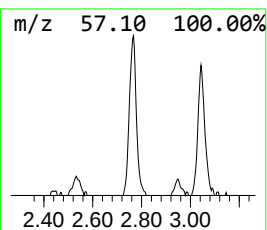
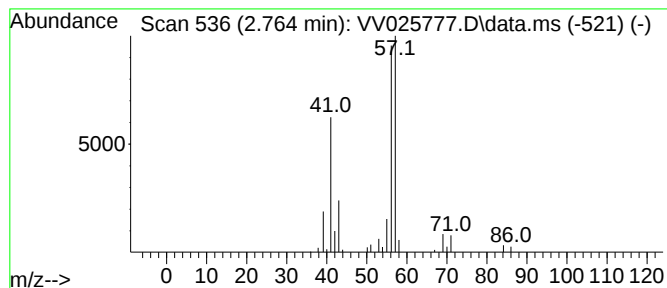
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MSVOA_V
ClientSampleId :
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.764	0.61 ug/L	44955	1,4-Difluorobenzene	5.619	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	72
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59
5	Cyclopropane, propyl-	84	C6H12	002415-72-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y2

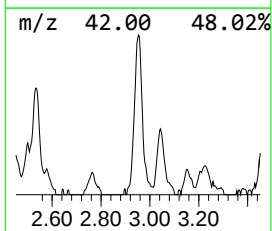
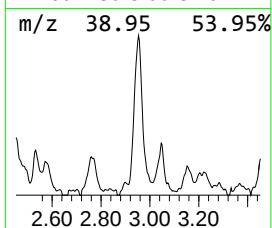
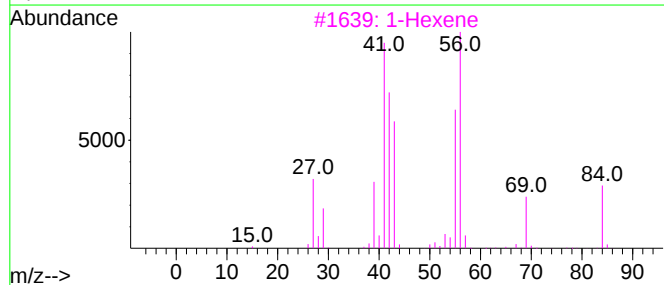
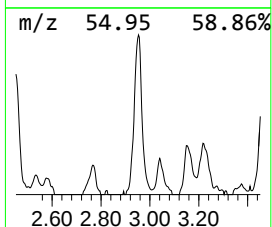
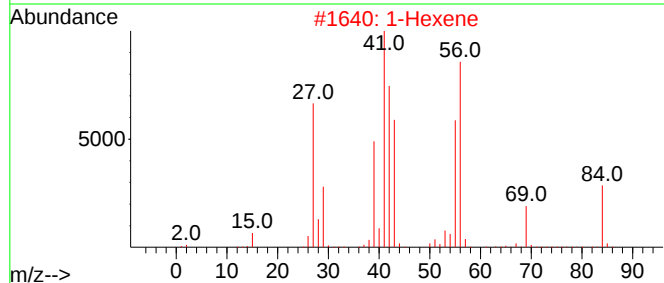
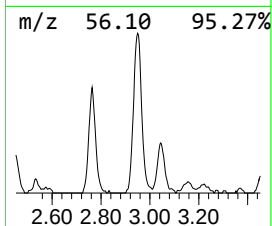
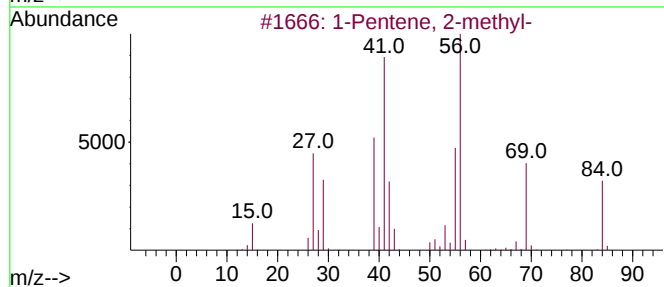
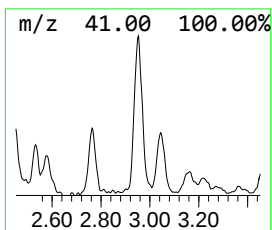
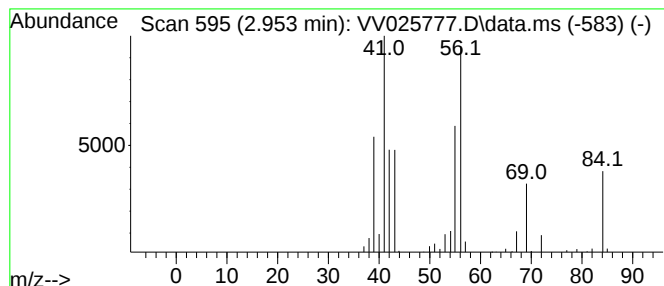
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Quant Title : TRACE VOA SFAM1.0

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	1.39 ug/L	102933	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	74
2		1-Hexene	84	C6H12	000592-41-6	68
3		1-Hexene	84	C6H12	000592-41-6	64
4		1-Hexene	84	C6H12	000592-41-6	58
5		Cyclohexane	84	C6H12	000110-82-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y2

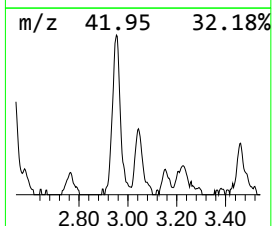
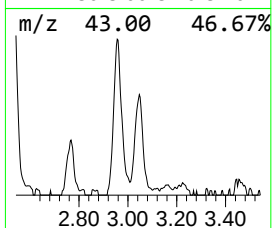
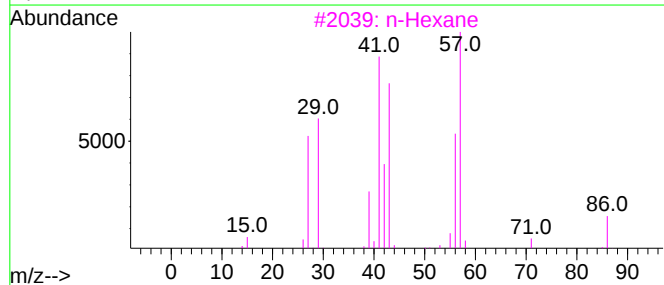
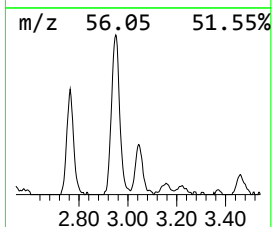
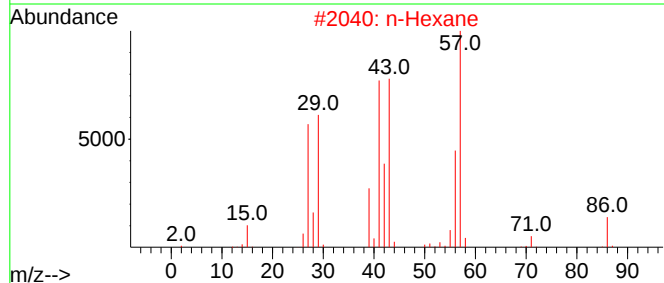
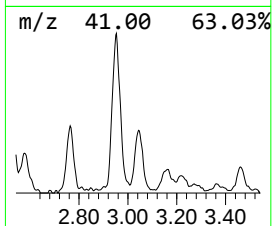
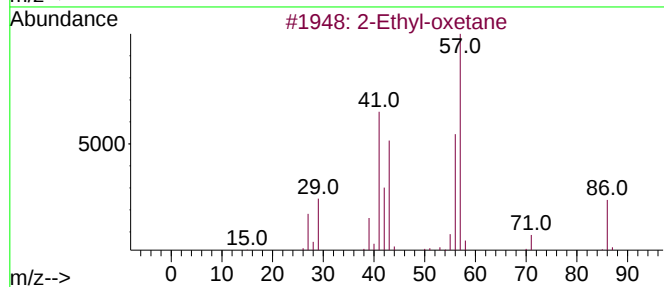
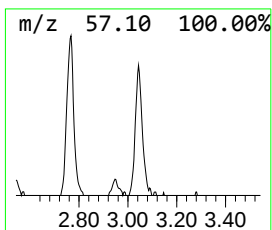
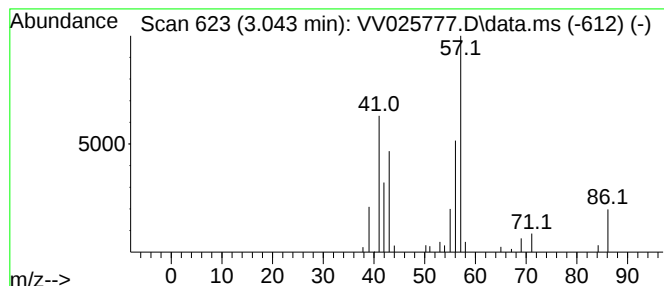
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 2-Ethyl-oxetane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.043	0.60 ug/L	44609	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Ethyl-oxetane	86	C5H10O	1010386-40-2	90
2		n-Hexane	86	C6H14	000110-54-3	78
3		n-Hexane	86	C6H14	000110-54-3	78
4		n-Hexane	86	C6H14	000110-54-3	56
5		1-Butanol	74	C4H10O	000071-36-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y2

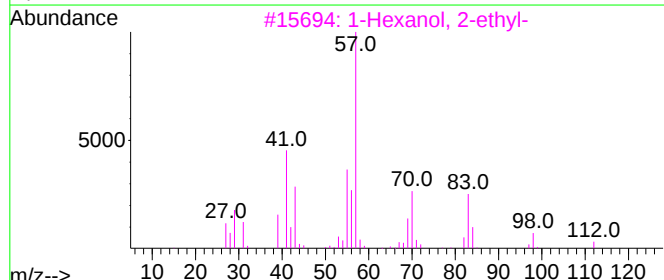
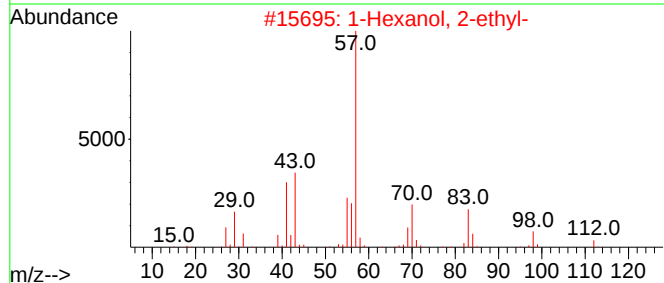
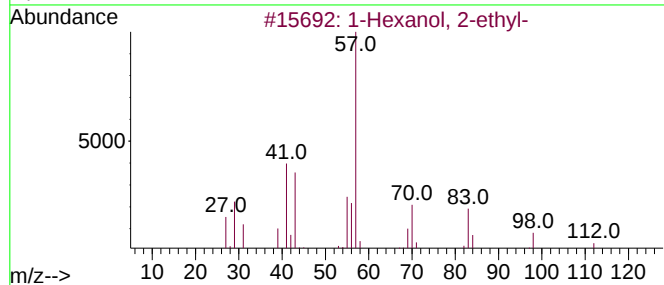
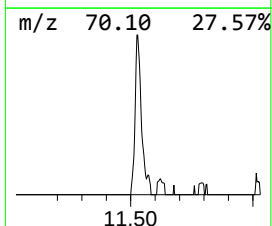
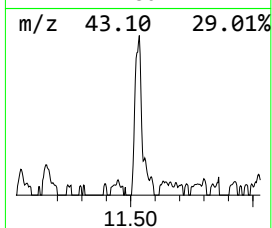
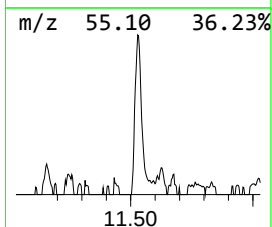
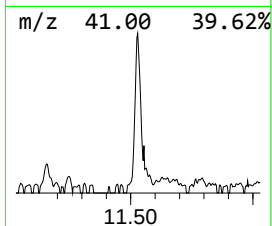
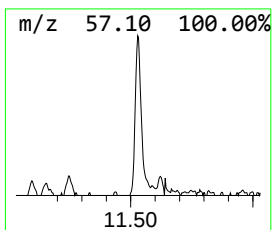
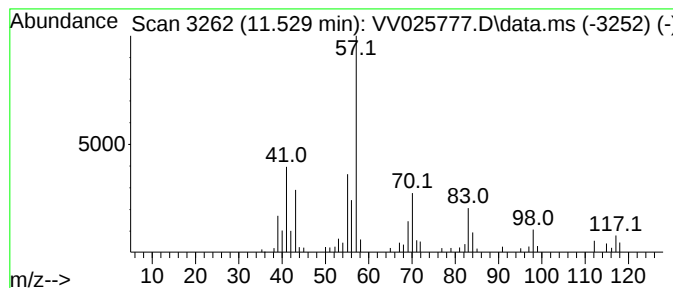
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 1-Hexanol, 2-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	0.75 ug/L	55958	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	78
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
4			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	64
5			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050222\
Data File : VV025777.D
Acq On : 02 May 2022 19:39
Operator : SY/MD
Sample : N2625-10
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EW6Y2

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	15.4	ug/L	1134870	1	5.619	369122	5.0
(DEL) Alkane: S...	1.420	1.9	ug/L	143163	1	5.619	369122	5.0
(DEL) Alkane: S...	1.635	2.8	ug/L	202875	1	5.619	369122	5.0
(DEL) Alkane: S...	1.770	3.7	ug/L	272062	1	5.619	369122	5.0
(DEL) Alkane: S...	1.815	5.2	ug/L	379929	1	5.619	369122	5.0
(DEL) Alkane: S...	1.982	1.6	ug/L	121406	1	5.619	369122	5.0
(DEL) Alkane: S...	2.445	0.9	ug/L	63812	1	5.619	369122	5.0
(DEL) Alkane: S...	2.764	0.6	ug/L	44955	1	5.619	369122	5.0
(DEL) Alkane: S...	2.953	1.4	ug/L	102933	1	5.619	369122	5.0
2-Ethyl-oxetane	3.043	0.6	ug/L	44609	1	5.619	369122	5.0
1-Hexanol, 2-et...	11.529	0.8	ug/L	55958	3	11.249	370867	5.0