

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102524\
 Data File : VU061220.D
 Acq On : 25 Oct 2024 12:18
 Operator : MD/SY
 Sample : P4533-06
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GCNW7

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_U\Method\SFAMUTR102324WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU061220.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.489	56	62	68	rBV	13283	13314	2.19%	0.162%
2	1.596	82	95	106	rBV2	125520	146187	24.01%	1.784%
3	1.911	179	193	209	rBV2	64242	105592	17.35%	1.288%
4	1.991	209	218	233	rVB	60657	84593	13.90%	1.032%
5	2.560	384	395	418	rBV	131905	237129	38.95%	2.893%
6	3.033	526	542	559	rBV3	26673	59642	9.80%	0.728%
7	3.129	559	572	600	rVB2	31692	70307	11.55%	0.858%
8	3.348	628	640	656	rVB2	5069	11584	1.90%	0.141%
9	3.695	737	748	761	rBV5	3060	6349	1.04%	0.077%
10	3.862	788	800	815	rBV2	5376	12906	2.12%	0.157%
11	3.978	822	836	848	rBV6	3147	7252	1.19%	0.088%
12	4.136	874	885	909	rBV2	7708	20858	3.43%	0.255%
13	4.370	945	958	959	rBV4	4079	6113	1.00%	0.075%
14	4.425	969	975	989	rVB2	6467	12000	1.97%	0.146%
15	4.521	992	1005	1018	rVB5	6342	14354	2.36%	0.175%
16	4.637	1020	1041	1073	rBV	74999	231439	38.02%	2.824%
17	4.865	1097	1112	1126	rBV3	5006	16417	2.70%	0.200%
18	5.055	1156	1171	1183	rBV	133515	306523	50.35%	3.740%
19	5.116	1185	1190	1208	rVB4	36633	73978	12.15%	0.903%
20	5.222	1212	1223	1227	rBV6	4248	8715	1.43%	0.106%
21	5.380	1257	1272	1285	rBV	42054	95876	15.75%	1.170%
22	5.454	1287	1295	1316	rVB9	6467	20140	3.31%	0.246%
23	5.627	1336	1349	1360	rBV3	25754	55461	9.11%	0.677%
24	5.718	1360	1377	1394	rBV2	235957	601772	98.85%	7.343%
25	5.885	1416	1429	1440	rBV2	138907	327517	53.80%	3.996%
26	5.959	1442	1452	1473	rVB	181451	423168	69.51%	5.163%
27	6.055	1474	1482	1490	rBV2	17396	32065	5.27%	0.391%
28	6.242	1528	1540	1565	rBV	215347	459625	75.50%	5.608%
29	6.505	1611	1622	1631	rBV	77016	154321	25.35%	1.883%
30	6.589	1643	1648	1664	rVB3	25444	45908	7.54%	0.560%
31	6.682	1664	1677	1688	rVV	142664	279679	45.94%	3.413%
32	6.756	1689	1700	1726	rVV4	71656	217340	35.70%	2.652%
33	7.004	1771	1777	1790	rVB4	3702	7958	1.31%	0.097%
34	7.152	1808	1823	1842	rBV6	20869	54625	8.97%	0.667%
35	7.251	1843	1854	1866	rVB6	5598	14956	2.46%	0.182%
36	7.389	1876	1897	1904	rBV7	6383	18129	2.98%	0.221%

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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_U\Method\SFAMUTR102324WMA.M
 Title : TRACE VOA SFAM1.0

37	7.431	1906	1910	1928	rVB4	6615	11706	1.92%	0.143%
38	7.560	1931	1950	1972	rBV	82140	179986	29.57%	2.196%
39	7.721	1988	2000	2002	rBV5	4866	7680	1.26%	0.094%
40	7.753	2002	2010	2023	rVB3	12092	22400	3.68%	0.273%
41	7.891	2028	2053	2080	rBV	308076	608758	100.00%	7.428%
42	8.020	2084	2093	2103	rBV6	9057	17552	2.88%	0.214%
43	8.106	2117	2120	2129	rVB5	6362	8316	1.37%	0.101%
44	8.174	2129	2141	2158	rVB2	45724	82987	13.63%	1.013%
45	8.258	2158	2167	2178	rVB4	6282	10599	1.74%	0.129%
46	8.348	2180	2195	2202	rBV4	5648	13410	2.20%	0.164%
47	8.402	2203	2212	2223	rBV2	32211	52641	8.65%	0.642%
48	8.627	2272	2282	2297	rVV	294309	565087	92.83%	6.895%
49	8.698	2298	2304	2314	rVB2	42198	71409	11.73%	0.871%
50	8.775	2323	2328	2347	rVB3	16442	31226	5.13%	0.381%
51	8.875	2350	2359	2368	rVV3	12002	21244	3.49%	0.259%
52	8.933	2368	2377	2390	rVB3	11507	23029	3.78%	0.281%
53	9.071	2409	2420	2438	rVB4	21595	40587	6.67%	0.495%
54	9.177	2438	2453	2463	rBV7	4259	9616	1.58%	0.117%
55	9.254	2464	2477	2487	rVV3	10216	21272	3.49%	0.260%
56	9.319	2487	2497	2507	rVV10	8958	18263	3.00%	0.223%
57	9.409	2507	2525	2551	rVV	288763	560985	92.15%	6.845%
58	9.521	2554	2560	2569	rVV5	14121	22235	3.65%	0.271%
59	9.589	2569	2581	2595	rVV10	6576	16212	2.66%	0.198%
60	9.682	2595	2610	2625	rVB9	28979	71816	11.80%	0.876%
61	9.804	2640	2648	2657	rBV9	3786	7834	1.29%	0.096%
62	10.013	2699	2713	2719	rBV8	4513	10217	1.68%	0.125%
63	10.058	2719	2727	2737	rVB5	8707	14906	2.45%	0.182%
64	10.521	2863	2871	2877	rBV4	6866	10217	1.68%	0.125%
65	10.682	2909	2921	2931	rVV6	10424	19145	3.14%	0.234%
66	10.746	2931	2941	2964	rVB	102269	176649	29.02%	2.155%
67	11.409	3137	3147	3159	rVB2	26386	42698	7.01%	0.521%
68	11.534	3174	3186	3194	rBV8	5317	10408	1.71%	0.127%
69	11.598	3194	3206	3211	rBV6	7156	12386	2.03%	0.151%
70	11.637	3212	3218	3231	rVB	13492	20669	3.40%	0.252%
71	11.804	3258	3270	3292	rBV	288302	479445	78.76%	5.850%
72	12.090	3345	3359	3370	rBV2	13421	23782	3.91%	0.290%
73	12.187	3374	3389	3409	rBV	275227	473321	77.75%	5.775%
74	12.293	3413	3422	3432	rVB10	4692	9492	1.56%	0.116%
75	12.856	3583	3597	3608	rBV2	16161	26696	4.39%	0.326%
76	13.110	3666	3676	3685	rBV2	14803	21735	3.57%	0.265%

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Instrument :
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ClientSampleId :
GCNW7

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_U\Method\SFAMUTR102324WMA.M
Title : TRACE VOA SFAM1.0

77	13.319	3731	3741	3755	rVB	25575	39781	6.53%	0.485%
78	13.402	3755	3767	3775	rBV8	4275	8619	1.42%	0.105%
79	13.524	3794	3805	3812	rBV7	4578	7346	1.21%	0.090%
80	13.785	3877	3886	3893	rBV3	7295	11281	1.85%	0.138%
81	13.881	3904	3916	3930	rVB4	9712	17673	2.90%	0.216%
82	14.277	4028	4039	4047	rBV4	5670	8462	1.39%	0.103%

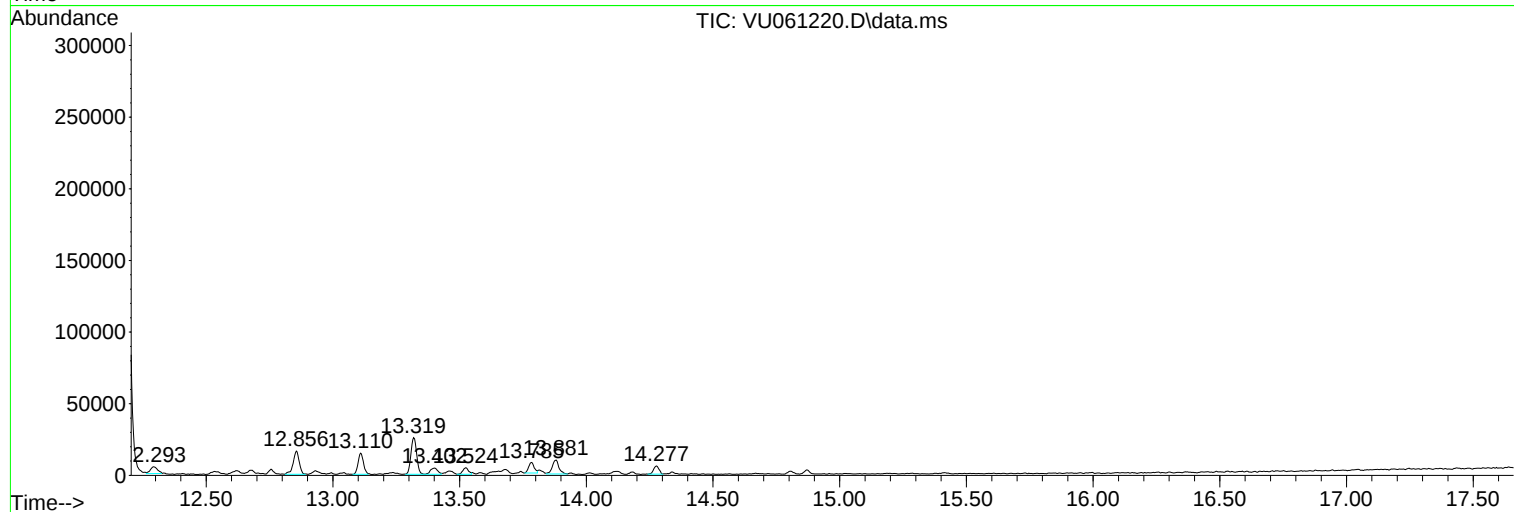
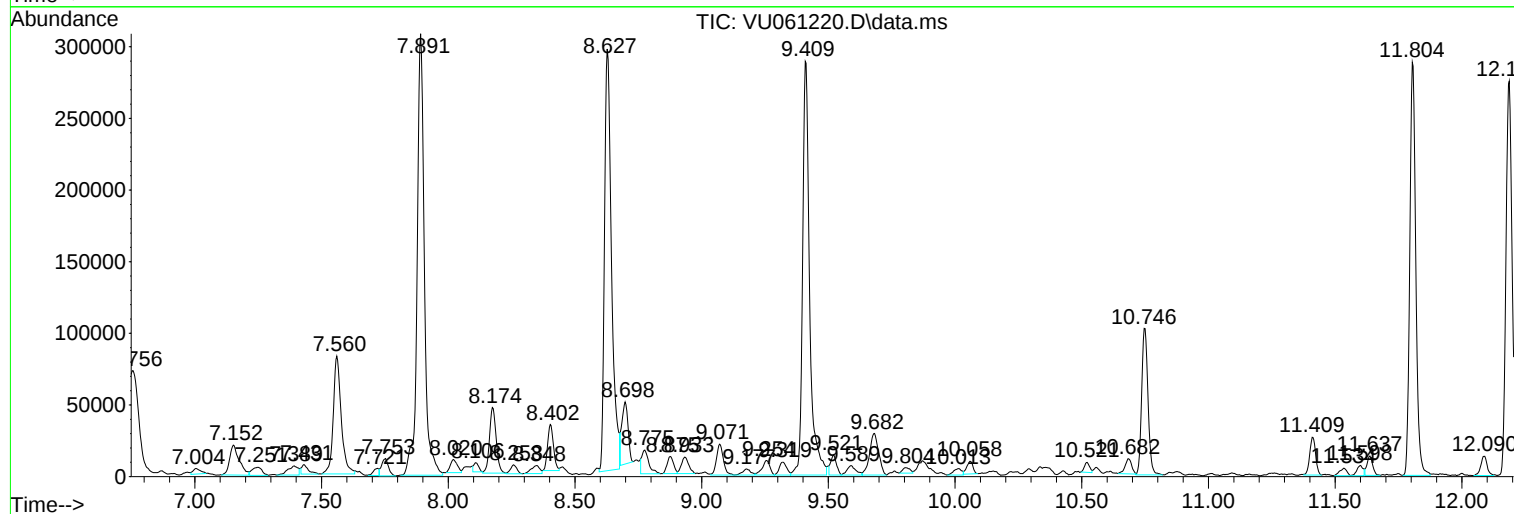
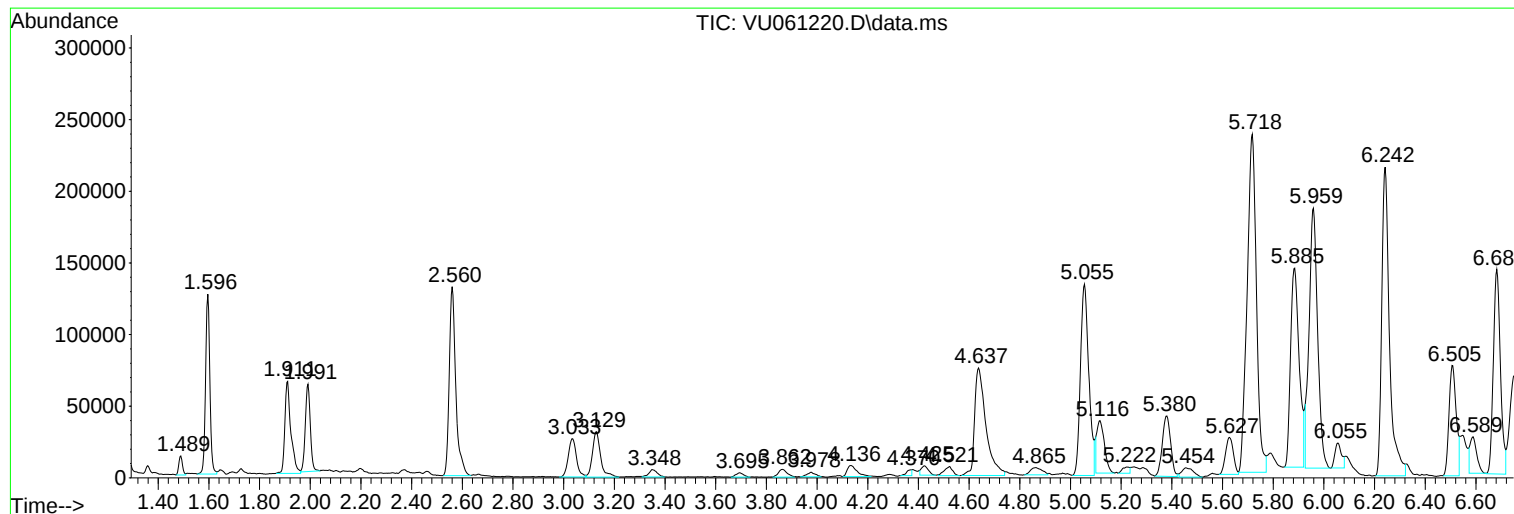
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Instrument :
MSVOA_U
ClientSampleId :
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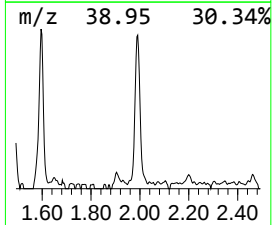
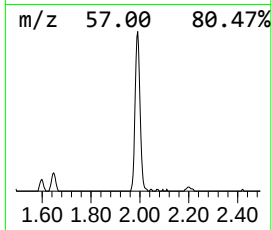
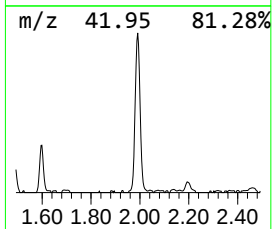
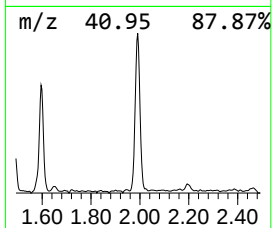
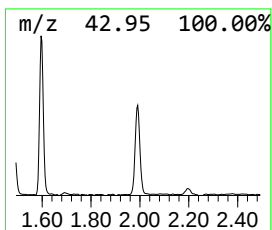
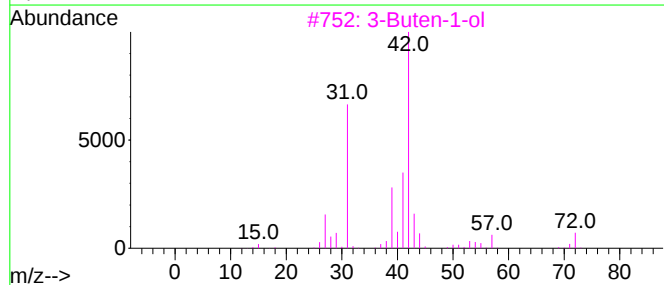
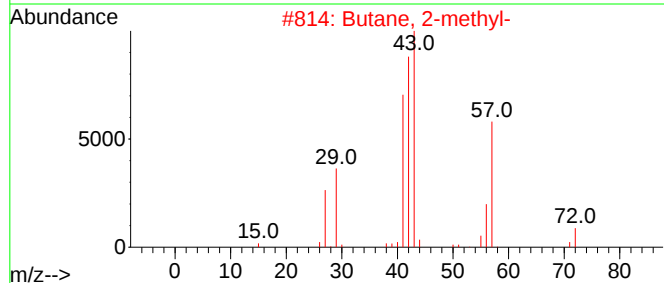
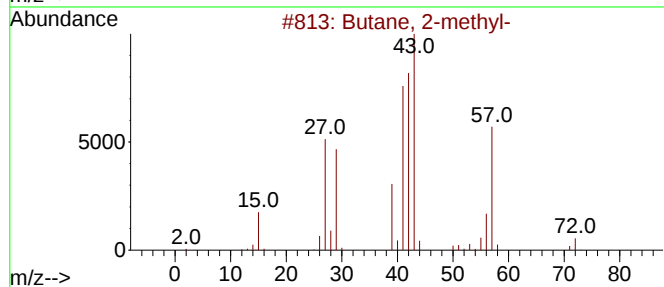
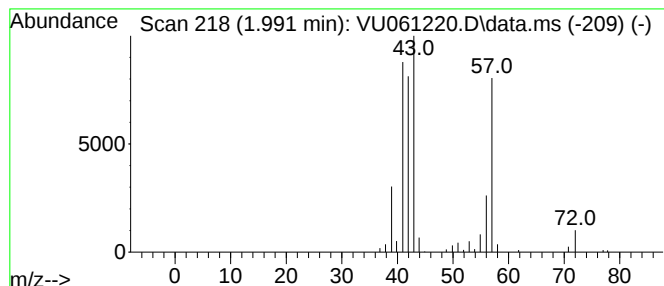
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.991	0.92 ug/L	84593	1,4-Difluorobenzene	6.242

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	80
3		3-Buten-1-ol	72	C4H8O	000627-27-0	43
4		3-Buten-1-ol	72	C4H8O	000627-27-0	37
5		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	36



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Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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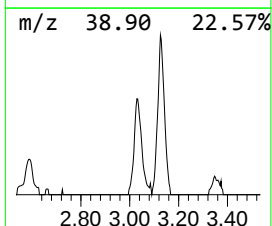
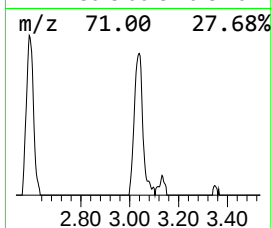
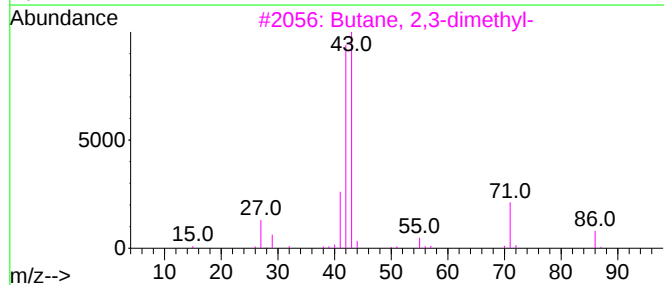
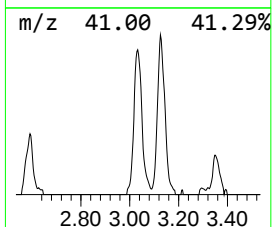
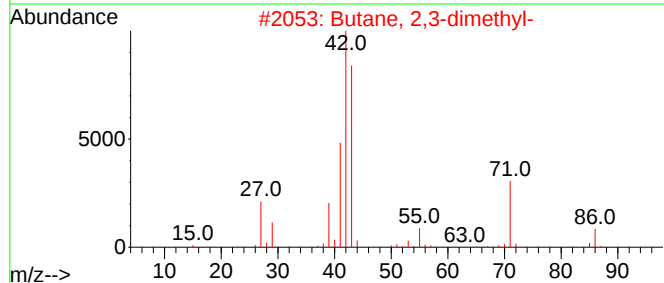
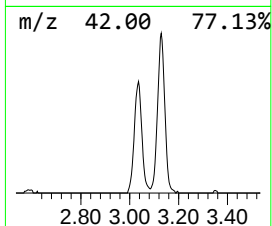
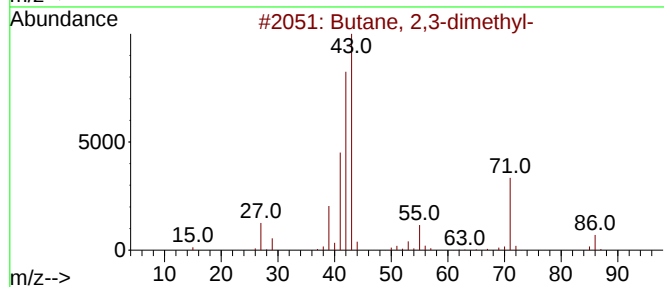
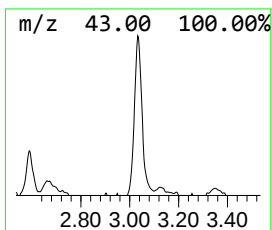
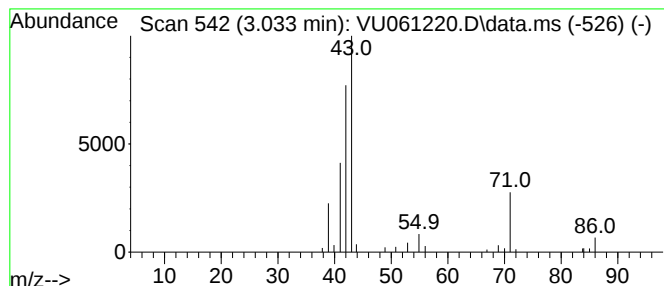
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.033	0.65 ug/L	59642	1,4-Difluorobenzene	6.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	83
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	78
4			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
5			Pentane, 2-methyl-	86	C6H14	000107-83-5	64



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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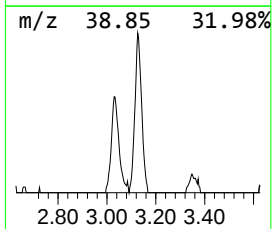
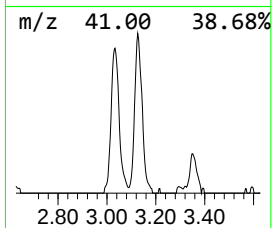
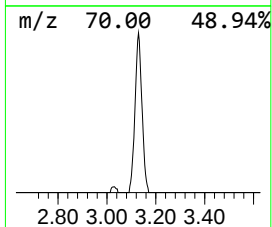
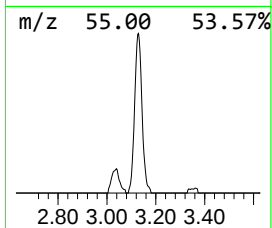
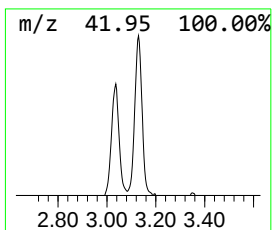
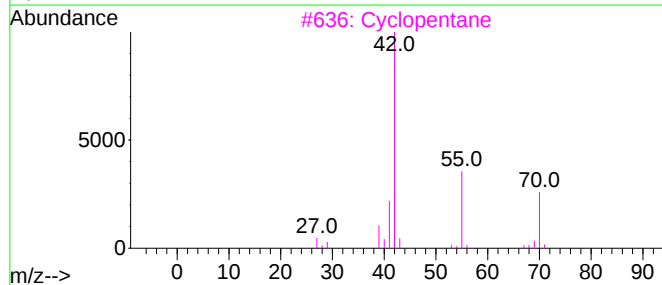
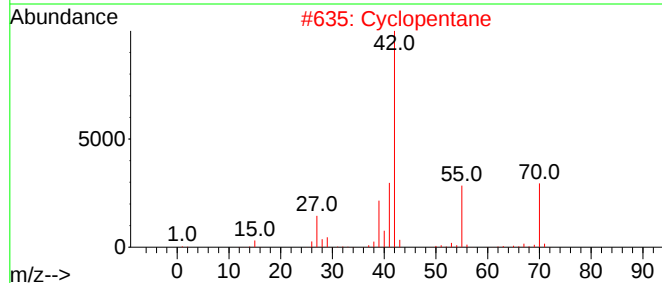
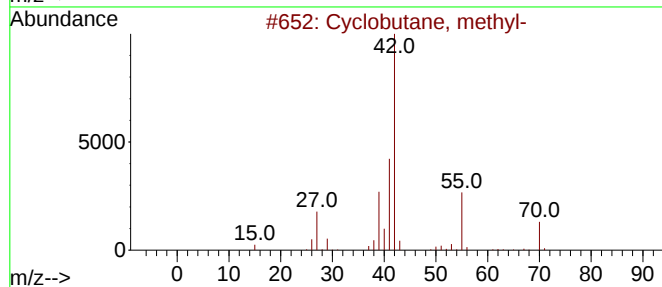
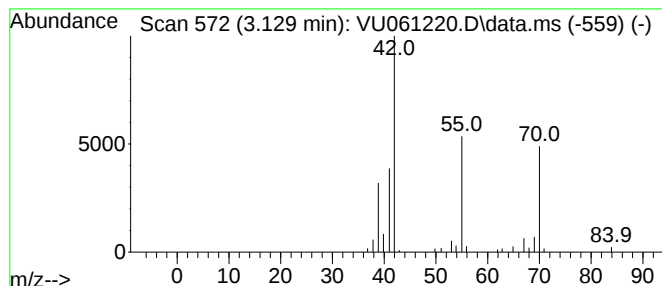
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Cyclic3.129 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.129	0.76 ug/L	70307	1,4-Difluorobenzene	6.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutane, methyl-	70	C5H10	000598-61-8	86
2			Cyclopentane	70	C5H10	000287-92-3	86
3			Cyclopentane	70	C5H10	000287-92-3	72
4			1-Pentene	70	C5H10	000109-67-1	72
5			1-Pentene	70	C5H10	000109-67-1	64



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Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

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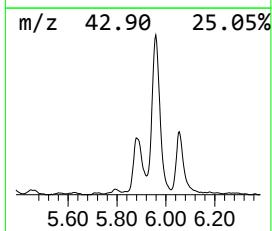
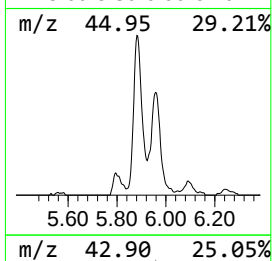
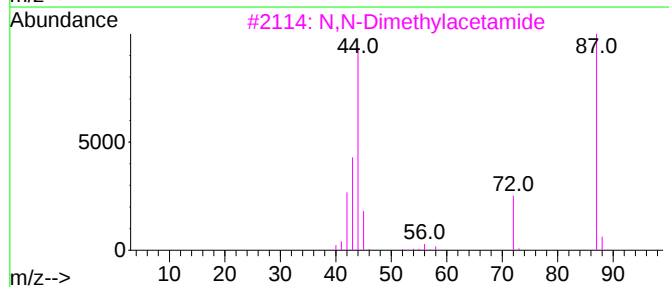
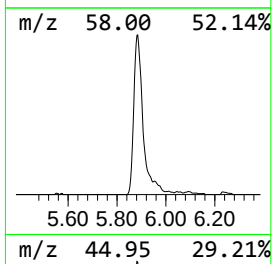
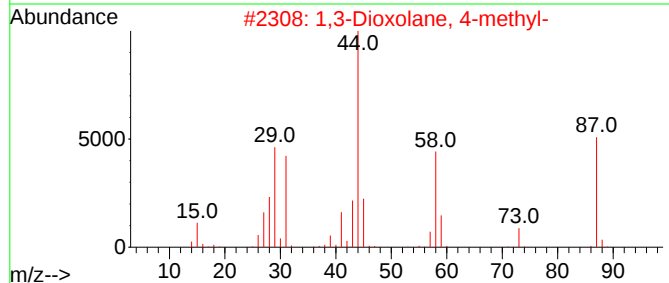
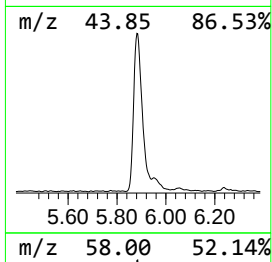
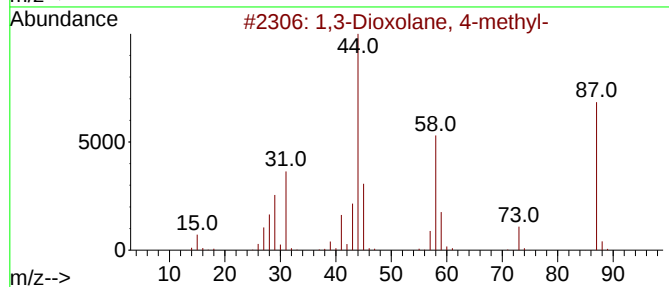
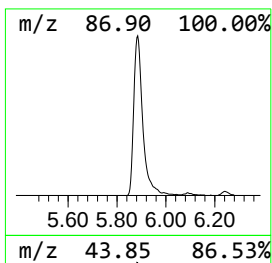
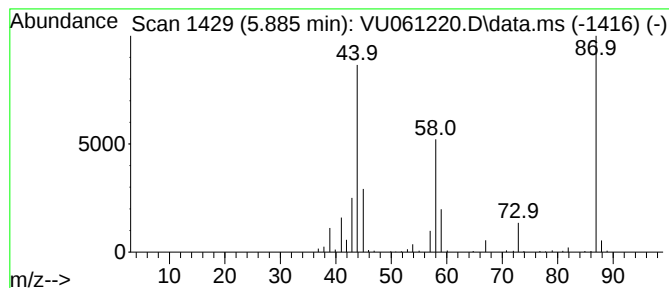
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 1,3-Dioxolane, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.885	3.56 ug/L	327517	1,4-Difluorobenzene	6.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 4-methyl-	88	C4H8O2	001072-47-5	91
2			1,3-Dioxolane, 4-methyl-	88	C4H8O2	001072-47-5	78
3			N,N-Dimethylacetamide	87	C4H9NO	000127-19-5	40
4			Diazene, butyl[1-(2,2-dimethylhy...	172	C8H20N4	061940-95-2	36
5			1,2-Ethanediamine, N,N'-dimethyl-	88	C4H12N2	000110-70-3	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102524\
Data File : VU061220.D
Acq On : 25 Oct 2024 12:18
Operator : MD/SY
Sample : P4533-06
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GCNW7

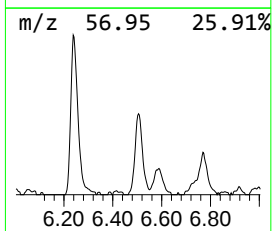
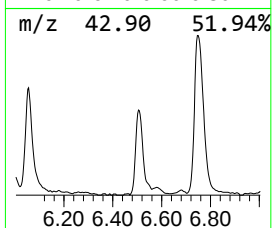
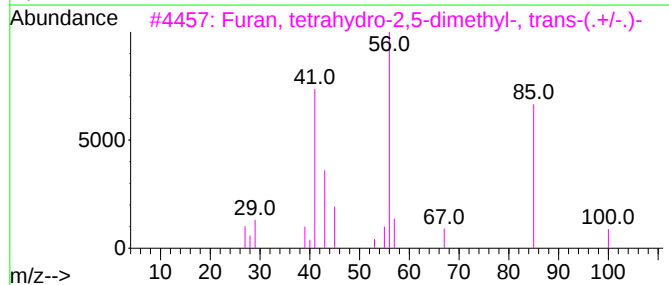
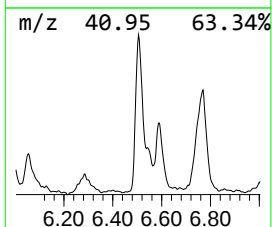
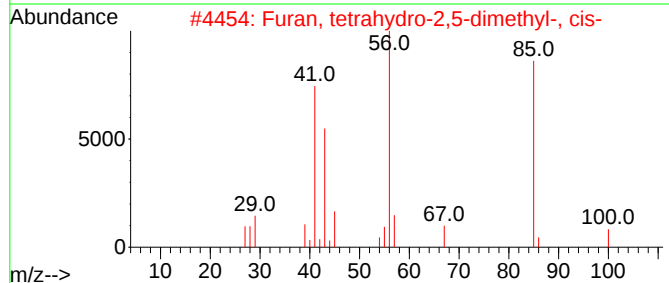
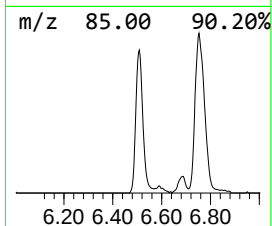
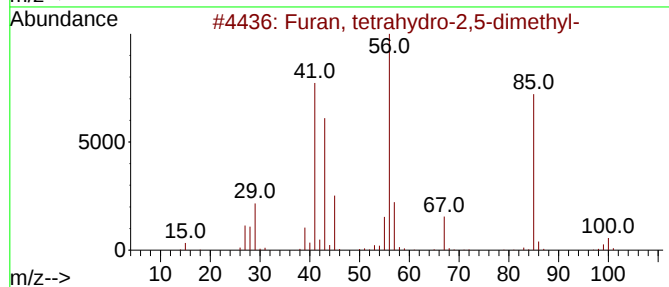
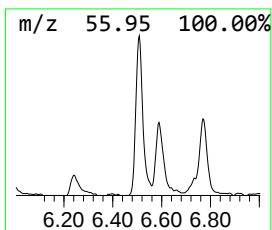
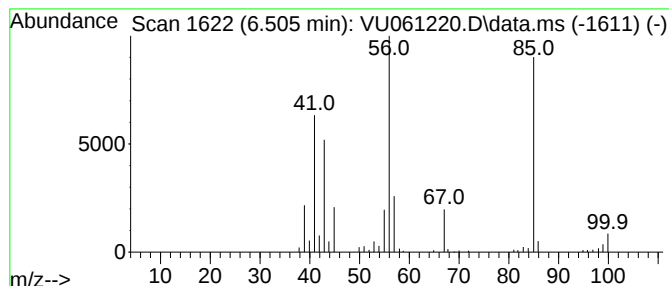
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Furan, tetrahydro-2,5-dimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.505	1.68 ug/L	154321	1,4-Difluorobenzene	6.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	91
2			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	002144-41-4	86
3			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	038484-59-2	72
4			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	64
5			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102524\
Data File : VU061220.D
Acq On : 25 Oct 2024 12:18
Operator : MD/SY
Sample : P4533-06
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GCNW7

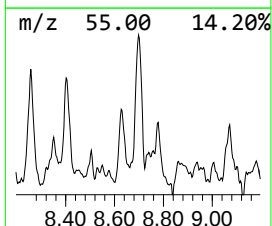
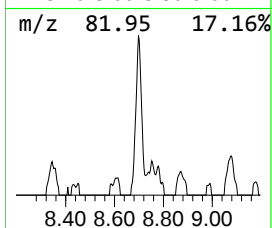
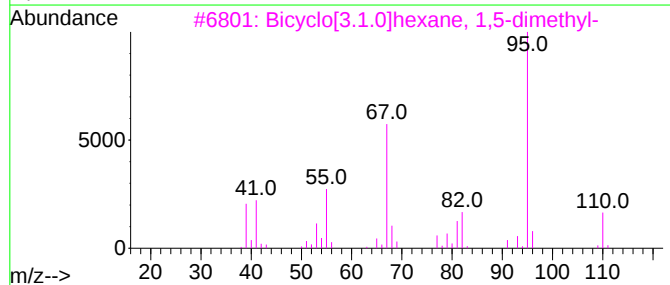
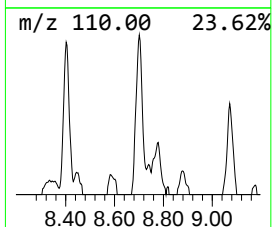
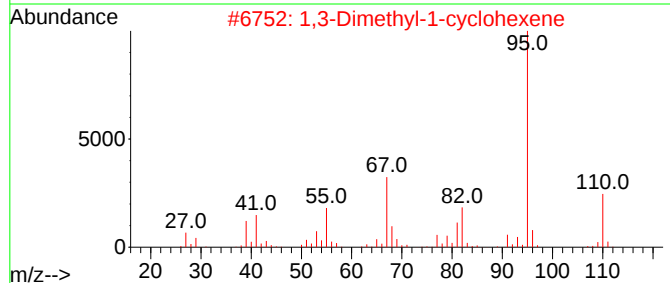
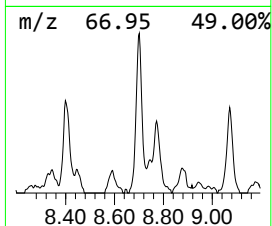
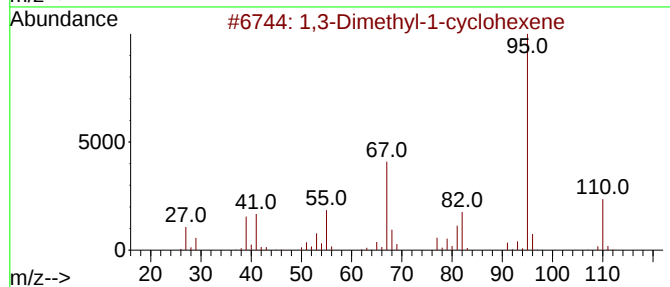
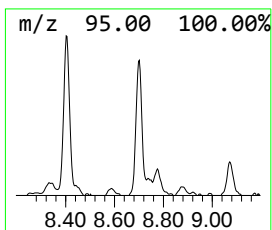
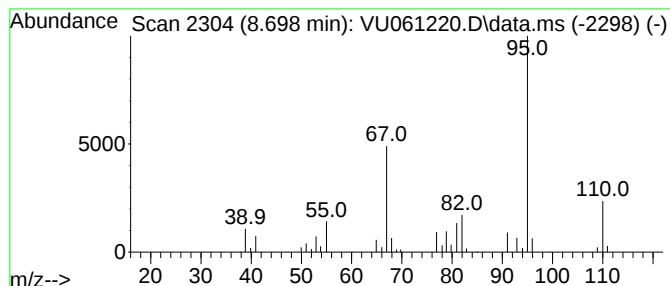
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 1,3-Dimethyl-1-cyclohexene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.698	0.64 ug/L	71409	Chlorobenzene-d5	9.412

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	91
2			1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	91
3			Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1010142-17-5	91
4			Cyclohexene, 3,5-dimethyl-	110	C8H14	000823-17-6	87
5			5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102524\
Data File : VU061220.D
Acq On : 25 Oct 2024 12:18
Operator : MD/SY
Sample : P4533-06
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GCNW7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102324WMA.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
(DEL) Alkane: S...	1.991	0.9	ug/L	84593	1	6.242	459625	5.0
(DEL) Alkane: S...	3.033	0.7	ug/L	59642	1	6.242	459625	5.0
(DEL) Alkane: C...	3.129	0.8	ug/L	70307	1	6.242	459625	5.0
1,3-Dioxolane, ...	5.885	3.6	ug/L	327517	1	6.242	459625	5.0
Furan, tetrahyd...	6.505	1.7	ug/L	154321	1	6.242	459625	5.0
1,3-Dimethyl-1-...	8.698	0.6	ug/L	71409	2	9.412	560985	5.0