

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
 Data File : VV034036.D
 Acq On : 07 Feb 2024 17:19
 Operator : SY/MD
 Sample : P1383-16
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 E0YB9

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\SFAMVTR012524WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VV034036.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	8	14	31	rBV	545179	552242	75.14%	8.964%
2	1.191	41	47	63	rVB	63358	56156	7.64%	0.912%
3	1.272	63	72	86	rBV2	340790	507001	68.98%	8.229%
4	1.326	86	89	94	rVV	12728	12196	1.66%	0.198%
5	1.359	94	99	102	rVV	9002	10975	1.49%	0.178%
6	1.384	102	107	119	rVB	51753	59132	8.05%	0.960%
7	1.532	142	153	165	rVB2	53147	75577	10.28%	1.227%
8	1.596	165	173	187	rBV	56448	70378	9.58%	1.142%
9	1.728	205	214	220	rBV	78828	99794	13.58%	1.620%
10	1.770	220	227	245	rVB3	85312	156898	21.35%	2.547%
11	1.857	247	254	264	rVB	8563	11008	1.50%	0.179%
12	1.934	267	278	286	rBV3	33226	48440	6.59%	0.786%
13	1.976	286	291	302	rVB	8279	10817	1.47%	0.176%
14	2.059	306	317	328	rBV	84562	126306	17.19%	2.050%
15	2.294	382	390	404	rVB	10782	16983	2.31%	0.276%
16	2.387	405	419	426	rBV5	8498	19656	2.67%	0.319%
17	2.471	437	445	452	rBV3	7157	12496	1.70%	0.203%
18	2.513	452	458	470	rVB3	8679	14933	2.03%	0.242%
19	2.696	500	515	530	rVB3	8721	18934	2.58%	0.307%
20	2.879	560	572	589	rVV6	22052	49836	6.78%	0.809%
21	2.969	589	600	615	rVB2	12837	25763	3.51%	0.418%
22	3.079	622	634	643	rBV6	4462	9519	1.30%	0.155%
23	3.375	711	726	739	rBV5	5753	14360	1.95%	0.233%
24	3.667	801	817	834	rBV7	8728	25394	3.46%	0.412%
25	3.821	850	865	897	rBV3	31568	140766	19.15%	2.285%
26	4.005	918	922	940	rVB5	3187	7382	1.00%	0.120%
27	4.252	983	999	1023	rBV2	56847	148762	20.24%	2.415%
28	4.580	1087	1101	1117	rBV5	7335	18682	2.54%	0.303%
29	4.825	1162	1177	1195	rBV7	2994	8672	1.18%	0.141%
30	4.960	1204	1219	1252	rBV2	129489	366372	49.85%	5.947%
31	5.236	1293	1305	1315	rBV7	4726	10396	1.41%	0.169%
32	5.300	1317	1325	1346	rVB7	3985	11411	1.55%	0.185%
33	5.416	1350	1361	1372	rBV5	4296	9991	1.36%	0.162%
34	5.535	1385	1398	1433	rBV	118961	250876	34.13%	4.072%
35	5.847	1485	1495	1508	rBV	5823	14422	1.96%	0.234%
36	5.992	1528	1540	1553	rBV	71428	151033	20.55%	2.452%

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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\SFAMVTR012524WMA.M
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37	6.230	1604	1614	1623	rBV4	5594	11974	1.63%	0.194%
38	6.304	1628	1637	1683	rVB	94589	197272	26.84%	3.202%
39	6.918	1817	1828	1853	rBV2	34242	66653	9.07%	1.082%
40	7.246	1919	1930	1942	rVV	158634	280132	38.11%	4.547%
41	7.316	1942	1952	1979	rVB	419334	734969	100.00%	11.930%
42	7.561	2019	2028	2049	rVB2	18899	38763	5.27%	0.629%
43	7.876	2113	2126	2136	rBV3	8216	15731	2.14%	0.255%
44	8.027	2164	2173	2210	rBV	157326	315347	42.91%	5.119%
45	8.181	2212	2221	2236	rVV2	30942	59542	8.10%	0.966%
46	8.249	2236	2242	2262	rVB	12166	24417	3.32%	0.396%
47	8.786	2398	2409	2428	rBV	183882	301242	40.99%	4.890%
48	8.953	2451	2461	2473	rBV2	20551	33354	4.54%	0.541%
49	9.075	2491	2499	2516	rBV	27961	47309	6.44%	0.768%
50	9.493	2617	2629	2659	rBV2	59338	115530	15.72%	1.875%
51	10.156	2823	2835	2851	rBV	51423	82716	11.25%	1.343%
52	10.326	2881	2888	2900	rVB2	10101	16655	2.27%	0.270%
53	10.390	2900	2908	2923	rBV2	6055	12861	1.75%	0.209%
54	10.480	2923	2936	2944	rVB6	3626	7865	1.07%	0.128%
55	10.673	2985	2996	3004	rBV5	4609	9189	1.25%	0.149%
56	10.857	3044	3053	3063	rBV	15229	22726	3.09%	0.369%
57	11.098	3121	3128	3137	rBV4	6703	10696	1.46%	0.174%
58	11.188	3146	3156	3175	rBV	197499	306938	41.76%	4.982%
59	11.278	3177	3184	3193	rVB	5286	7488	1.02%	0.122%
60	11.490	3237	3250	3262	rBV6	8252	20796	2.83%	0.338%
61	11.561	3262	3272	3298	rVB	160574	258698	35.20%	4.199%
62	12.291	3490	3499	3514	rBV4	11405	18368	2.50%	0.298%

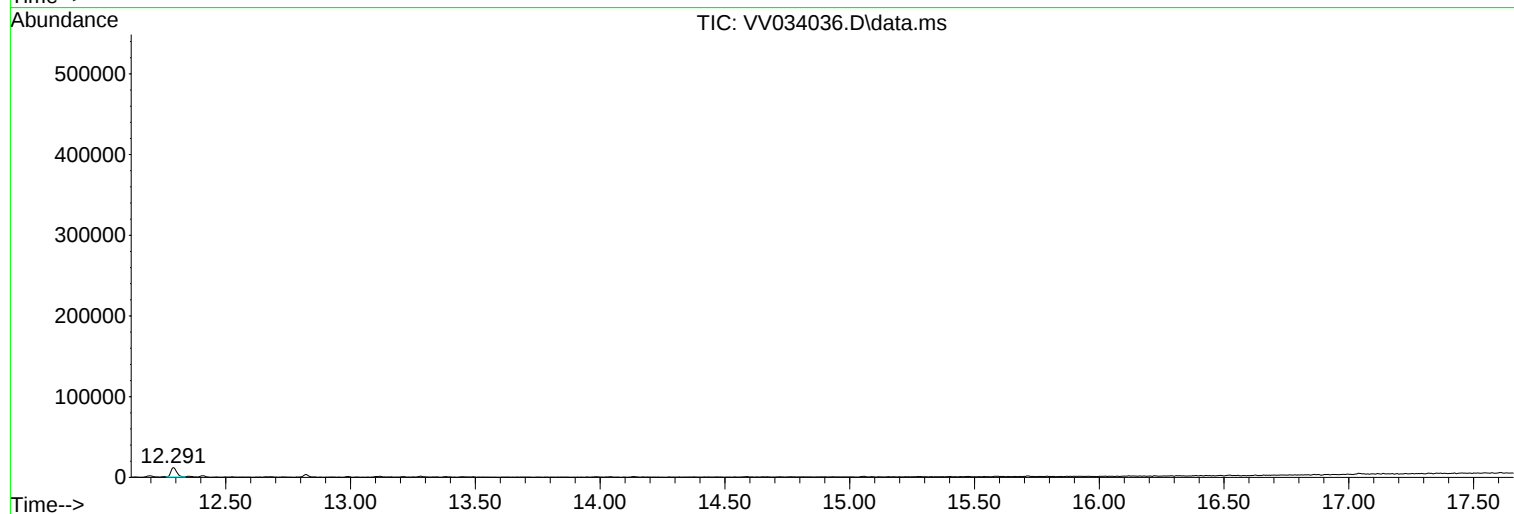
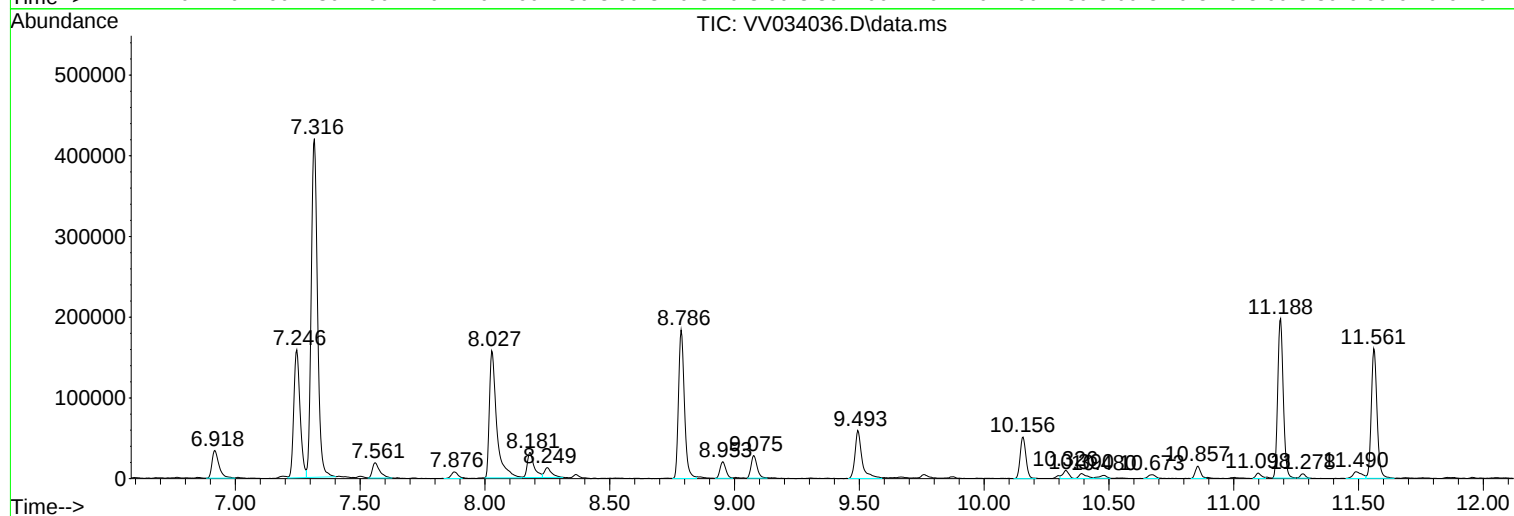
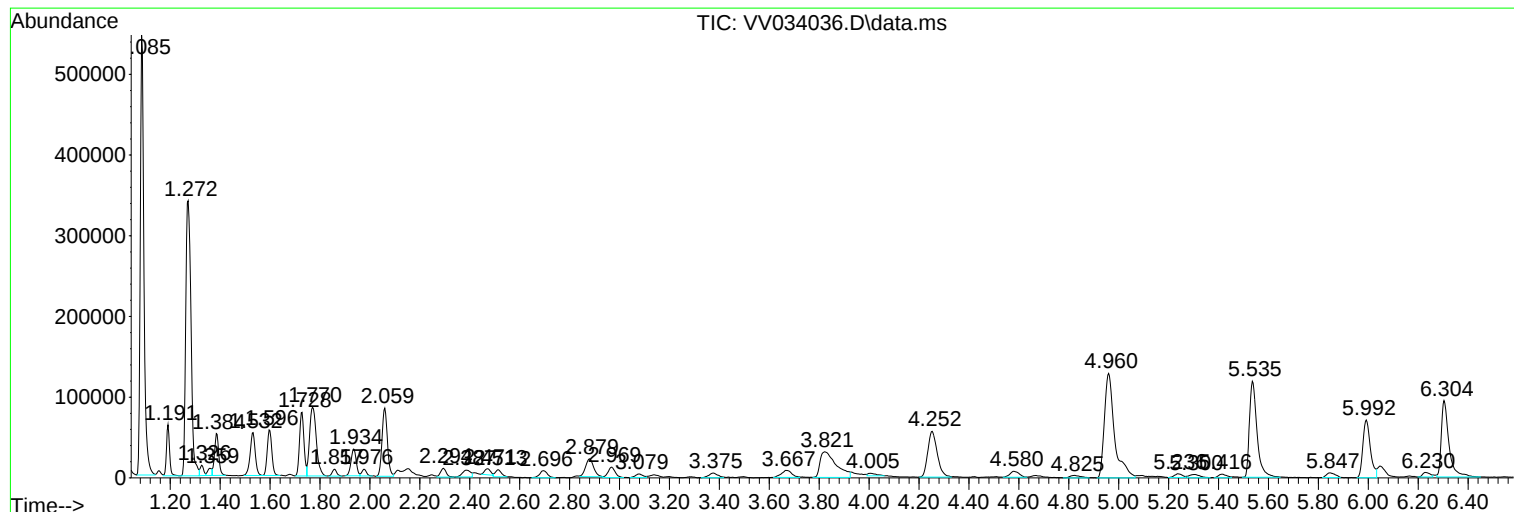
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR012524WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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MSVOA_V
ClientSampleId :
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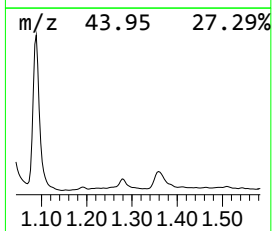
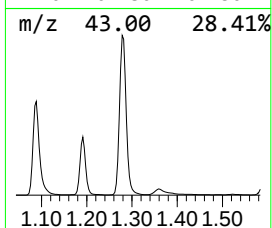
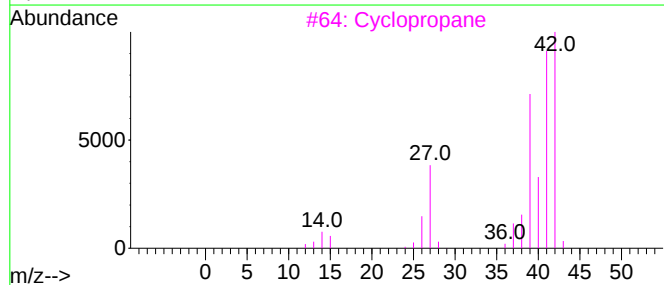
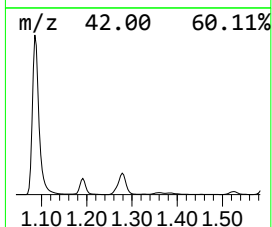
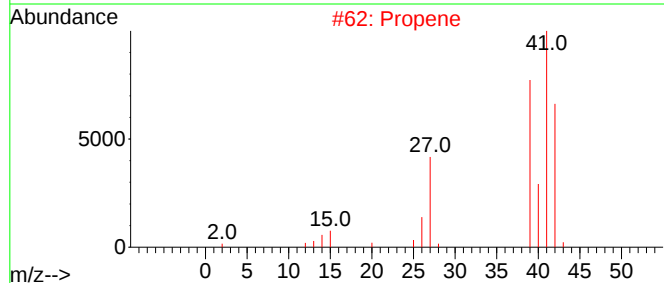
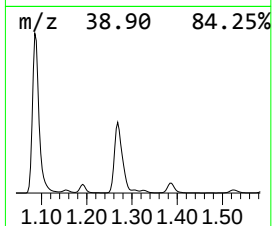
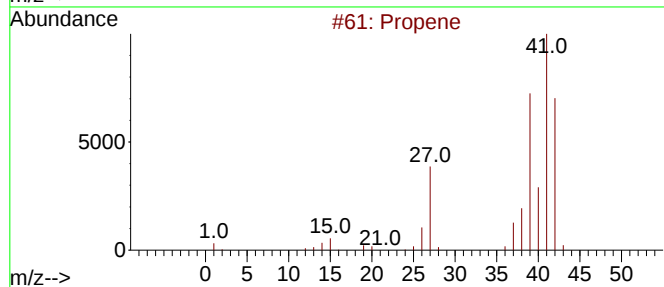
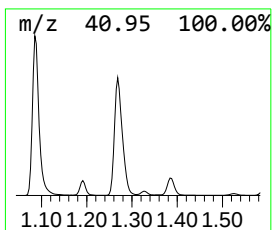
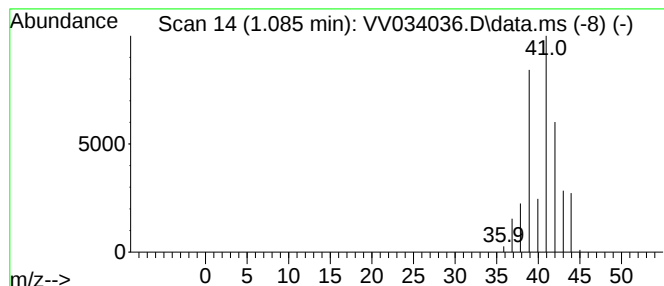
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.085	11.01 ug/L	552242	1,4-Difluorobenzene	5.535

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	7
4			Cyclopropane	42	C3H6	000075-19-4	7
5			Cyclopropene	40	C3H4	002781-85-3	5



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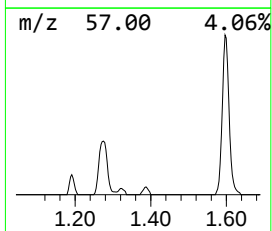
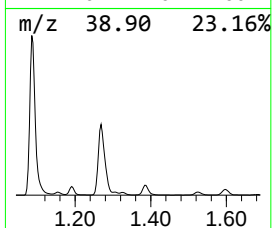
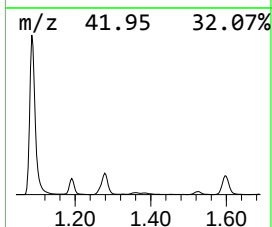
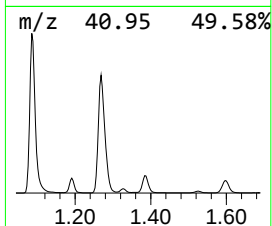
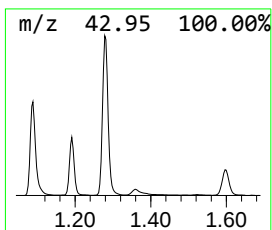
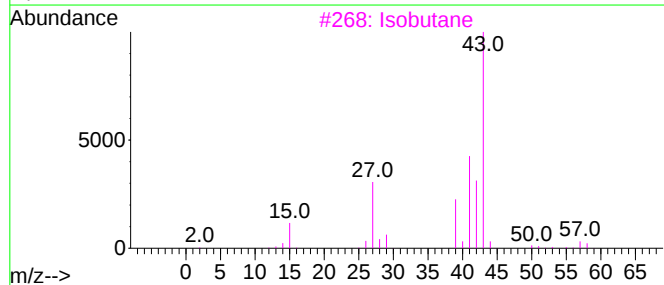
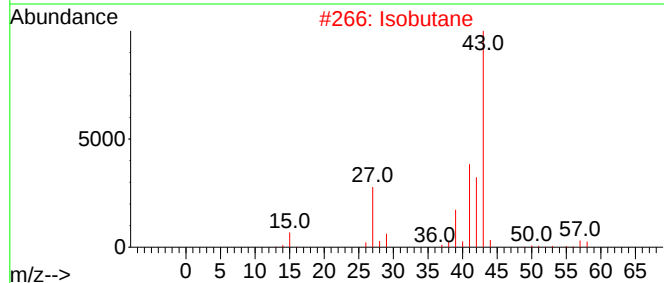
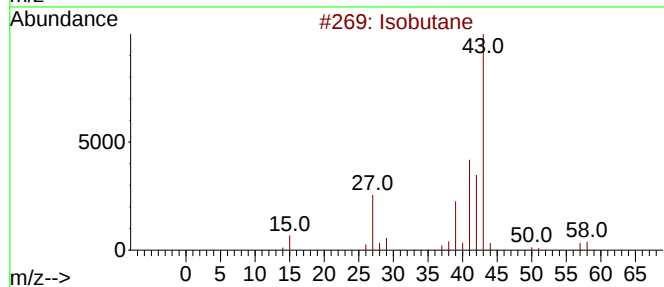
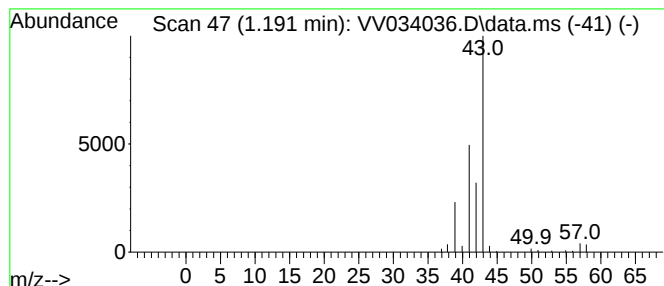
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR012524WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.191	1.12 ug/L	56156	1,4-Difluorobenzene	5.535

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	64
4			Isobutane	58	C4H10	000075-28-5	64
5			Isobutane	58	C4H10	000075-28-5	59



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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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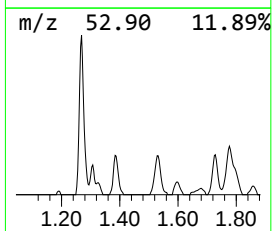
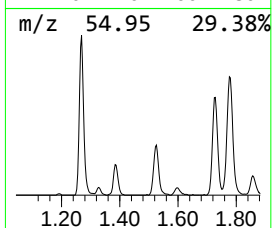
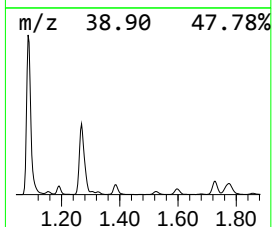
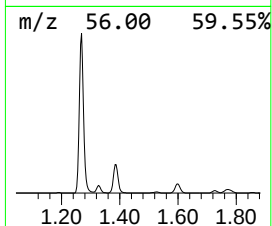
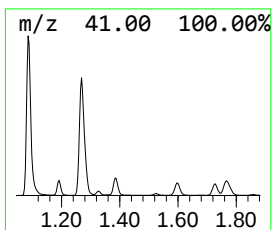
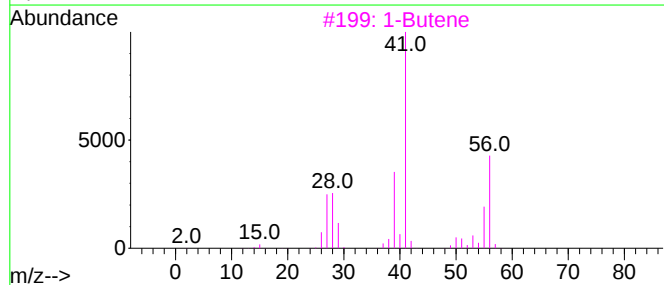
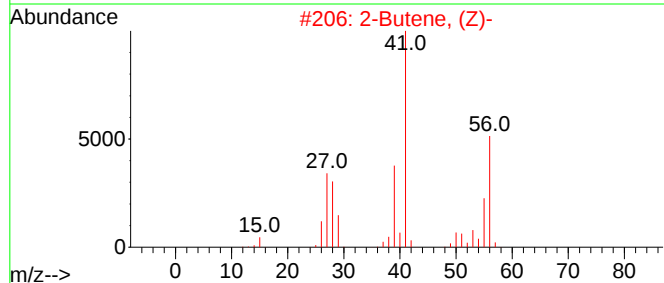
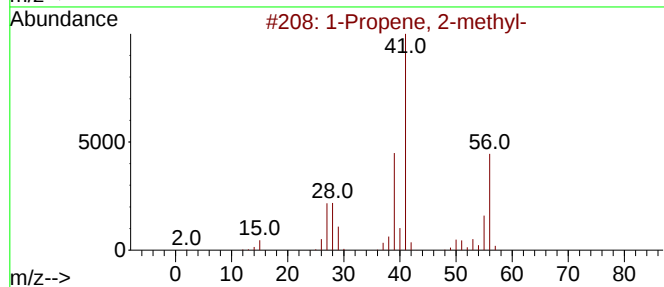
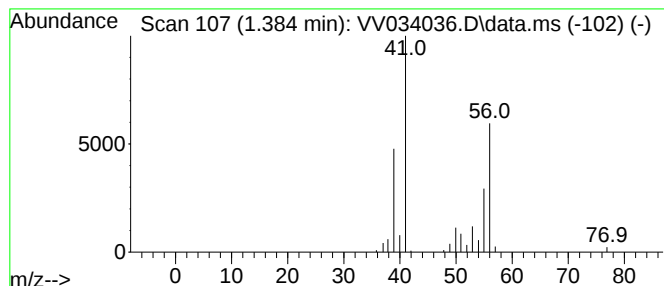
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR012524WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.384	1.18 ug/L	59132	1,4-Difluorobenzene	5.535

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 2-methyl-	56	C4H8	000115-11-7	86	
2	2-Butene, (Z)-	56	C4H8	000590-18-1	83	
3	1-Butene	56	C4H8	000106-98-9	80	
4	2-Butene, (E)-	56	C4H8	000624-64-6	80	
5	2-Butene, (Z)-	56	C4H8	000590-18-1	80	



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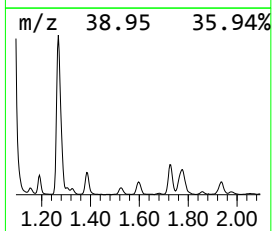
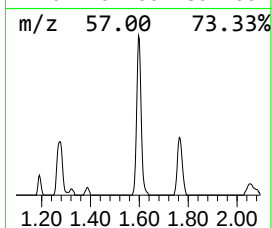
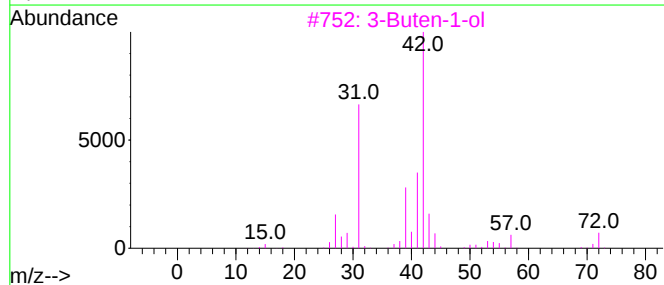
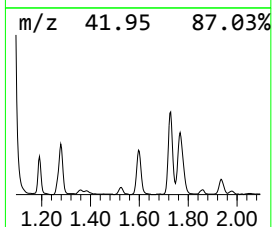
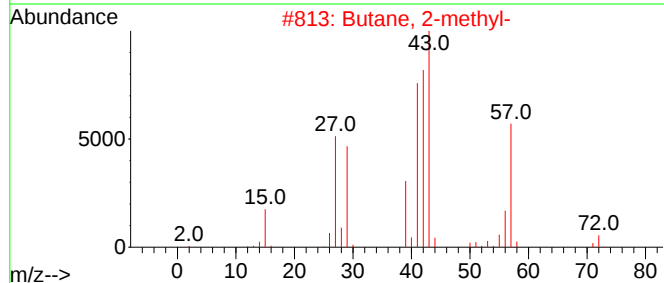
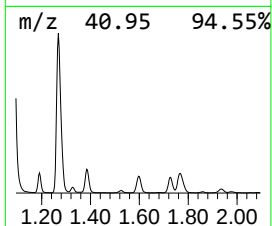
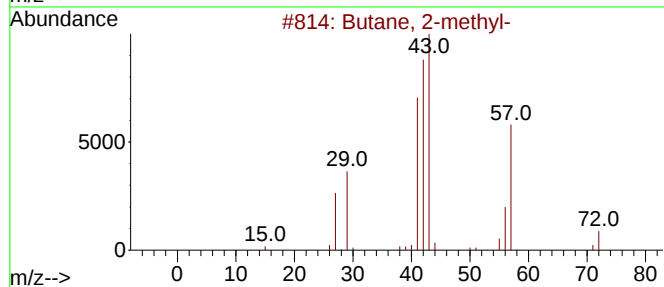
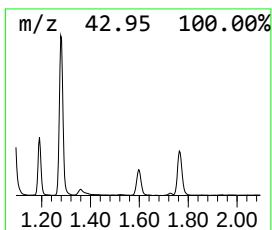
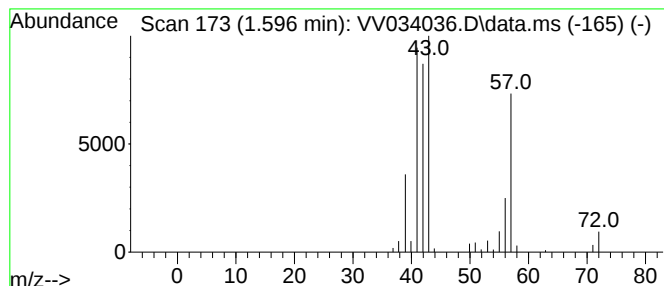
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.596	1.40 ug/L	70378	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	86
2			Butane, 2-methyl-	72	C5H12	000078-78-4	86
3			3-Buten-1-ol	72	C4H8O	000627-27-0	43
4			Oxirane, trimethyl-	86	C5H10O	005076-19-7	33
5			Pentane	72	C5H12	000109-66-0	28



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ClientSampleId :
E0YB9

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Quant Title : TRACE VOA SFAM1.0

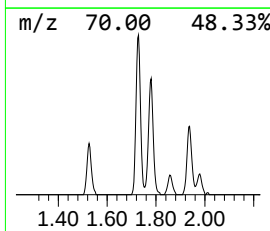
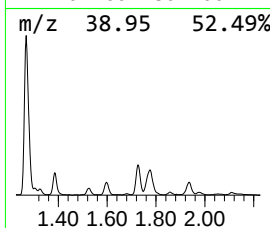
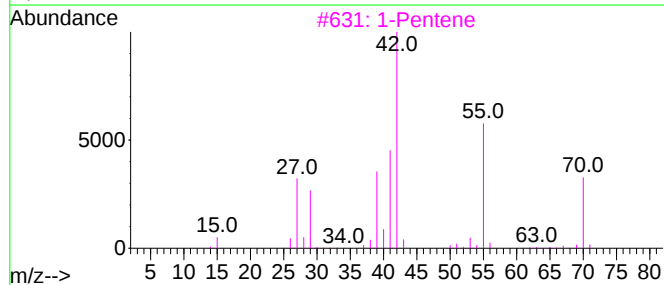
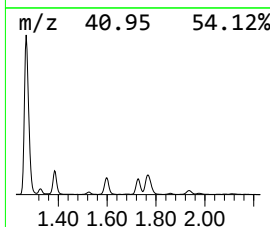
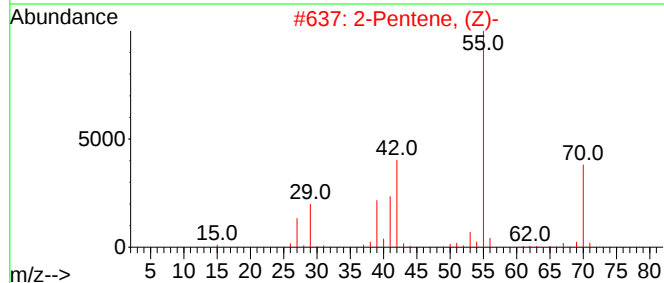
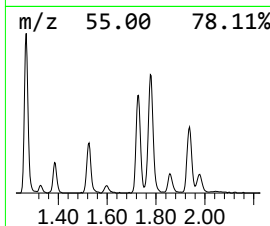
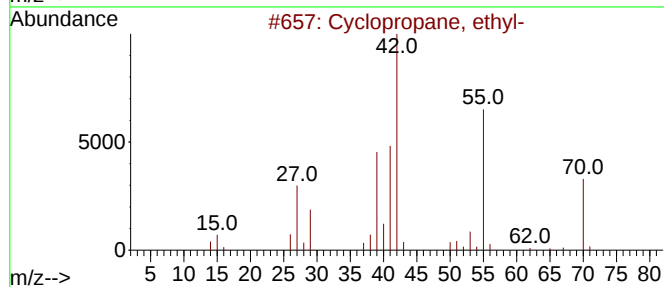
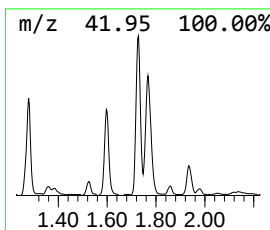
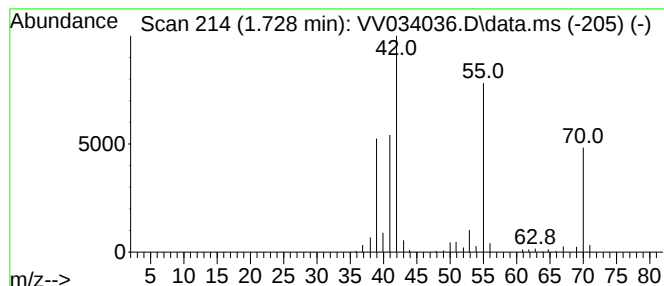
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TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic1.728 Concentration Rank 4

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1.728	1.99 ug/L	99794	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	87
3			1-Pentene	70	C5H10	000109-67-1	76
4			Cyclopropane, ethyl-	70	C5H10	001191-96-4	76
5			Cyclobutane, methyl-	70	C5H10	000598-61-8	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
E0YB9

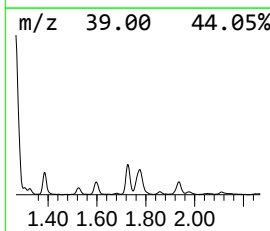
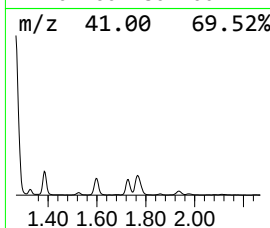
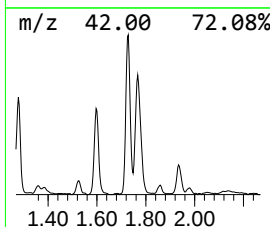
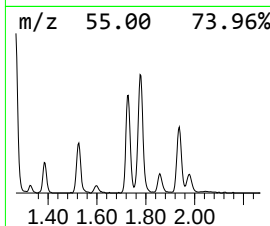
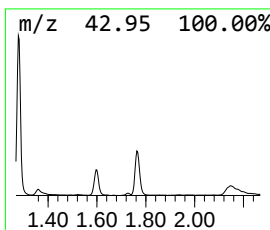
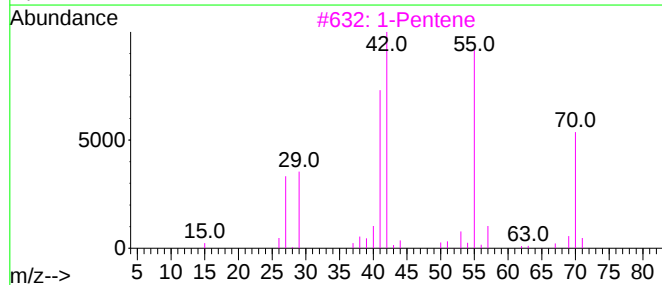
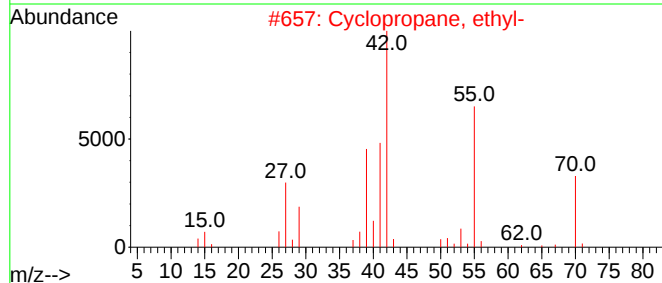
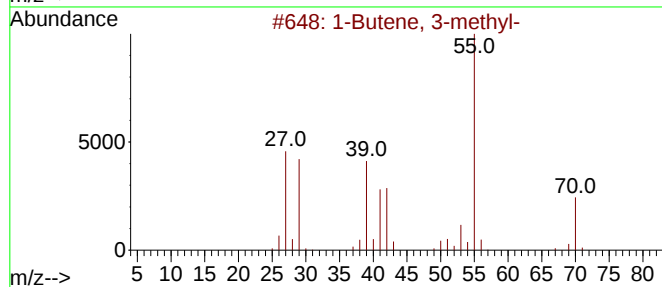
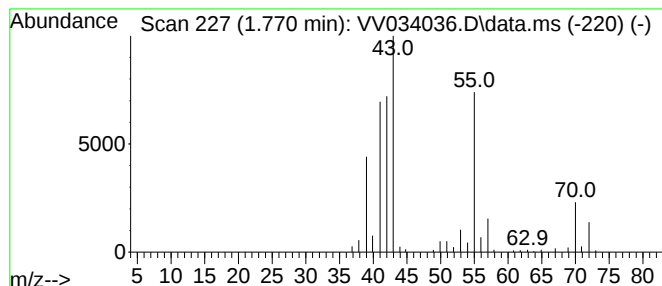
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.770	3.13 ug/L	156898	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butene, 3-methyl-	70	C5H10	000563-45-1	52
2			Cyclopropane, ethyl-	70	C5H10	001191-96-4	52
3			1-Pentene	70	C5H10	000109-67-1	50
4			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	47
5			2-Pentene, (E)-	70	C5H10	000646-04-8	41



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
E0YB9

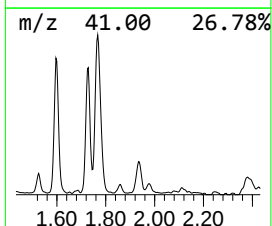
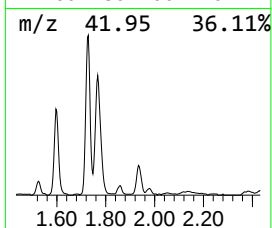
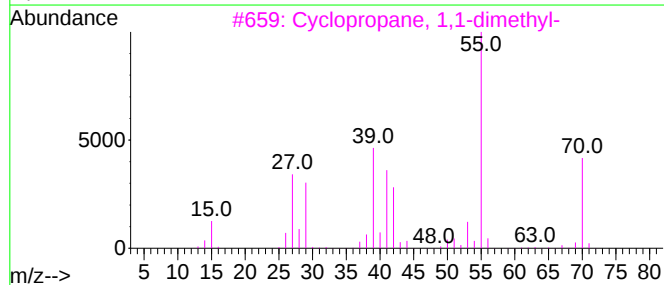
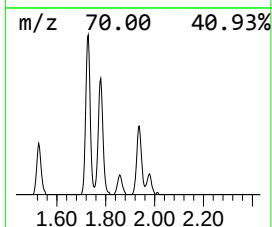
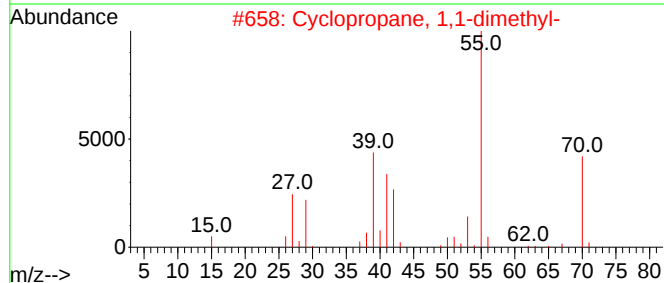
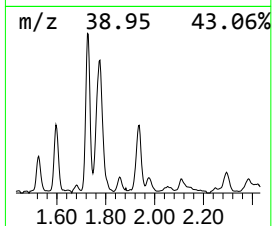
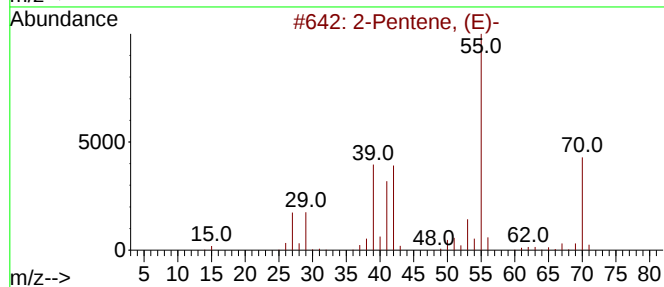
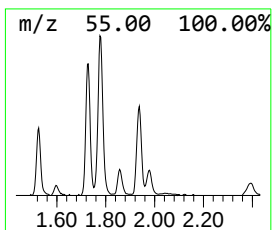
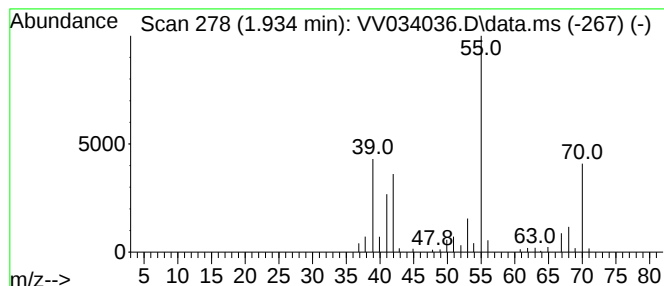
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Quant Title : TRACE VOA SFAM1.0

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.
1.934	0.97 ug/L	48440	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (E)-			70	C5H10	000646-04-8	87
2	Cyclopropane, 1,1-dimethyl-			70	C5H10	001630-94-0	87
3	Cyclopropane, 1,1-dimethyl-			70	C5H10	001630-94-0	83
4	Cyclopropane, 1,1-dimethyl-			70	C5H10	001630-94-0	72
5	Cyclopropane, 1,2-dimethyl-, trans-			70	C5H10	002402-06-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

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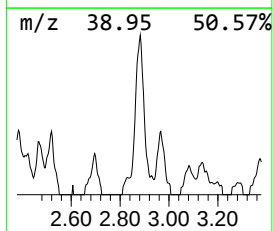
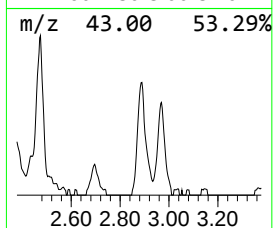
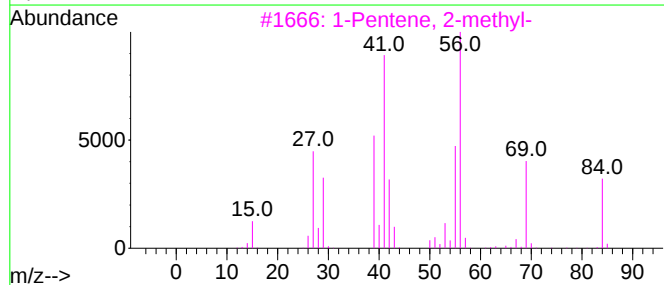
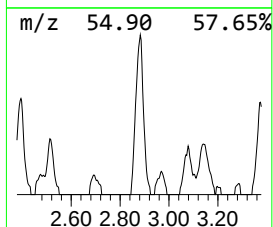
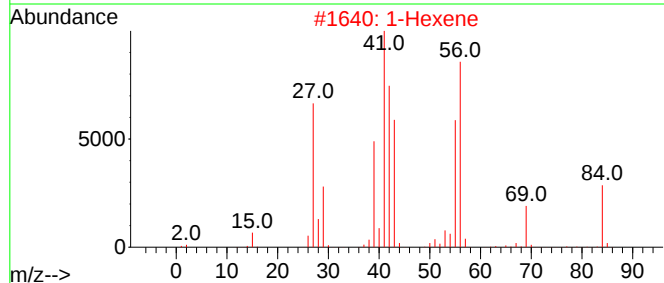
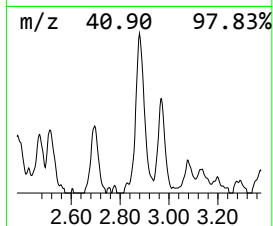
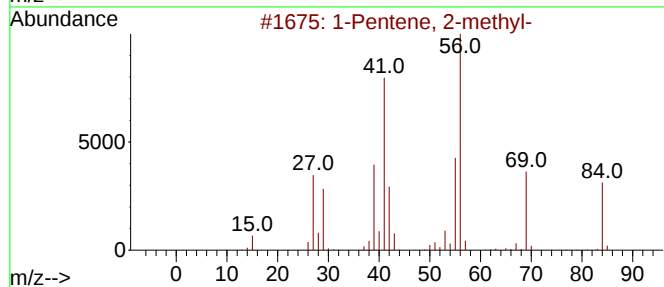
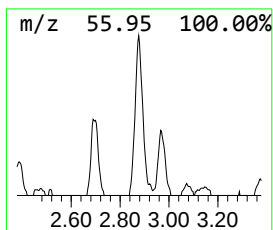
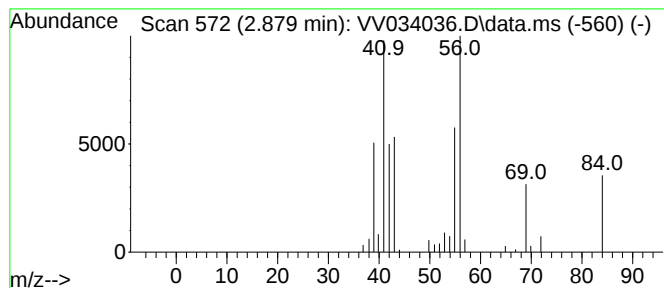
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.879	0.99 ug/L	49836	1,4-Difluorobenzene	5.535

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	83	
2	1-Hexene	84	C6H12	000592-41-6	81	
3	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72	
4	1-Hexene	84	C6H12	000592-41-6	64	
5	Formic acid, hexyl ester	130	C7H14O2	000629-33-4	50	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

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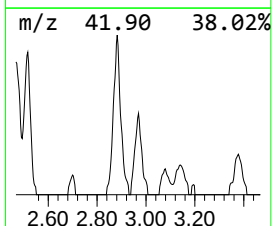
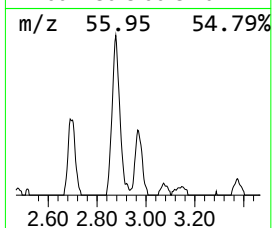
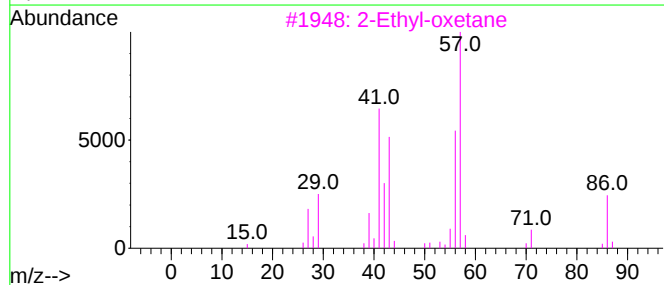
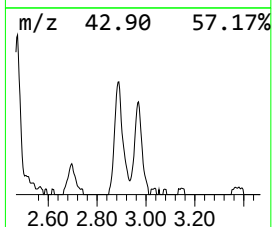
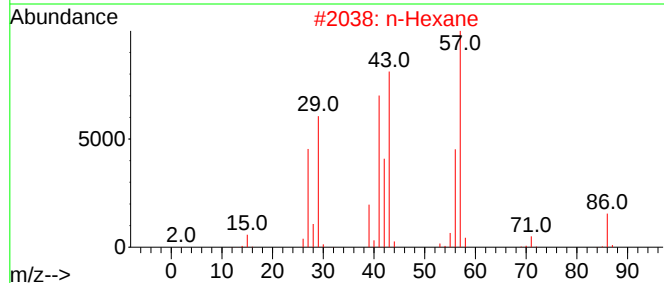
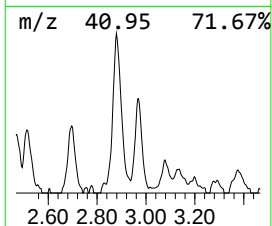
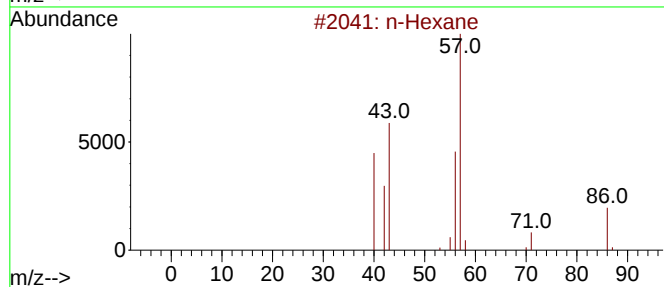
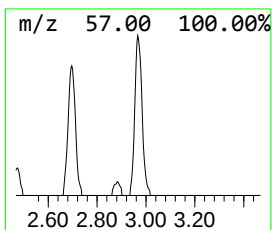
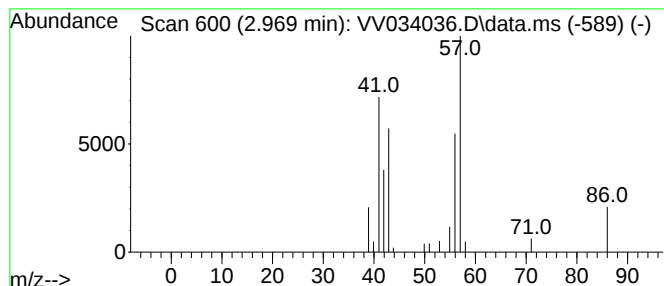
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.969	0.51 ug/L	25763	1,4-Difluorobenzene	5.535

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexane	86	C6H14	000110-54-3	80
2		n-Hexane	86	C6H14	000110-54-3	72
3		2-Ethyl-oxetane	86	C5H10O	1010386-40-2	72
4		n-Hexane	86	C6H14	000110-54-3	64
5		n-Hexane	86	C6H14	000110-54-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
E0YB9

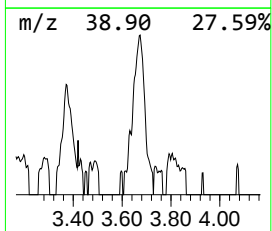
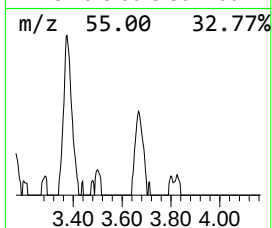
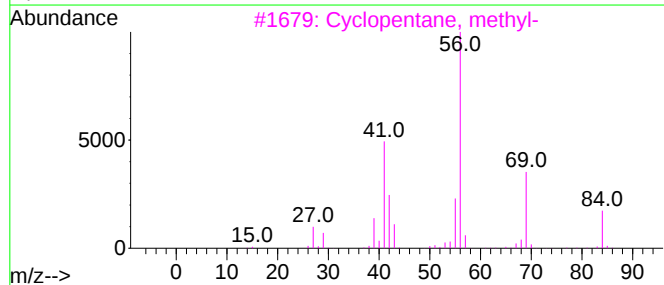
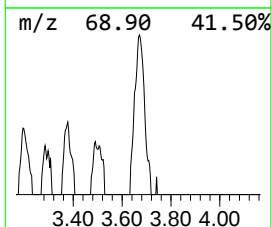
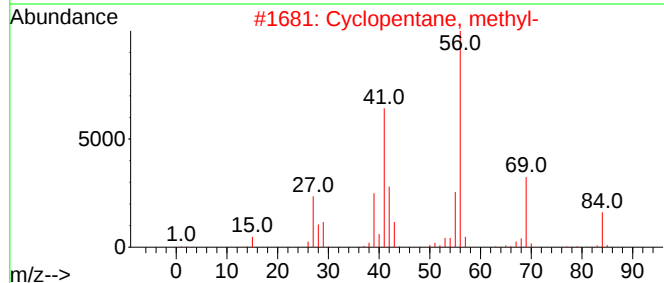
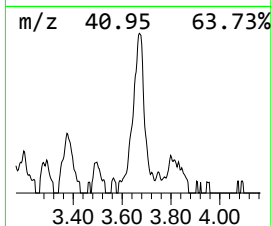
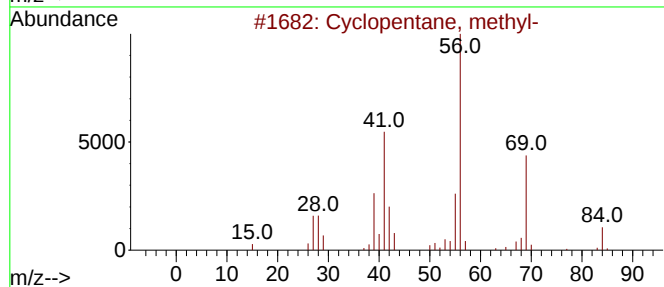
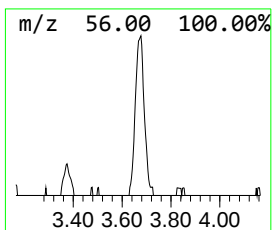
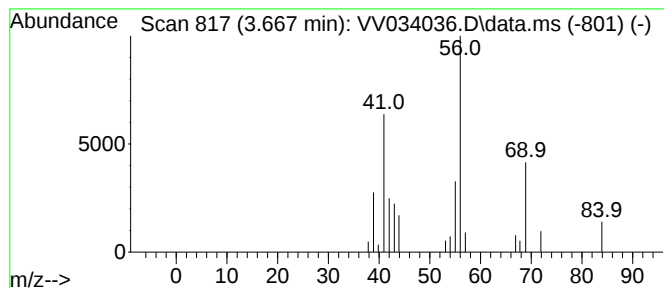
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 (DEL) Alkane: Cyclic3.667 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.667	0.51 ug/L	25394	1,4-Difluorobenzene	5.535

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	83
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	83
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	80
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

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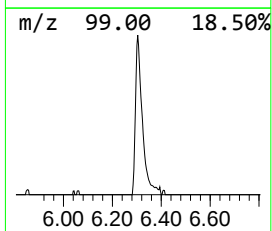
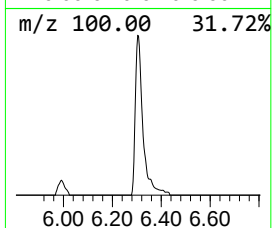
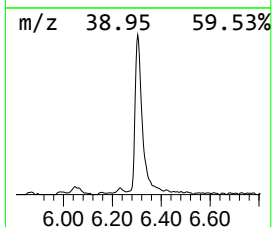
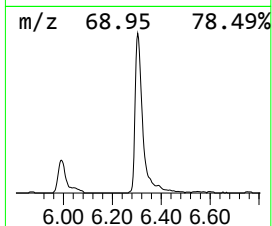
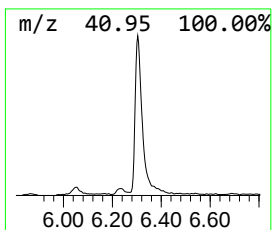
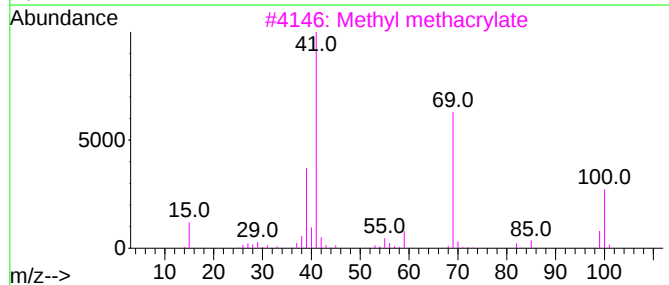
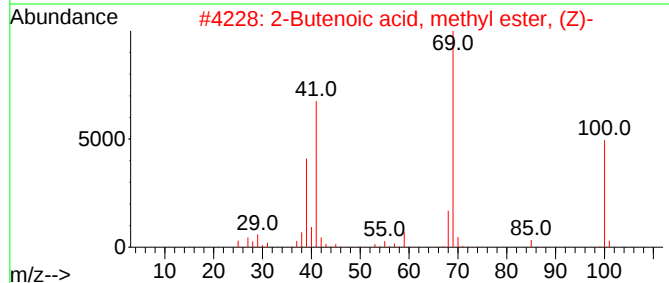
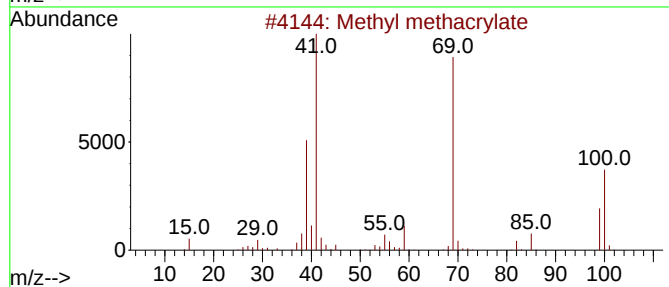
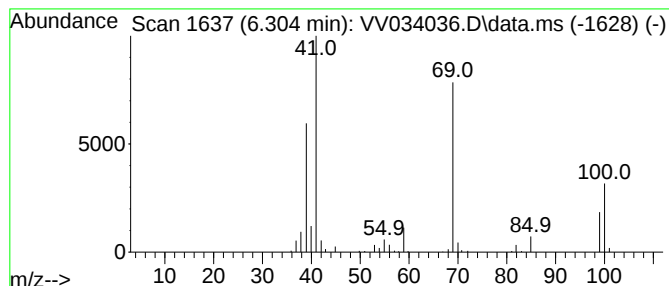
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Methyl methacrylate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.304	3.93 ug/L	197272	1,4-Difluorobenzene	5.535

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl methacrylate	100	C5H8O2	000080-62-6	94
2			2-Butenoic acid, methyl ester, (Z)-	100	C5H8O2	004358-59-2	86
3			Methyl methacrylate	100	C5H8O2	000080-62-6	76
4			Methyl methacrylate	100	C5H8O2	000080-62-6	76
5			2-Propenoic acid, 2-methyl-, 2-h...	144	C7H12O3	000923-26-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
E0YB9

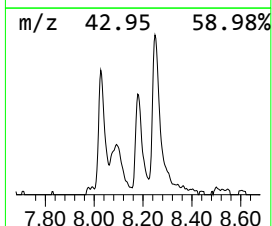
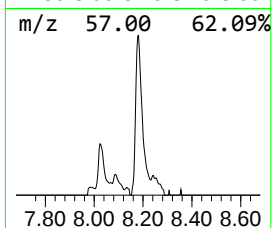
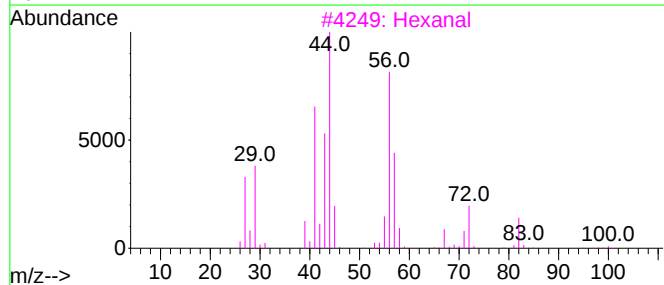
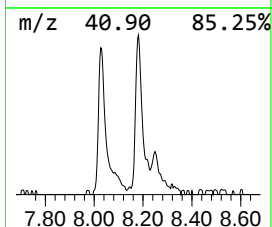
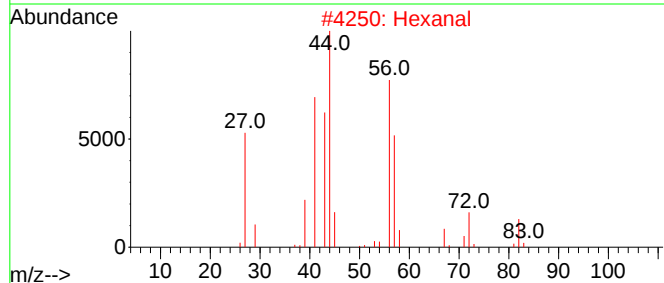
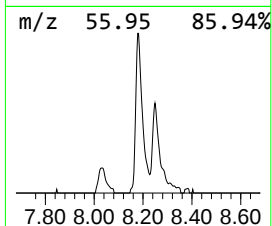
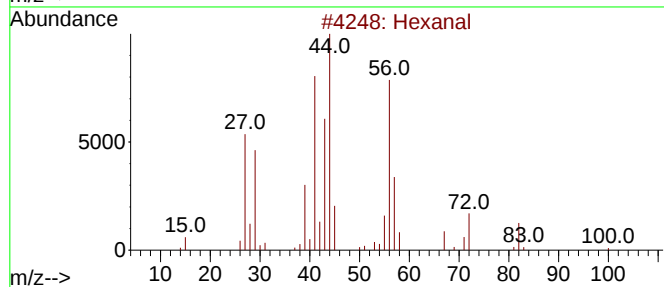
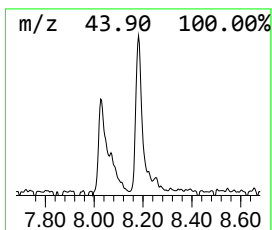
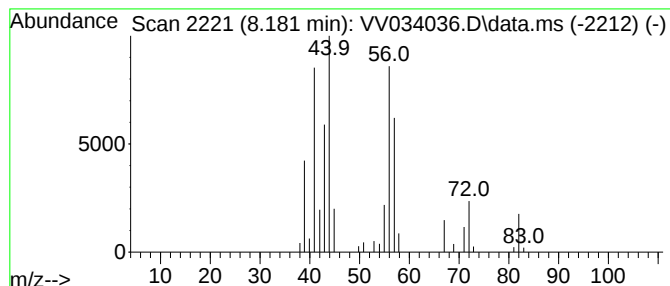
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Hexanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.181	0.99 ug/L	59542	Chlorobenzene-d5	8.786

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal			100	C6H12O	000066-25-1	72
2	Hexanal			100	C6H12O	000066-25-1	64
3	Hexanal			100	C6H12O	000066-25-1	64
4	Hexanal			100	C6H12O	000066-25-1	64
5	Hexanal			100	C6H12O	000066-25-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020724\
Data File : VV034036.D
Acq On : 07 Feb 2024 17:19
Operator : SY/MD
Sample : P1383-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
E0YB9

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR012524WMA.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...	1.085	11.0	ug/L	552242	1	5.535	250876	5.0
(DEL) Alkane: S...	1.191	1.1	ug/L	56156	1	5.535	250876	5.0
(DEL) Alkane: S...	1.384	1.2	ug/L	59132	1	5.535	250876	5.0
(DEL) Alkane: S...	1.596	1.4	ug/L	70378	1	5.535	250876	5.0
(DEL) Alkane: C...	1.728	2.0	ug/L	99794	1	5.535	250876	5.0
(DEL) Alkane: S...	1.770	3.1	ug/L	156898	1	5.535	250876	5.0
(DEL) Alkane: S...	1.934	1.0	ug/L	48440	1	5.535	250876	5.0
(DEL) Alkane: S...	2.879	1.0	ug/L	49836	1	5.535	250876	5.0
(DEL) Alkane: S...	2.969	0.5	ug/L	25763	1	5.535	250876	5.0
(DEL) Alkane: C...	3.667	0.5	ug/L	25394	1	5.535	250876	5.0
Methyl methacry...	6.304	3.9	ug/L	197272	1	5.535	250876	5.0
Hexanal	8.181	1.0	ug/L	59542	2	8.786	301242	5.0