

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
 Data File : VV023378.D
 Acq On : 11 Nov 2021 11:06
 Operator : SY/MD
 Sample : M4558-09DL 100X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GB871DL

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\SFAMVTR110421WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV023378.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.214	47	54	64	rVB	22699	20146	1.66%	0.381%
2	1.307	74	83	93	rVB	159176	158673	13.11%	2.998%
3	1.568	156	164	173	rBV	47579	54345	4.49%	1.027%
4	1.632	175	184	202	rVB	164995	208166	17.20%	3.934%
5	1.809	229	239	253	rVB3	37151	62238	5.14%	1.176%
6	1.902	258	268	279	rVB	18910	25402	2.10%	0.480%
7	2.024	298	306	318	rVB	45226	63753	5.27%	1.205%
8	2.108	321	332	349	rVB	92596	138167	11.42%	2.611%
9	2.497	436	453	460	rBV8	28405	71067	5.87%	1.343%
10	2.577	471	478	493	rVB3	27025	50904	4.21%	0.962%
11	2.761	523	535	548	rBV2	21245	42167	3.48%	0.797%
12	3.764	831	847	863	rBV3	29411	74139	6.13%	1.401%
13	3.909	882	892	918	rBV2	29507	100446	8.30%	1.898%
14	4.352	1012	1030	1057	rBV	61566	171816	14.20%	3.247%
15	4.677	1115	1131	1147	rBV4	24213	63080	5.21%	1.192%
16	4.761	1148	1157	1178	rVB5	10977	29733	2.46%	0.562%
17	5.050	1232	1247	1253	rBV2	158490	353333	29.19%	6.677%
18	5.098	1253	1262	1285	rVB	557838	1210326	100.00%	22.871%
19	5.230	1296	1303	1320	rVB7	13701	37481	3.10%	0.708%
20	5.323	1323	1332	1345	rVB7	6727	14047	1.16%	0.265%
21	5.619	1411	1424	1450	rBV	136812	294546	24.34%	5.566%
22	6.072	1552	1565	1578	rBV	83860	177247	14.64%	3.349%
23	6.654	1733	1746	1765	rBV3	8763	19035	1.57%	0.360%
24	6.777	1770	1784	1798	rBV3	16342	37131	3.07%	0.702%
25	6.989	1840	1850	1870	rBV	41119	77571	6.41%	1.466%
26	7.317	1941	1952	1971	rBV	178425	312284	25.80%	5.901%
27	7.625	2040	2048	2066	rBV	23481	42210	3.49%	0.798%
28	8.092	2184	2193	2224	rBV	137276	274874	22.71%	5.194%
29	8.854	2419	2430	2454	rBV	206780	338770	27.99%	6.402%
30	10.217	2844	2854	2874	rVB2	56245	88446	7.31%	1.671%
31	11.249	3165	3175	3199	rBV	211484	329811	27.25%	6.232%
32	11.529	3253	3262	3276	rBV	29242	52344	4.32%	0.989%
33	11.625	3280	3292	3315	rVV	180501	284545	23.51%	5.377%
34	12.114	3435	3444	3459	rVB2	8276	13621	1.13%	0.257%

Sum of corrected areas: 5291864

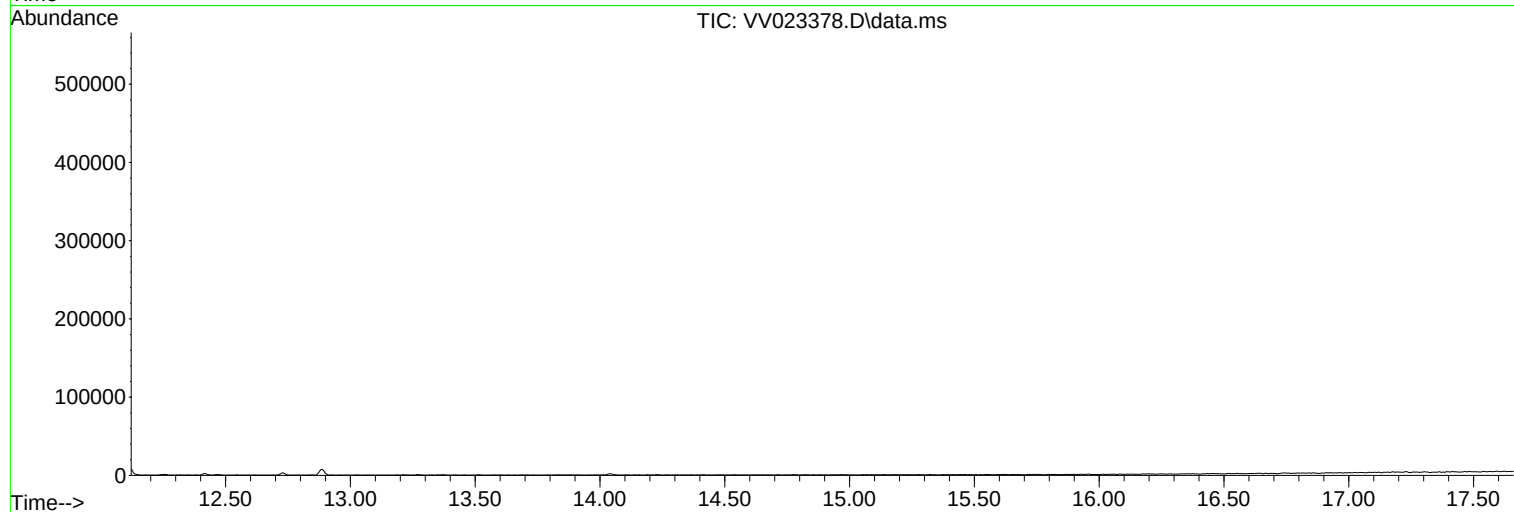
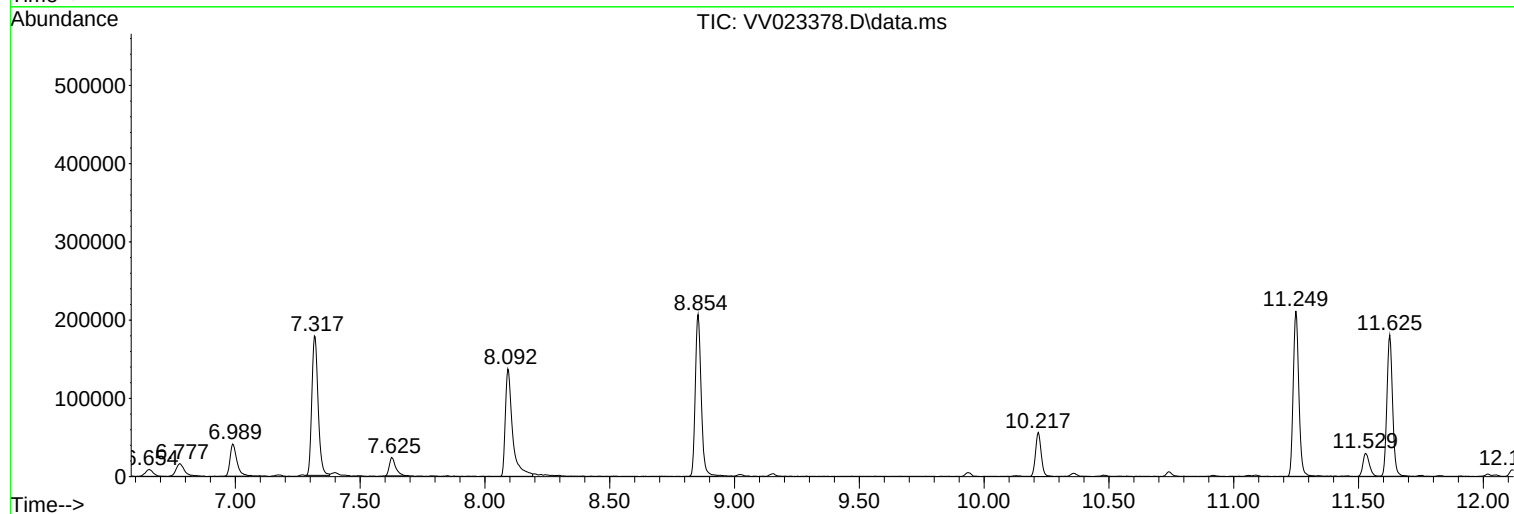
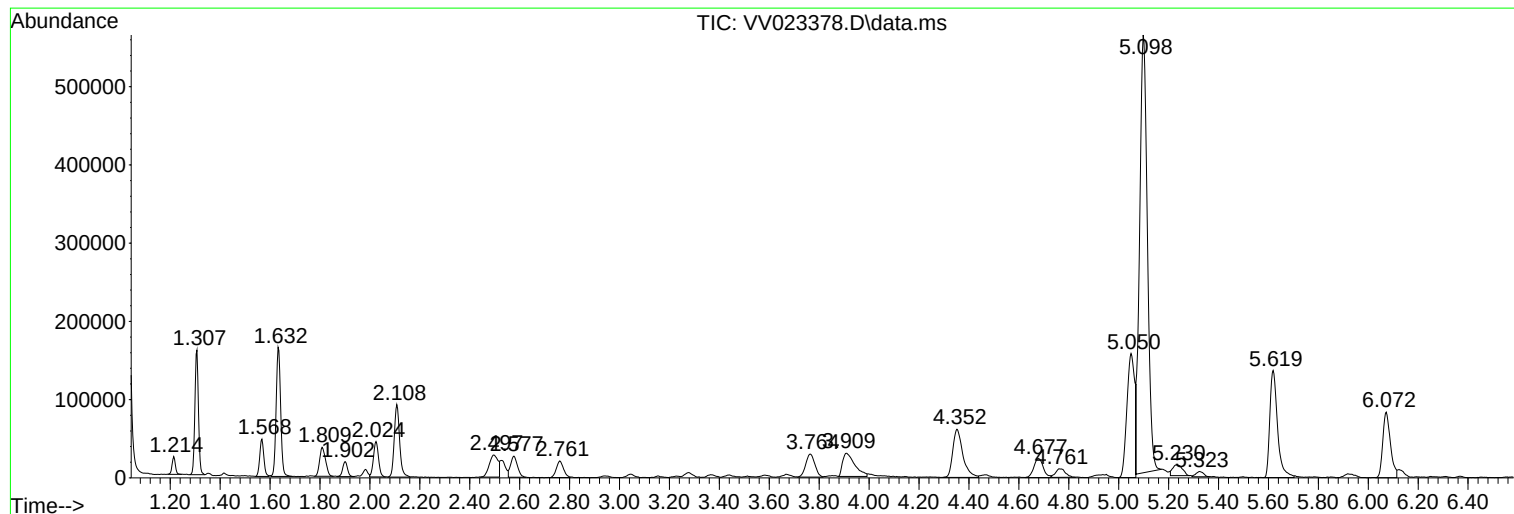
Instrument :
MSVOA_V
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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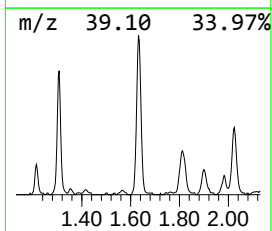
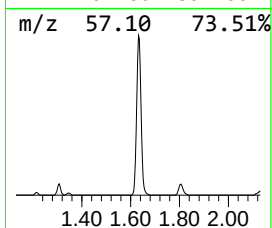
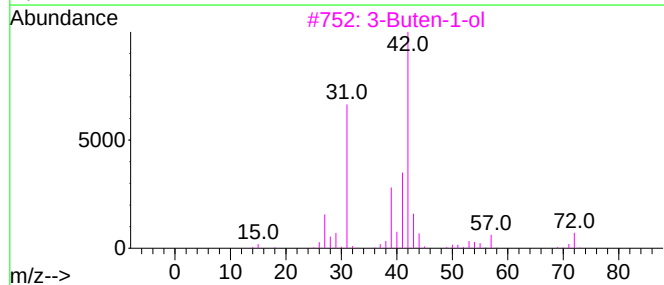
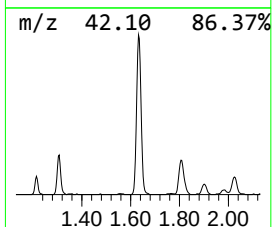
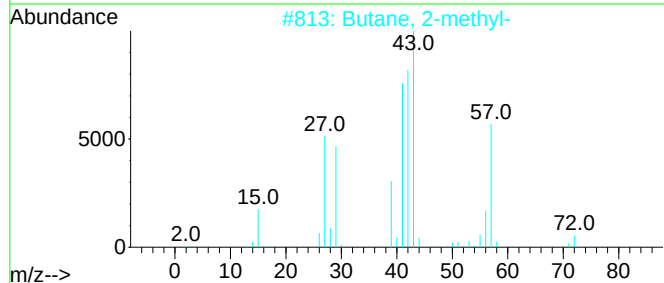
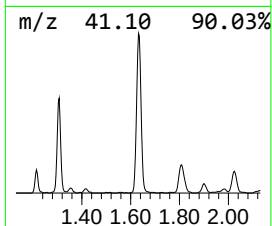
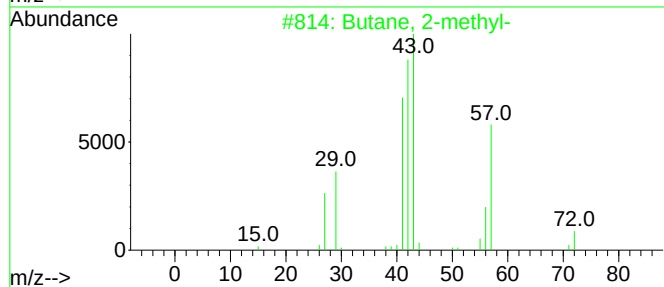
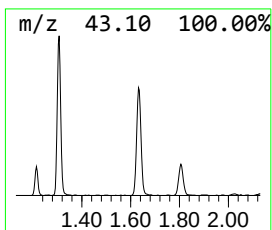
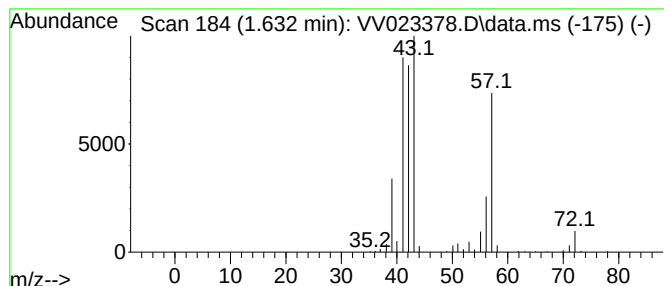
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Quant Title : TRACE VOA SFAM1.0

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.
1.632	3.53 ug/L	208166	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-			72	C5H12	000078-78-4	86
2	Butane, 2-methyl-			72	C5H12	000078-78-4	80
3	3-Buten-1-ol			72	C4H8O	000627-27-0	43
4	1-Propanol, 2-methyl-			74	C4H10O	000078-83-1	36
5	Pentane			72	C5H12	000109-66-0	28



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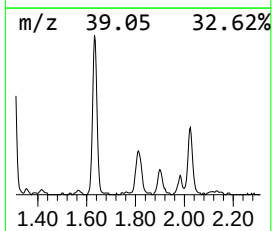
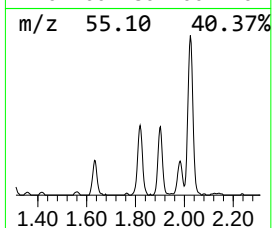
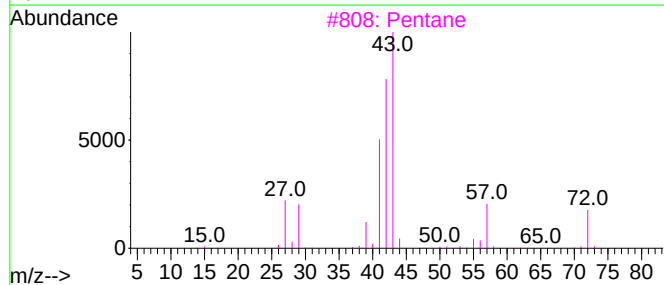
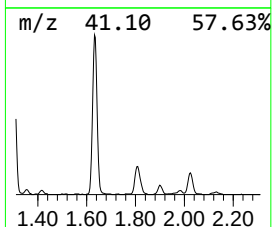
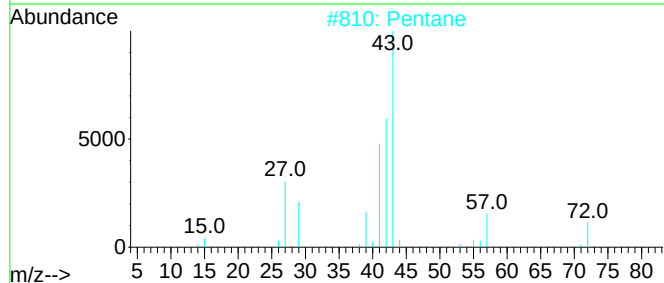
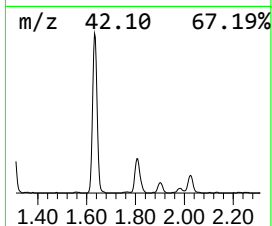
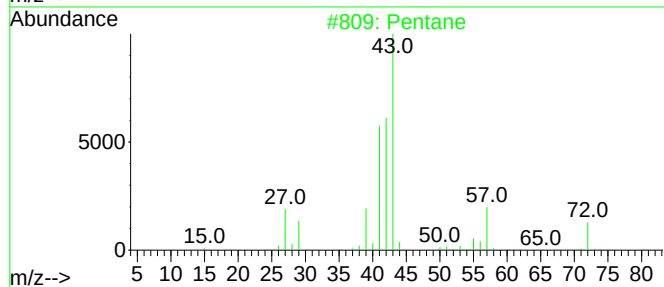
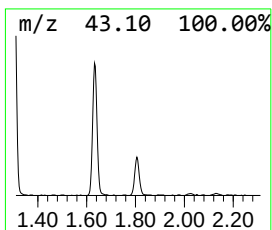
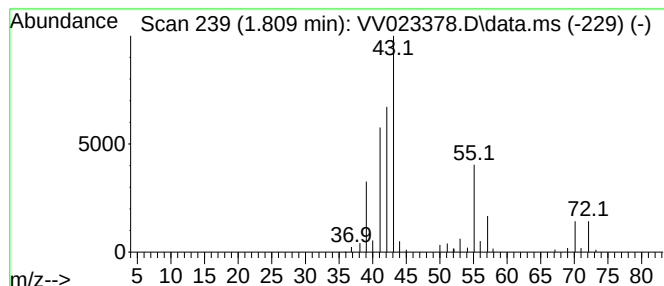
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.809	1.06 ug/L	62238	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	64
3	Pentane			72	C5H12	000109-66-0	64
4	Pentane			72	C5H12	000109-66-0	59
5	Pentane			72	C5H12	000109-66-0	59



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ALS Vial : 65 Sample Multiplier: 1

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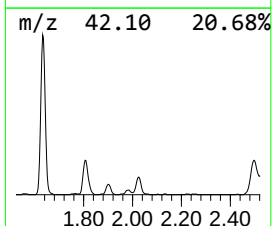
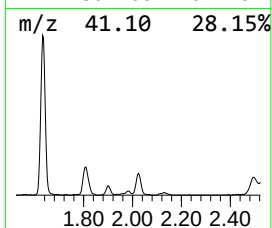
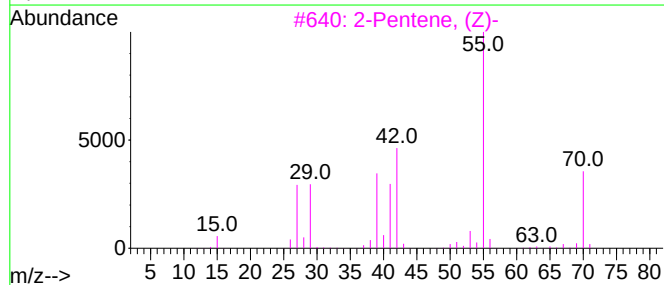
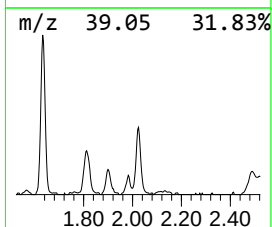
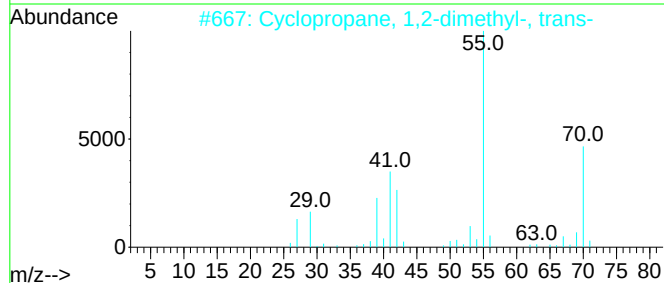
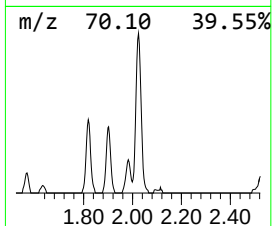
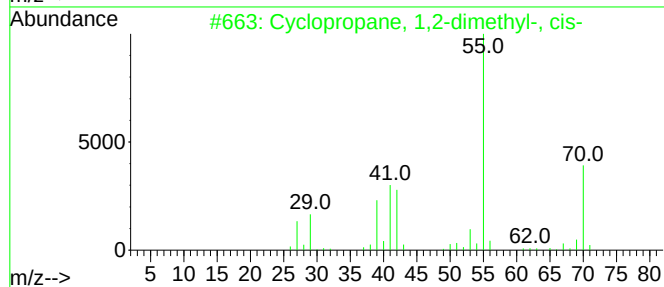
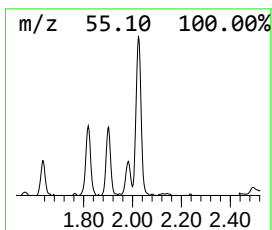
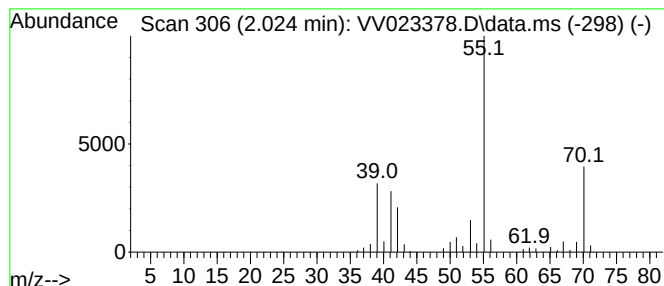
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Cyclic2.024 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.024	1.08 ug/L	63753	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	86
2			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	86
3			2-Pentene, (Z)-	70	C5H10	000627-20-3	80
4			2-Butene, 2-methyl-	70	C5H10	000513-35-9	72
5			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	72



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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

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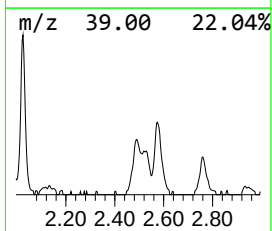
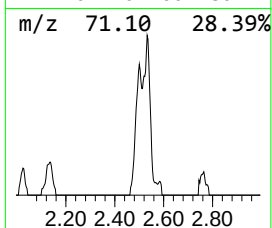
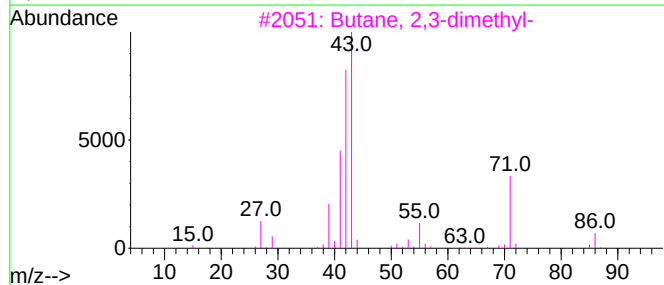
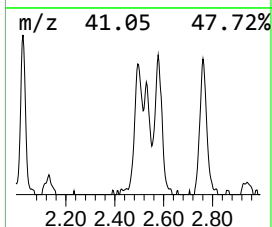
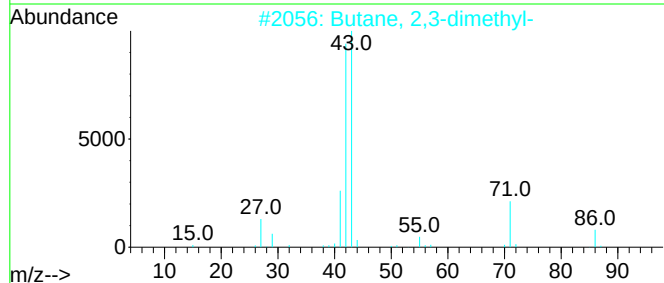
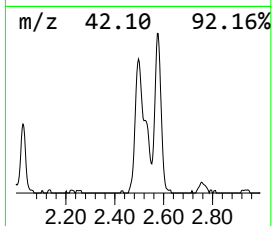
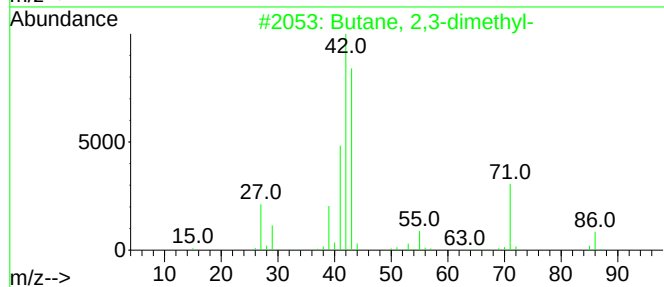
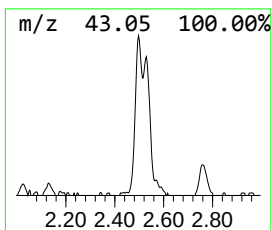
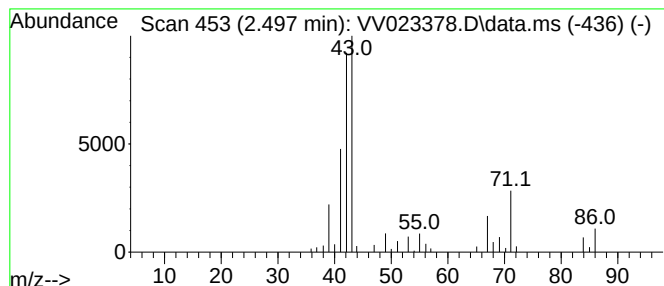
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.497	1.21 ug/L	71067	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	83
2			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	80
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	58
4			1,2,4,5-Tetrazine, 1,4-dihydro-3...	112	C4H8N4	037454-64-1	25
5			2-Oxetanone, 4-methylene-	84	C4H4O2	000674-82-8	25



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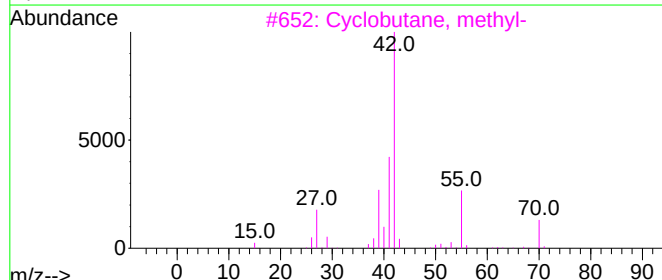
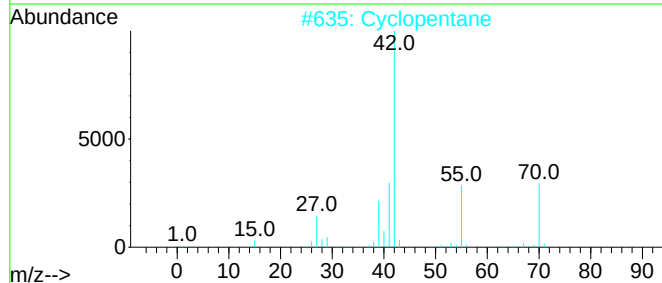
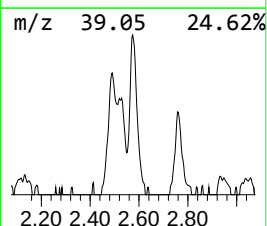
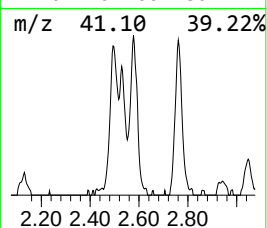
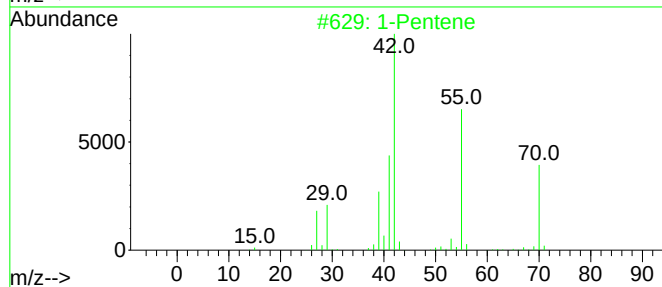
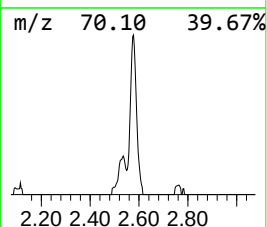
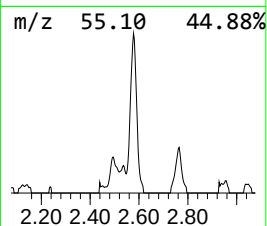
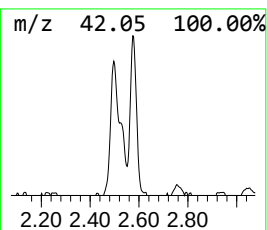
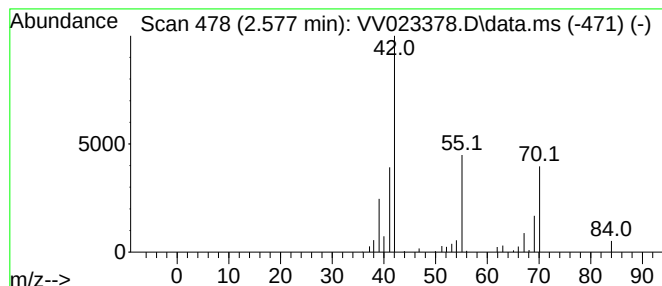
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.577	0.86 ug/L	50904	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene	70	C5H10	000109-67-1	64
2	Cyclopentane	70	C5H10	000287-92-3	64
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	64
4	3-Buten-1-ol	72	C4H8O	000627-27-0	23
5	Cyclobutanone	70	C4H6O	001191-95-3	9



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GB871DL

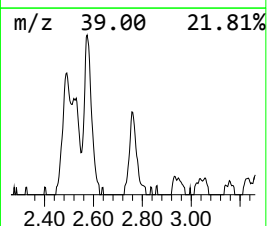
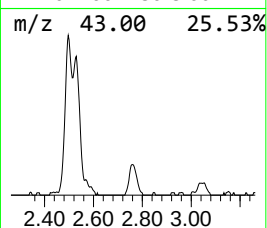
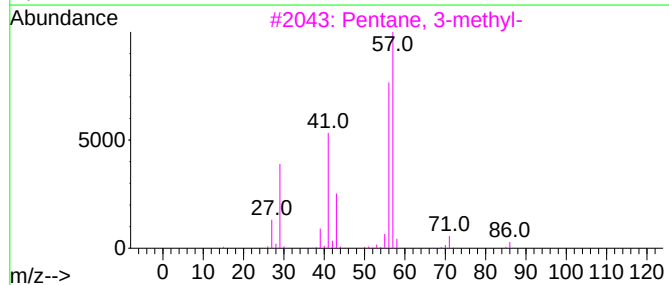
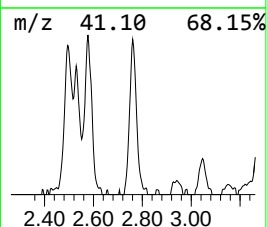
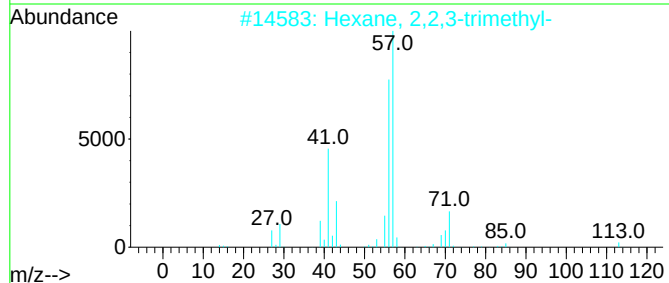
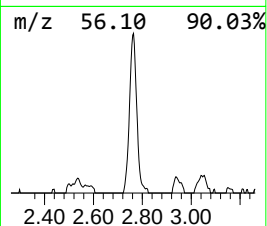
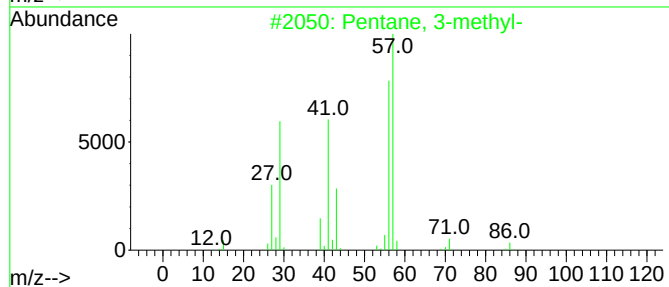
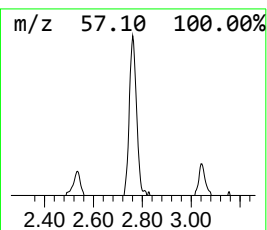
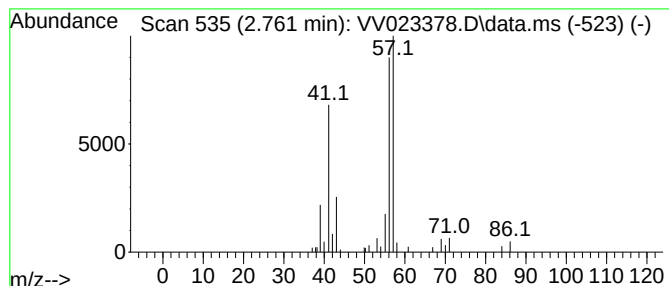
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.761	0.72 ug/L	42167	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	80
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	72
3			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
4			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
5			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
Data File : VV023378.D
Acq On : 11 Nov 2021 11:06
Operator : SY/MD
Sample : M4558-09DL 100X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GB871DL

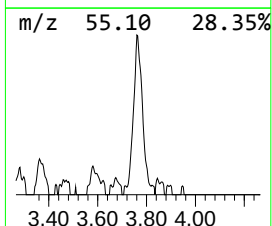
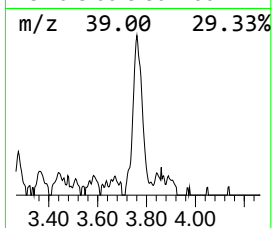
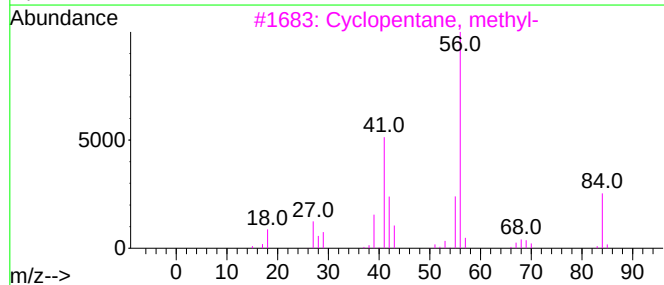
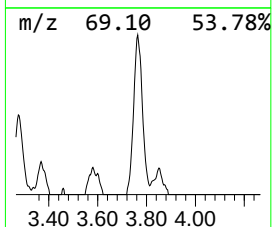
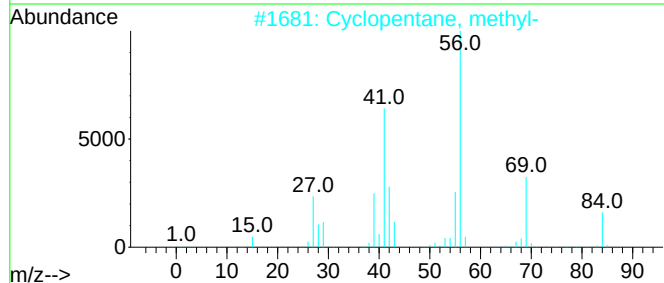
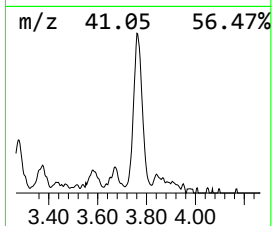
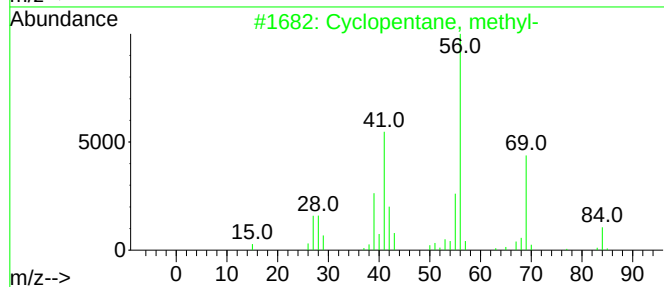
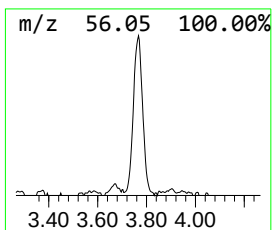
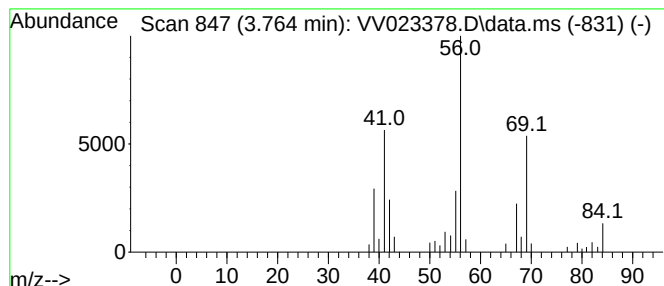
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Cyclic3.764 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.764	1.26 ug/L	74139	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	81
2			Cyclopentane, methyl-	84	C6H12	000096-37-7	64
3			Cyclopentane, methyl-	84	C6H12	000096-37-7	53
4			Cyclopentane, methyl-	84	C6H12	000096-37-7	53
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
Data File : VV023378.D
Acq On : 11 Nov 2021 11:06
Operator : SY/MD
Sample : M4558-09DL 100X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

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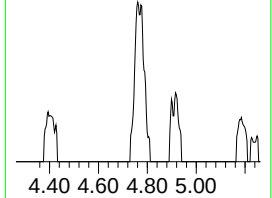
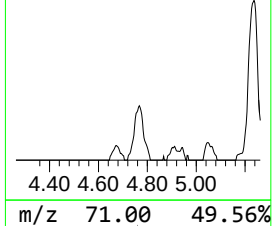
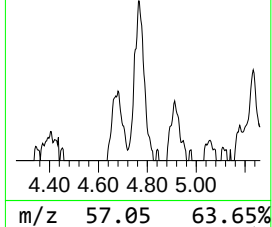
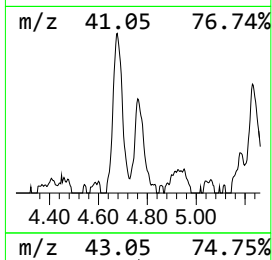
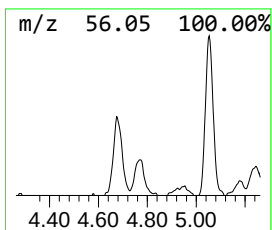
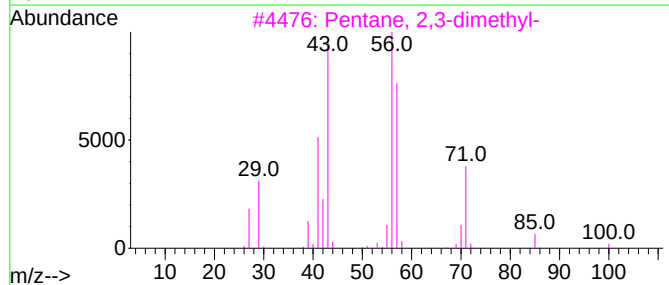
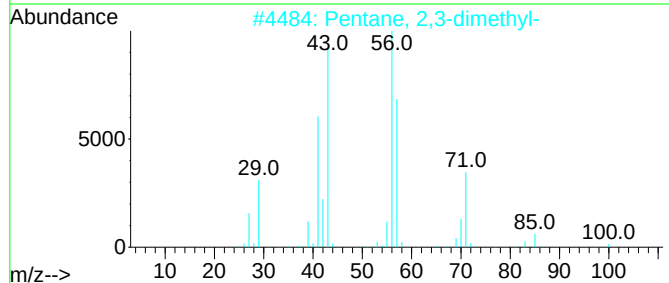
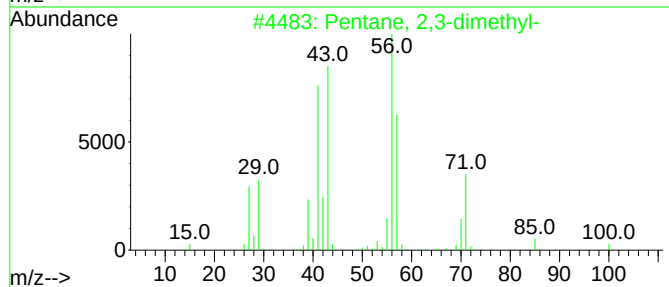
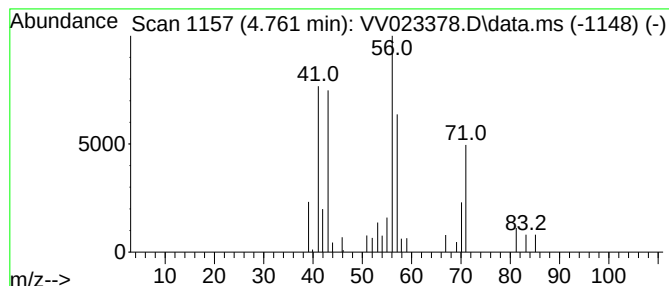
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.761	0.50 ug/L	29733	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	53
2			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	50
3			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	42
4			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	42
5			1-Hexene, 5-methyl-	98	C7H14	003524-73-0	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
Data File : VV023378.D
Acq On : 11 Nov 2021 11:06
Operator : SY/MD
Sample : M4558-09DL 100X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

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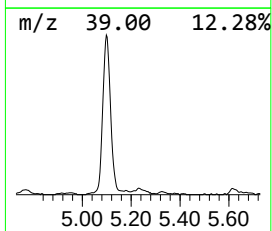
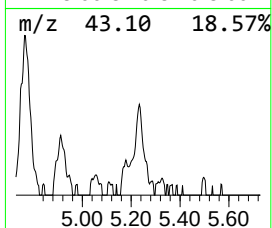
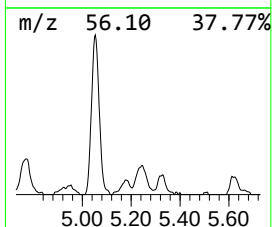
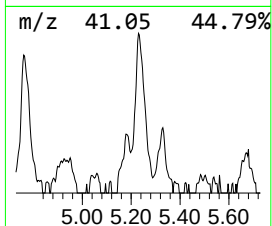
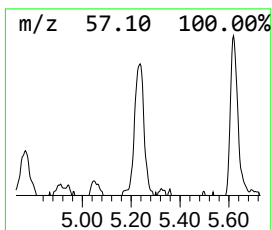
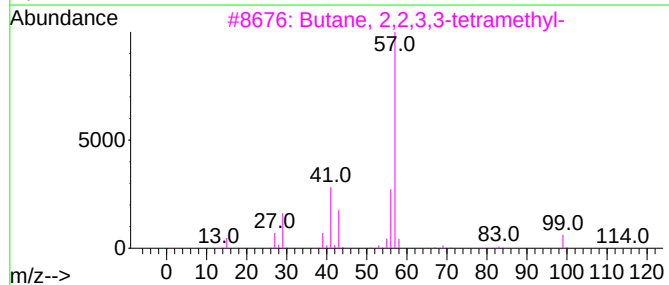
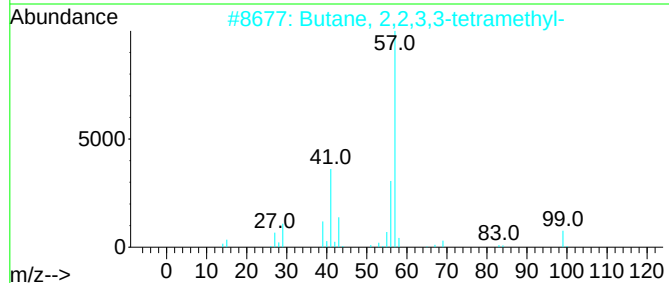
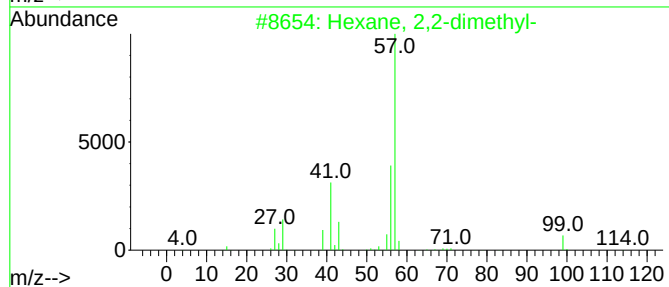
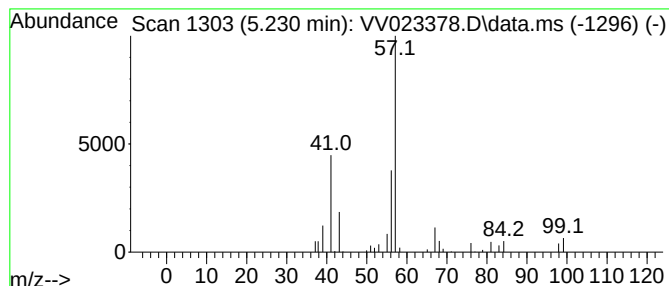
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.230	0.64 ug/L	37481	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	32
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	25
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	23
4		2-Propenal	56	C3H4O	000107-02-8	9
5		Methane, isocyanato-	57	C2H3NO	000624-83-9	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
Data File : VV023378.D
Acq On : 11 Nov 2021 11:06
Operator : SY/MD
Sample : M4558-09DL 100X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

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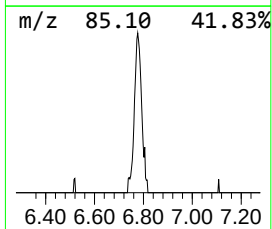
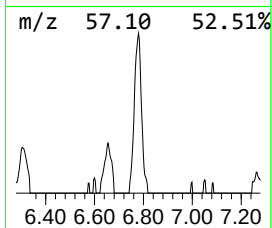
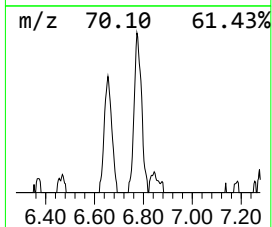
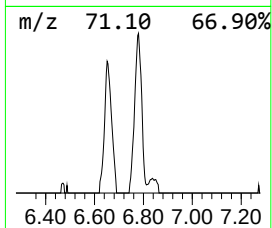
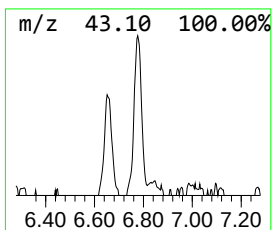
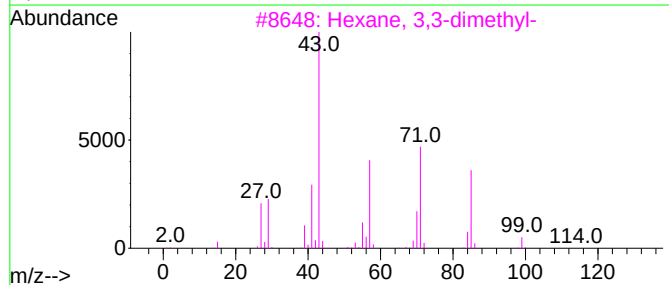
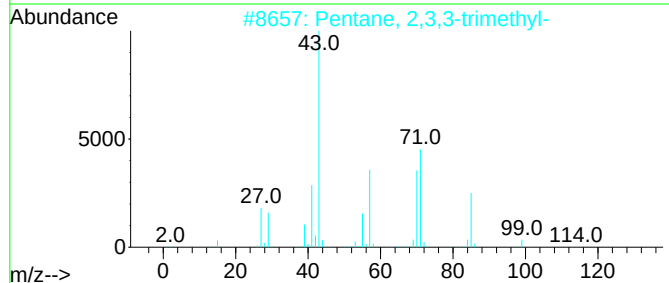
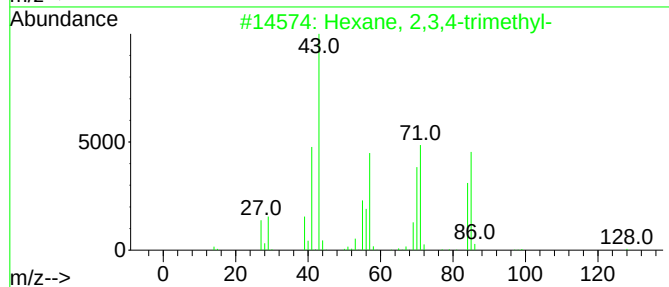
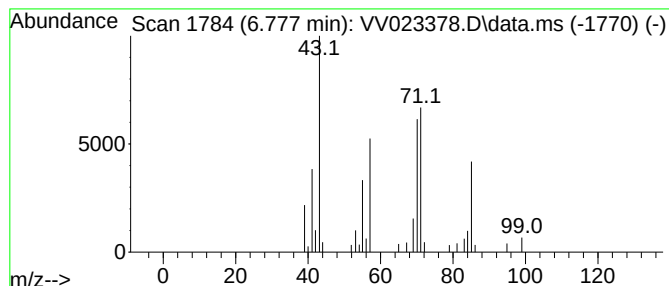
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.777	0.63 ug/L	37131	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	72
2		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
Data File : VV023378.D
Acq On : 11 Nov 2021 11:06
Operator : SY/MD
Sample : M4558-09DL 100X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

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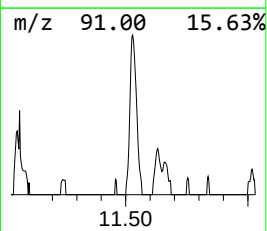
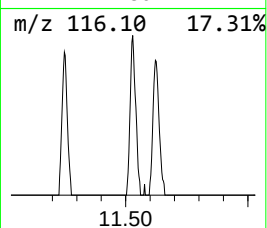
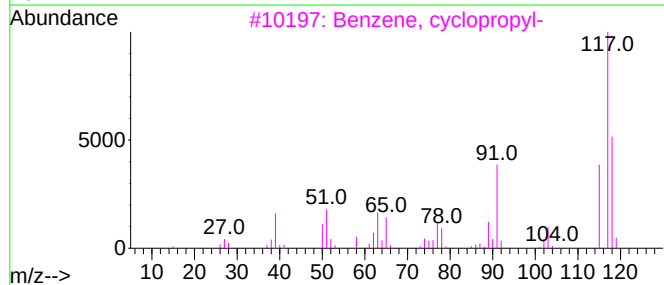
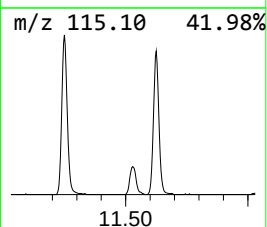
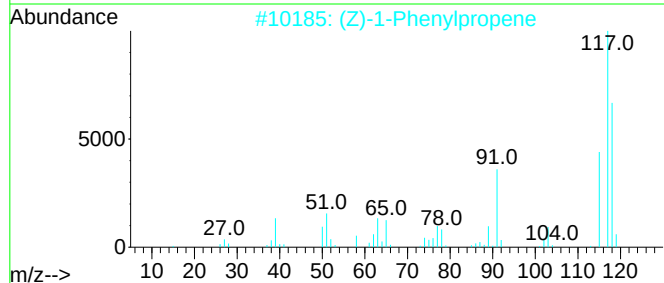
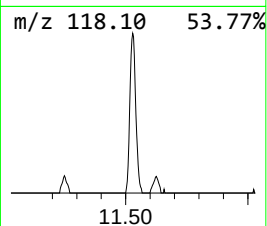
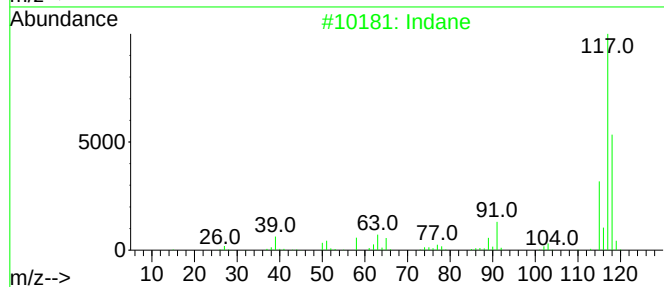
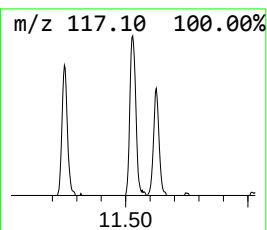
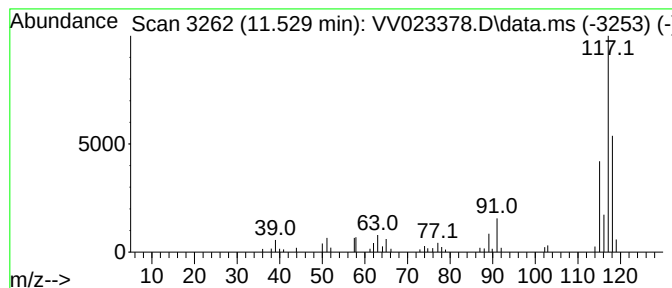
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	0.79 ug/L	52344	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	76
2			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	70
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5			3-Bromo-1-phenyl-1-propene	196	C9H9Br	004392-24-9	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\
Data File : VV023378.D
Acq On : 11 Nov 2021 11:06
Operator : SY/MD
Sample : M4558-09DL 100X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 65 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GB871DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

					--Internal Standard--			
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(DEL) Alkane: S...	1.809	1.1	ug/L	62238	1	5.619	294546	5.0
(DEL) Alkane: C...	2.024	1.1	ug/L	63753	1	5.619	294546	5.0
(DEL) Alkane: S...	2.497	1.2	ug/L	71067	1	5.619	294546	5.0
(DEL) Alkane: S...	2.577	0.9	ug/L	50904	1	5.619	294546	5.0
(DEL) Alkane: S...	2.761	0.7	ug/L	42167	1	5.619	294546	5.0
(DEL) Alkane: C...	3.764	1.3	ug/L	74139	1	5.619	294546	5.0
(DEL) Alkane: S...	4.761	0.5	ug/L	29733	1	5.619	294546	5.0
(DEL) Alkane: S...	5.230	0.6	ug/L	37481	1	5.619	294546	5.0
(DEL) Alkane: S...	6.777	0.6	ug/L	37131	1	5.619	294546	5.0
Indane	11.529	0.8	ug/L	52344	3	11.249	329811	5.0