

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
 Data File : VV032601.D
 Acq On : 24 Oct 2023 19:43
 Operator : SY/MD
 Sample : 04992-05
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0AJ1

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\SFAMVTR092923WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV032601.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.085	7	14	31	rBV	1792660	1701468	100.00%	20.677%
2	1.192	40	47	63	rBV	126538	141488	8.32%	1.719%
3	1.272	63	72	86	rBV2	706779	986881	58.00%	11.993%
4	1.388	101	108	123	rVB2	99187	146481	8.61%	1.780%
5	1.529	138	152	164	rVB3	94879	134953	7.93%	1.640%
6	1.597	165	173	182	rBV	95621	119980	7.05%	1.458%
7	1.725	203	213	220	rBV	220409	288157	16.94%	3.502%
8	1.764	220	225	241	rVB	109249	152793	8.98%	1.857%
9	1.854	243	253	261	rBV2	30832	59549	3.50%	0.724%
10	1.938	271	279	287	rVB2	31840	44811	2.63%	0.545%
11	2.060	306	317	328	rBV	70465	108604	6.38%	1.320%
12	2.179	347	354	364	rVB3	9202	17920	1.05%	0.218%
13	2.388	402	419	432	rBV5	31829	81183	4.77%	0.987%
14	2.696	501	515	531	rVB4	28632	64450	3.79%	0.783%
15	2.883	562	573	588	rVB2	38630	77764	4.57%	0.945%
16	2.970	590	600	624	rVB2	17226	39910	2.35%	0.485%
17	3.368	712	724	735	rBV7	7605	19166	1.13%	0.233%
18	3.802	841	859	893	rBV3	80441	287862	16.92%	3.498%
19	3.970	900	911	944	rVB	175849	455770	26.79%	5.539%
20	4.253	982	999	1017	rBV2	61818	164562	9.67%	2.000%
21	4.365	1024	1034	1053	rVB4	11240	31902	1.87%	0.388%
22	4.658	1114	1125	1147	rVB8	8010	22549	1.33%	0.274%
23	4.960	1204	1219	1235	rBV4	134102	350944	20.63%	4.265%
24	5.535	1386	1398	1415	rBV	130535	277448	16.31%	3.372%
25	5.992	1525	1540	1552	rBV2	83755	179686	10.56%	2.184%
26	6.918	1819	1828	1854	rBV2	34845	76087	4.47%	0.925%
27	7.246	1919	1930	1948	rBV	139180	267011	15.69%	3.245%
28	7.323	1948	1954	1968	rVB	15192	28255	1.66%	0.343%
29	7.558	2018	2027	2046	rBV	21172	46589	2.74%	0.566%
30	8.027	2163	2173	2212	rBV	275216	559666	32.89%	6.801%
31	8.786	2398	2409	2437	rBV	212131	378252	22.23%	4.597%
32	9.075	2492	2499	2517	rVB3	11049	20029	1.18%	0.243%
33	10.156	2823	2835	2852	rBV	86456	141023	8.29%	1.714%
34	10.387	2899	2907	2922	rBV	13195	27158	1.60%	0.330%
35	10.857	3045	3053	3063	rVB2	10829	17277	1.02%	0.210%
36	11.188	3146	3156	3174	rVV	221565	348040	20.46%	4.230%

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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

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37	11.278	3174	3184	3197	rVB	24715	40530	2.38%	0.493%
38	11.564	3263	3273	3291	rVB	179753	304111	17.87%	3.696%
39	13.445	3849	3858	3868	rBV3	11301	18396	1.08%	0.224%

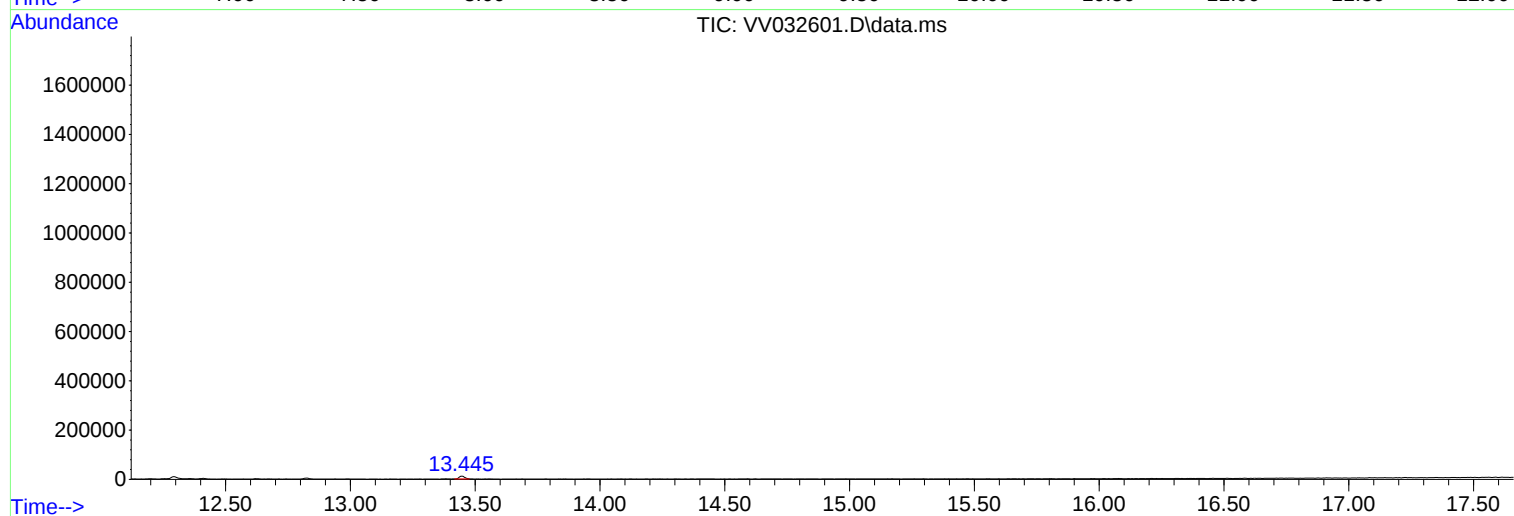
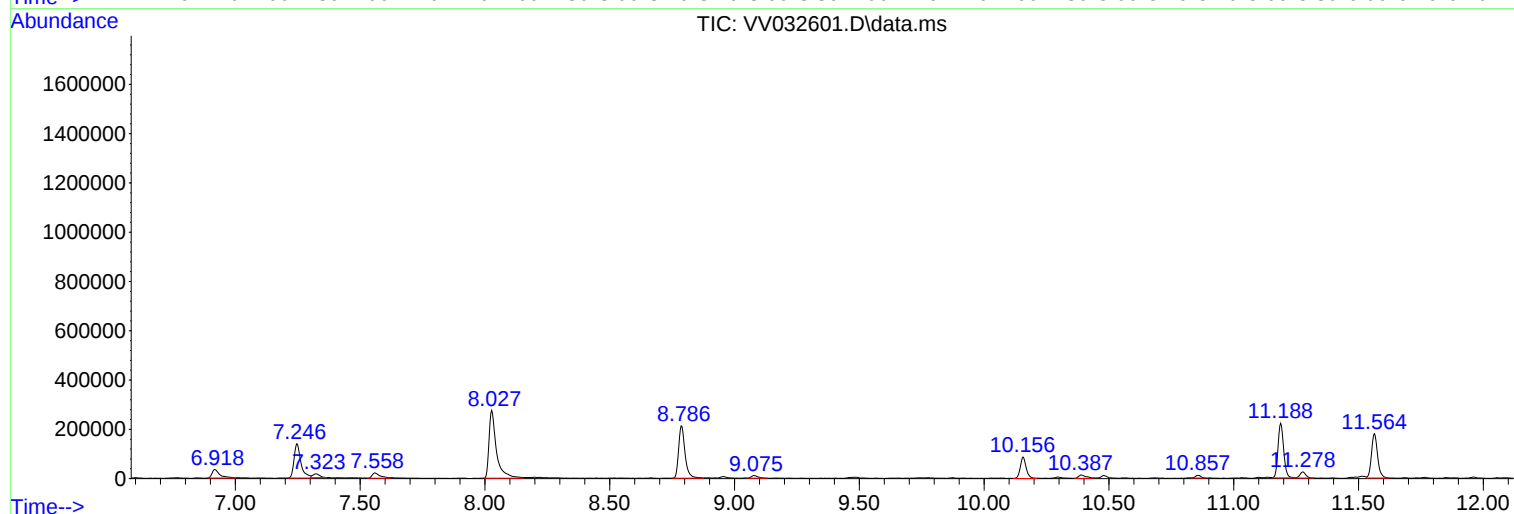
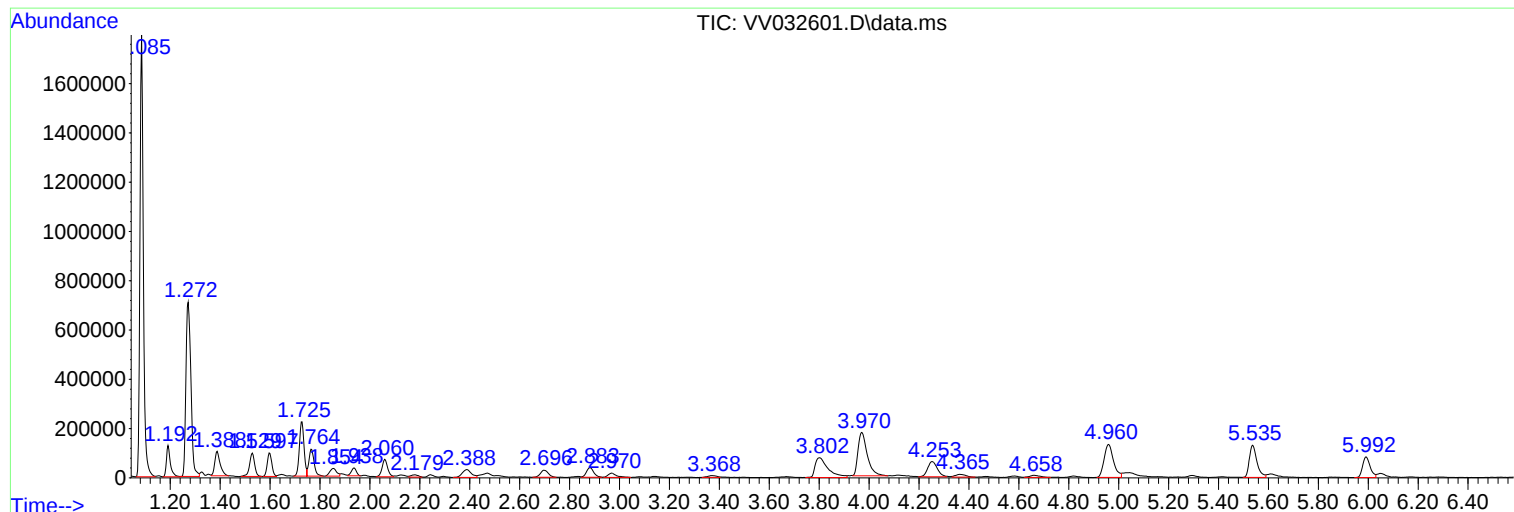
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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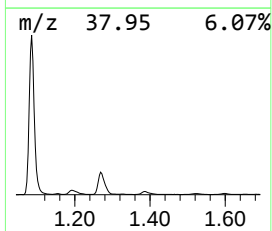
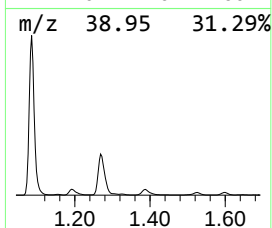
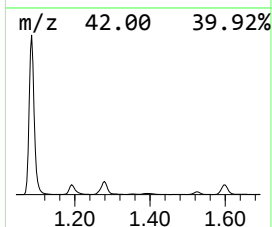
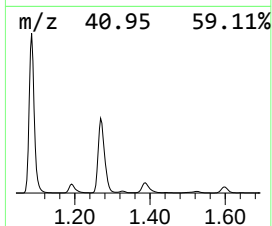
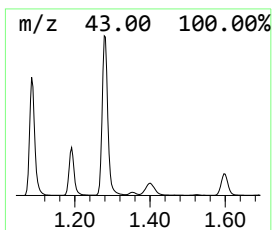
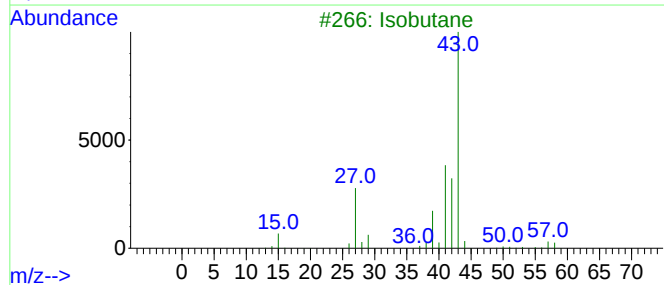
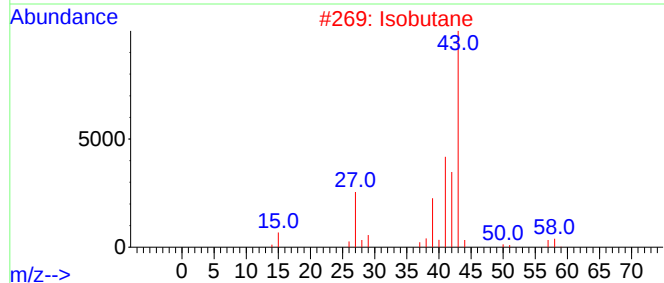
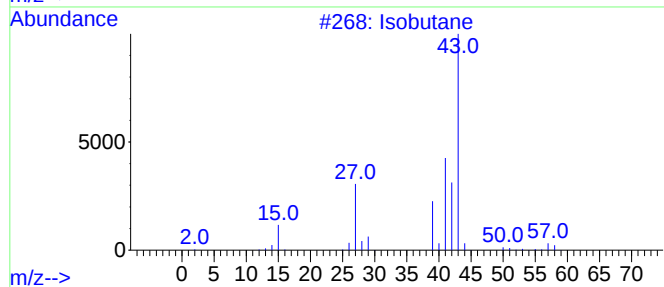
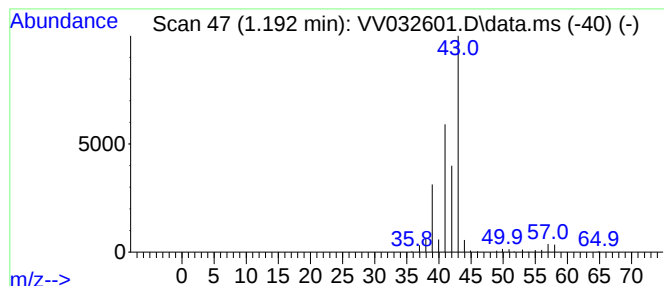
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.192	2.55 ug/L	141488	1,4-Difluorobenzene	5.535

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	50
5		Isobutane	58	C4H10	000075-28-5	36



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Instrument :
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ClientSampleId :
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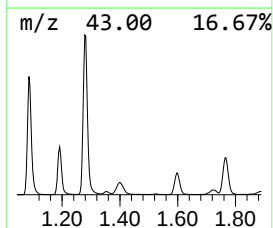
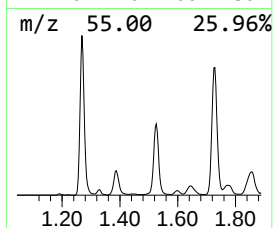
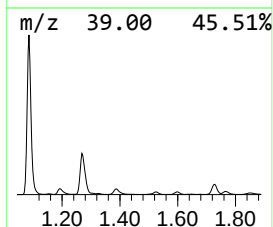
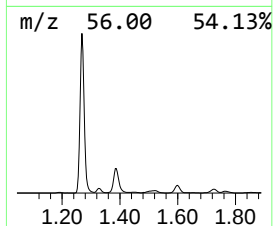
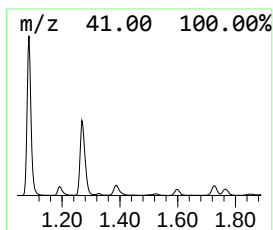
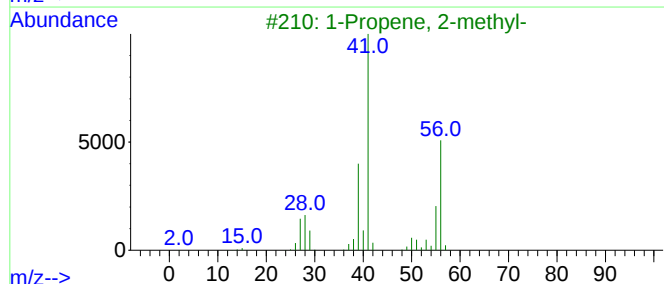
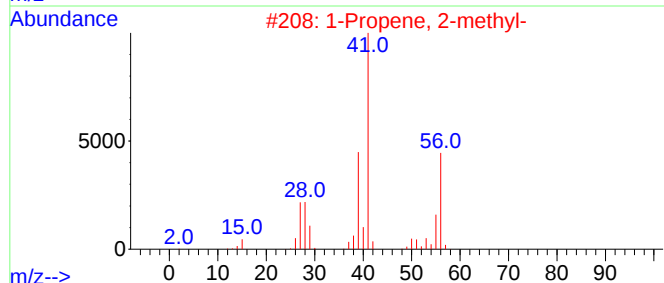
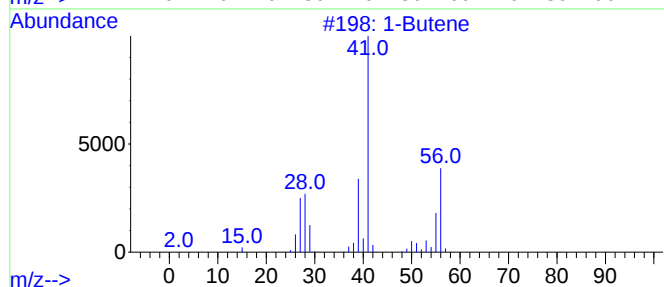
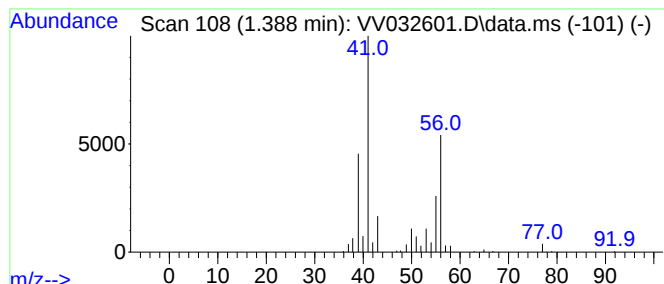
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.388	2.64 ug/L	146481	1,4-Difluorobenzene	5.535

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



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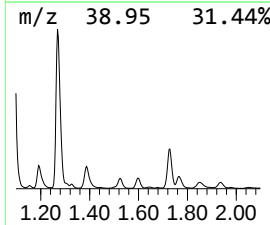
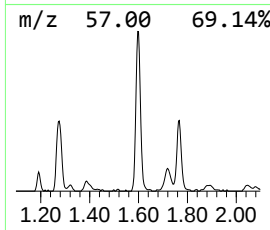
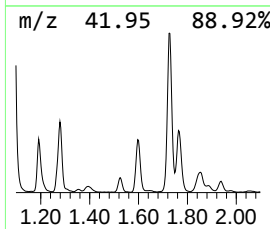
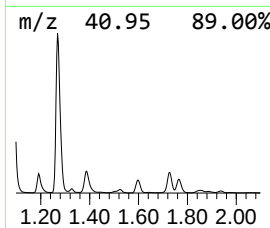
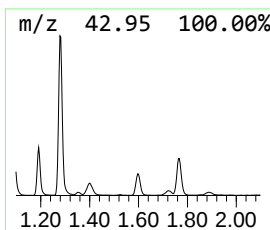
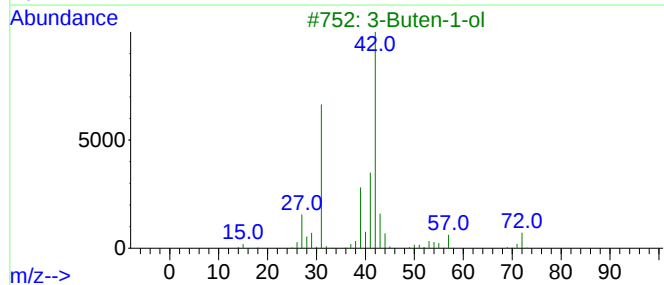
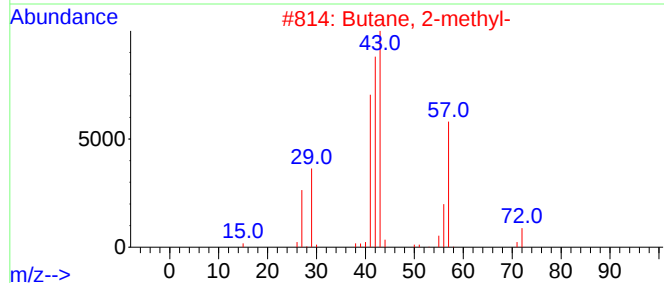
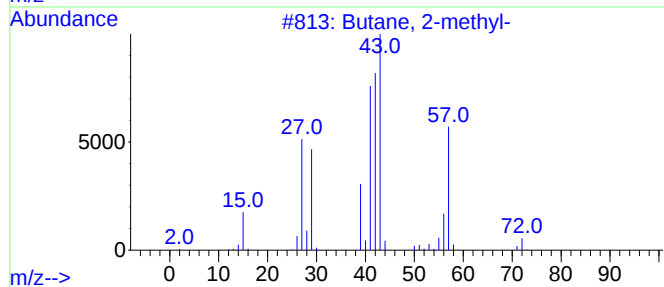
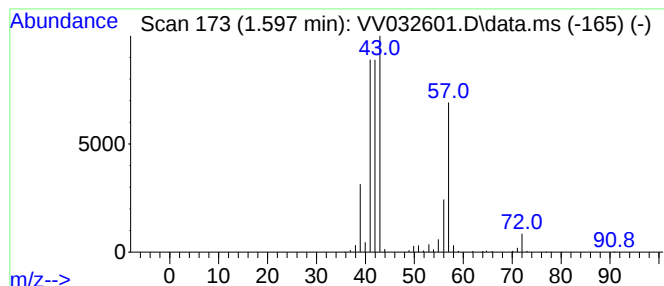
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.597	2.16 ug/L	119980	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	90
2			Butane, 2-methyl-	72	C5H12	000078-78-4	80
3			3-Buten-1-ol	72	C4H8O	000627-27-0	47
4			Pentane	72	C5H12	000109-66-0	36
5			1-Propene, 2-methyl-	56	C4H8	000115-11-7	9



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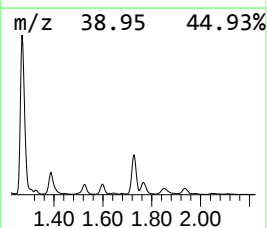
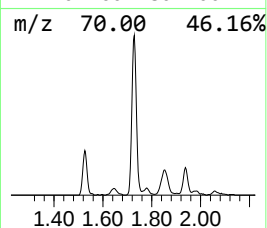
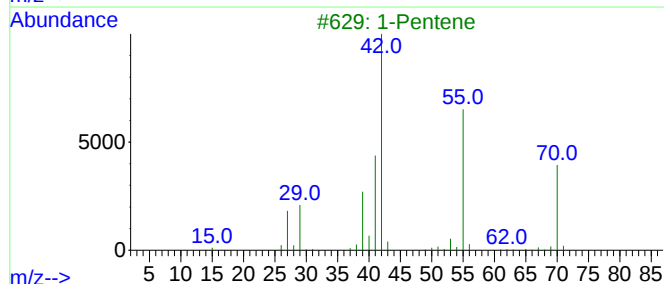
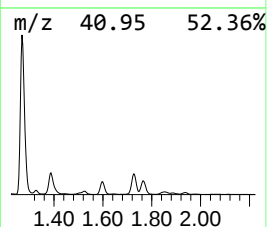
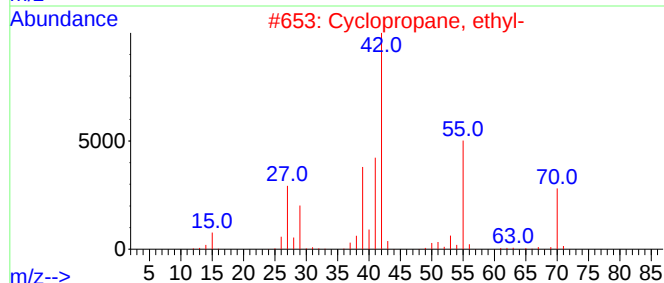
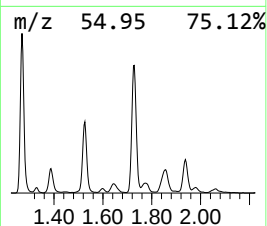
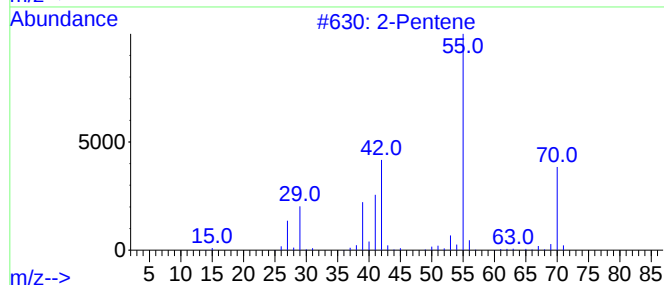
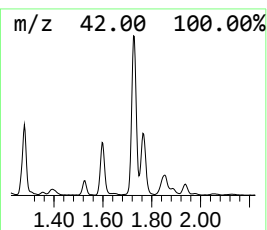
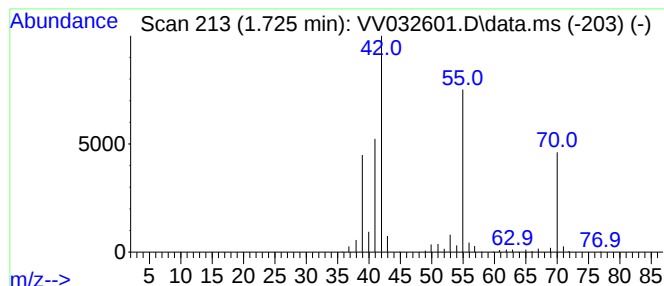
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.725	5.19 ug/L	288157	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentene	70	C5H10	000109-68-2	83
2			Cyclopropane, ethyl-	70	C5H10	001191-96-4	76
3			1-Pentene	70	C5H10	000109-67-1	76
4			Cyclobutane, methyl-	70	C5H10	000598-61-8	72
5			Cyclopropane, ethyl-	70	C5H10	001191-96-4	72



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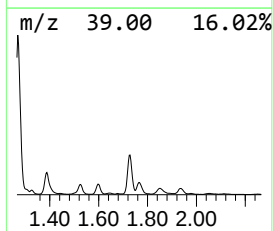
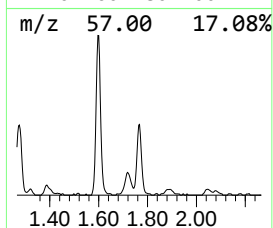
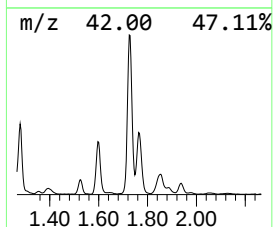
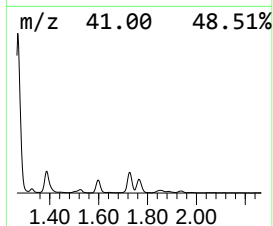
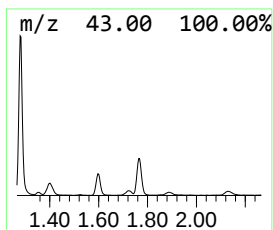
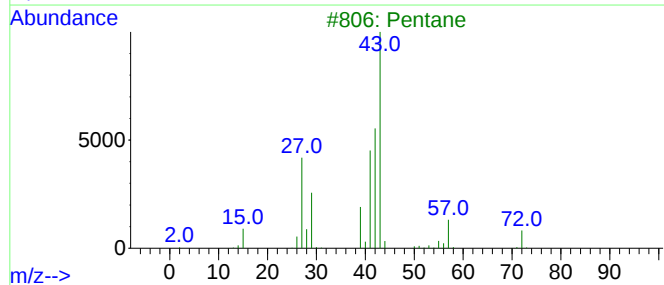
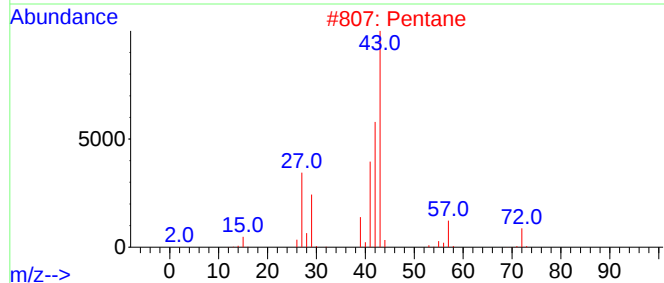
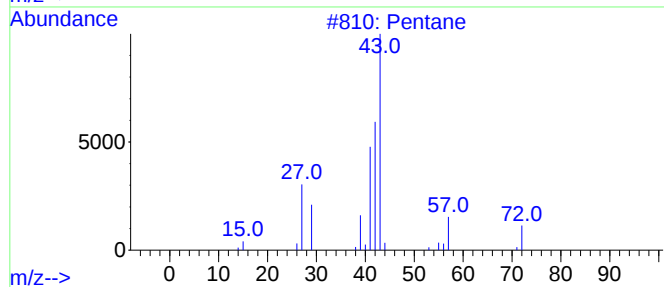
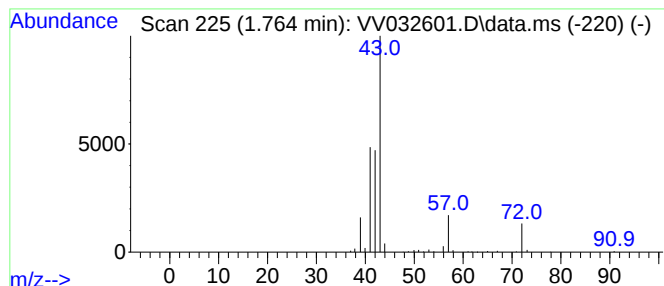
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.764	2.75 ug/L	152793	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane			72	C5H12	000109-66-0	86
2	Pentane			72	C5H12	000109-66-0	80
3	Pentane			72	C5H12	000109-66-0	78
4	Pentane			72	C5H12	000109-66-0	72
5	Pentane			72	C5H12	000109-66-0	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AJ1

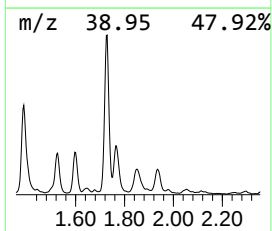
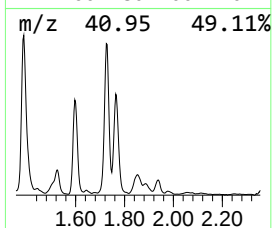
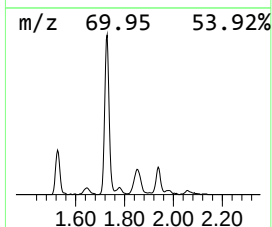
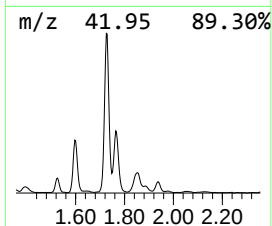
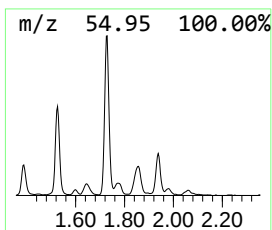
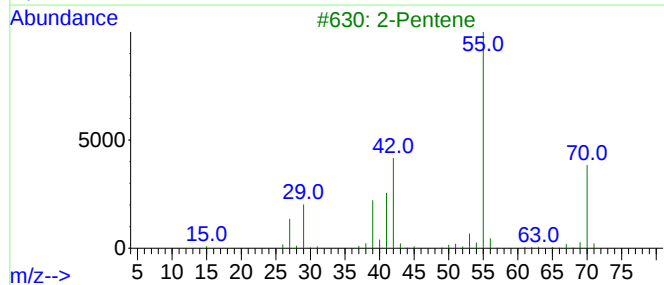
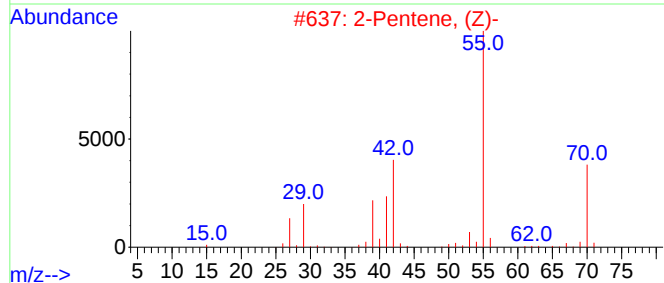
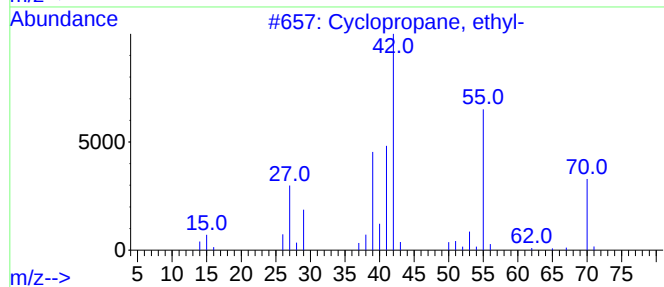
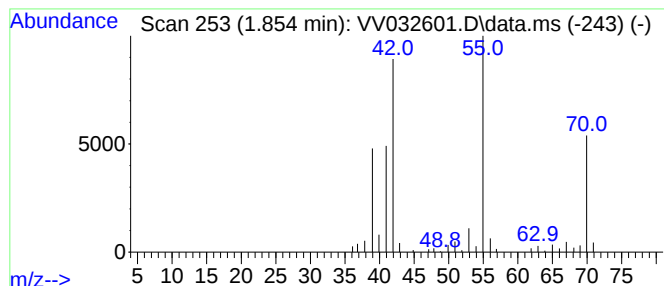
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Cyclic1.854 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.854	1.07 ug/L	59549	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, ethyl-	70	C5H10	001191-96-4	90
2			2-Pentene, (Z)-	70	C5H10	000627-20-3	83
3			2-Pentene	70	C5H10	000109-68-2	80
4			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	68
5			1-Butanol, 3-methyl-	88	C5H12O	000123-51-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AJ1

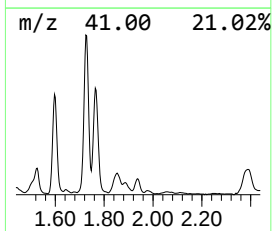
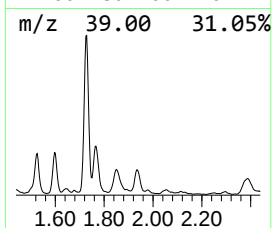
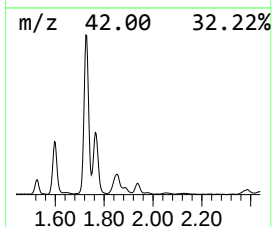
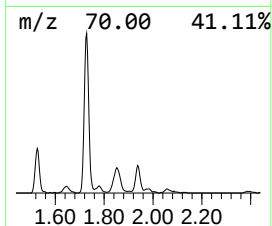
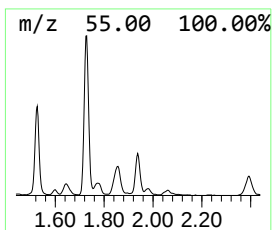
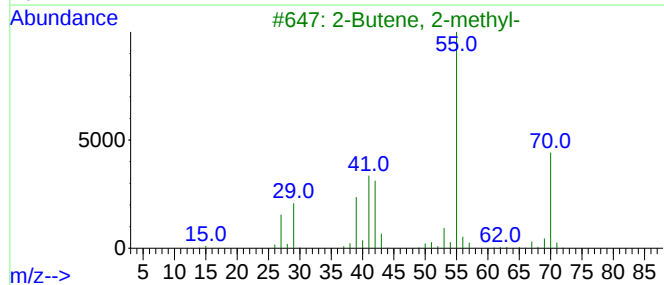
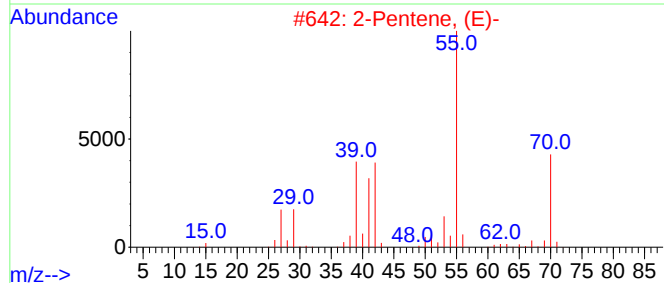
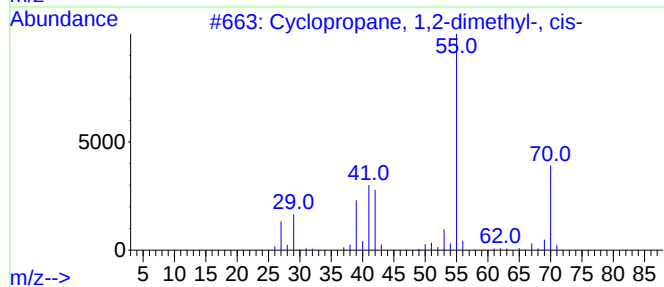
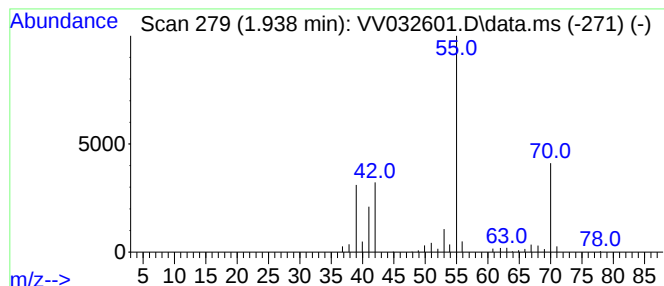
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Cyclic1.938 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.938	0.81 ug/L	44811	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	90
2			2-Pentene, (E)-	70	C5H10	000646-04-8	87
3			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86
4			2-Methyl-1-butene	70	C5H10	000563-46-2	86
5			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AJ1

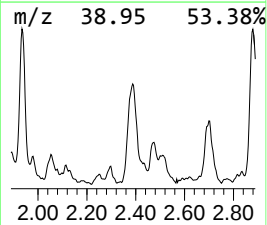
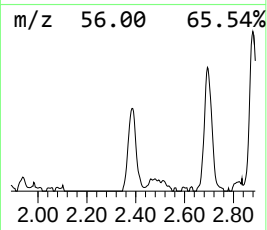
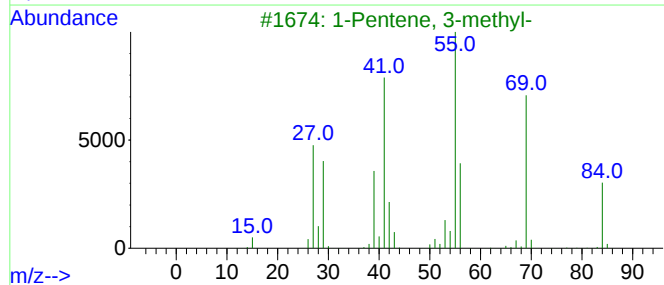
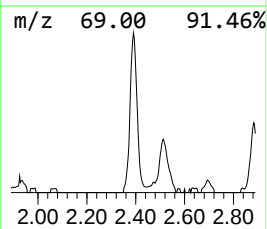
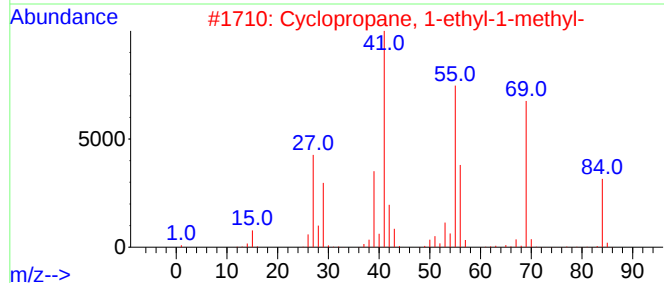
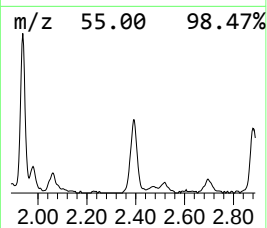
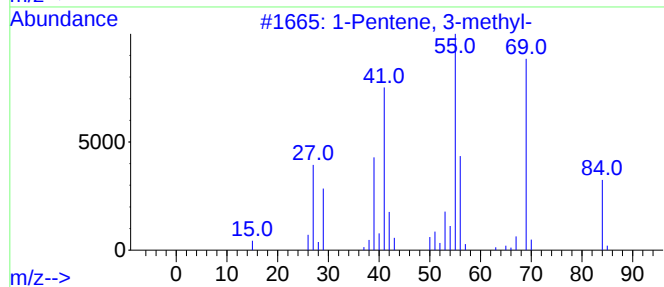
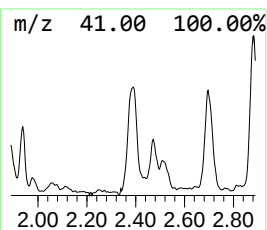
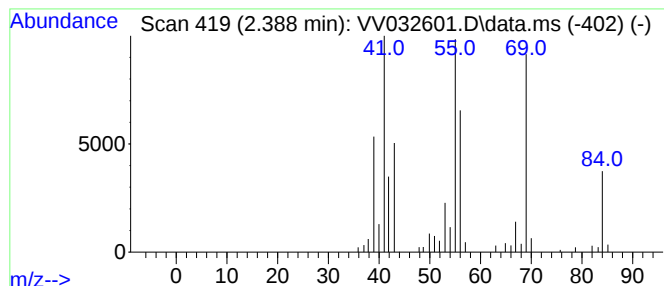
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.388	1.46 ug/L	81183	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 3-methyl-	84	C6H12	000760-20-3	87
2			Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	70
3			1-Pentene, 3-methyl-	84	C6H12	000760-20-3	70
4			Pentane, 3-methylene-	84	C6H12	000760-21-4	62
5			Cyclopropane, 1-ethyl-2-methyl-,...	84	C6H12	019781-68-1	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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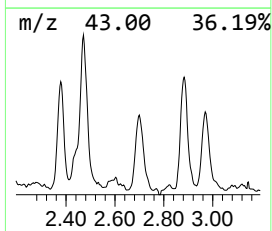
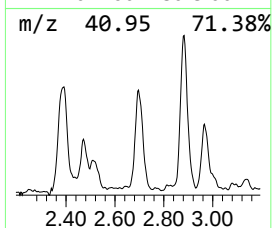
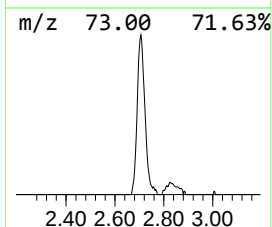
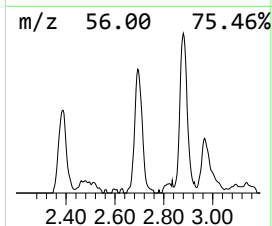
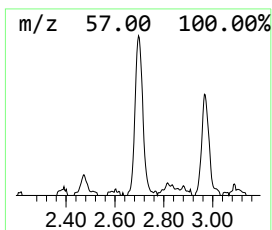
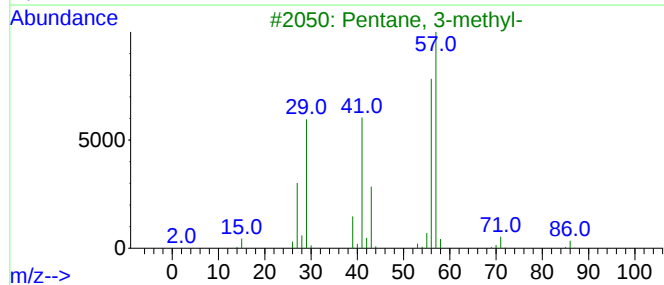
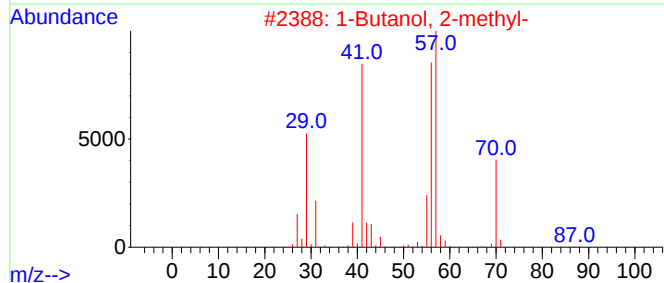
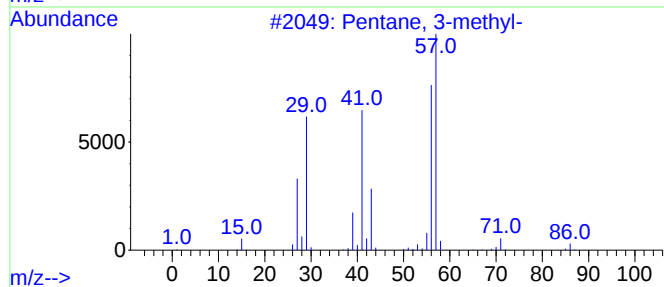
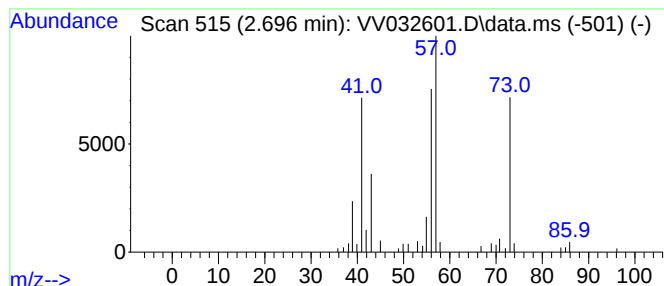
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.696	1.16 ug/L	64450	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
2			1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	59
3			Pentane, 3-methyl-	86	C6H14	000096-14-0	59
4			Pentane, 3-methyl-	86	C6H14	000096-14-0	59
5			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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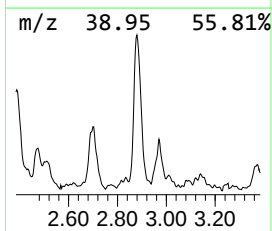
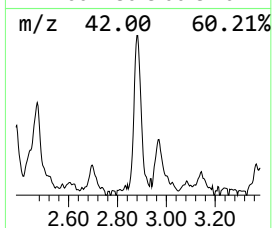
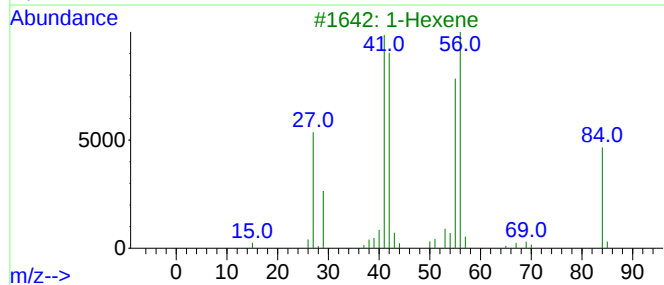
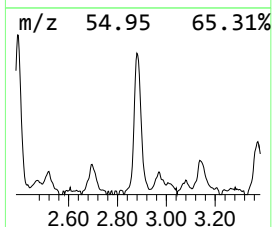
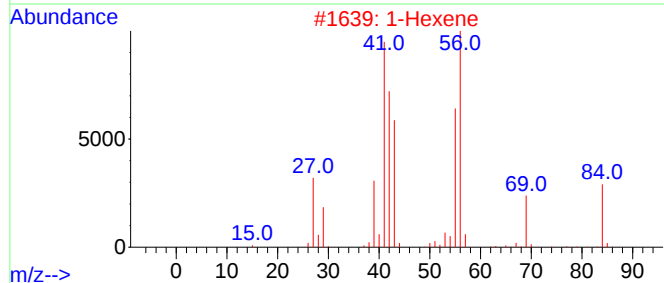
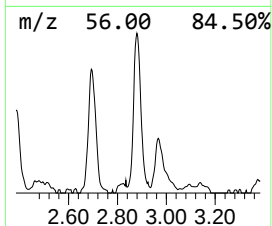
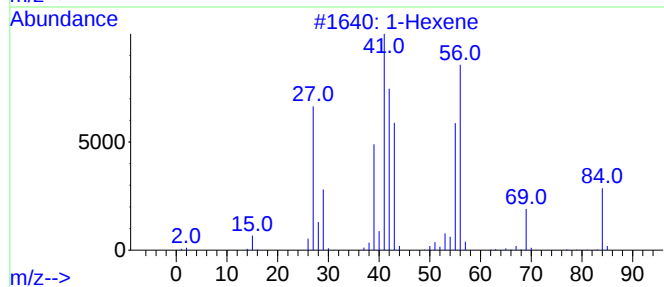
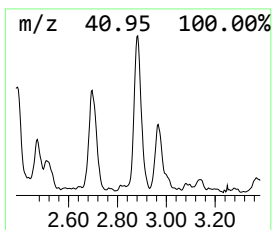
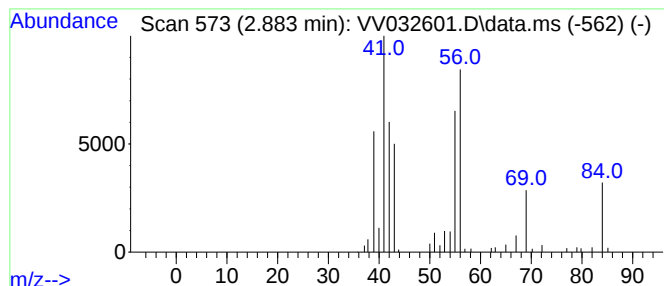
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.883	1.40 ug/L	77764	1,4-Difluorobenzene	5.535

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene		84	C6H12	000592-41-6	91
2	1-Hexene		84	C6H12	000592-41-6	81
3	1-Hexene		84	C6H12	000592-41-6	64
4	2-Hexene, (Z)-		84	C6H12	007688-21-3	58
5	1-Pentene, 2-methyl-		84	C6H12	000763-29-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
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Acq On : 24 Oct 2023 19:43
Operator : SY/MD
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ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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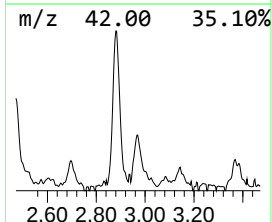
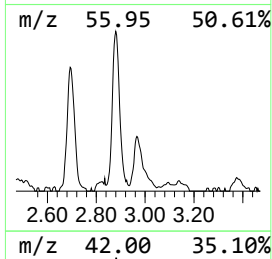
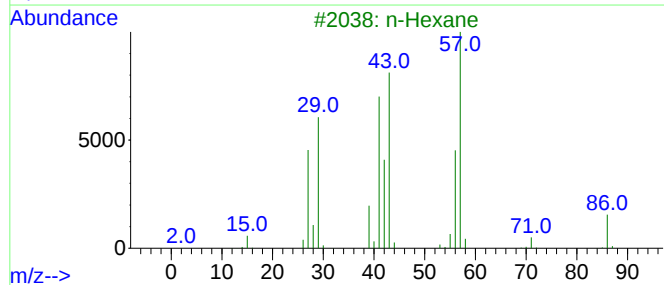
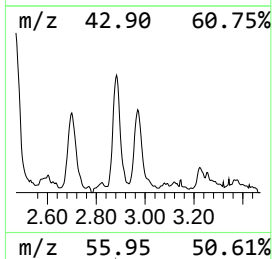
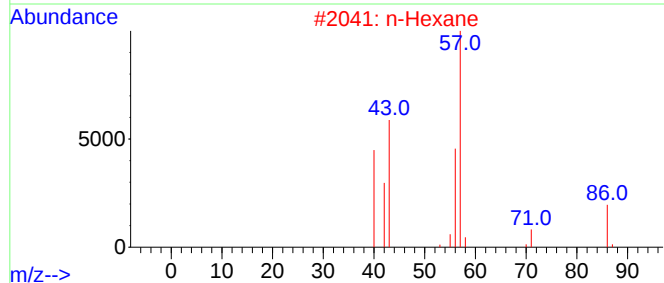
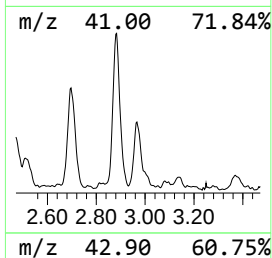
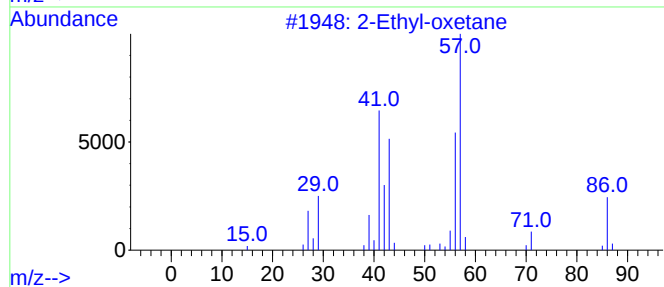
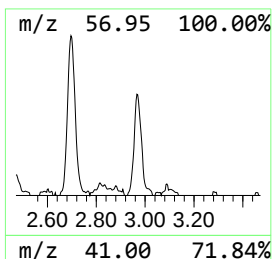
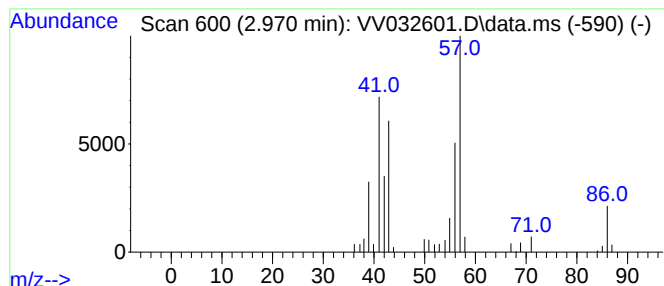
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 2-Ethyl-oxetane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.970	0.72 ug/L	39910	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethyl-oxetane	86	C5H10O	1010386-40-2	87
2			n-Hexane	86	C6H14	000110-54-3	80
3			n-Hexane	86	C6H14	000110-54-3	59
4			n-Hexane	86	C6H14	000110-54-3	59
5			n-Hexane	86	C6H14	000110-54-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
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Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AJ1

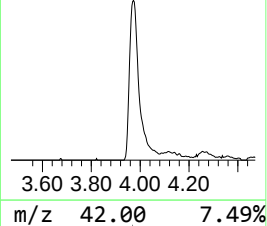
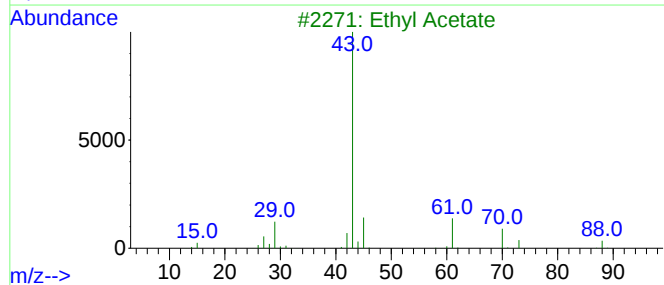
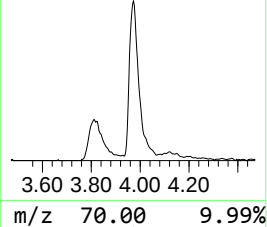
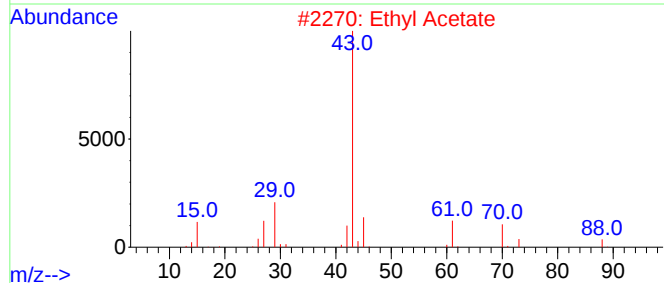
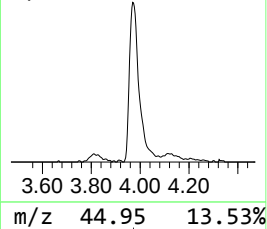
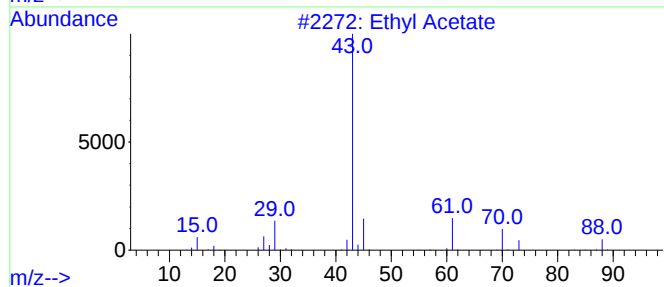
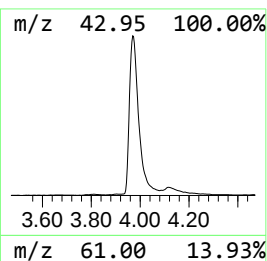
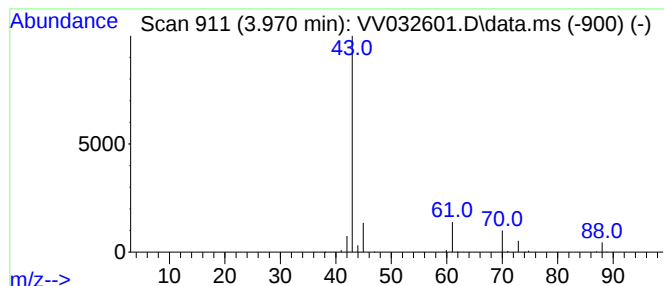
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Ethyl Acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.970	8.21 ug/L	455770	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl Acetate	88	C4H8O2	000141-78-6	86
2			Ethyl Acetate	88	C4H8O2	000141-78-6	86
3			Ethyl Acetate	88	C4H8O2	000141-78-6	86
4			Ethyl Acetate	88	C4H8O2	000141-78-6	86
5			CH3C(O)CH2CH2OH	88	C4H8O2	000590-90-9	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AJ1

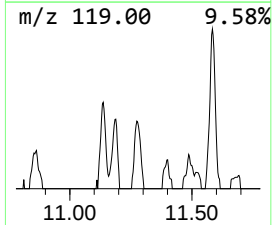
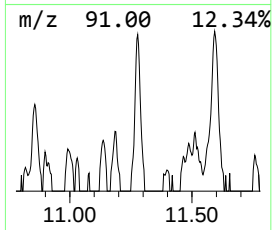
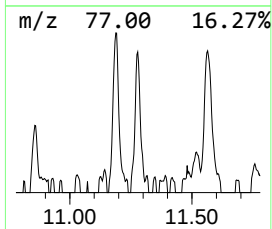
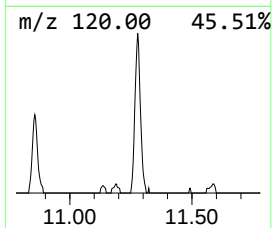
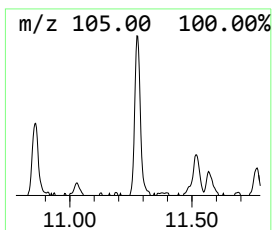
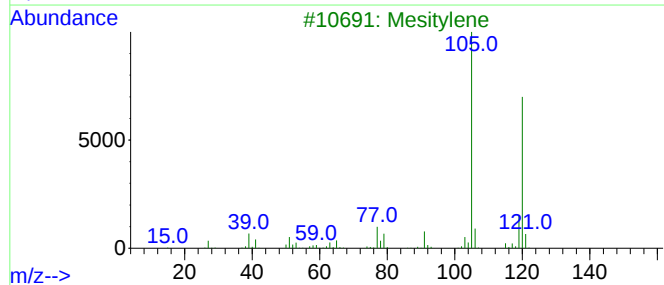
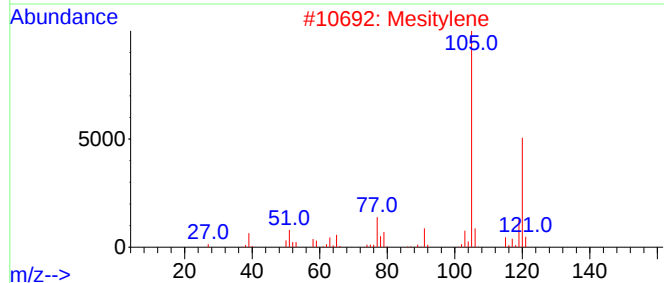
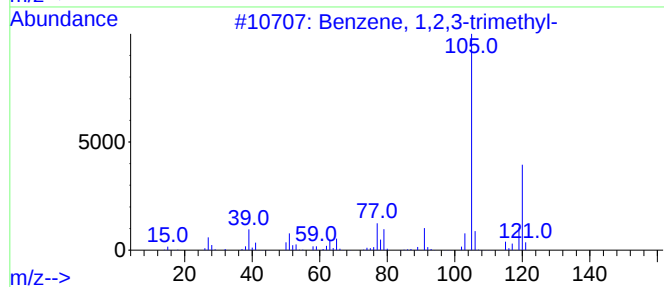
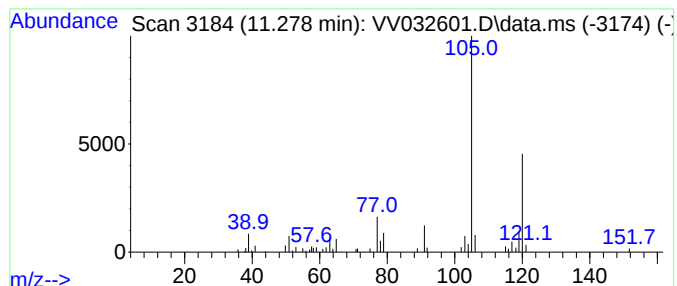
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.278	0.58 ug/L	40530	1,4-Dichlorobenzene-d4	11.188

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2			Mesitylene	120	C9H12	000108-67-8	95
3			Mesitylene	120	C9H12	000108-67-8	94
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102423\
Data File : VV032601.D
Acq On : 24 Oct 2023 19:43
Operator : SY/MD
Sample : 04992-05
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AJ1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.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...	1.192	2.5	ug/L	141488	1	5.535	277448	5.0
(DEL) Alkane: S...	1.388	2.6	ug/L	146481	1	5.535	277448	5.0
(DEL) Alkane: S...	1.597	2.2	ug/L	119980	1	5.535	277448	5.0
(DEL) Alkane: S...	1.725	5.2	ug/L	288157	1	5.535	277448	5.0
(DEL) Alkane: S...	1.764	2.8	ug/L	152793	1	5.535	277448	5.0
(DEL) Alkane: C...	1.854	1.1	ug/L	59549	1	5.535	277448	5.0
(DEL) Alkane: C...	1.938	0.8	ug/L	44811	1	5.535	277448	5.0
(DEL) Alkane: S...	2.388	1.5	ug/L	81183	1	5.535	277448	5.0
(DEL) Alkane: S...	2.696	1.2	ug/L	64450	1	5.535	277448	5.0
(DEL) Alkane: S...	2.883	1.4	ug/L	77764	1	5.535	277448	5.0
2-Ethyl-oxetane	2.970	0.7	ug/L	39910	1	5.535	277448	5.0
Ethyl Acetate	3.970	8.2	ug/L	455770	1	5.535	277448	5.0
Benzene, 1,2,3-...	11.278	0.6	ug/L	40530	3	11.188	348040	5.0