

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
 Data File : VV025542.D
 Acq On : 14 Apr 2022 13:43
 Operator : SY/MD
 Sample : N2356-10DL 50X
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GBDS8DL

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: VV025542.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.217	49	55	68	rBV	32206	32575	2.18%	0.464%
2	1.310	76	84	94	rBV	319023	311002	20.78%	4.427%
3	1.571	157	165	173	rBV	89370	97767	6.53%	1.392%
4	1.638	176	186	199	rVV	397637	494215	33.02%	7.034%
5	1.809	231	239	254	rVB2	54610	79151	5.29%	1.127%
6	1.905	261	269	281	rVB	24433	33022	2.21%	0.470%
7	1.986	281	294	299	rVV	11219	18008	1.20%	0.256%
8	2.027	299	307	321	rVV	67075	96523	6.45%	1.374%
9	2.111	321	333	357	rVB	163143	255155	17.05%	3.632%
10	2.500	429	454	461	rBV4	44255	105401	7.04%	1.500%
11	2.580	472	479	497	rVB2	35875	66009	4.41%	0.940%
12	2.764	523	536	556	rBV2	36426	74819	5.00%	1.065%
13	3.670	807	818	833	rVB3	8017	19077	1.27%	0.272%
14	3.767	833	848	867	rBV2	38077	94720	6.33%	1.348%
15	3.905	880	891	923	rVB4	41606	137077	9.16%	1.951%
16	4.352	1013	1030	1061	rBV2	80567	212500	14.20%	3.025%
17	4.680	1114	1132	1146	rBV3	32388	82276	5.50%	1.171%
18	4.767	1147	1159	1182	rVB3	26645	68818	4.60%	0.979%
19	5.050	1232	1247	1252	rBV2	193444	399314	26.68%	5.683%
20	5.098	1253	1262	1281	rVB	693877	1496586	100.00%	21.301%
21	5.233	1296	1304	1323	rVB5	32329	87625	5.85%	1.247%
22	5.323	1323	1332	1346	rVB5	11010	24837	1.66%	0.354%
23	5.619	1412	1424	1446	rBV	149210	309277	20.67%	4.402%
24	6.072	1552	1565	1596	rBV2	100397	240359	16.06%	3.421%
25	6.654	1730	1746	1760	rBV	23836	49142	3.28%	0.699%
26	6.776	1769	1784	1799	rBV2	41525	89030	5.95%	1.267%
27	6.989	1839	1850	1866	rBV2	41996	79544	5.32%	1.132%
28	7.317	1942	1952	1970	rBV	215091	373889	24.98%	5.322%
29	7.625	2040	2048	2070	rBV2	22461	43897	2.93%	0.625%
30	8.092	2183	2193	2223	rBV	143659	296521	19.81%	4.220%
31	8.854	2419	2430	2445	rBV	231631	377129	25.20%	5.368%
32	9.146	2508	2521	2534	rBV2	11615	20518	1.37%	0.292%
33	10.217	2842	2854	2868	rBV	60284	96088	6.42%	1.368%
34	11.249	3163	3175	3197	rBV	219233	343757	22.97%	4.893%
35	11.529	3251	3262	3280	rBV2	40140	70801	4.73%	1.008%
36	11.622	3280	3291	3314	rVV	192095	311991	20.85%	4.441%

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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

37	12.114	3434	3444	3457	rBV2	14085	22265	1.49%	0.317%
38	12.882	3672	3683	3693	rVB2	9961	15181	1.01%	0.216%

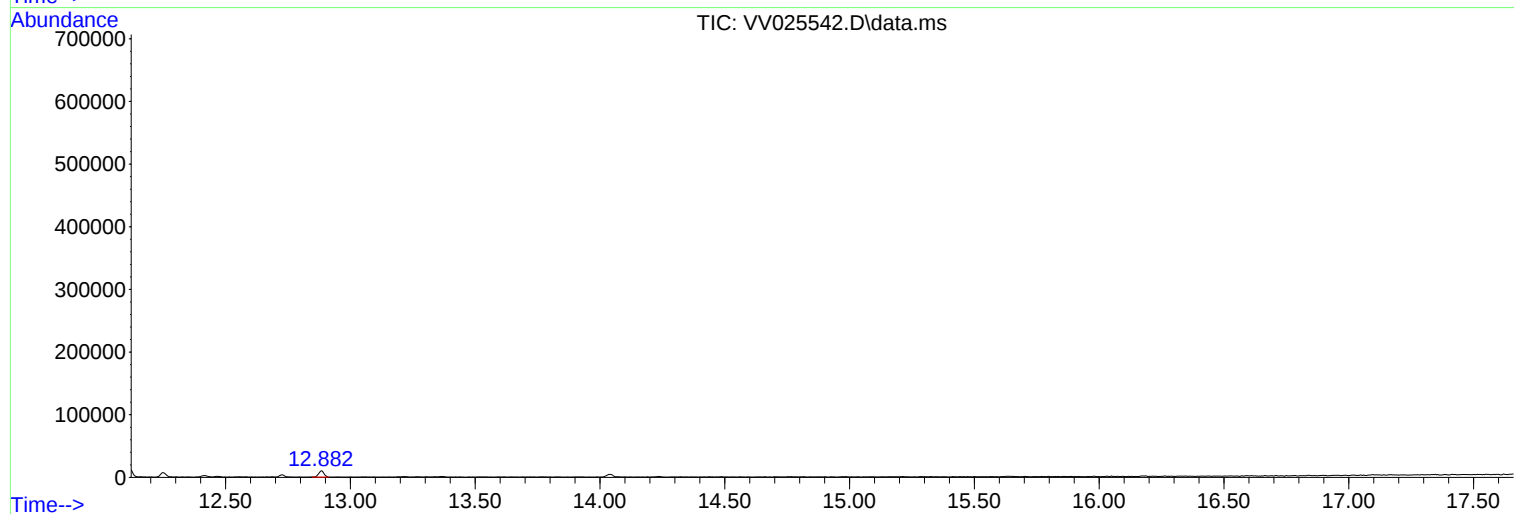
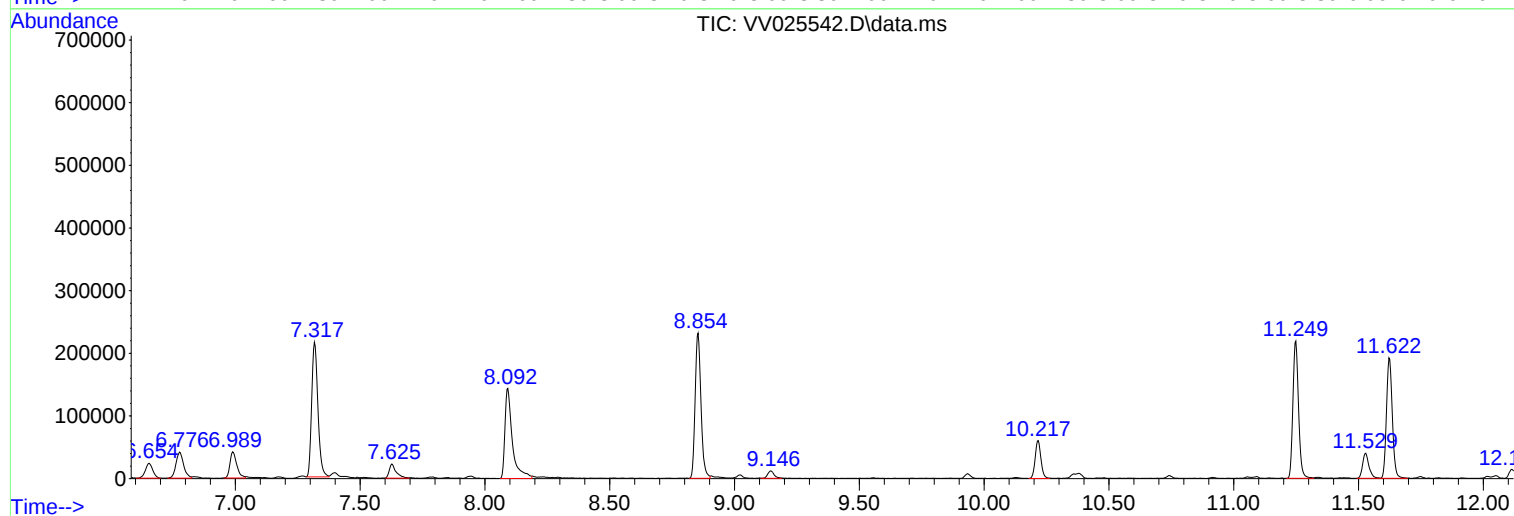
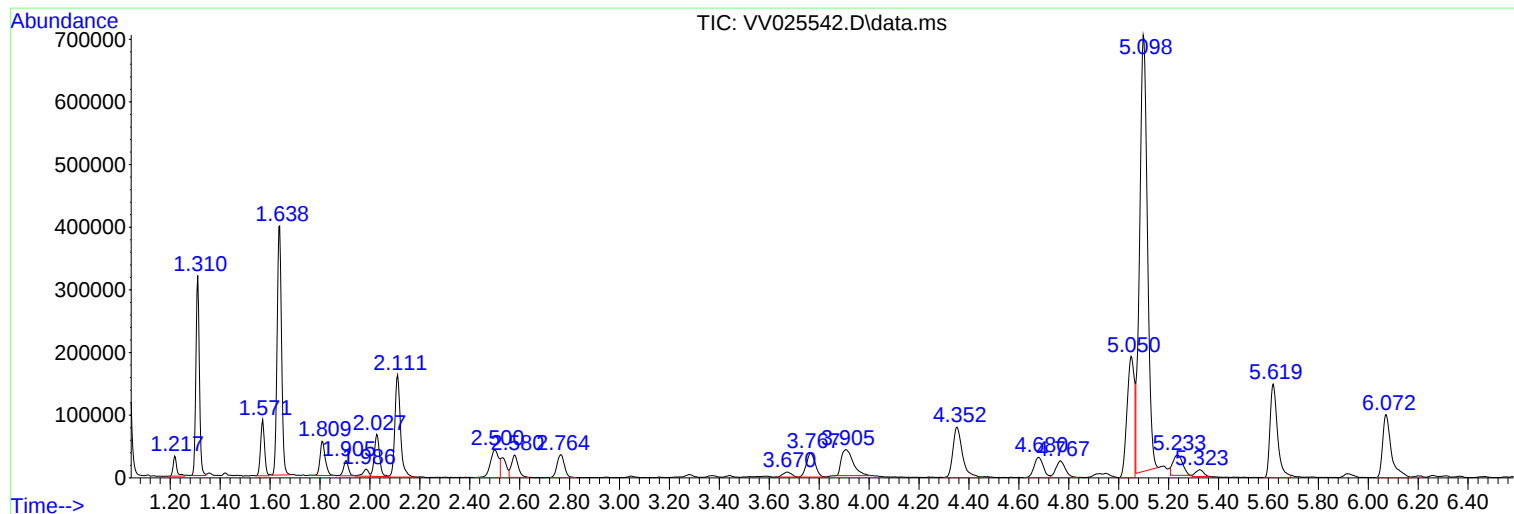
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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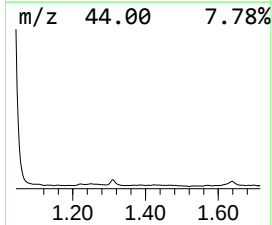
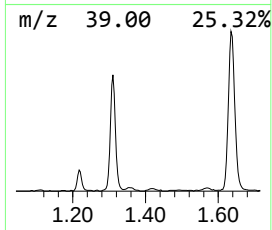
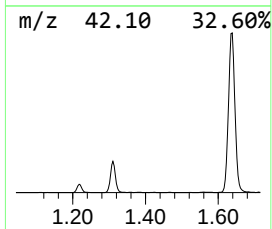
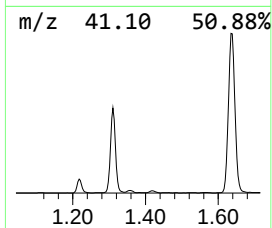
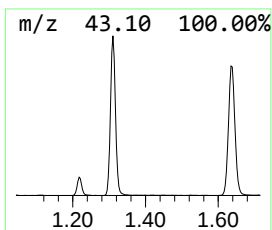
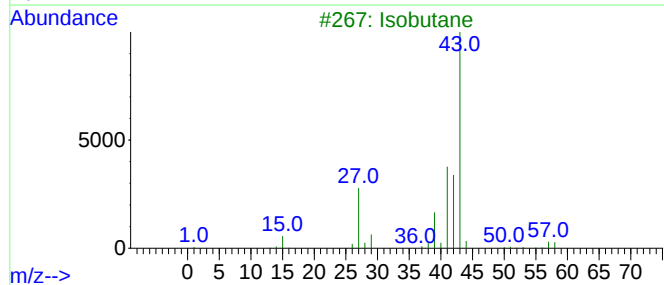
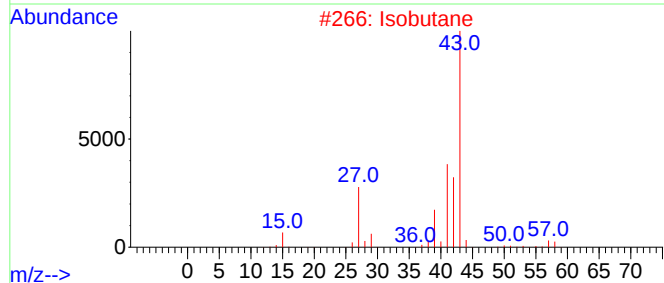
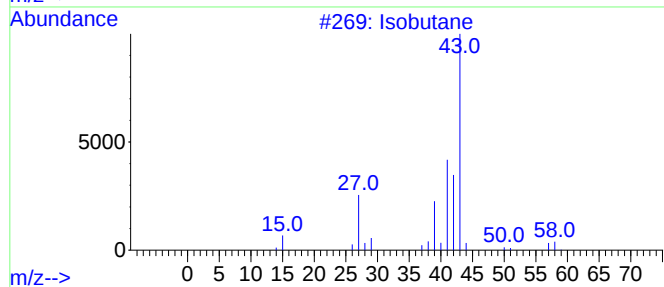
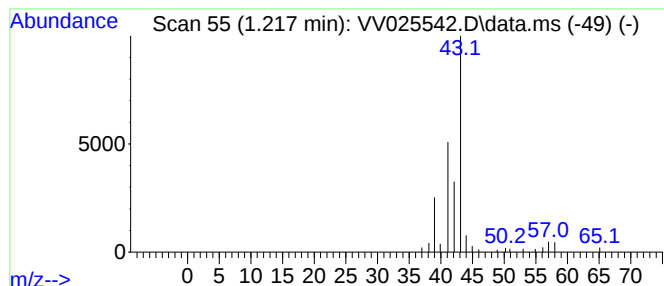
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.217	0.53 ug/L	32575	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutane	58	C4H10	000075-28-5	64
2			Isobutane	58	C4H10	000075-28-5	64
3			Isobutane	58	C4H10	000075-28-5	50
4			Isobutane	58	C4H10	000075-28-5	45
5			1-[3-(4-Bromophenyl)-2-thioureid...	560	C21H25BrN2O9S	094273-12-8	12



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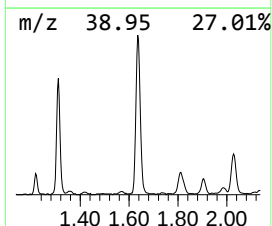
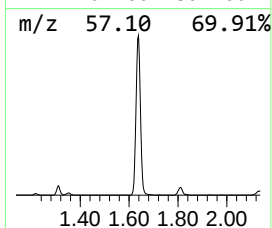
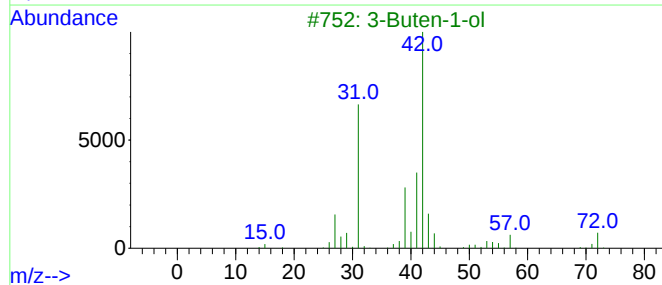
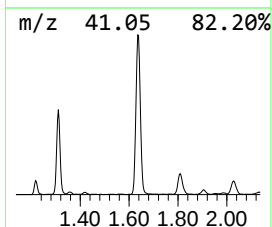
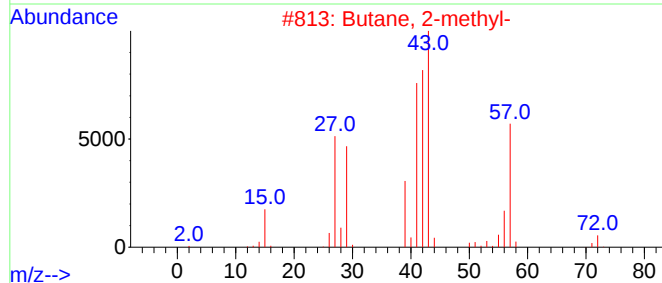
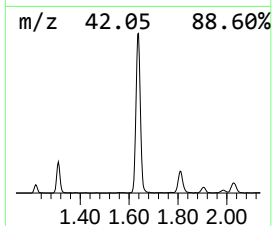
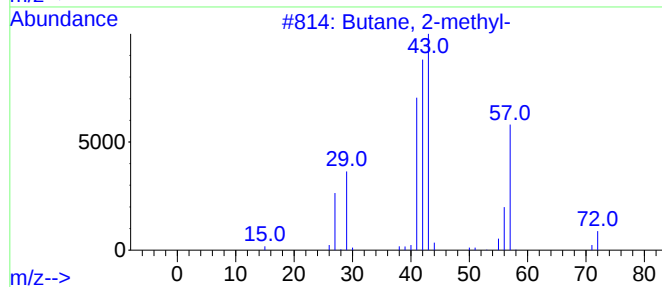
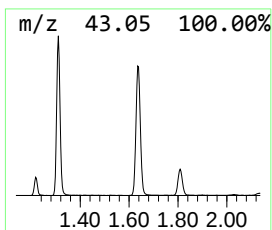
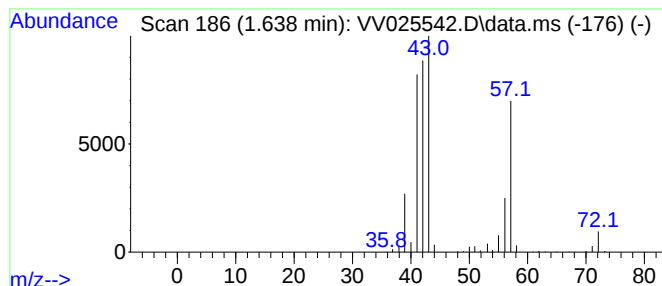
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.638	7.99 ug/L	494215	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	86
3	3-Buten-1-ol			72	C4H8O	000627-27-0	47
4	Pentane			72	C5H12	000109-66-0	38
5	Methane, isocyanato-			57	C2H3NO	000624-83-9	9



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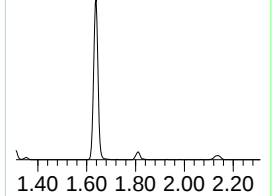
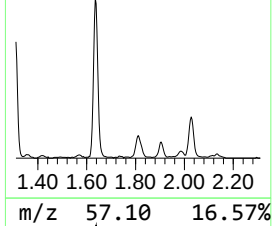
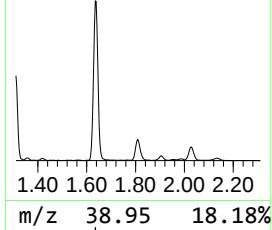
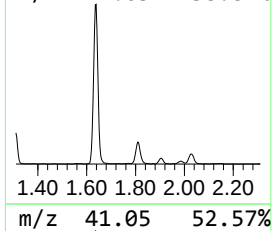
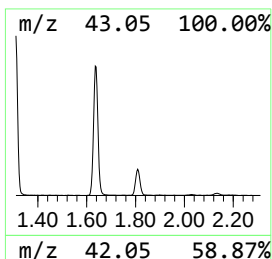
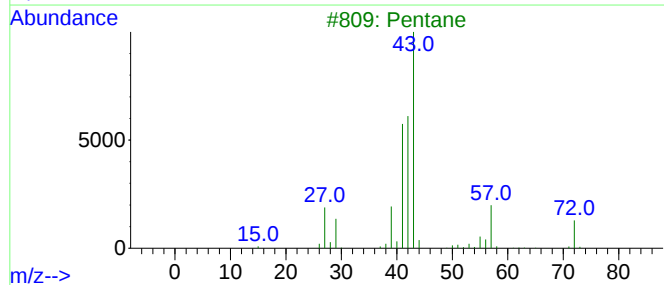
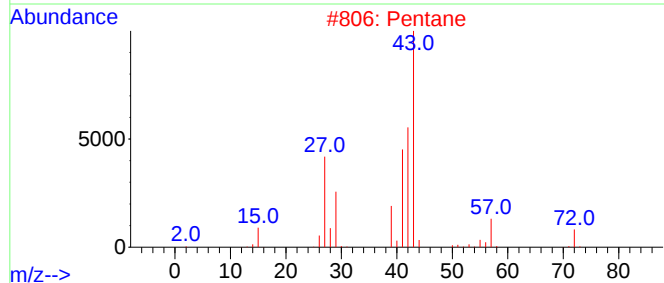
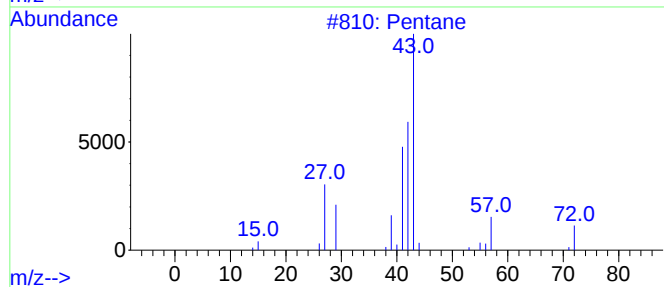
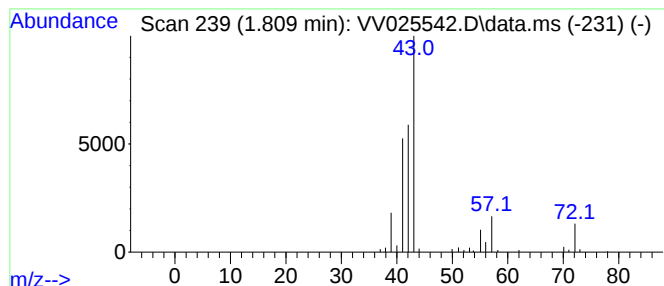
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.809	1.28 ug/L	79151	1,4-Difluorobenzene	5.619

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



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

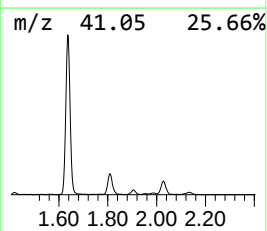
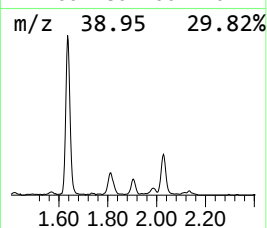
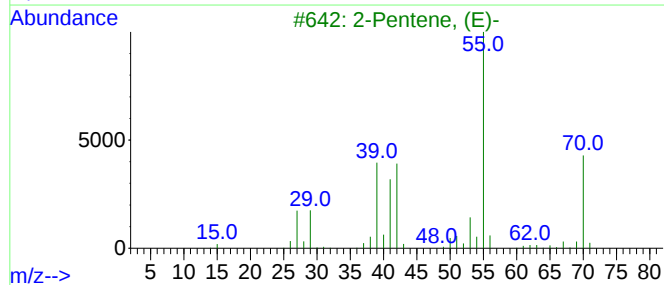
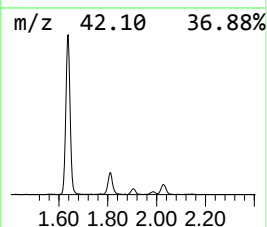
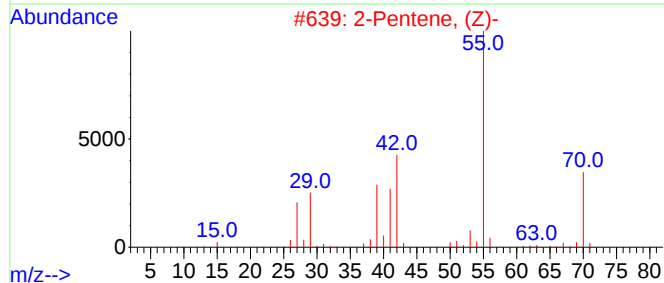
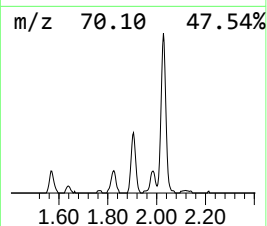
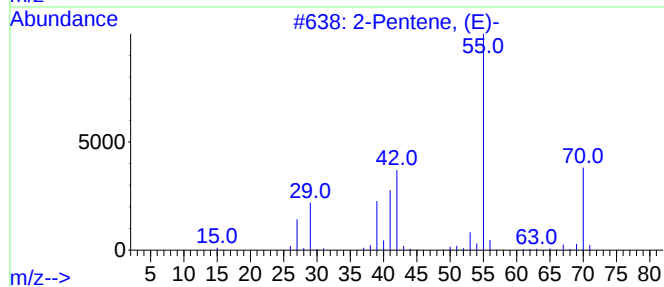
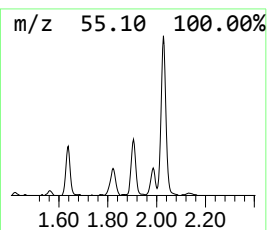
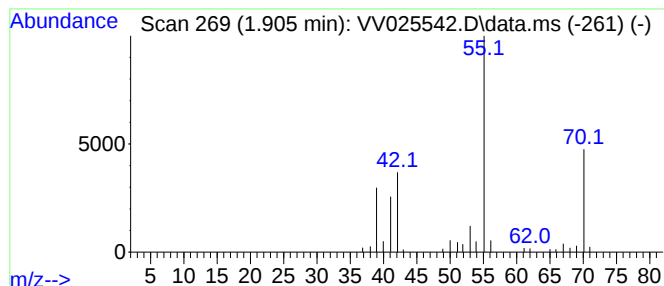
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.905	0.53 ug/L	33022	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (E)-		70	C5H10	000646-04-8	90
2	2-Pentene, (Z)-		70	C5H10	000627-20-3	87
3	2-Pentene, (E)-		70	C5H10	000646-04-8	87
4	Cyclopropane, 1,2-dimethyl-, trans-		70	C5H10	002402-06-4	87
5	2-Butene, 2-methyl-		70	C5H10	000513-35-9	86



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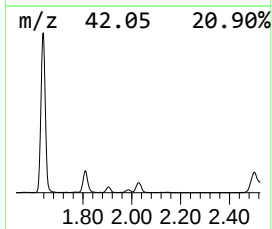
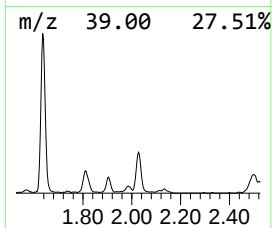
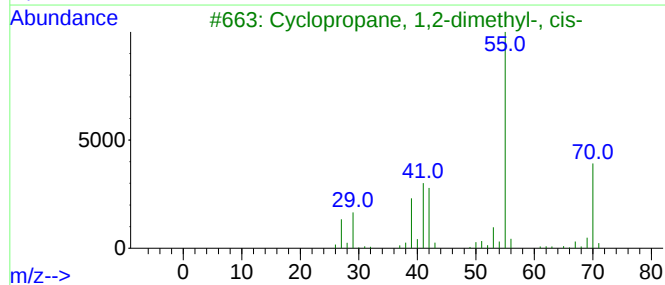
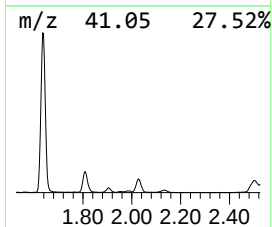
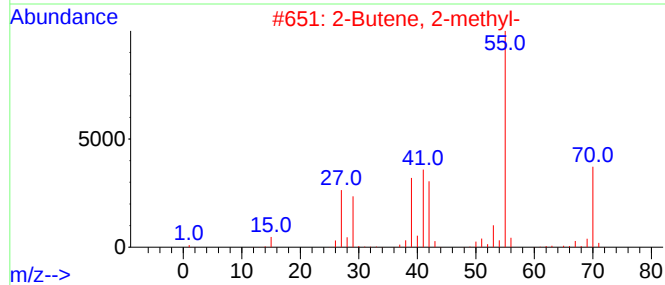
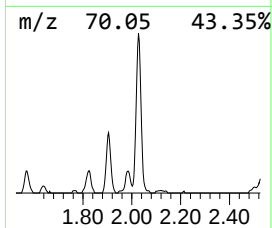
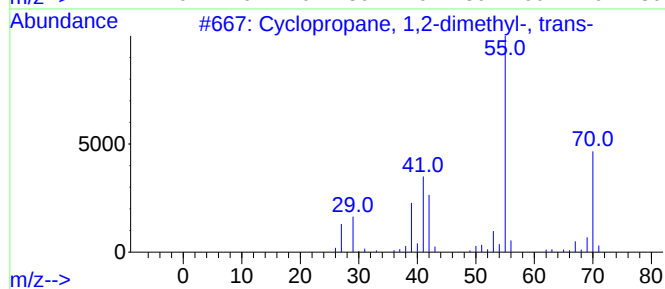
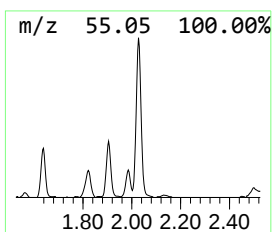
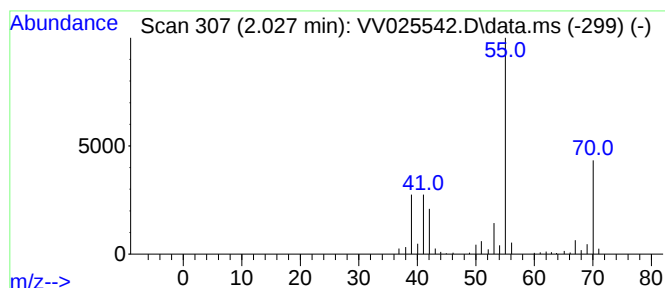
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Cyclic2.027 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.027	1.56 ug/L	96523	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	91
2		2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
3		Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	87
4		2-Pentene, (E)-	70	C5H10	000646-04-8	86
5		1-Butene, 3-methyl-	70	C5H10	000563-45-1	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

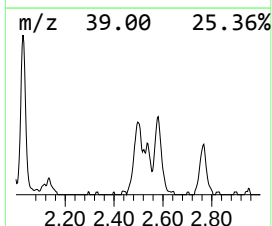
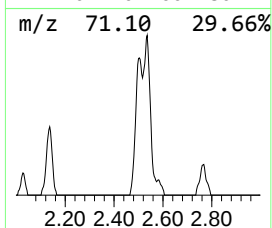
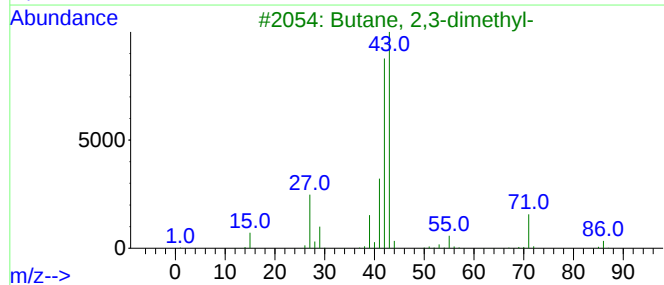
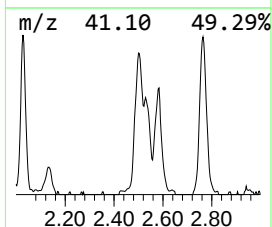
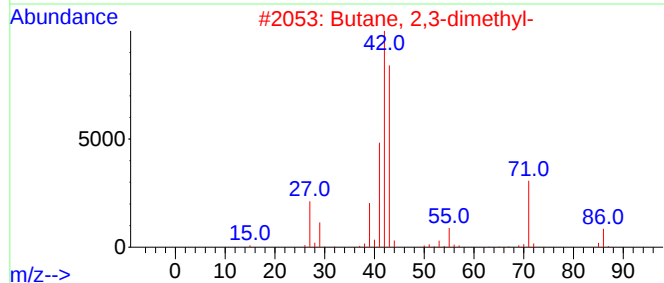
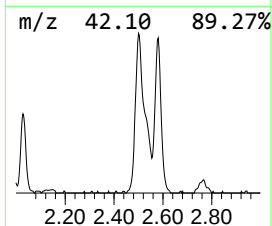
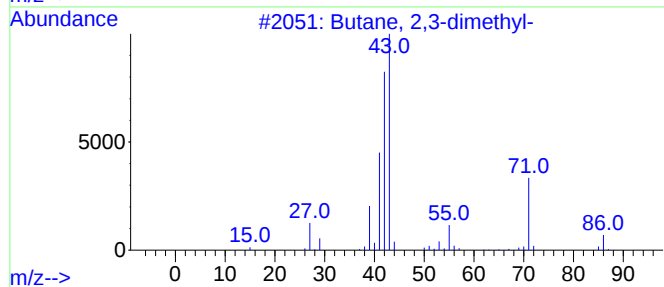
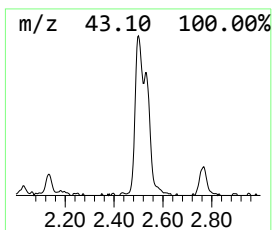
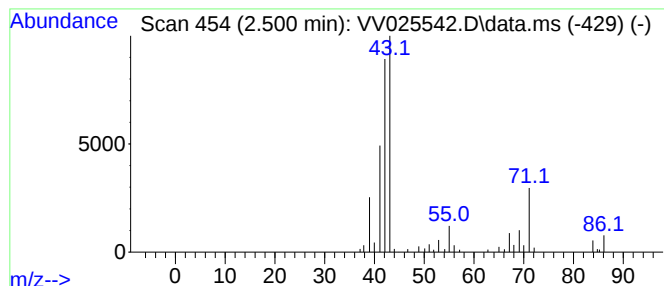
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.500	1.70 ug/L	105401	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	83
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
4			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
5			1,2,3-Trimethyldiaziridine	86	C4H10N2	113604-56-1	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

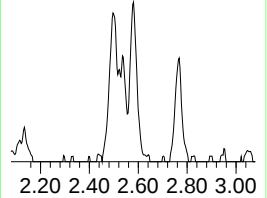
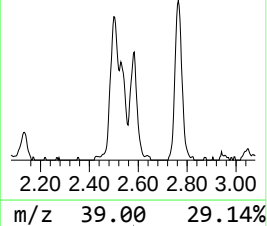
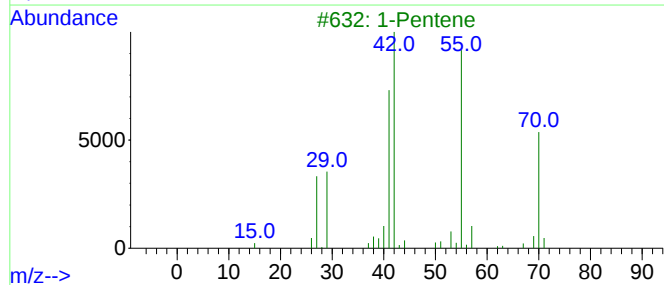
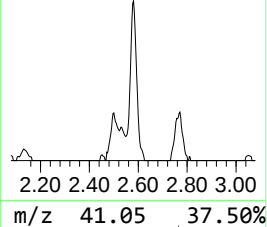
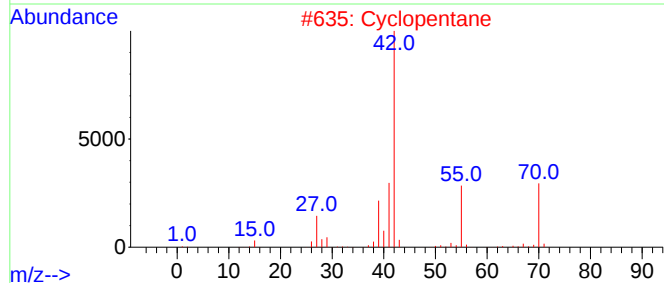
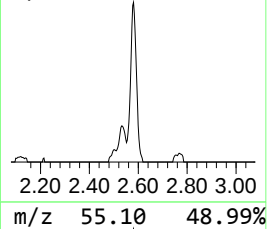
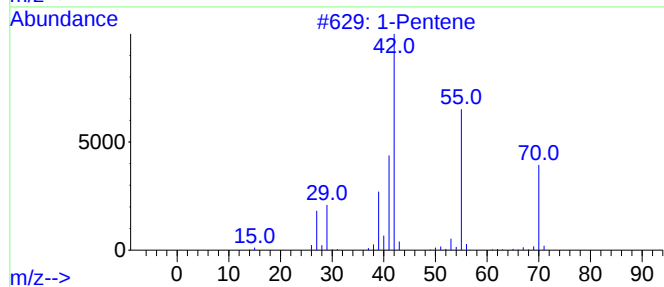
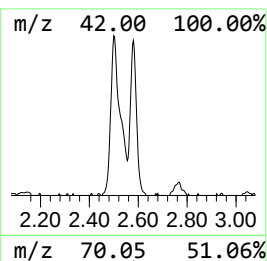
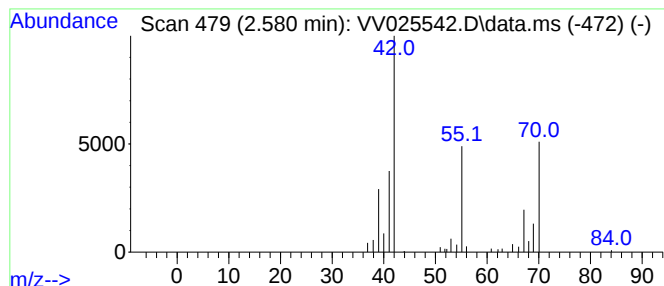
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.580	1.07 ug/L	66009	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene	70	C5H10	000109-67-1	59
2		Cyclopentane	70	C5H10	000287-92-3	59
3		1-Pentene	70	C5H10	000109-67-1	53
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	47
5		Cyclopropane, ethyl-	70	C5H10	001191-96-4	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

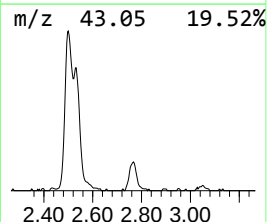
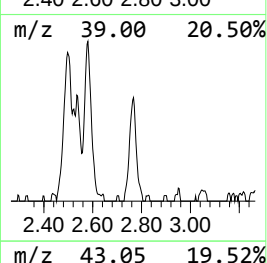
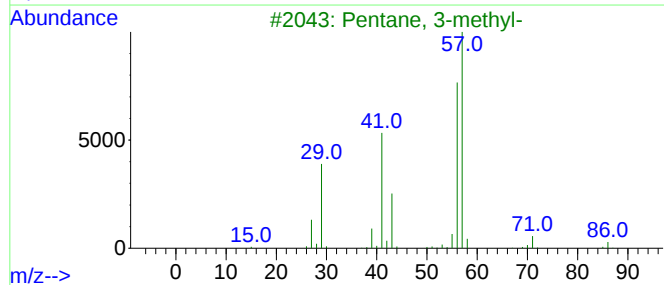
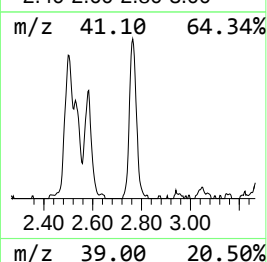
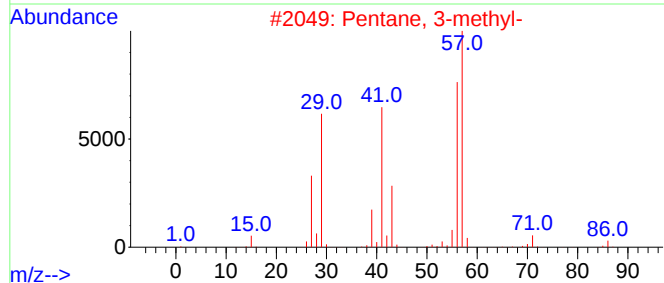
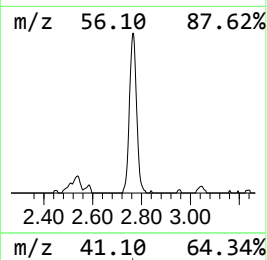
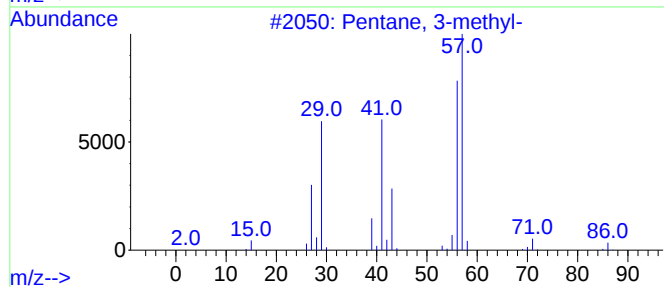
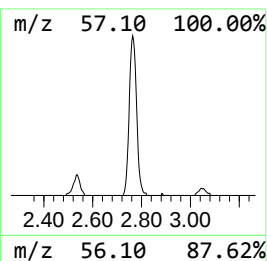
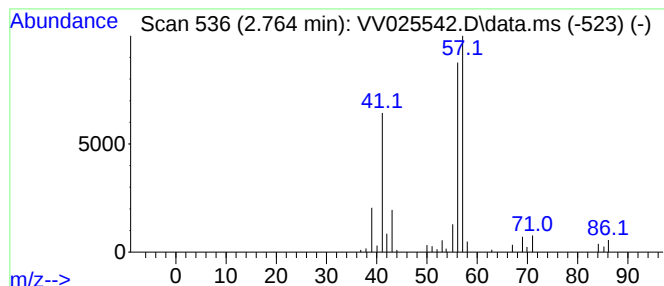
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Quant Title : TRACE VOA SFAM1.0

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.
2.764	1.21 ug/L	74819	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		Pentane, 3-methyl-	86	C6H14	000096-14-0	80
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	80
4		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
5		Cyclopentane, methyl-	84	C6H12	000096-37-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

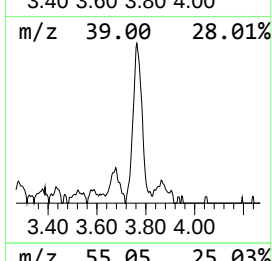
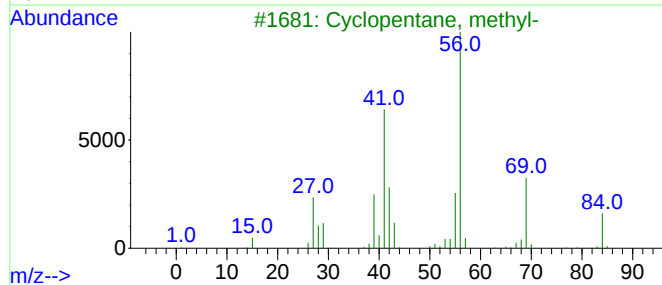
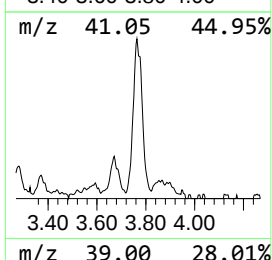
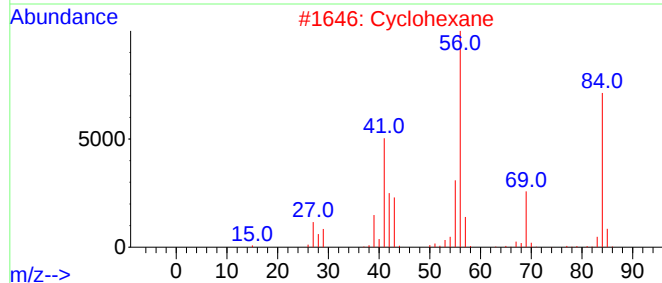
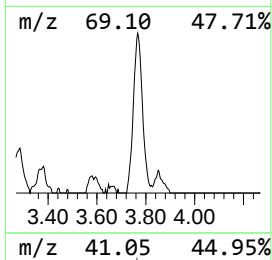
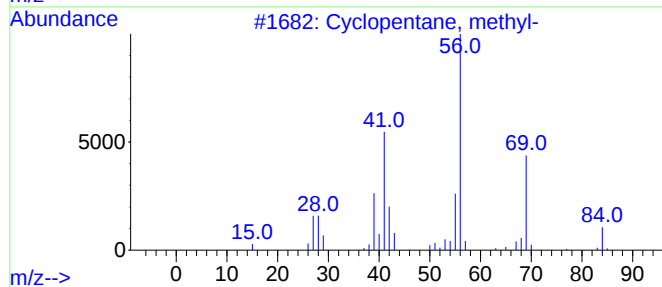
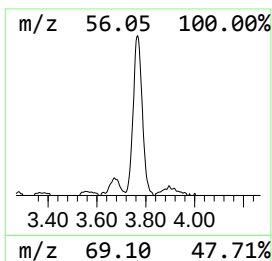
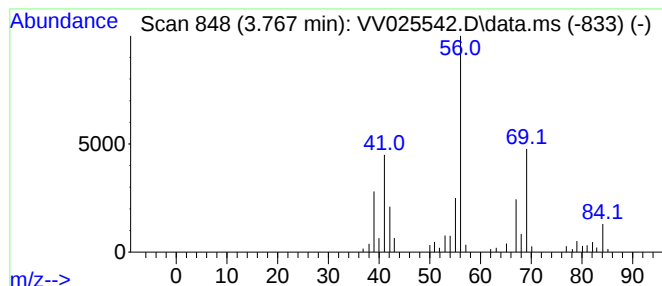
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 (DEL) Alkane: Cyclic3.767 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.767	1.53 ug/L	94720	1,4-Difluorobenzene	5.619

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	64
3			Cyclopentane, methyl-	84	C6H12	000096-37-7	64
4			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

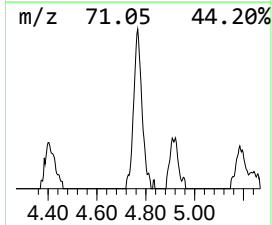
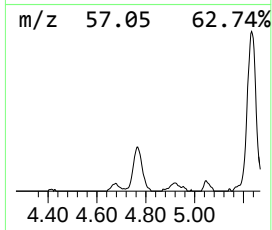
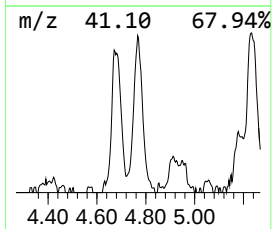
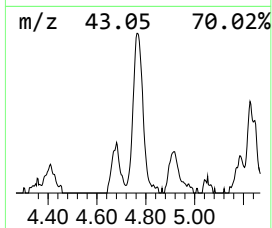
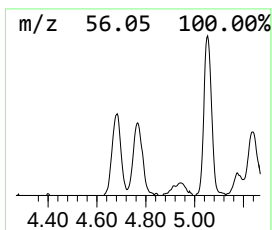
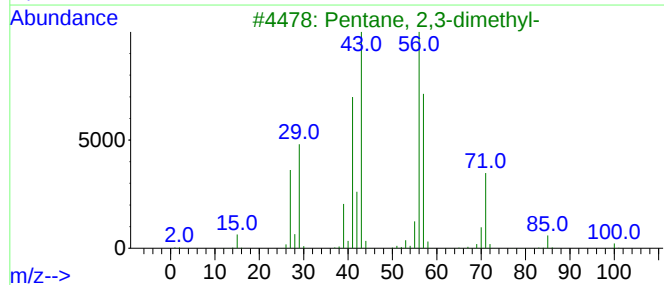
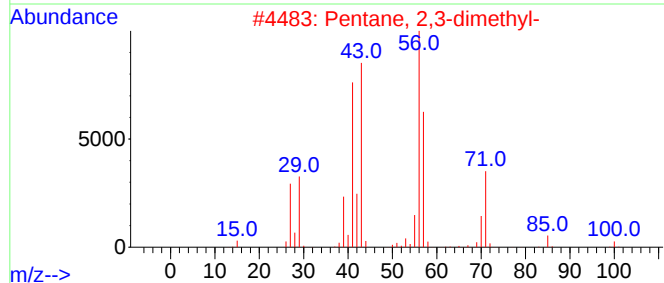
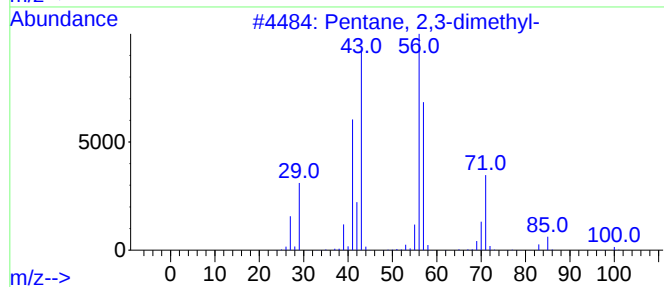
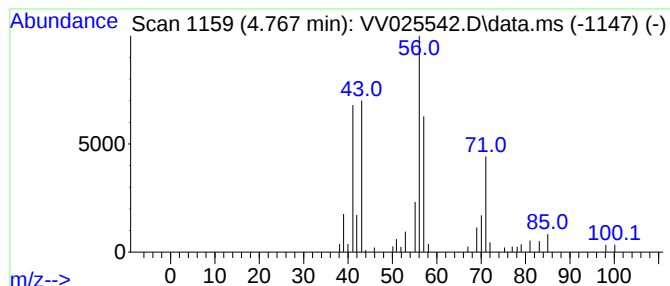
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.767	1.11 ug/L	68818	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	90
2		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	86
3		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	72
4		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	59
5		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GBDS8DL

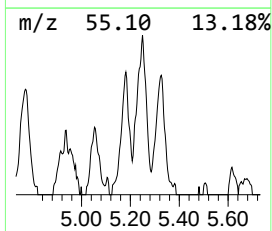
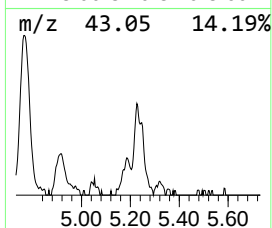
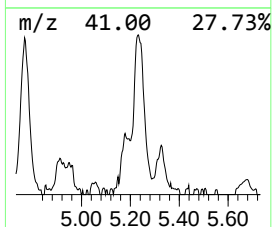
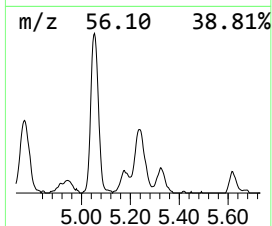
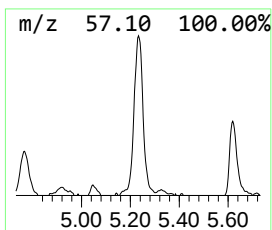
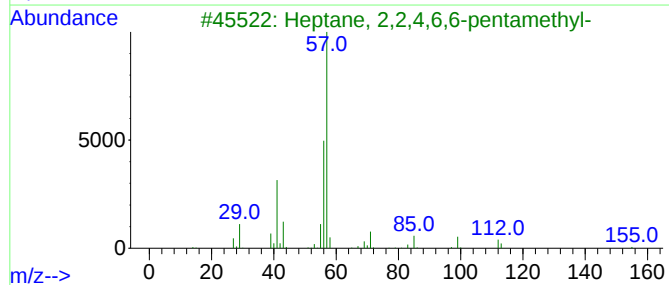
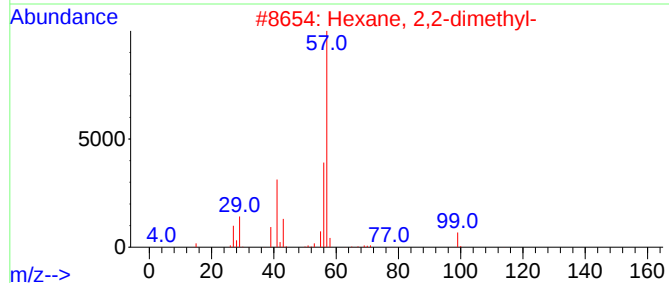
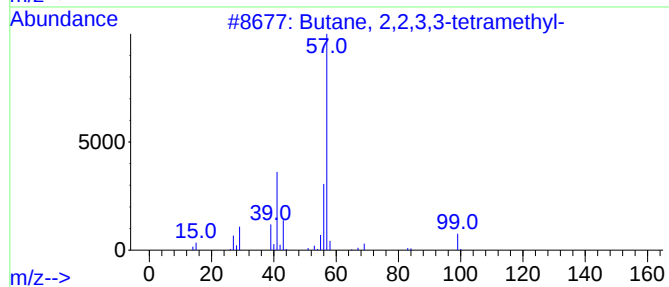
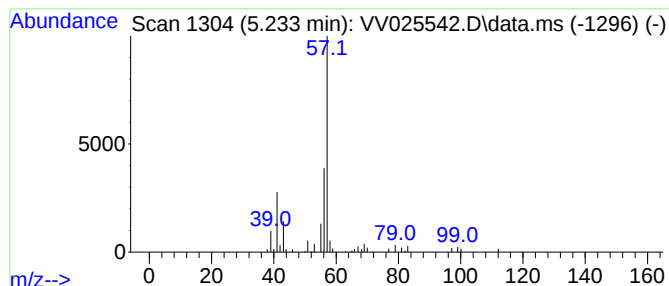
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.233	1.42 ug/L	87625	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	64
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	53
3		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	50
4		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	45
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

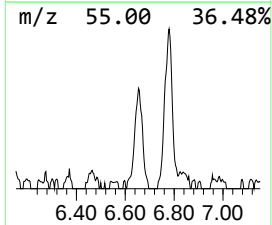
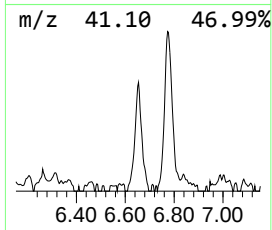
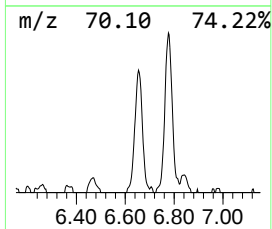
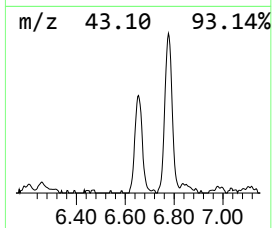
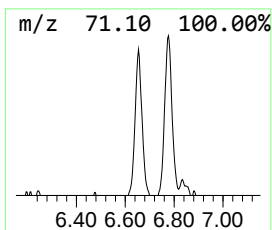
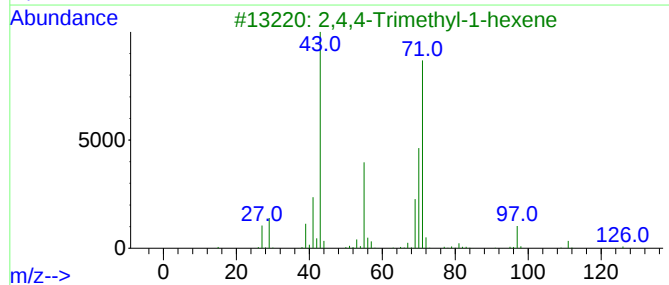
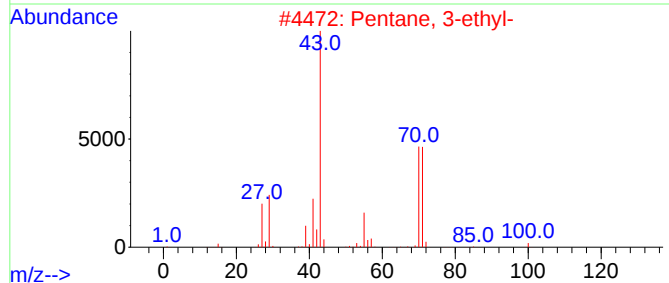
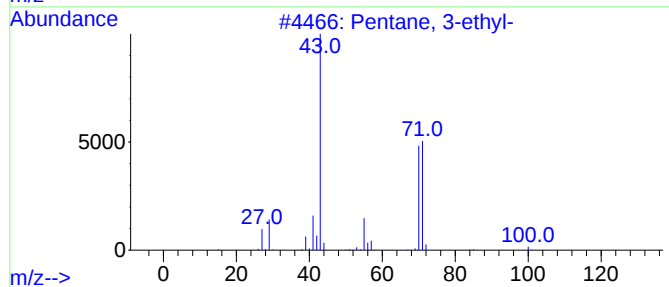
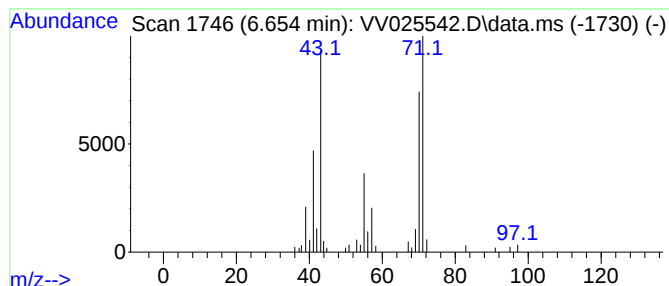
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Quant Title : TRACE VOA SFAM1.0

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	0.79 ug/L	49142	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
2		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
3		2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
4		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

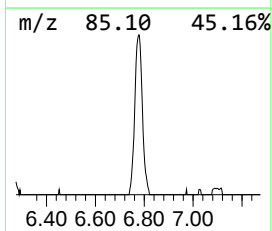
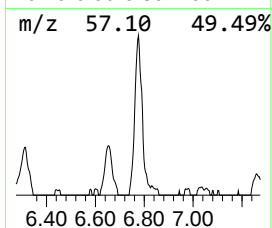
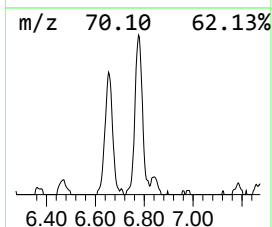
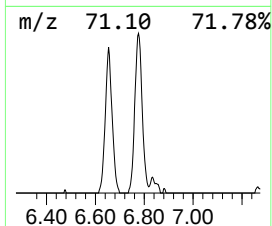
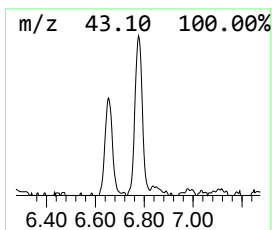
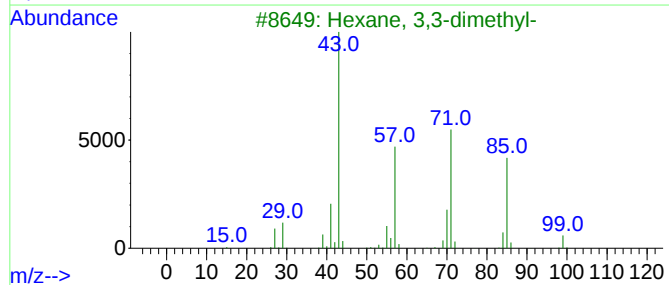
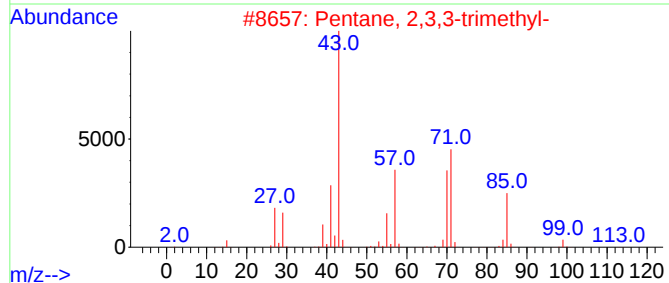
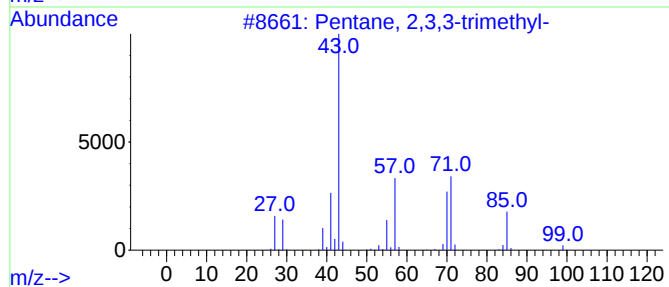
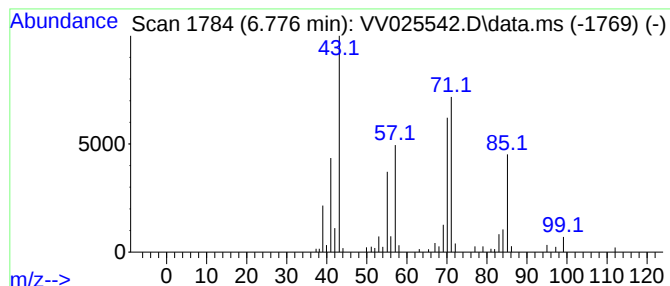
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.776	1.44 ug/L	89030	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	78
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	64
3			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4			Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	59
5			Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
Data File : VV025542.D
Acq On : 14 Apr 2022 13:43
Operator : SY/MD
Sample : N2356-10DL 50X
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS8DL

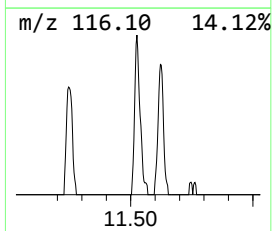
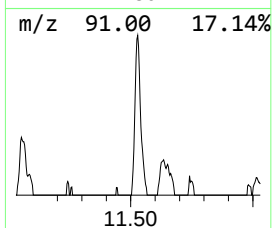
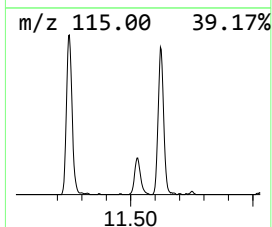
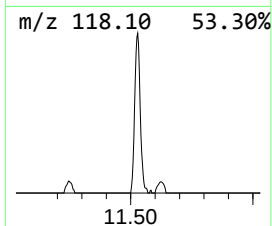
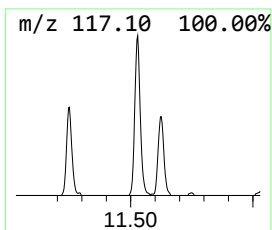
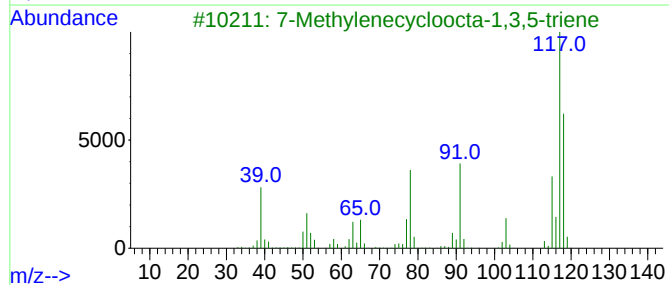
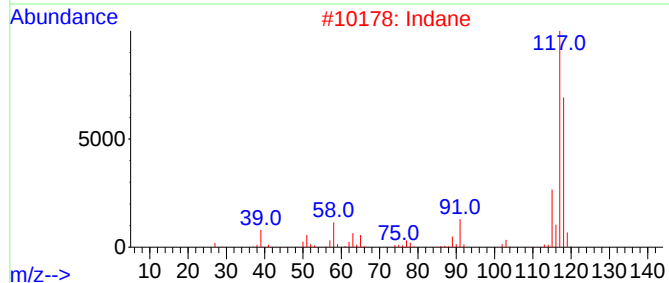
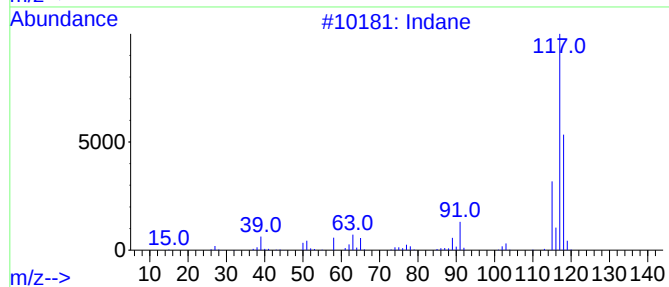
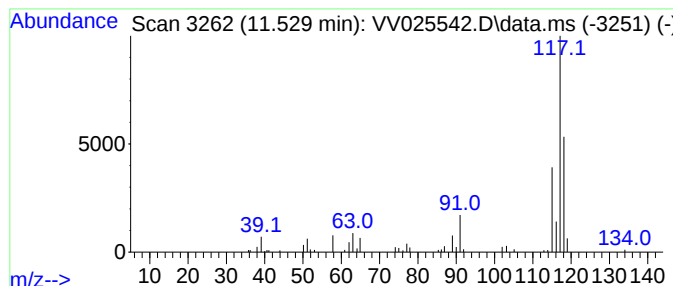
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Indane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	1.03 ug/L	70801	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	76
3			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	68
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	62
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041422\
 Data File : VV025542.D
 Acq On : 14 Apr 2022 13:43
 Operator : SY/MD
 Sample : N2356-10DL 50X
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GBDS8DL

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...	1.217	0.5	ug/L	32575	1	5.619	309277	5.0	
(DEL) Alkane: S...	1.638	8.0	ug/L	494215	1	5.619	309277	5.0	
(DEL) Alkane: S...	1.809	1.3	ug/L	79151	1	5.619	309277	5.0	
(DEL) Alkane: S...	1.905	0.5	ug/L	33022	1	5.619	309277	5.0	
(DEL) Alkane: C...	2.027	1.6	ug/L	96523	1	5.619	309277	5.0	
(DEL) Alkane: S...	2.500	1.7	ug/L	105401	1	5.619	309277	5.0	
(DEL) Alkane: S...	2.580	1.1	ug/L	66009	1	5.619	309277	5.0	
(DEL) Alkane: S...	2.764	1.2	ug/L	74819	1	5.619	309277	5.0	
(DEL) Alkane: C...	3.767	1.5	ug/L	94720	1	5.619	309277	5.0	
(DEL) Alkane: S...	4.767	1.1	ug/L	68818	1	5.619	309277	5.0	
(DEL) Alkane: S...	5.233	1.4	ug/L	87625	1	5.619	309277	5.0	
(DEL) Alkane: S...	6.654	0.8	ug/L	49142	1	5.619	309277	5.0	
(DEL) Alkane: S...	6.776	1.4	ug/L	89030	1	5.619	309277	5.0	
Indane	11.529	1.0	ug/L	70801	3	11.249	343757	5.0	