

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
 Data File : VV025473.D
 Acq On : 12 Apr 2022 14:45
 Operator : SY/MD
 Sample : N2372-01
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BFY47

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: VV025473.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.111	15	22	37	rVB	230247	210251	14.59%	2.196%
2	1.217	48	55	60	rBV2	25821	23849	1.65%	0.249%
3	1.307	73	83	90	rBV2	201889	293999	20.40%	3.071%
4	1.420	113	118	132	rVB2	19972	26622	1.85%	0.278%
5	1.571	156	165	177	rVB	86730	102028	7.08%	1.066%
6	1.638	178	186	195	rBV	16334	21409	1.49%	0.224%
7	1.770	219	227	232	rBV	38899	47910	3.32%	0.500%
8	1.809	232	239	259	rVB	118794	165460	11.48%	1.728%
9	1.983	282	293	302	rBV4	16591	26071	1.81%	0.272%
10	2.111	321	333	352	rVB4	197130	344505	23.90%	3.598%
11	2.770	524	538	552	rVB5	9483	22769	1.58%	0.238%
12	3.195	655	670	689	rBV	67600	145420	10.09%	1.519%
13	3.915	879	894	927	rBV2	109011	322165	22.35%	3.365%
14	4.352	1014	1030	1057	rBV2	92711	247596	17.18%	2.586%
15	4.613	1096	1111	1125	rVB4	6140	14686	1.02%	0.153%
16	5.050	1231	1247	1283	rBV2	190293	489505	33.96%	5.112%
17	5.619	1411	1424	1448	rBV	185095	377494	26.19%	3.943%
18	5.918	1504	1517	1533	rBV5	67889	145718	10.11%	1.522%
19	6.072	1551	1565	1590	rBV	112115	232682	16.14%	2.430%
20	6.288	1622	1632	1650	rBV2	33804	72219	5.01%	0.754%
21	6.989	1839	1850	1873	rBV2	49571	94607	6.56%	0.988%
22	7.317	1941	1952	1969	rBV	226389	400127	27.76%	4.179%
23	7.394	1969	1976	1989	rVB	16249	31434	2.18%	0.328%
24	7.625	2038	2048	2066	rBV2	25573	48111	3.34%	0.502%
25	7.976	2146	2157	2176	rVB	117856	201492	13.98%	2.104%
26	8.092	2183	2193	2225	rBV	161326	355169	24.64%	3.709%
27	8.236	2227	2238	2273	rVB	522036	869398	60.32%	9.080%
28	8.850	2417	2429	2453	rBV	294414	489658	33.97%	5.114%
29	9.805	2713	2726	2749	rBV2	98247	167245	11.60%	1.747%
30	10.217	2843	2854	2871	rVB	71915	115108	7.99%	1.202%
31	10.715	2999	3009	3018	rBV	82016	126902	8.80%	1.325%
32	10.757	3018	3022	3035	rVB6	14594	24342	1.69%	0.254%
33	11.146	3133	3143	3163	rBV	291163	450876	31.28%	4.709%
34	11.246	3164	3174	3194	rVB	337700	524150	36.37%	5.474%
35	11.622	3279	3291	3307	rBV	218596	350756	24.34%	3.663%
36	12.008	3401	3411	3429	rBV3	38922	73785	5.12%	0.771%

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 Title : TRACE VOA SFAM1.0

37	12.243	3474	3484	3502	rBV6	8516	21044	1.46%	0.220%
38	12.339	3502	3514	3547	rBV	979946	1441336	100.00%	15.053%
39	13.133	3754	3761	3775	rVB7	15075	24823	1.72%	0.259%
40	13.323	3811	3820	3838	rBV8	8532	18037	1.25%	0.188%
41	13.432	3845	3854	3863	rBV8	14041	22387	1.55%	0.234%
42	14.172	4073	4084	4107	rBV2	149368	255655	17.74%	2.670%
43	14.567	4197	4207	4221	rBV4	9733	18571	1.29%	0.194%
44	14.805	4270	4281	4303	rBV3	25828	51361	3.56%	0.536%
45	15.143	4379	4386	4405	rVB5	31526	51563	3.58%	0.539%
46	16.178	4700	4708	4720	rBV9	9758	14604	1.01%	0.153%

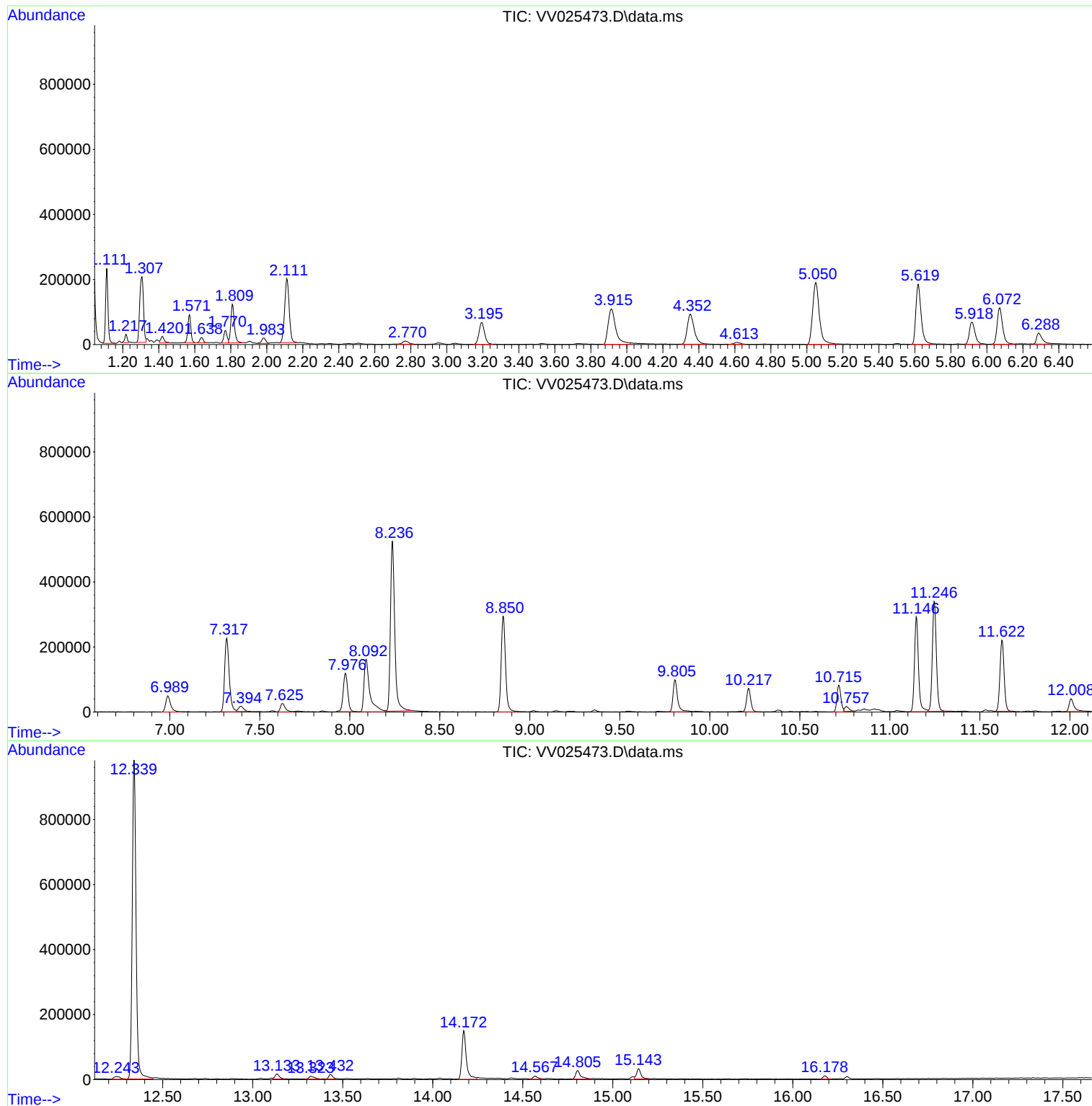
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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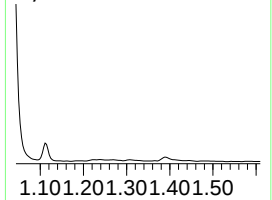
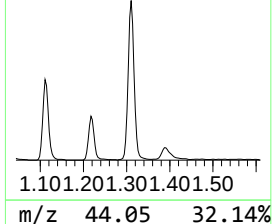
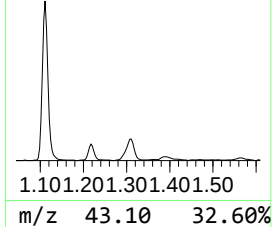
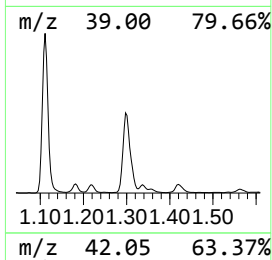
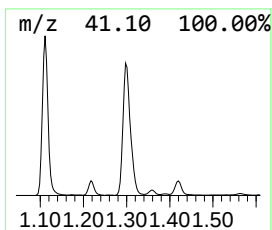
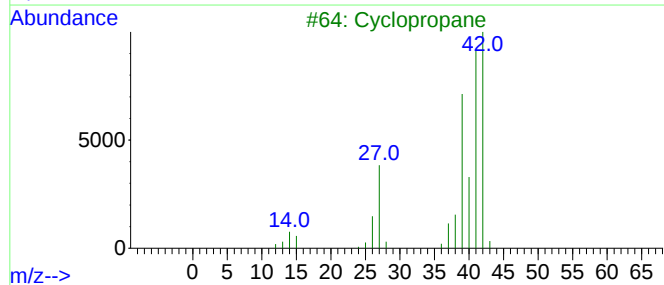
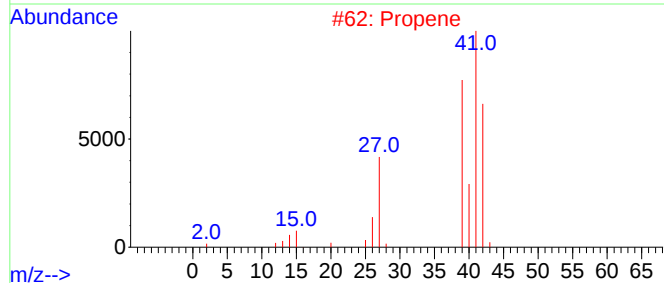
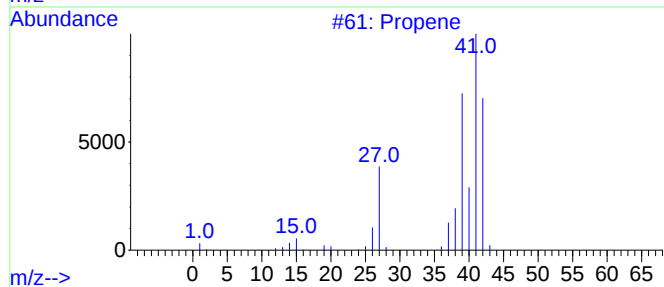
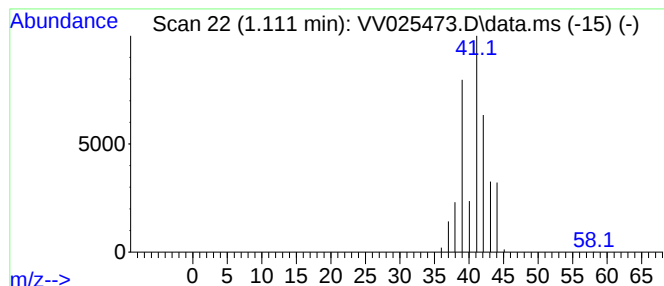
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.111	2.78 ug/L	210251	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	90
2			Propene	42	C3H6	000115-07-1	42
3			Cyclopropane	42	C3H6	000075-19-4	9
4			Cyclopropene	40	C3H4	002781-85-3	9
5			Cyclopropane	42	C3H6	000075-19-4	7



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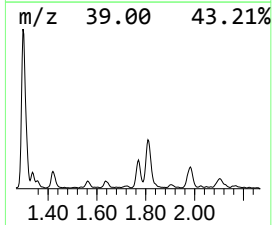
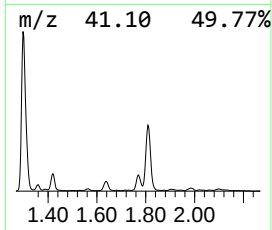
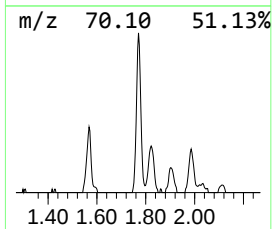
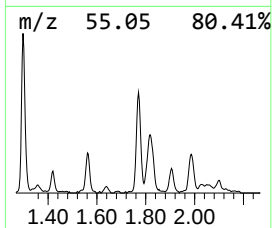
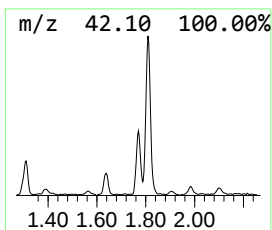
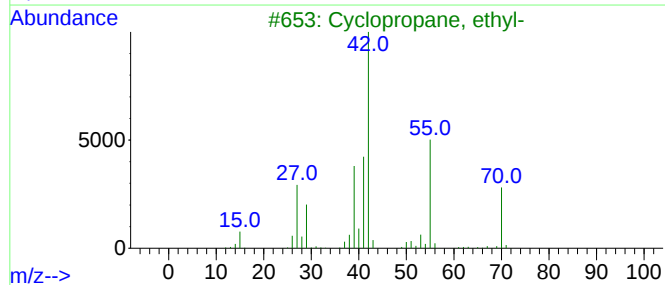
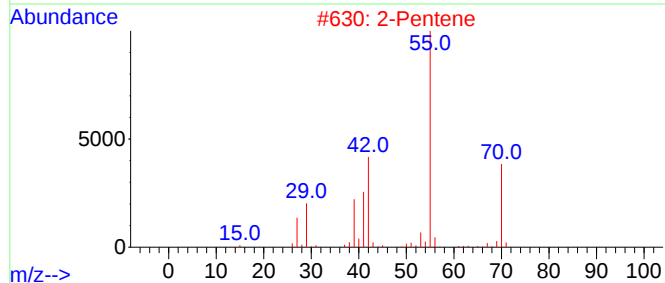
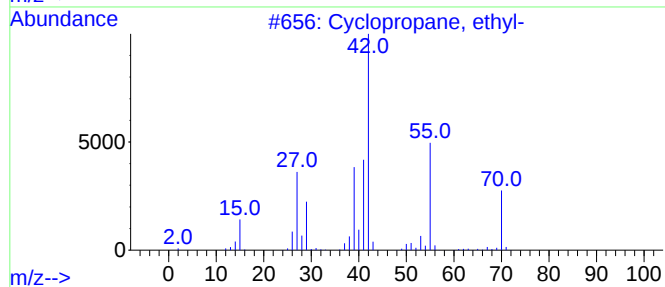
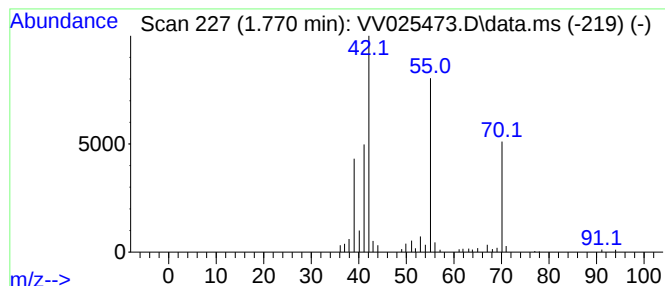
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Cyclic1.770 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.770	0.63 ug/L	47910	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, ethyl-	70	C5H10	001191-96-4	83
2			2-Pentene	70	C5H10	000109-68-2	72
3			Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
4			1-Pentene	70	C5H10	000109-67-1	58
5			1-Pentene	70	C5H10	000109-67-1	58



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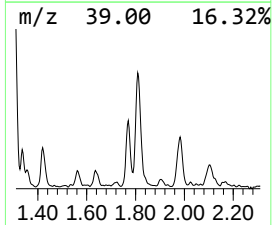
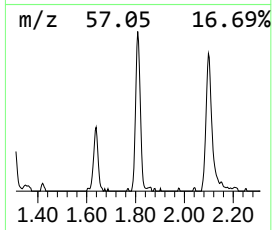
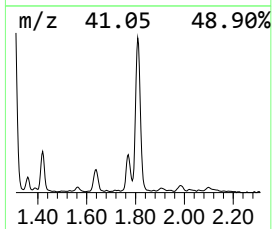
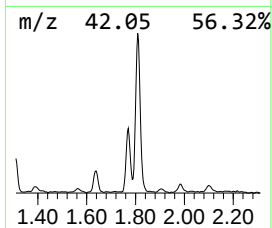
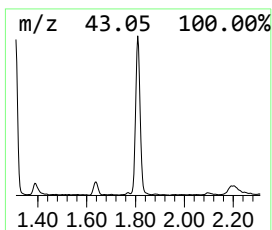
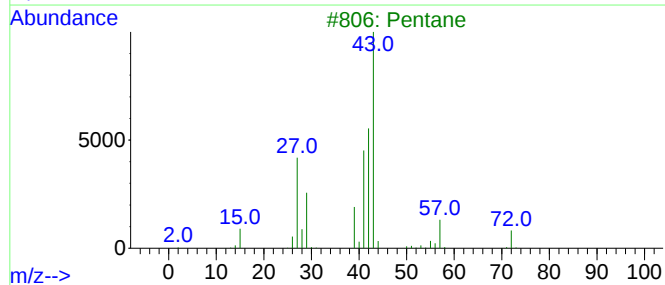
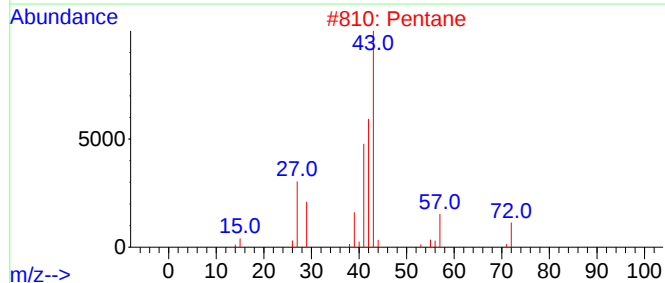
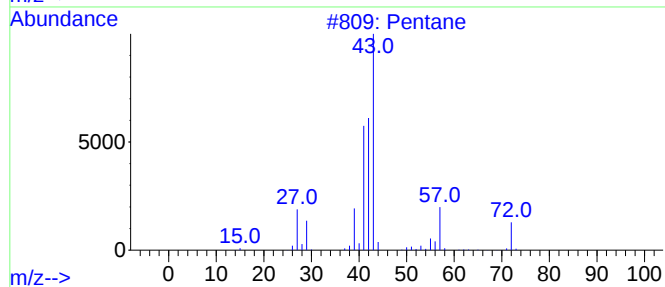
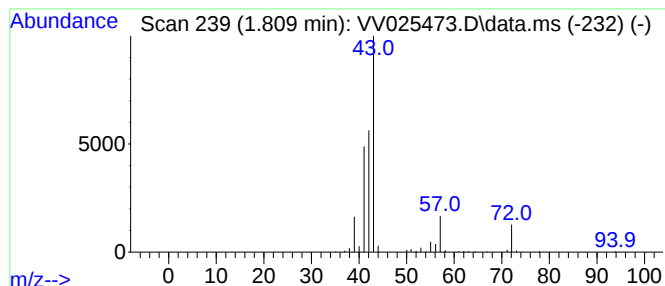
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.809	2.19 ug/L	165460	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane			72	C5H12	000109-66-0	91
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	80



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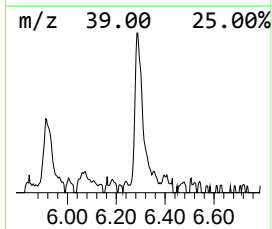
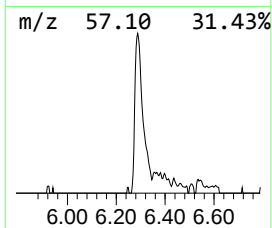
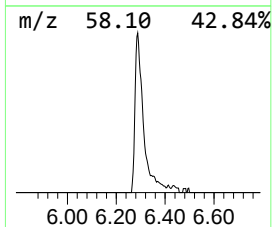
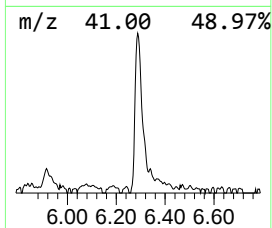
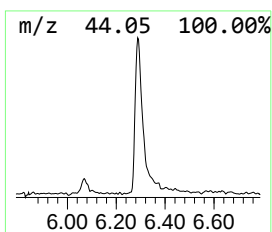
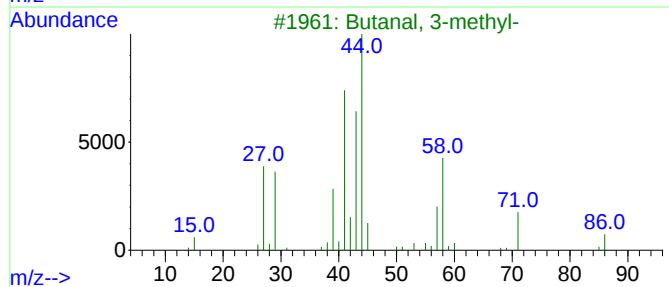
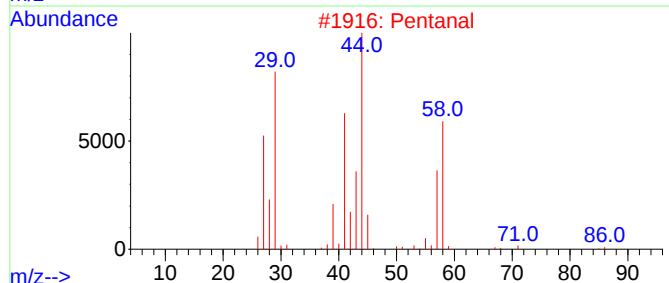
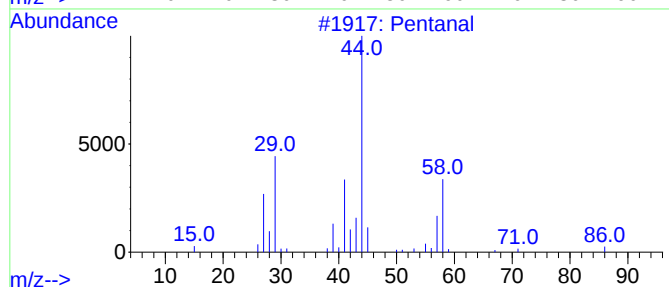
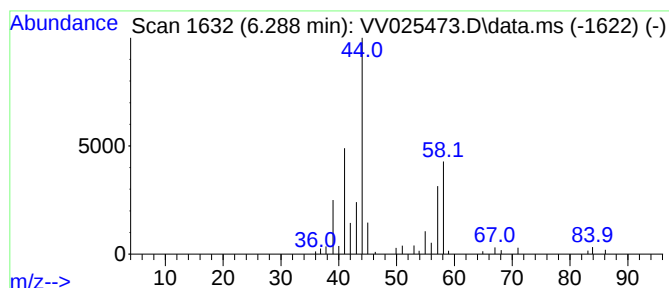
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Peak Number 4 Pentanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.288	0.96 ug/L	72219	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanal			86	C5H10O	000110-62-3	90
2	Pentanal			86	C5H10O	000110-62-3	78
3	Butanal, 3-methyl-			86	C5H10O	000590-86-3	74
4	Pentanal			86	C5H10O	000110-62-3	72
5	Pentanal			86	C5H10O	000110-62-3	64



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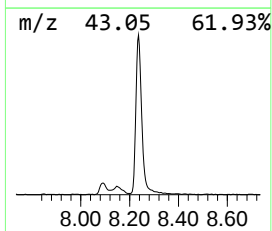
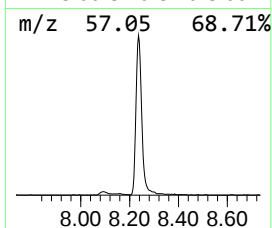
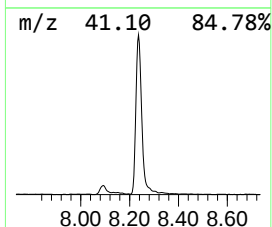
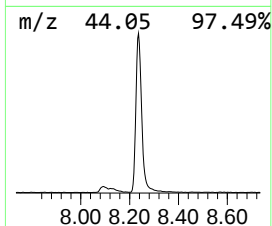
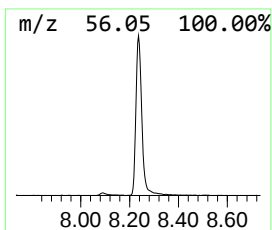
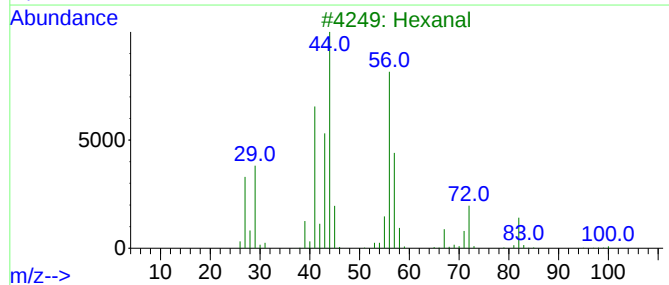
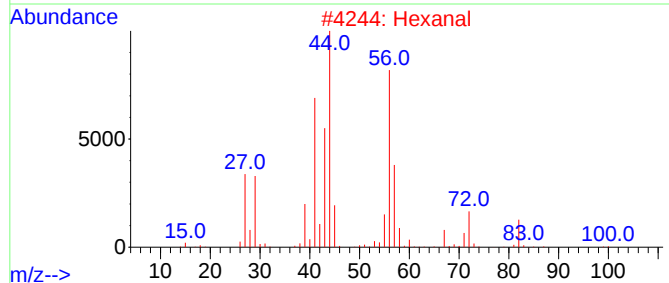
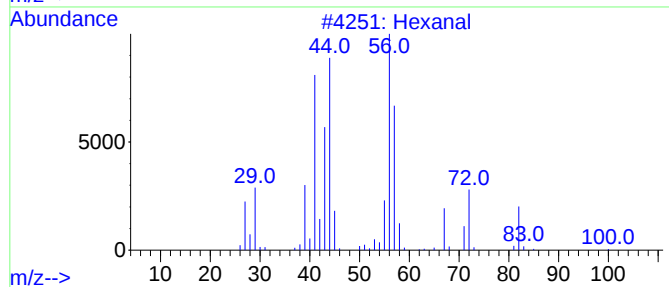
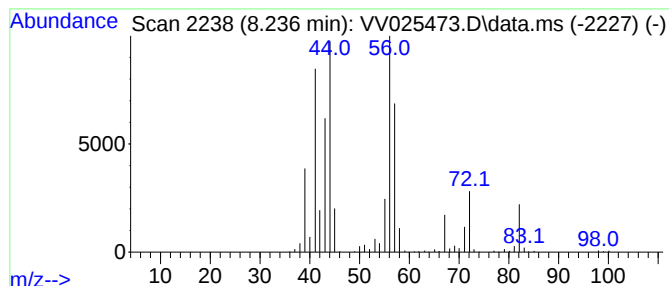
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Hexanal Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.236	8.88 ug/L	869398	Chlorobenzene-d5	8.854

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal			100	C6H12O	000066-25-1	90
2	Hexanal			100	C6H12O	000066-25-1	81
3	Hexanal			100	C6H12O	000066-25-1	74
4	Hexanal			100	C6H12O	000066-25-1	59
5	Hexanal			100	C6H12O	000066-25-1	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFY47

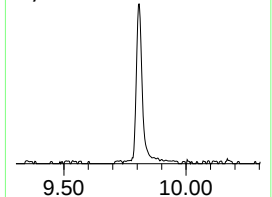
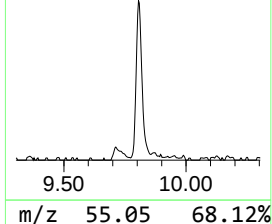
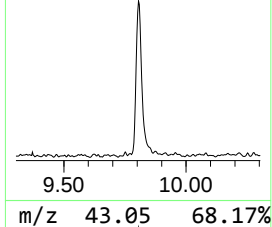
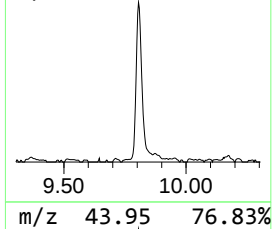
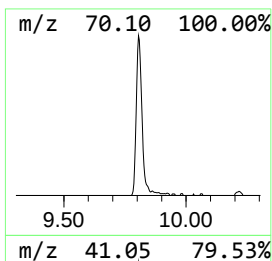
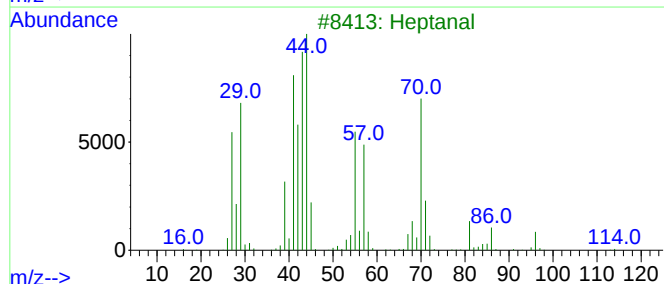
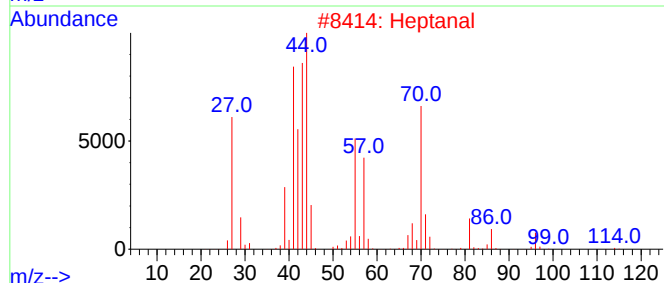
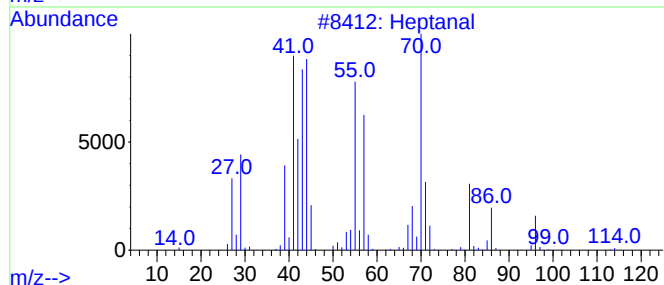
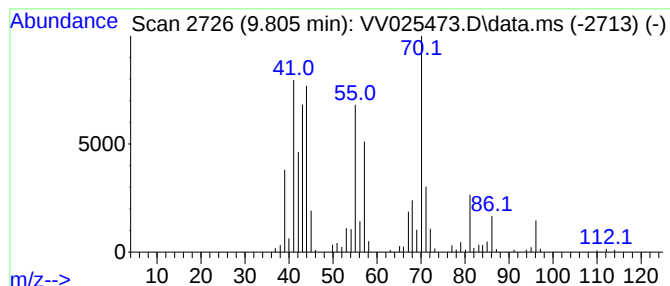
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Heptanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.805	1.71 ug/L	167245	Chlorobenzene-d5	8.854

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanal			114	C7H14O	000111-71-7	96
2	Heptanal			114	C7H14O	000111-71-7	95
3	Heptanal			114	C7H14O	000111-71-7	95
4	Heptanal			114	C7H14O	000111-71-7	93
5	Heptanal			114	C7H14O	000111-71-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFY47

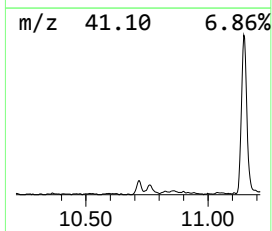
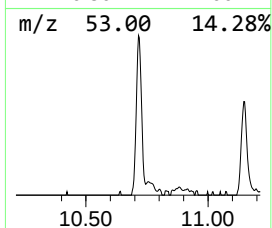
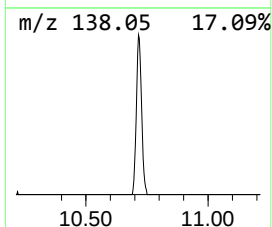
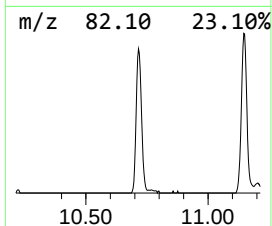
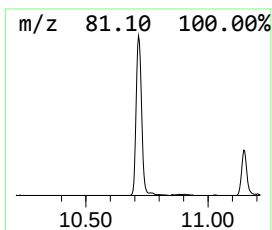
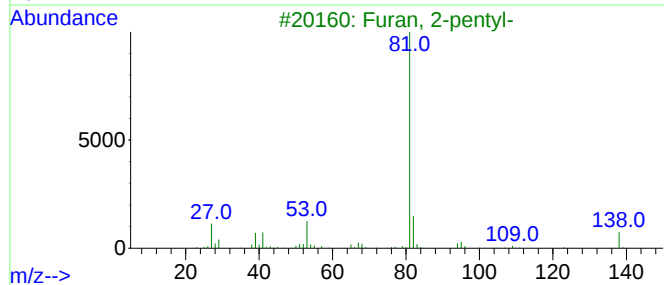
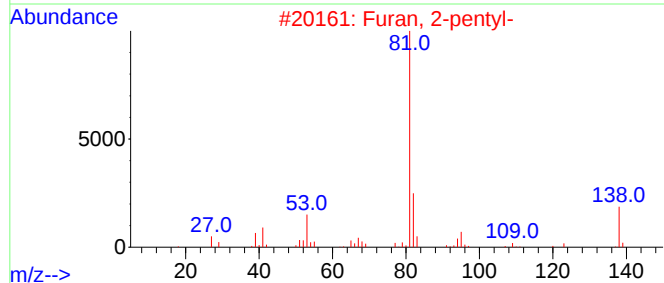
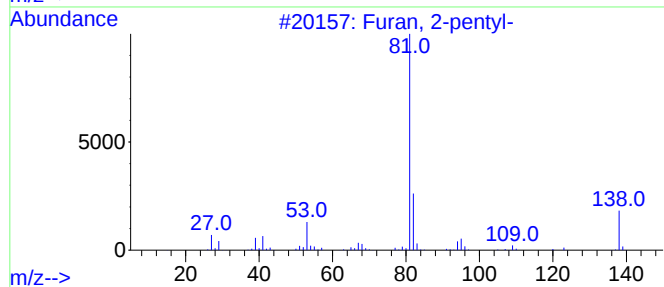
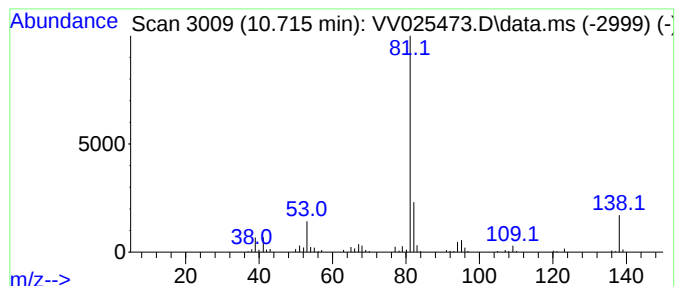
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Furan, 2-pentyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.715	1.21 ug/L	126902	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, 2-pentyl-	138	C9H14O	003777-69-3	91
2			Furan, 2-pentyl-	138	C9H14O	003777-69-3	90
3			Furan, 2-pentyl-	138	C9H14O	003777-69-3	72
4			2-[(2-Furylmethyl)amino]-2-methy...	169	C9H15NO2	889949-94-4	45
5			2,4-Nonadienal, (E,E)-	138	C9H14O	005910-87-2	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFY47

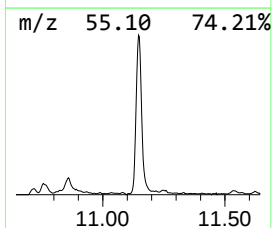
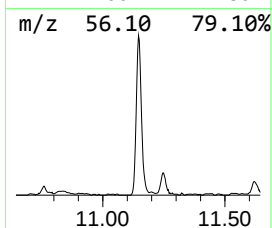
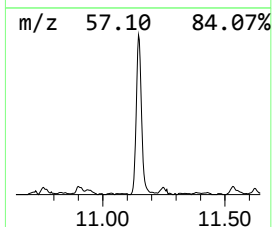
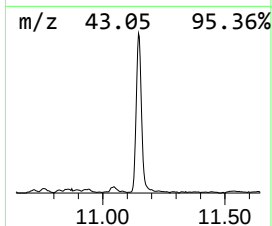
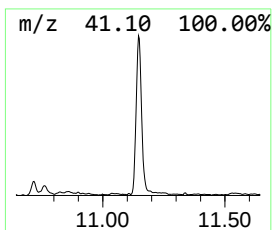
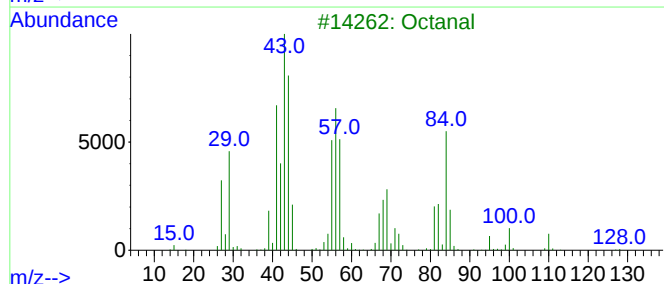
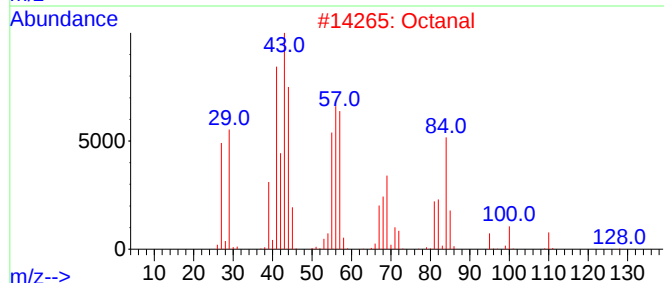
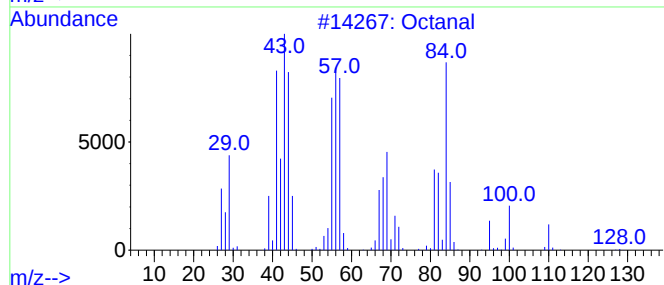
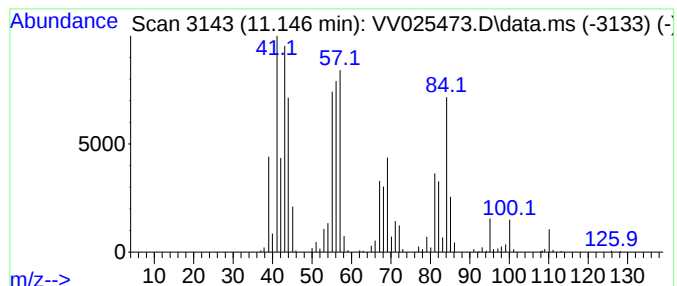
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Octanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.146	4.30 ug/L	450876	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanal			128	C8H16O	000124-13-0	97
2	Octanal			128	C8H16O	000124-13-0	94
3	Octanal			128	C8H16O	000124-13-0	93
4	Octanal			128	C8H16O	000124-13-0	91
5	Octanal			128	C8H16O	000124-13-0	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

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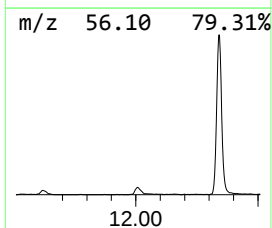
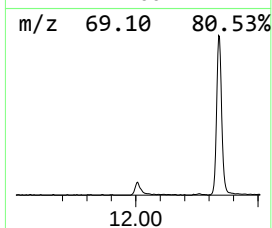
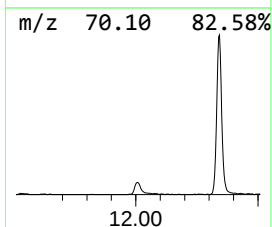
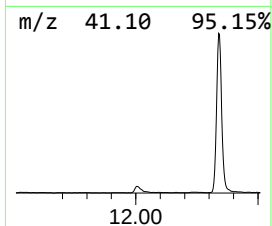
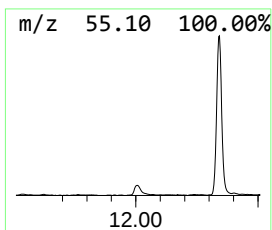
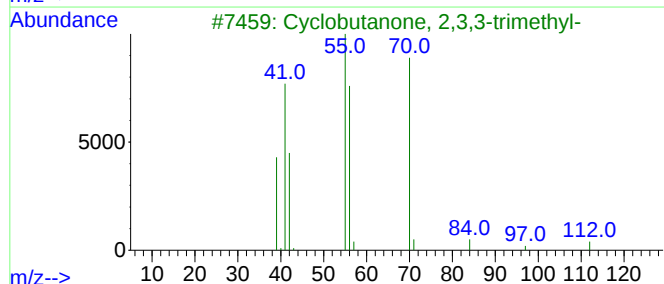
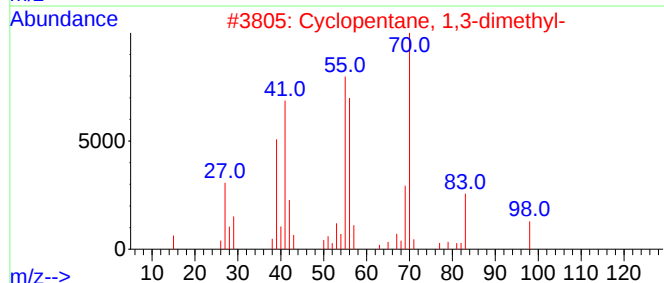
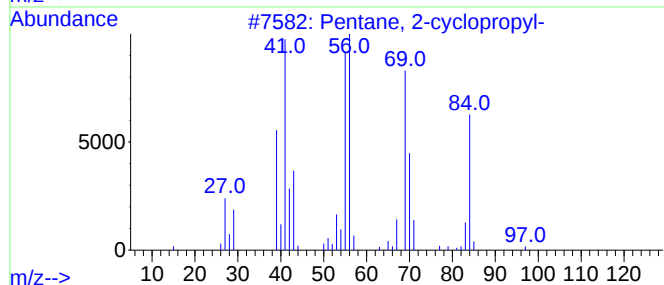
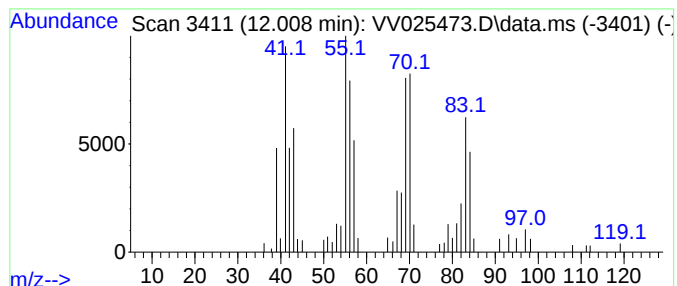
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 (DEL) Alkane: Cyclic12.008 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.008	0.70 ug/L	73785	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-cyclopropyl-	112	C8H16	005458-16-2	49
2			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	38
3			Cyclobutanone, 2,3,3-trimethyl-	112	C7H12O	028290-01-9	38
4			Pentane, 3-methylene-	84	C6H12	000760-21-4	30
5			Pentane, 3-methylene-	84	C6H12	000760-21-4	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

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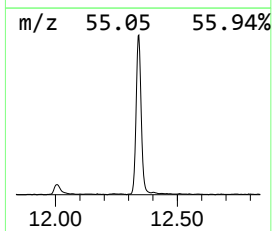
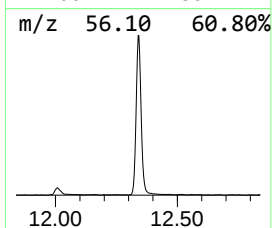
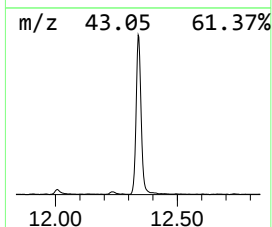
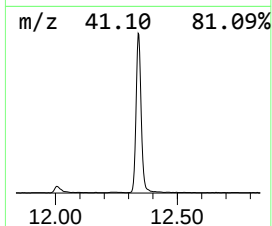
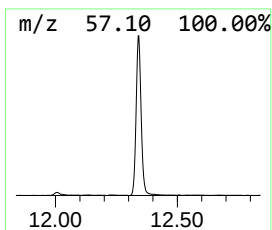
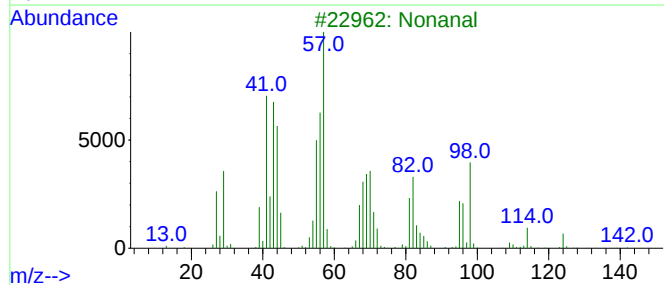
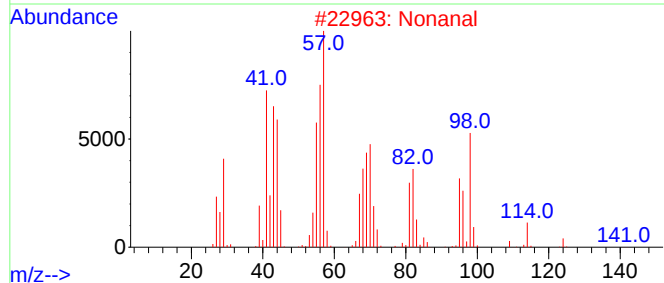
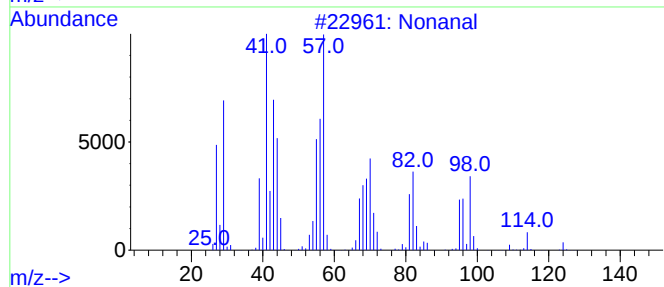
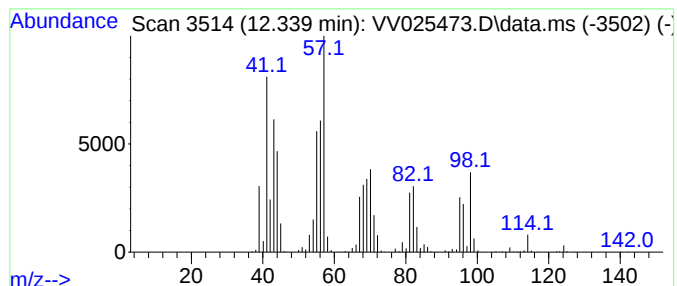
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Nonanal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.339	13.75 ug/L	1441340	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanal	142	C9H18O	000124-19-6	91
2			Nonanal	142	C9H18O	000124-19-6	91
3			Nonanal	142	C9H18O	000124-19-6	91
4			Nonanal	142	C9H18O	000124-19-6	90
5			Heptane, 2-chloro-	134	C7H15Cl	001001-89-4	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFY47

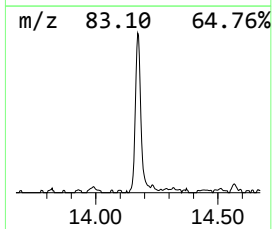
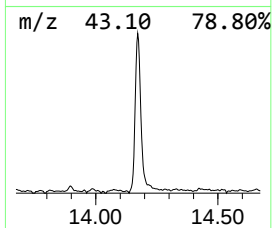
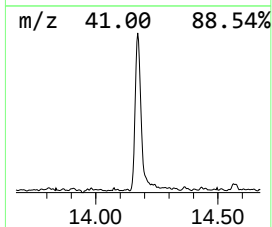
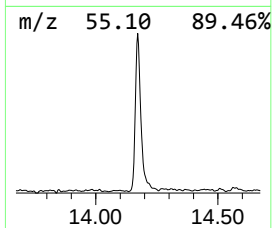
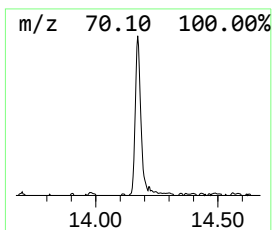
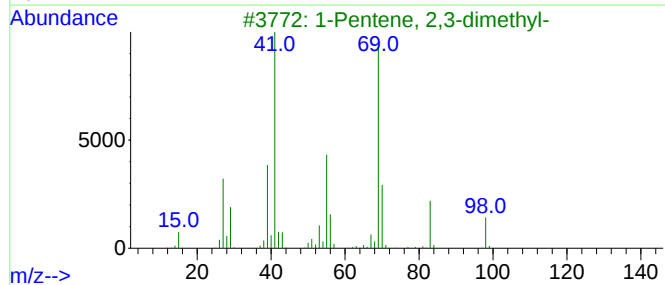
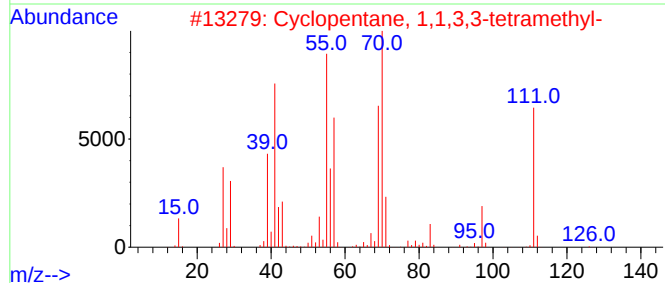
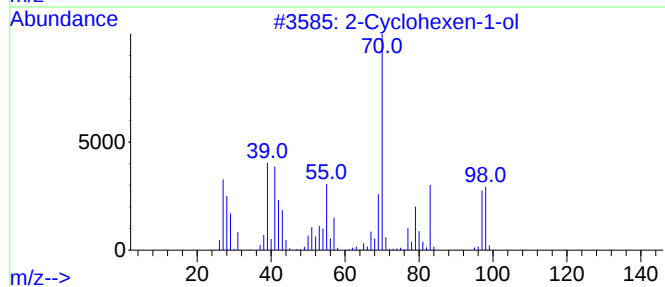
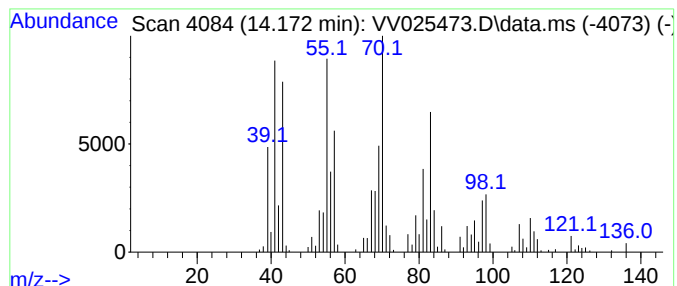
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 2-Cyclohexen-1-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.172	2.44 ug/L	255655	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-ol	98	C6H10O	000822-67-3	53
2			Cyclopentane, 1,1,3,3-tetramethyl-	126	C9H18	050876-33-0	46
3			1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	38
4			n-Heptadecanol-1	256	C17H36O	001454-85-9	38
5			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025473.D
Acq On : 12 Apr 2022 14:45
Operator : SY/MD
Sample : N2372-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFY47

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: C...	1.770	0.6	ug/L	47910	1	5.619	377494	5.0
(DEL) Alkane: S...	1.809	2.2	ug/L	165460	1	5.619	377494	5.0
Pentanal	6.288	1.0	ug/L	72219	1	5.619	377494	5.0
Hexanal	8.236	8.9	ug/L	869398	2	8.854	489658	5.0
Heptanal	9.805	1.7	ug/L	167245	2	8.854	489658	5.0
Furan, 2-pentyl-	10.715	1.2	ug/L	126902	3	11.246	524150	5.0
Octanal	11.146	4.3	ug/L	450876	3	11.246	524150	5.0
(DEL) Alkane: C...	12.008	0.7	ug/L	73785	3	11.246	524150	5.0
Nonanal	12.339	13.8	ug/L	1441340	3	11.246	524150	5.0
2-Cyclohexen-1-ol	14.172	2.4	ug/L	255655	3	11.246	524150	5.0