

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
 Data File : VV025541.D
 Acq On : 14 Apr 2022 13:19
 Operator : SY/MD
 Sample : N2356-09DL 50X
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GBDS7DL

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: VV025541.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.220	49	56	64	rBV	31976	30808	1.93%	0.439%
2	1.310	76	84	95	rBV	316379	314699	19.75%	4.482%
3	1.571	157	165	176	rBV	87841	100561	6.31%	1.432%
4	1.638	176	186	202	rVB	390530	483230	30.33%	6.883%
5	1.812	231	240	254	rVB2	52833	74469	4.67%	1.061%
6	1.905	262	269	280	rVB	23736	30264	1.90%	0.431%
7	2.027	300	307	320	rVV	55838	79989	5.02%	1.139%
8	2.111	323	333	358	rVB	157742	251204	15.77%	3.578%
9	2.503	440	455	461	rBV3	41770	91989	5.77%	1.310%
10	2.580	472	479	497	rVB	34742	66767	4.19%	0.951%
11	2.767	523	537	553	rBV	34452	69937	4.39%	0.996%
12	3.677	804	820	832	rBV4	7276	18614	1.17%	0.265%
13	3.767	832	848	868	rVV2	38229	94097	5.91%	1.340%
14	3.908	880	892	919	rBV2	44112	149678	9.39%	2.132%
15	4.352	1011	1030	1057	rBV2	81956	217828	13.67%	3.103%
16	4.680	1117	1132	1145	rBV4	32051	79213	4.97%	1.128%
17	4.767	1147	1159	1180	rVB5	23734	61208	3.84%	0.872%
18	5.053	1230	1248	1253	rBV2	196259	427455	26.83%	6.089%
19	5.101	1253	1263	1282	rVB	735831	1593255	100.00%	22.694%
20	5.236	1295	1305	1322	rVB2	30275	83922	5.27%	1.195%
21	5.330	1324	1334	1346	rVB9	10839	24898	1.56%	0.355%
22	5.619	1413	1424	1457	rBV	138627	290649	18.24%	4.140%
23	6.072	1547	1565	1595	rBV	106322	246653	15.48%	3.513%
24	6.654	1727	1746	1762	rBV2	20641	44291	2.78%	0.631%
25	6.776	1770	1784	1799	rBV3	37365	81906	5.14%	1.167%
26	6.989	1838	1850	1875	rBV2	44248	87282	5.48%	1.243%
27	7.317	1941	1952	1969	rBV	215002	373696	23.45%	5.323%
28	7.629	2036	2049	2066	rBV3	22521	45147	2.83%	0.643%
29	8.092	2183	2193	2232	rBV	153720	323936	20.33%	4.614%
30	8.854	2419	2430	2451	rBV	210044	347681	21.82%	4.952%
31	9.143	2510	2520	2535	rBV	11832	21910	1.38%	0.312%
32	10.214	2843	2853	2873	rVB	62342	100780	6.33%	1.435%
33	11.249	3164	3175	3194	rBV	199021	314292	19.73%	4.477%
34	11.529	3251	3262	3277	rBV	41754	69972	4.39%	0.997%
35	11.622	3280	3291	3312	rVV	190524	307493	19.30%	4.380%
36	12.114	3434	3444	3454	rBV	13962	20869	1.31%	0.297%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

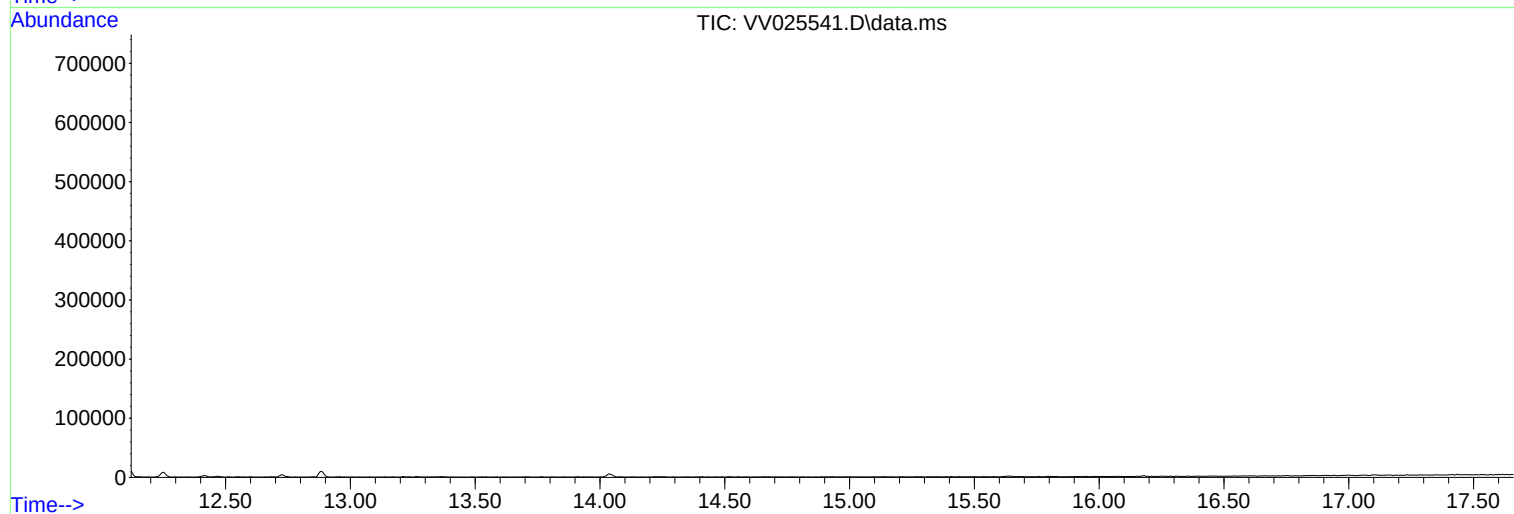
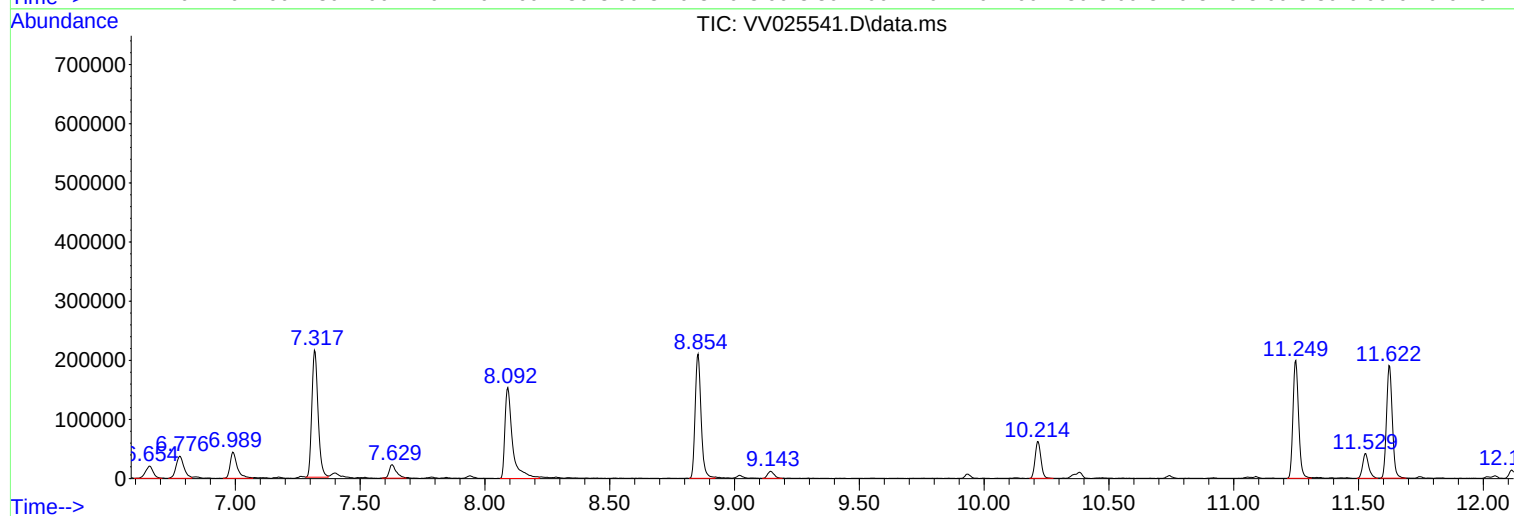
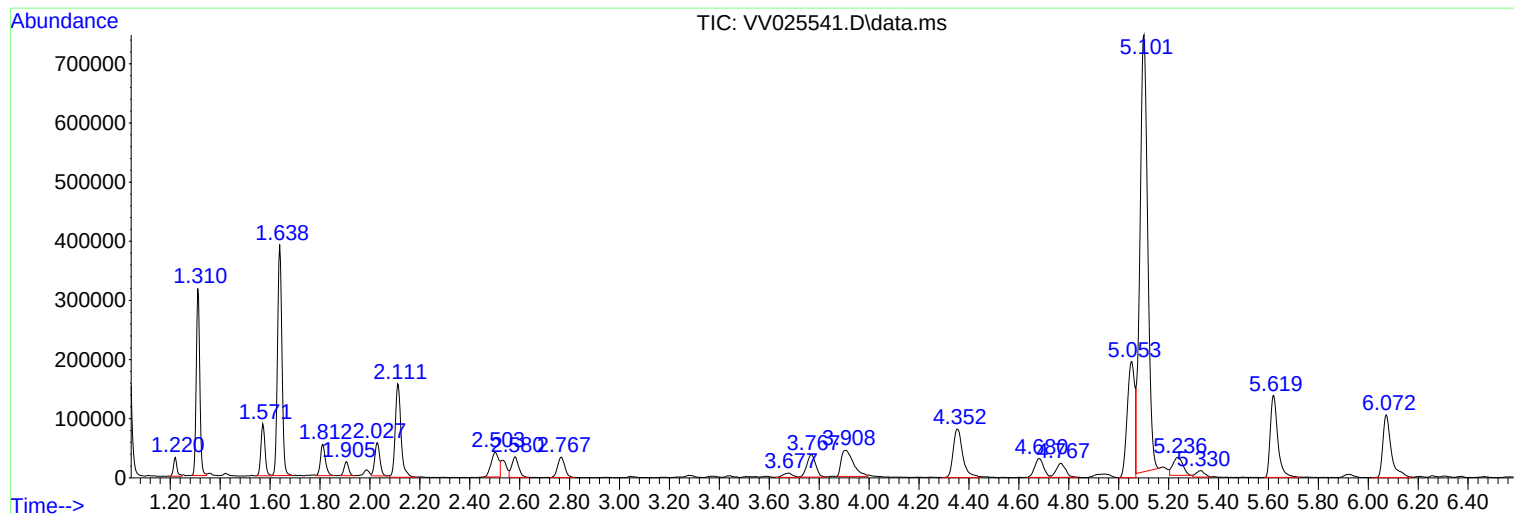
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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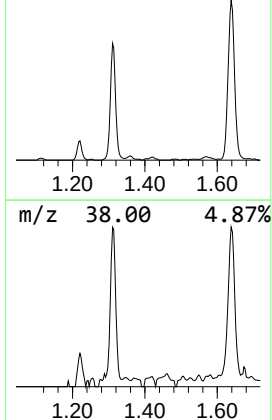
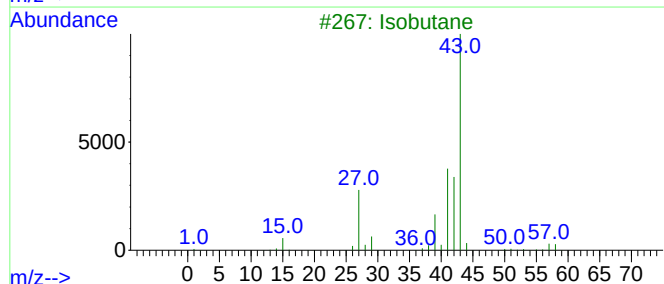
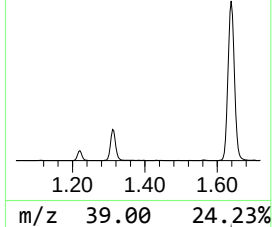
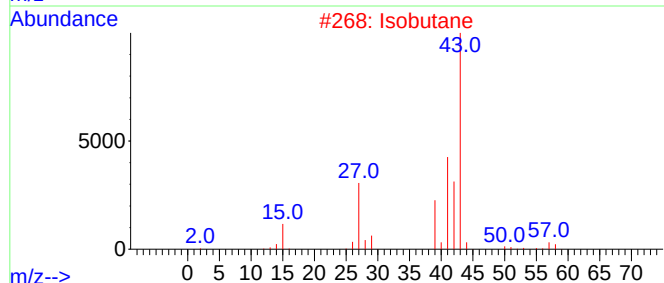
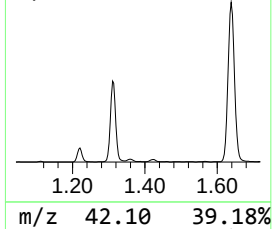
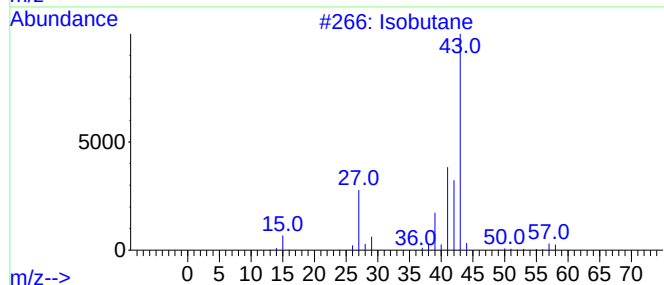
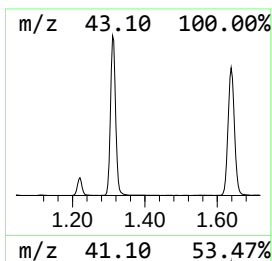
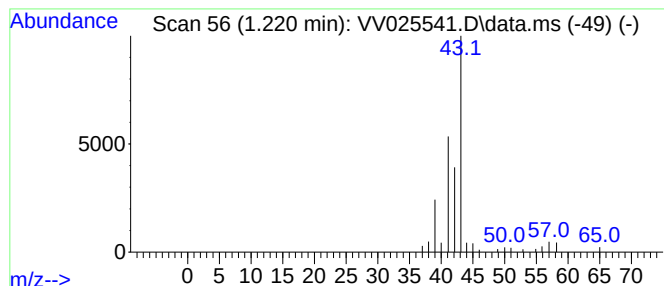
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.220	0.53 ug/L	30808	1,4-Difluorobenzene	5.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isobutane	58	C4H10	000075-28-5	72
2		Isobutane	58	C4H10	000075-28-5	72
3		Isobutane	58	C4H10	000075-28-5	72
4		Isobutane	58	C4H10	000075-28-5	72
5		1-[3-(4-Bromophenyl)-2-thioureid...	560	C21H25BrN2O9S	094273-12-8	10



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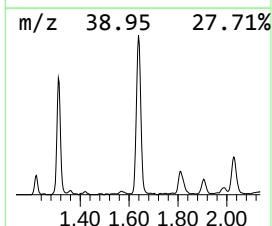
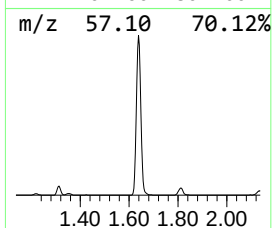
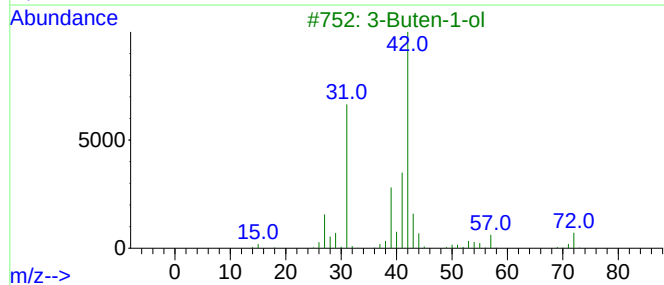
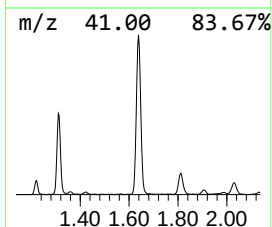
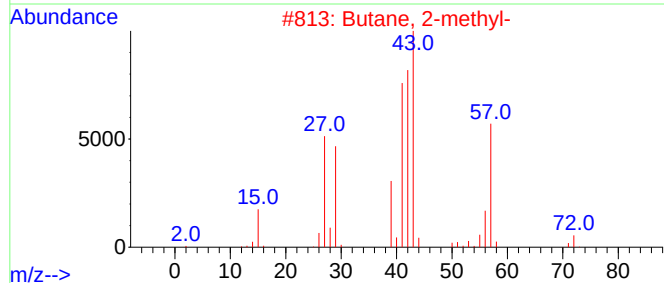
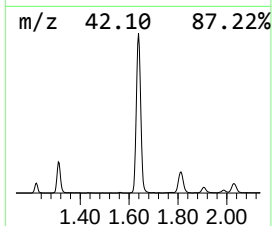
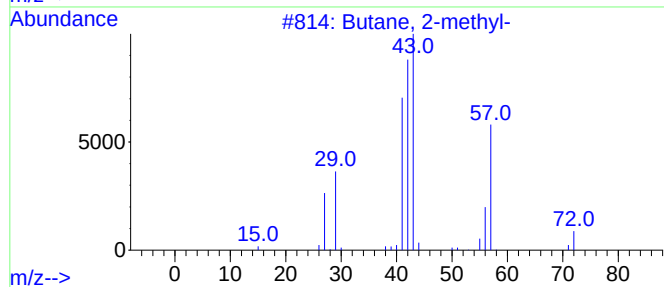
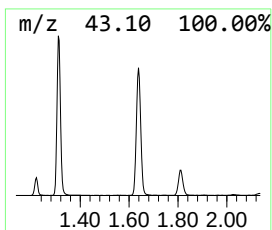
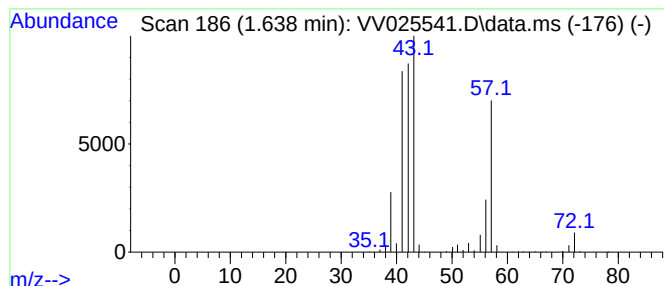
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.638	8.31 ug/L	483230	1,4-Difluorobenzene	5.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	87
2		Butane, 2-methyl-	72	C5H12	000078-78-4	86
3		3-Buten-1-ol	72	C4H8O	000627-27-0	47
4		Pentane	72	C5H12	000109-66-0	36
5		Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	27



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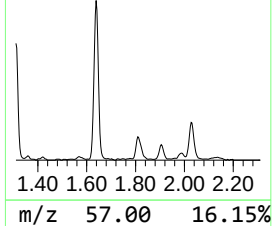
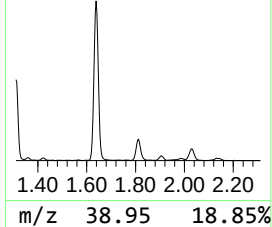
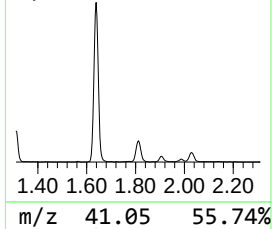
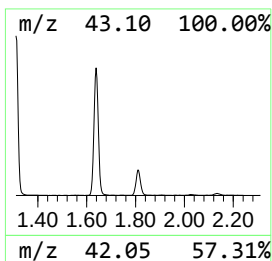
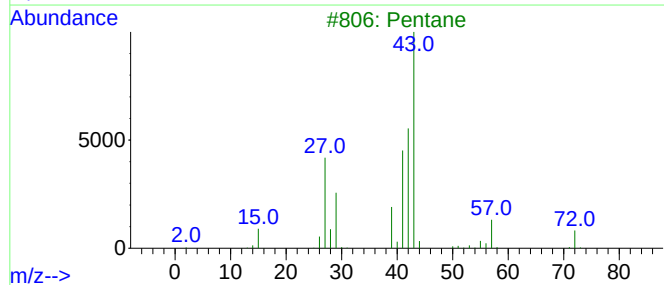
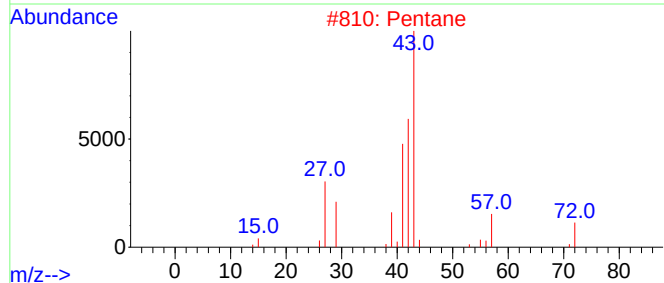
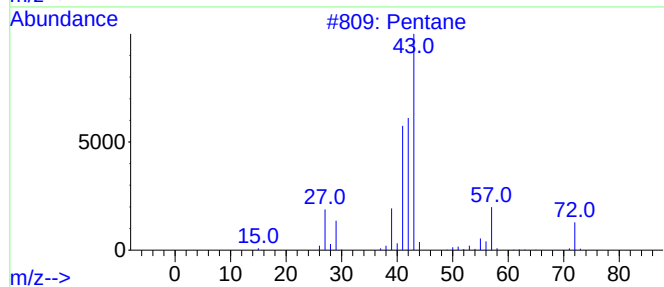
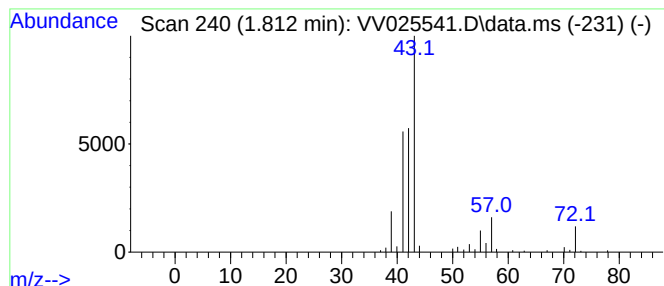
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.812	1.28 ug/L	74469	1,4-Difluorobenzene	5.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane			72	C5H12	000109-66-0	87
2	Pentane			72	C5H12	000109-66-0	86
3	Pentane			72	C5H12	000109-66-0	86
4	Pentane			72	C5H12	000109-66-0	80
5	Pentane			72	C5H12	000109-66-0	59



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ALS Vial : 11 Sample Multiplier: 1

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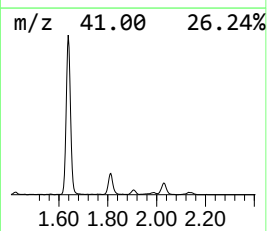
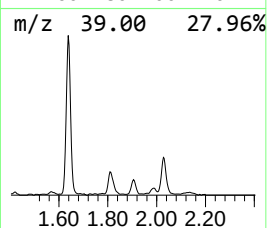
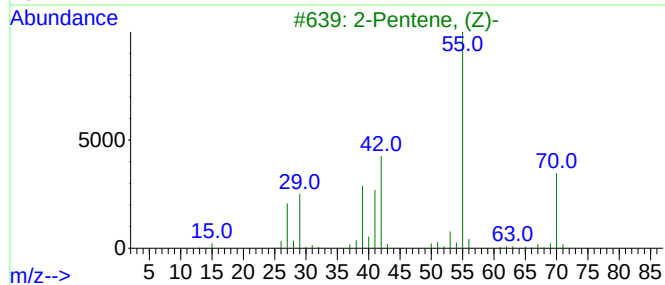
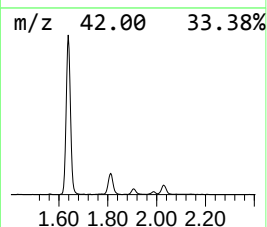
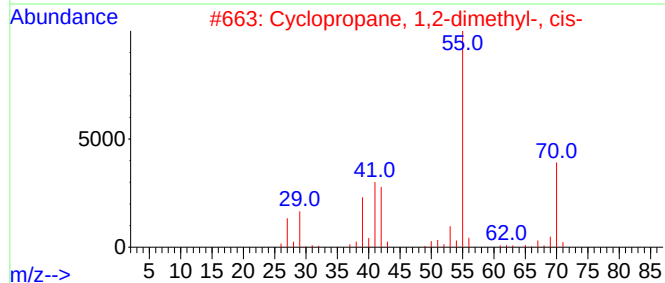
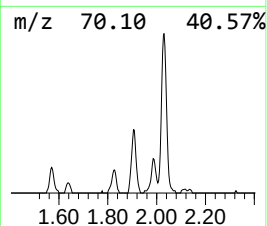
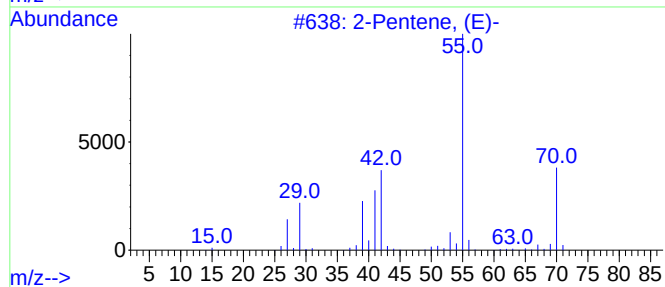
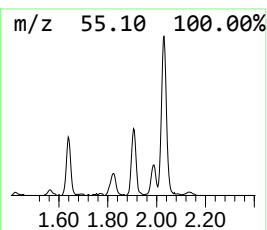
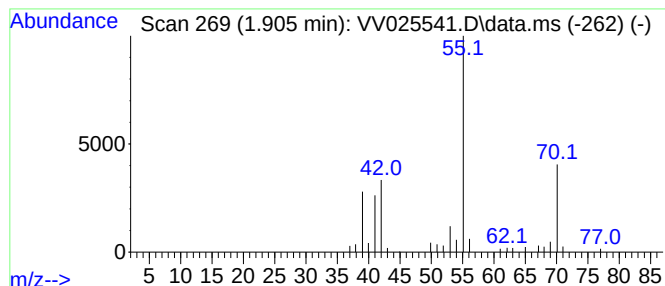
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.905	0.52 ug/L	30264	1,4-Difluorobenzene	5.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (E)-			70	C5H10	000646-04-8	91
2	Cyclopropane, 1,2-dimethyl-, cis-			70	C5H10	000930-18-7	91
3	2-Pentene, (Z)-			70	C5H10	000627-20-3	90
4	1-Butene, 3-methyl-			70	C5H10	000563-45-1	90
5	2-Methyl-1-butene			70	C5H10	000563-46-2	90



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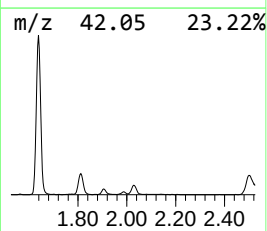
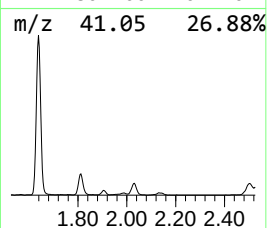
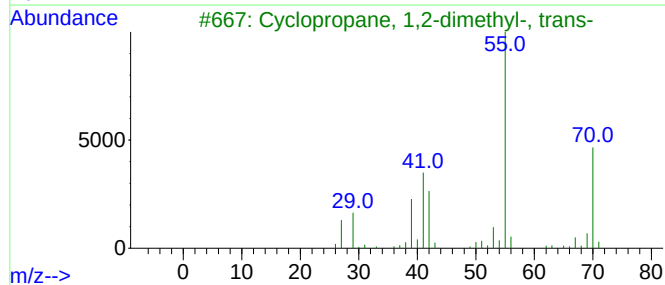
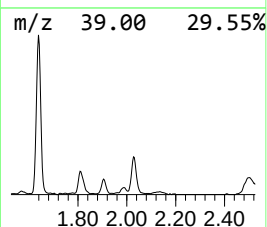
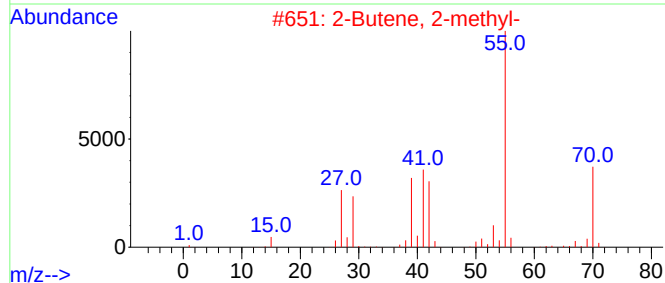
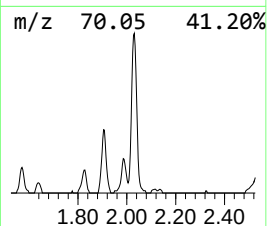
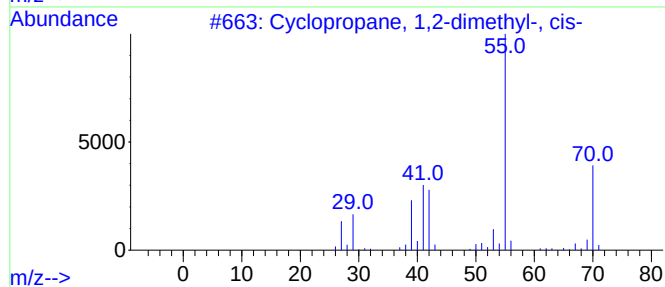
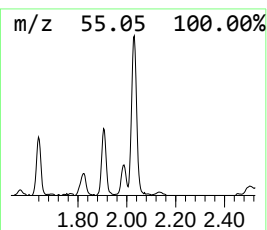
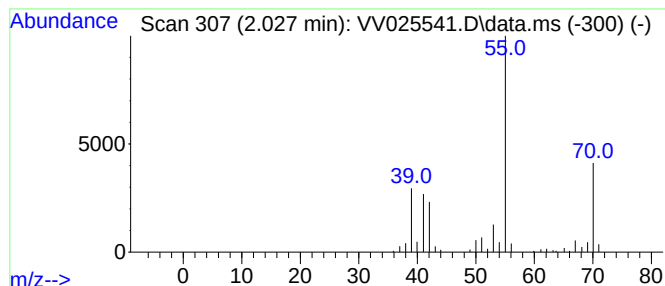
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Cyclic2.027 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.027	1.38 ug/L	79989	1,4-Difluorobenzene	5.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	87
2			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86
3			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	86
4			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86
5			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80



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Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS7DL

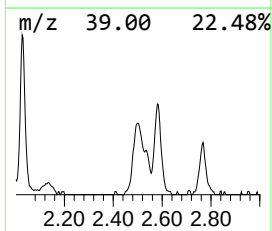
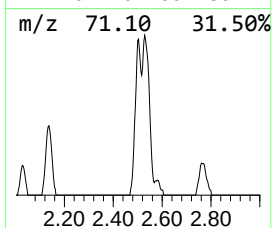
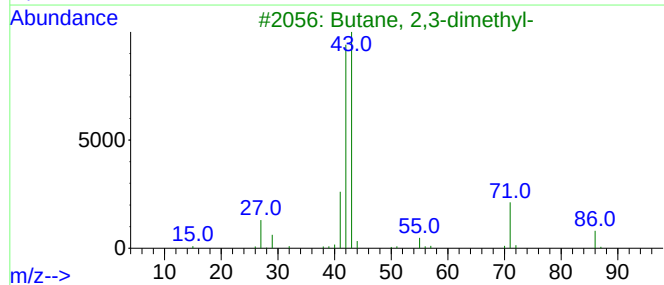
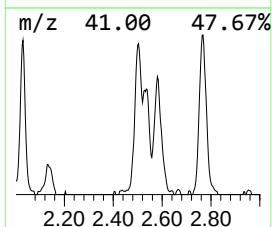
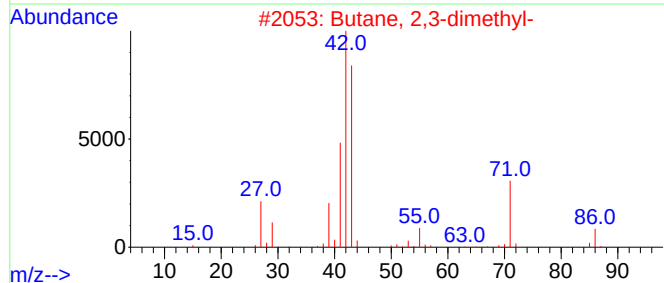
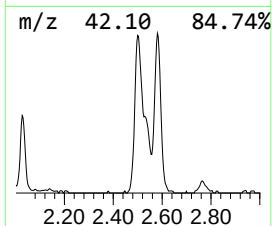
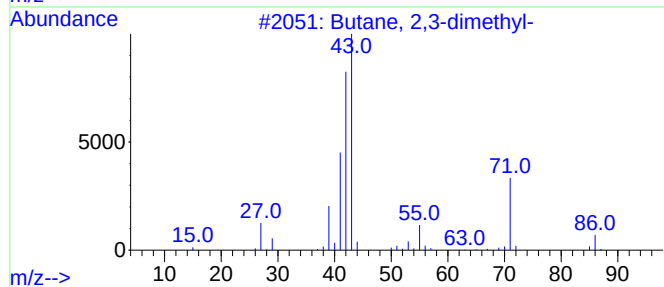
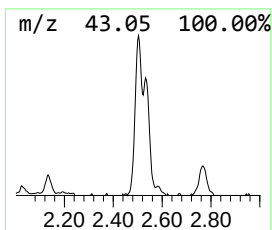
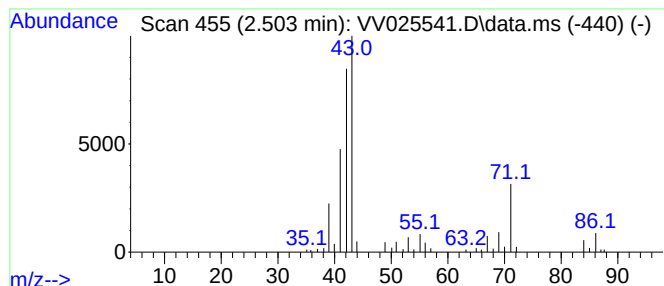
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.503	1.58 ug/L	91989	1,4-Difluorobenzene	5.622

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	74
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	64
4			Pentane, 2-methyl-	86	C6H14	000107-83-5	53
5			Pentane, 2-methyl-	86	C6H14	000107-83-5	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS7DL

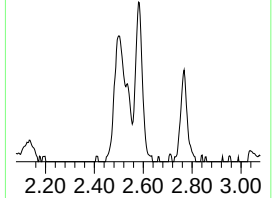
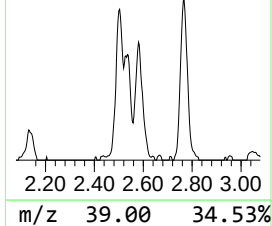
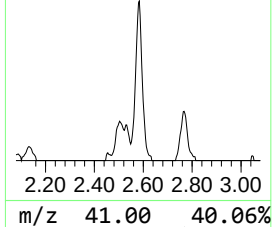
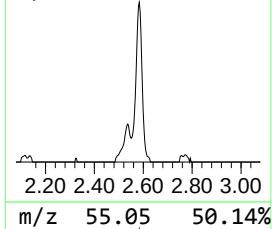
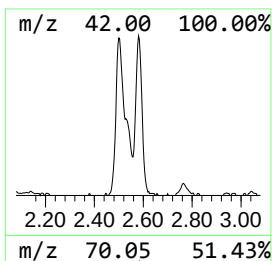
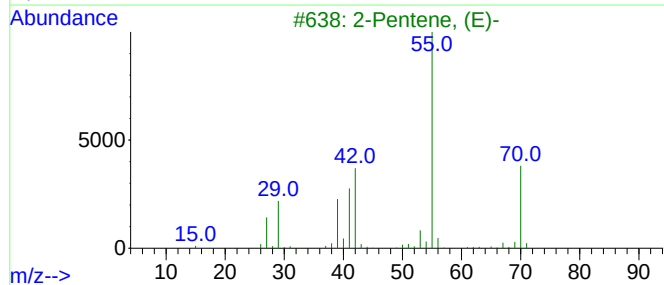
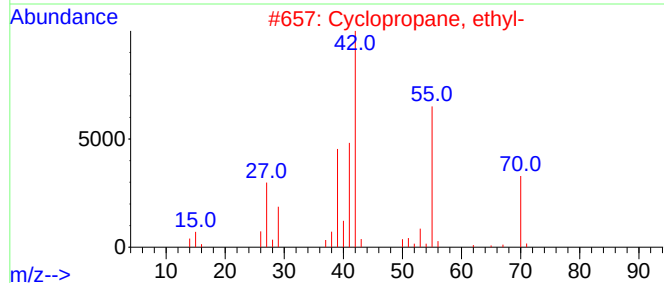
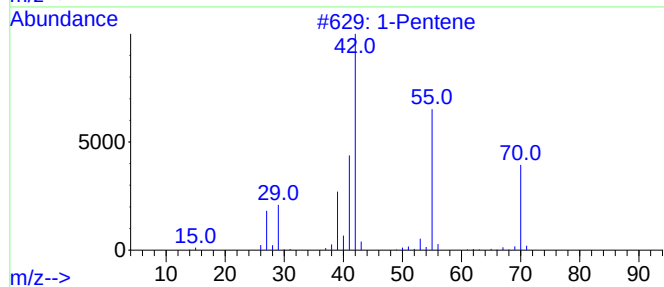
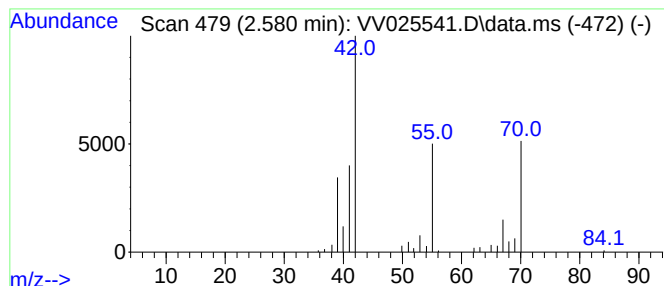
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.580	1.15 ug/L	66767	1,4-Difluorobenzene	5.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene	70	C5H10	000109-67-1	72
2		Cyclopropane, ethyl-	70	C5H10	001191-96-4	45
3		2-Pentene, (E)-	70	C5H10	000646-04-8	43
4		Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	43
5		2-Pentene	70	C5H10	000109-68-2	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS7DL

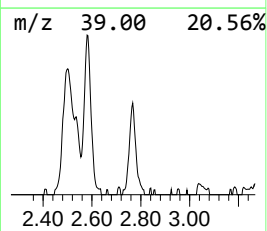
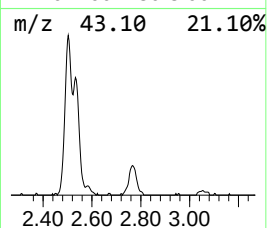
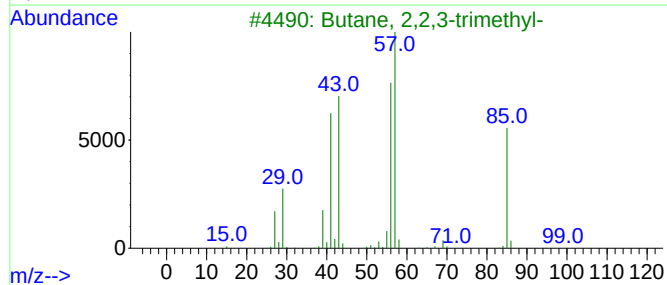
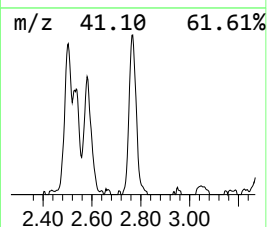
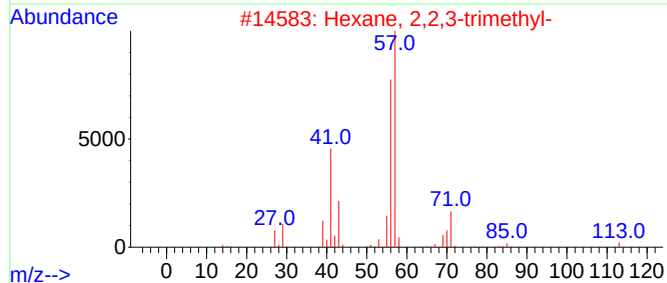
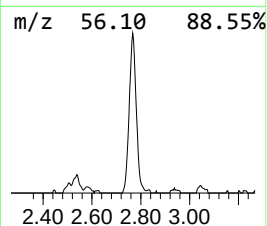
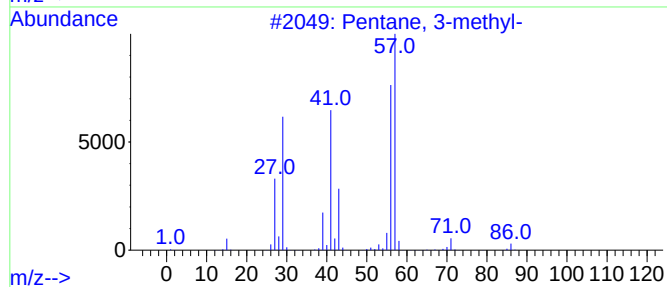
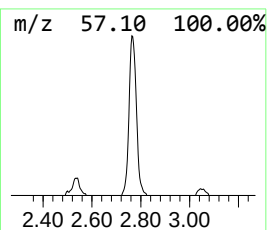
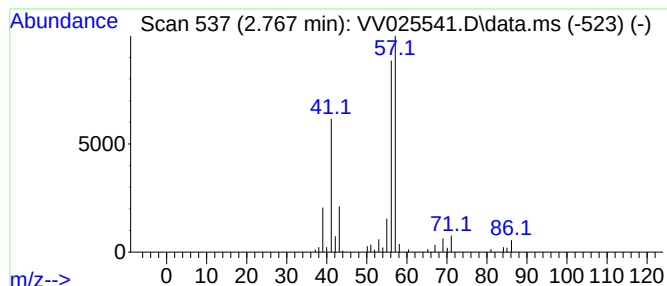
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.767	1.20 ug/L	69937	1,4-Difluorobenzene	5.622

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	78
3			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
4			Pentane, 3-methyl-	86	C6H14	000096-14-0	53
5			Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
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ALS Vial : 11 Sample Multiplier: 1

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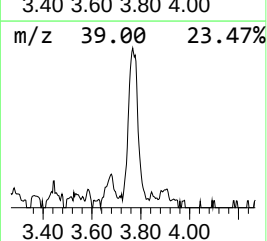
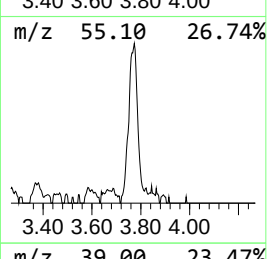
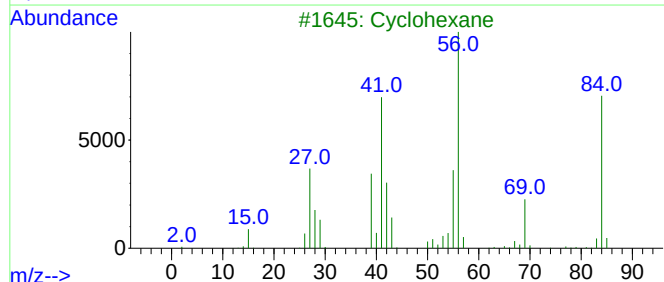
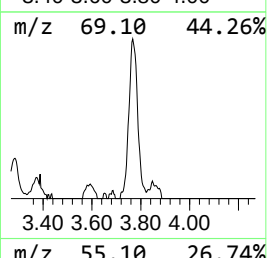
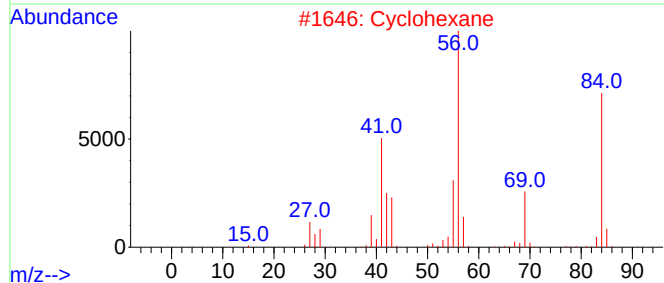
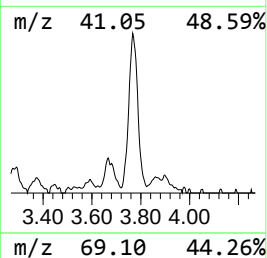
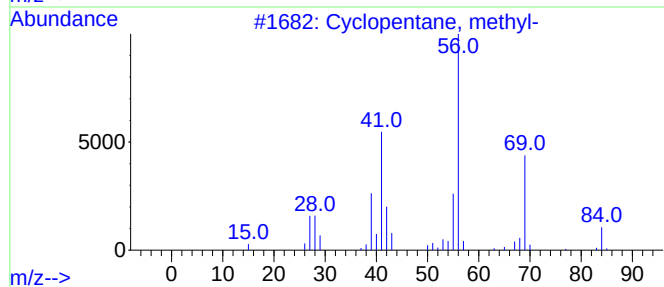
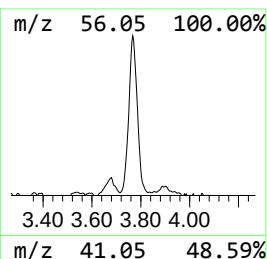
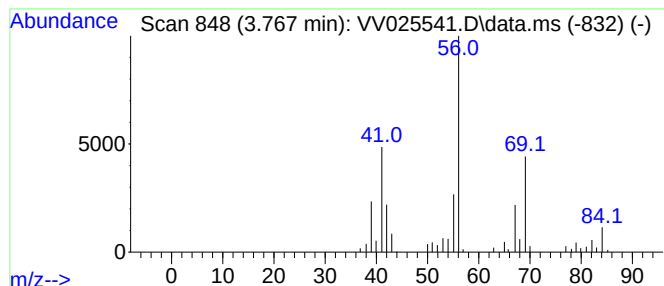
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 (DEL) Alkane: Cyclic3.767 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.767	1.62 ug/L	94097	1,4-Difluorobenzene	5.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	74
2			Cyclohexane	84	C6H12	000110-82-7	72
3			Cyclohexane	84	C6H12	000110-82-7	64
4			Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5			Cyclopentane, methyl-	84	C6H12	000096-37-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
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ALS Vial : 11 Sample Multiplier: 1

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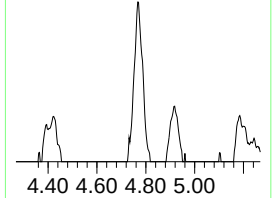
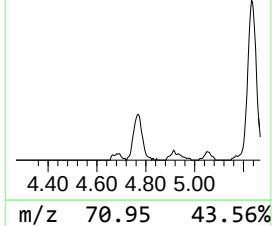
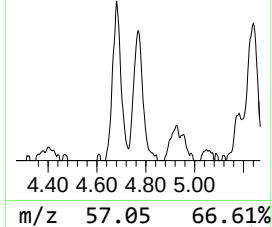
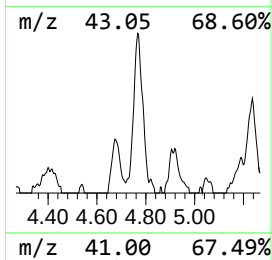
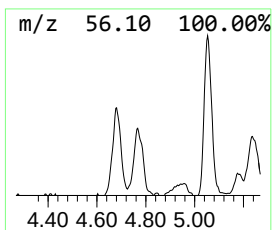
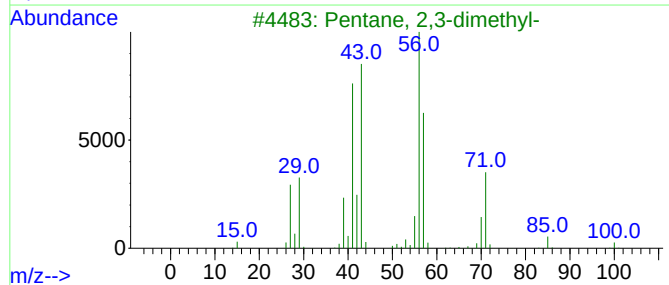
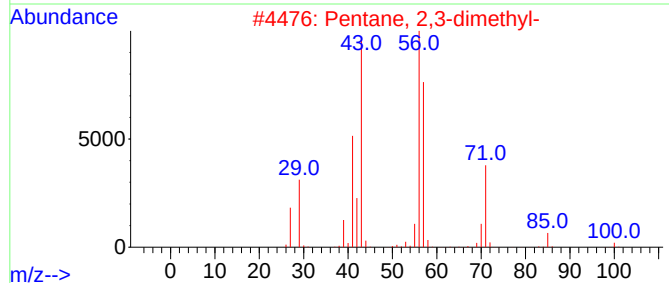
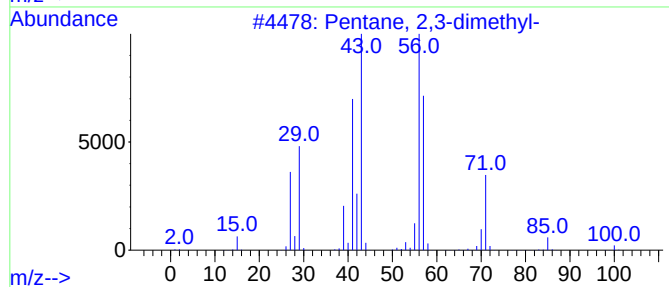
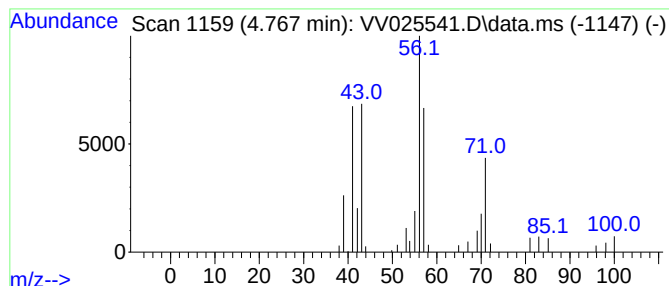
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.767	1.05 ug/L	61208	1,4-Difluorobenzene	5.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	78
2		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	64
3		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	64
4		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	56
5		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
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ALS Vial : 11 Sample Multiplier: 1

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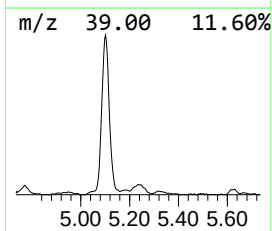
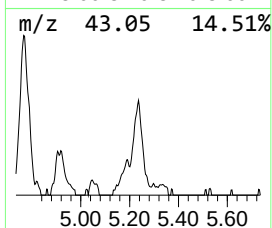
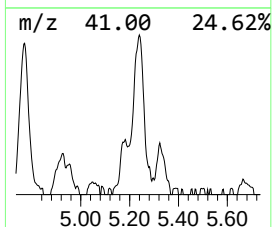
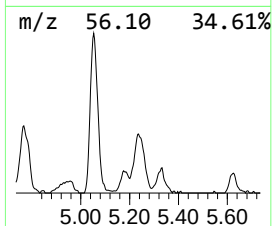
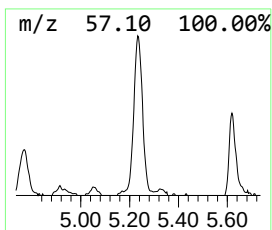
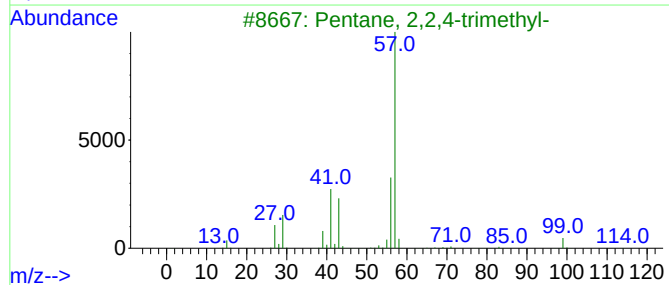
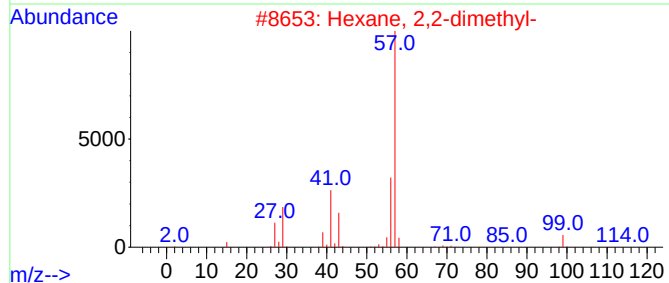
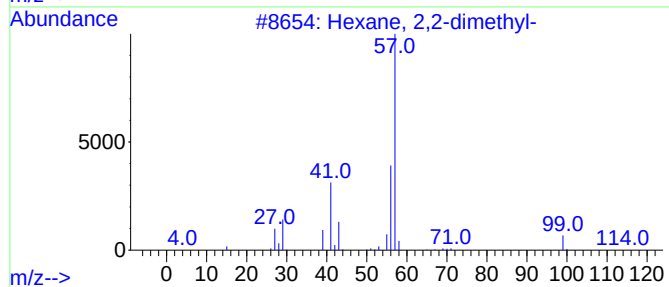
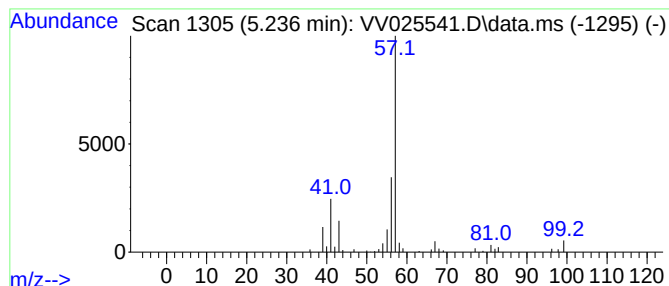
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.236	1.44 ug/L	83922	1,4-Difluorobenzene	5.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	45
4		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	45
5		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

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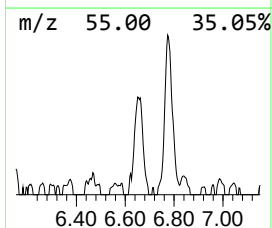
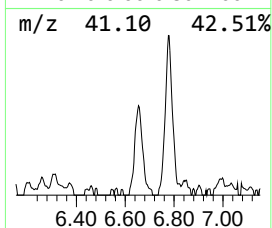
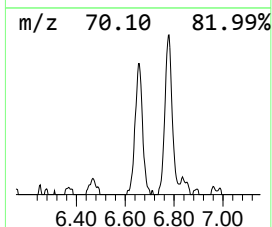
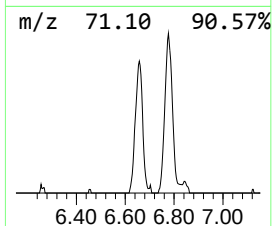
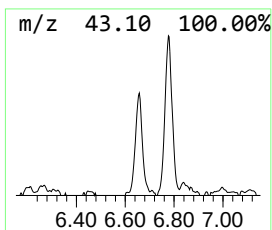
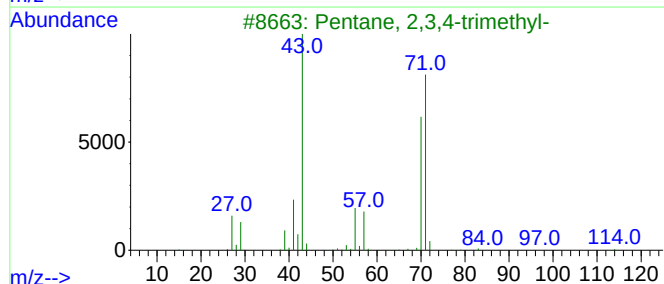
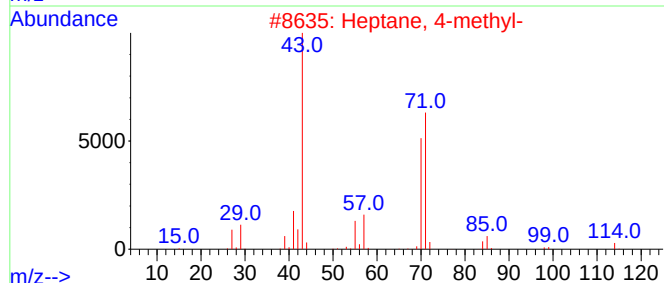
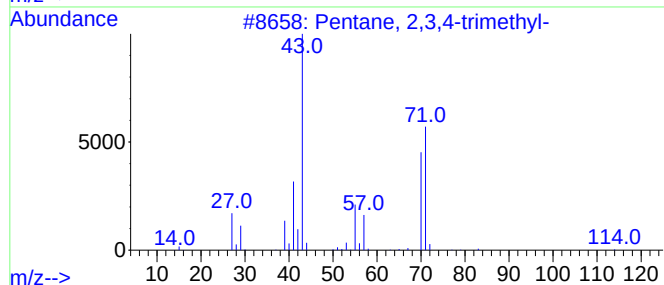
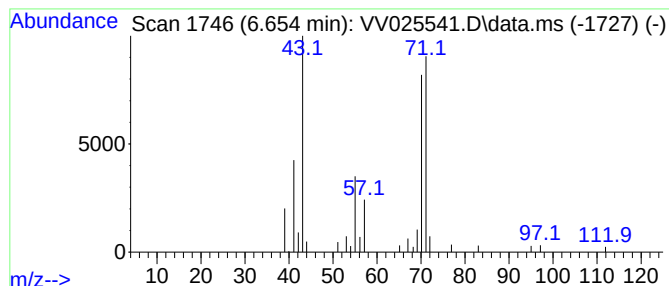
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.654	0.76 ug/L	44291	1,4-Difluorobenzene	5.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	86
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	78
3		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
4		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS7DL

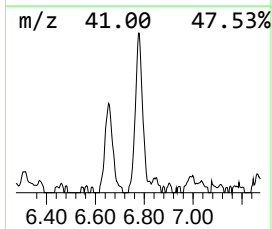
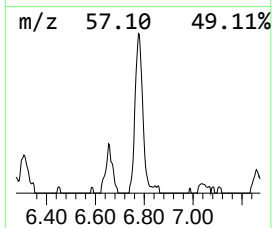
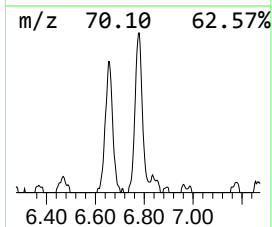
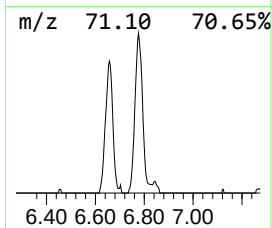
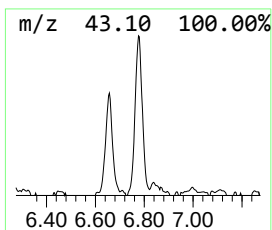
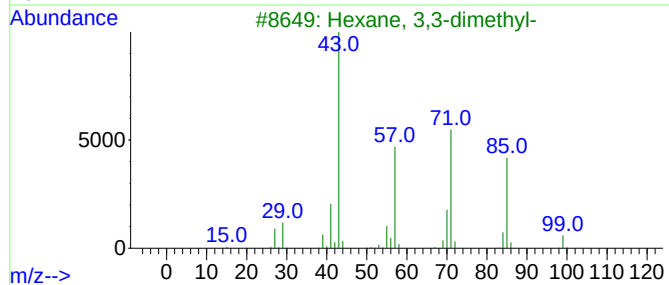
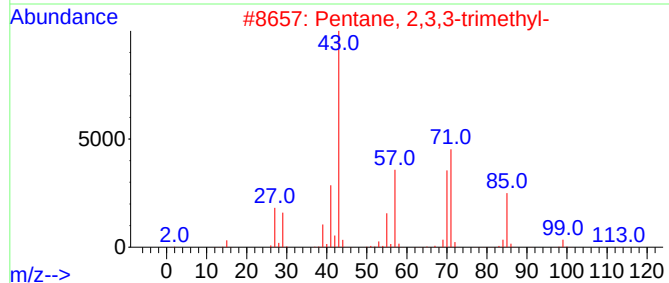
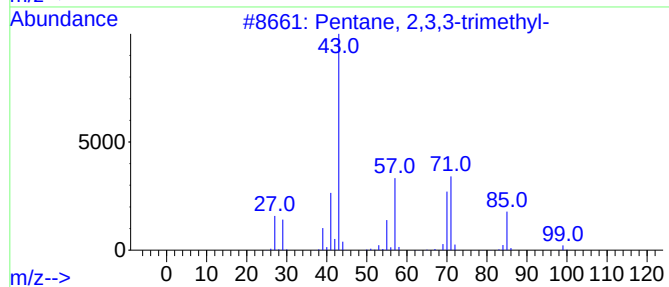
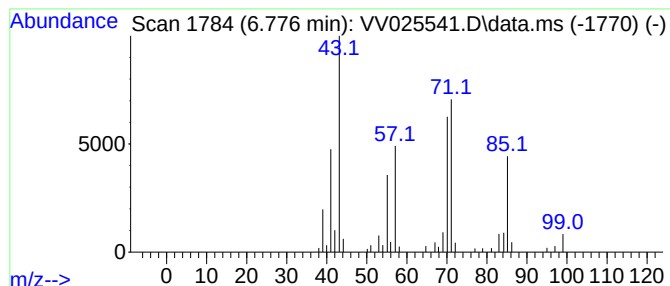
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.776	1.41 ug/L	81906	1,4-Difluorobenzene	5.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
3			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4			Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	59
5			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025541.D
Acq On : 14 Apr 2022 13:19
Operator : SY/MD
Sample : N2356-09DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS7DL

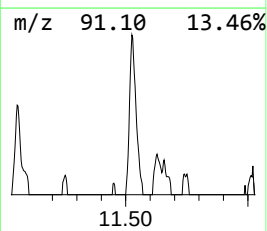
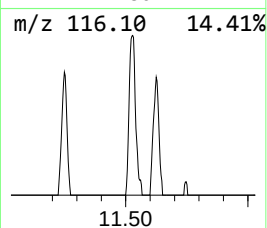
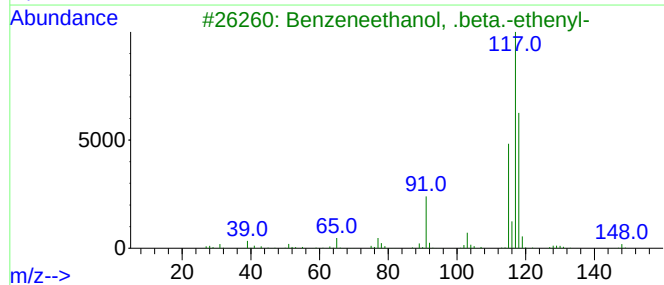
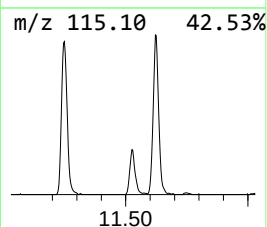
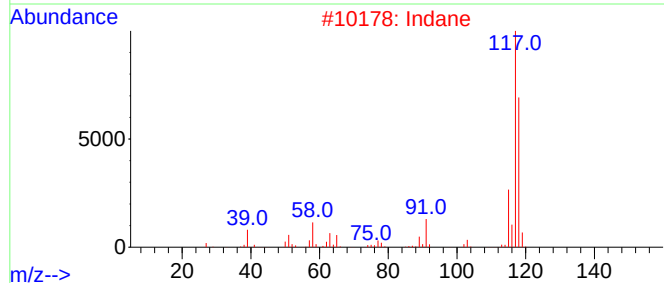
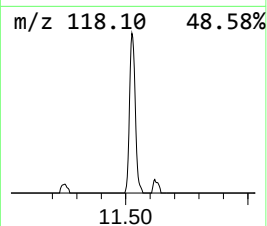
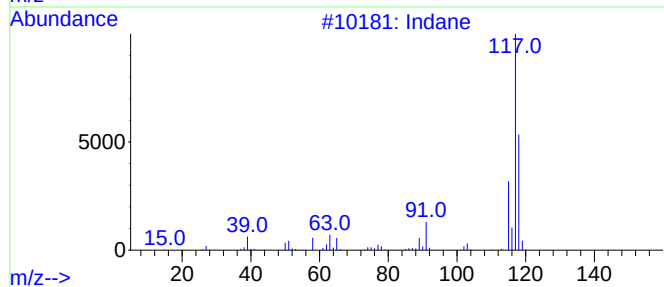
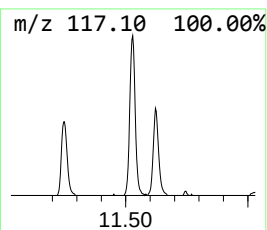
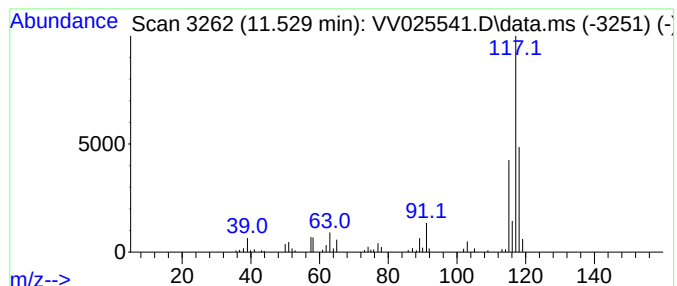
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Indane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	1.11 ug/L	69972	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	94
3			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	78
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
 Data File : VV025541.D
 Acq On : 14 Apr 2022 13:19
 Operator : SY/MD
 Sample : N2356-09DL 50X
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GBDS7DL

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...	1.220	0.5	ug/L	30808	1	5.622	290649	5.0	
(DEL) Alkane: S...	1.638	8.3	ug/L	483230	1	5.622	290649	5.0	
(DEL) Alkane: S...	1.812	1.3	ug/L	74469	1	5.622	290649	5.0	
(DEL) Alkane: S...	1.905	0.5	ug/L	30264	1	5.622	290649	5.0	
(DEL) Alkane: C...	2.027	1.4	ug/L	79989	1	5.622	290649	5.0	
(DEL) Alkane: S...	2.503	1.6	ug/L	91989	1	5.622	290649	5.0	
(DEL) Alkane: S...	2.580	1.1	ug/L	66767	1	5.622	290649	5.0	
(DEL) Alkane: S...	2.767	1.2	ug/L	69937	1	5.622	290649	5.0	
(DEL) Alkane: C...	3.767	1.6	ug/L	94097	1	5.622	290649	5.0	
(DEL) Alkane: S...	4.767	1.1	ug/L	61208	1	5.622	290649	5.0	
(DEL) Alkane: S...	5.236	1.4	ug/L	83922	1	5.622	290649	5.0	
(DEL) Alkane: S...	6.654	0.8	ug/L	44291	1	5.622	290649	5.0	
(DEL) Alkane: S...	6.776	1.4	ug/L	81906	1	5.622	290649	5.0	
Indane	11.529	1.1	ug/L	69972	3	11.249	314292	5.0	