

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
 Data File : VW022492.D
 Acq On : 11 Apr 2022 18:57
 Operator : SY/VA
 Sample : N2316-13
 Misc : 6.00g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNR3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

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_W\Method\SFAMWLM040722SMA.M
 Title : SFAM01.0

Signal : TIC: VW022492.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.958	4	9	20	rVB	643330	860491	43.59%	4.402%
2	2.214	46	51	54	rBV5	5326	9987	0.51%	0.051%
3	2.355	66	74	83	rBV	146067	265706	13.46%	1.359%
4	2.501	92	98	109	rBV	54443	132546	6.71%	0.678%
5	2.885	154	161	174	rVB	105074	220360	11.16%	1.127%
6	3.391	238	244	249	rBV5	3380	7239	0.37%	0.037%
7	3.848	310	319	324	rBV9	3822	9297	0.47%	0.048%
8	3.940	329	334	338	rBV7	3818	8512	0.43%	0.044%
9	4.025	338	348	359	rBV2	161697	502445	25.45%	2.570%
10	4.208	374	378	387	rVB10	5267	13709	0.69%	0.070%
11	4.391	400	408	410	rBV2	3651	7626	0.39%	0.039%
12	4.781	461	472	483	rBV4	10817	35850	1.82%	0.183%
13	5.013	486	510	527	rBV	265036	970610	49.17%	4.965%
14	5.183	530	538	547	rVB5	6778	21412	1.08%	0.110%
15	5.433	568	579	589	rBV8	6165	22896	1.16%	0.117%
16	5.561	594	600	608	rVB10	3956	11416	0.58%	0.058%
17	5.946	652	663	676	rVB5	14349	48418	2.45%	0.248%
18	6.214	698	707	717	rBV2	9929	34282	1.74%	0.175%
19	6.525	752	758	761	rBV7	2357	5061	0.26%	0.026%
20	6.622	768	774	780	rVB7	2832	6939	0.35%	0.035%
21	6.866	810	814	815	rVV4	2695	3649	0.18%	0.019%
22	6.982	824	833	841	rVV2	22841	72054	3.65%	0.369%
23	7.092	841	851	873	rVV4	90217	369728	18.73%	1.891%
24	7.646	930	942	959	rBV	283097	746622	37.82%	3.819%
25	7.951	981	992	1000	rBV4	44326	144805	7.34%	0.741%
26	8.031	1000	1005	1017	rVB6	9765	27227	1.38%	0.139%
27	8.159	1017	1026	1033	rBV3	27191	60208	3.05%	0.308%
28	8.274	1033	1045	1048	rBV	536880	1198014	60.69%	6.129%
29	8.323	1048	1053	1061	rVB3	848589	1973894	100.00%	10.098%
30	8.402	1061	1066	1072	rBV	47476	94818	4.80%	0.485%
31	8.506	1080	1083	1089	rVB8	4404	5609	0.28%	0.029%
32	8.567	1089	1093	1102	rVB8	10252	25864	1.31%	0.132%
33	8.695	1105	1114	1125	rVB2	28427	58230	2.95%	0.298%
34	8.841	1128	1138	1153	rBV	563501	1118625	56.67%	5.723%
35	8.988	1156	1162	1168	rBV2	21769	43767	2.22%	0.224%
36	9.079	1168	1177	1186	rBV3	102311	245394	12.43%	1.255%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
 Title : SFAM01.0

37	9.158	1188	1190	1196	rVB4	4961	6119	0.31%	0.031%
38	9.274	1199	1209	1215	rBV	395097	818526	41.47%	4.187%
39	9.335	1215	1219	1236	rVB2	60190	153885	7.80%	0.787%
40	9.518	1242	1249	1255	rVB3	9119	18961	0.96%	0.097%
41	9.610	1260	1264	1269	rVB4	3186	4768	0.24%	0.024%
42	9.677	1269	1275	1276	rBV5	3549	5646	0.29%	0.029%
43	9.756	1284	1288	1296	rVB9	4532	8272	0.42%	0.042%
44	9.908	1304	1313	1315	rBV7	4059	7983	0.40%	0.041%
45	9.963	1315	1322	1328	rBV	106919	202136	10.24%	1.034%
46	10.036	1328	1334	1341	rVB	210556	373056	18.90%	1.908%
47	10.207	1356	1362	1366	rBV6	6310	11548	0.59%	0.059%
48	10.244	1366	1368	1373	rVB6	3931	5376	0.27%	0.028%
49	10.323	1373	1381	1387	rBV	776365	1361295	68.96%	6.964%
50	10.384	1387	1391	1399	rVV2	166219	298032	15.10%	1.525%
51	10.469	1399	1405	1406	rVV5	8774	14060	0.71%	0.072%
52	10.506	1406	1411	1416	rVV6	11880	27753	1.41%	0.142%
53	10.579	1416	1423	1432	rVV2	147656	254661	12.90%	1.303%
54	10.664	1432	1437	1439	rVV4	7668	12665	0.64%	0.065%
55	10.713	1439	1445	1459	rVB3	53416	128680	6.52%	0.658%
56	10.859	1464	1469	1473	rVB7	4789	9005	0.46%	0.046%
57	10.926	1473	1480	1493	rBV2	334863	601466	30.47%	3.077%
58	11.060	1496	1502	1508	rVB8	7300	13209	0.67%	0.068%
59	11.176	1517	1521	1526	rVB6	5547	9624	0.49%	0.049%
60	11.292	1533	1540	1551	rVB2	30766	59087	2.99%	0.302%
61	11.627	1588	1595	1607	rBV2	834807	1712942	86.78%	8.763%
62	11.731	1607	1612	1615	rVV5	7095	14087	0.71%	0.072%
63	11.835	1619	1629	1638	rVV2	15067	35316	1.79%	0.181%
64	12.066	1662	1667	1673	rVV3	11442	20334	1.03%	0.104%
65	12.164	1677	1683	1690	rVB2	10831	20958	1.06%	0.107%
66	12.359	1705	1715	1719	rBV10	7059	15085	0.76%	0.077%
67	12.469	1728	1733	1737	rBV4	5632	9439	0.48%	0.048%
68	12.603	1748	1755	1760	rBV3	16158	29064	1.47%	0.149%
69	12.688	1760	1769	1775	rBV	338665	570077	28.88%	2.916%
70	12.749	1775	1779	1786	rVV3	14490	27426	1.39%	0.140%
71	12.859	1794	1797	1798	rBV3	3386	3399	0.17%	0.017%
72	12.889	1800	1802	1805	rVV4	5008	6532	0.33%	0.033%
73	12.938	1806	1810	1815	rVV7	7368	14499	0.73%	0.074%
74	13.188	1845	1851	1853	rBV6	6031	10978	0.56%	0.056%
75	13.249	1855	1861	1865	rVV8	8686	19860	1.01%	0.102%
76	13.359	1876	1879	1882	rBV4	2931	3573	0.18%	0.018%

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 Title : SFAM01.0

77	13.493	1894	1901	1904	rVV9	6329	13954	0.71%	0.071%
78	13.554	1904	1911	1923	rVV	879017	1457069	73.82%	7.454%
79	13.645	1923	1926	1933	rVB9	5673	9078	0.46%	0.046%
80	13.853	1952	1960	1970	rVB	713319	1262418	63.96%	6.458%
81	13.987	1978	1982	1984	rBV5	2898	3563	0.18%	0.018%
82	14.066	1989	1995	2000	rVV3	5830	11562	0.59%	0.059%
83	14.194	2012	2016	2019	rBV5	3849	5448	0.28%	0.028%
84	14.322	2031	2037	2039	rBV5	3499	5843	0.30%	0.030%
85	14.468	2055	2061	2067	rVV3	32444	61534	3.12%	0.315%
86	14.517	2067	2069	2074	rVB5	4446	5842	0.30%	0.030%
87	14.627	2083	2087	2090	rBV5	4100	6244	0.32%	0.032%
88	14.688	2092	2097	2099	rVV6	3924	6885	0.35%	0.035%
89	14.737	2099	2105	2112	rVV4	37167	75523	3.83%	0.386%
90	14.822	2115	2119	2122	rVV5	2703	5284	0.27%	0.027%
91	14.852	2122	2124	2130	rVB6	3259	4994	0.25%	0.026%
92	15.084	2158	2162	2163	rVV4	2442	3899	0.20%	0.020%
93	15.127	2163	2169	2176	rVB	65636	116640	5.91%	0.597%
94	15.487	2221	2228	2232	rBV	21399	39926	2.02%	0.204%
95	15.547	2232	2238	2245	rBV2	53277	95417	4.83%	0.488%
96	15.852	2283	2288	2292	rBV8	3032	4980	0.25%	0.025%
97	16.108	2325	2330	2331	rBV5	2615	3836	0.19%	0.020%
98	16.444	2378	2385	2387	rBV7	5492	12724	0.64%	0.065%
99	16.492	2389	2393	2401	rVB7	8955	18640	0.94%	0.095%
100	16.730	2424	2432	2438	rVB7	5668	13667	0.69%	0.070%

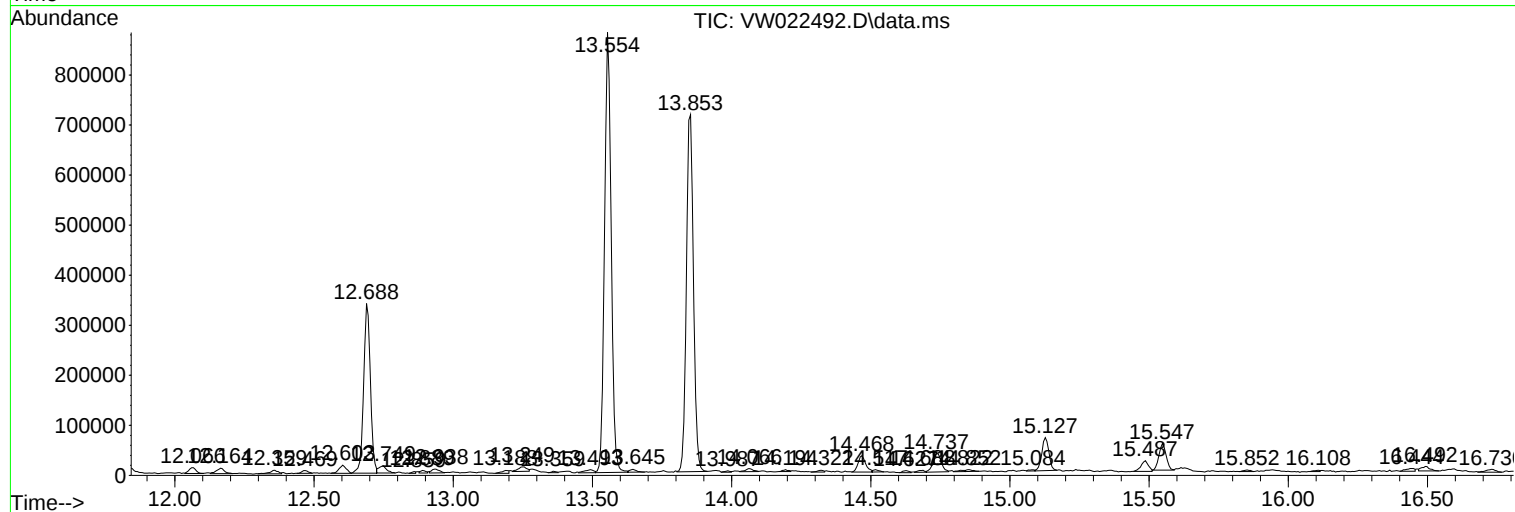
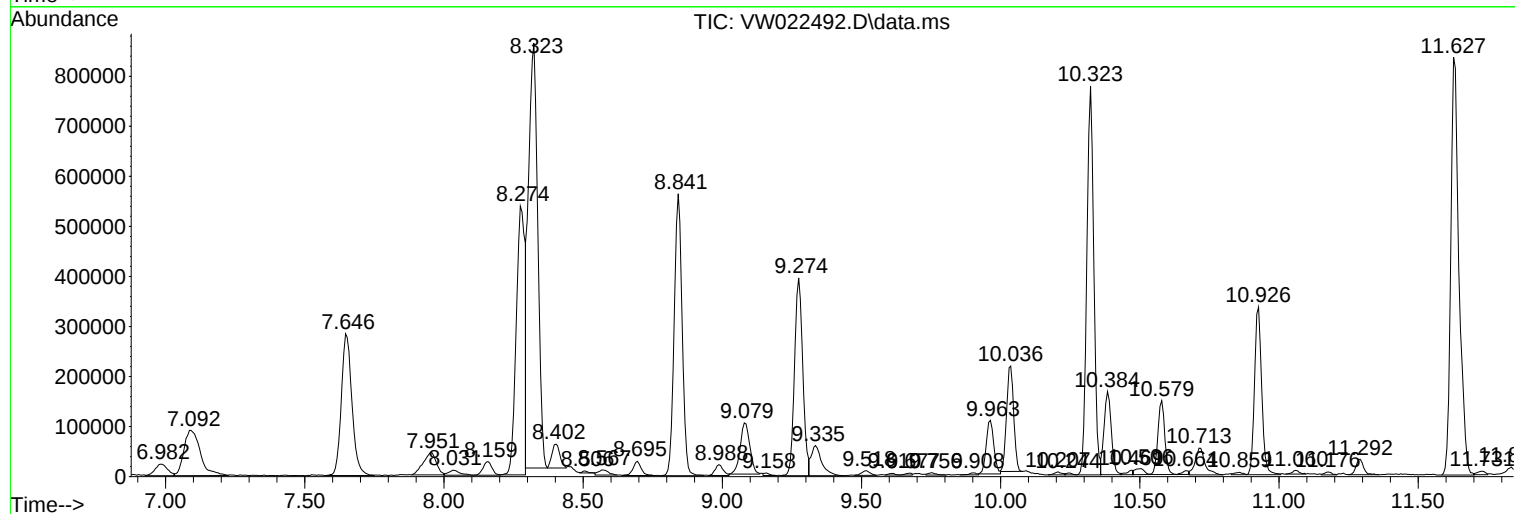
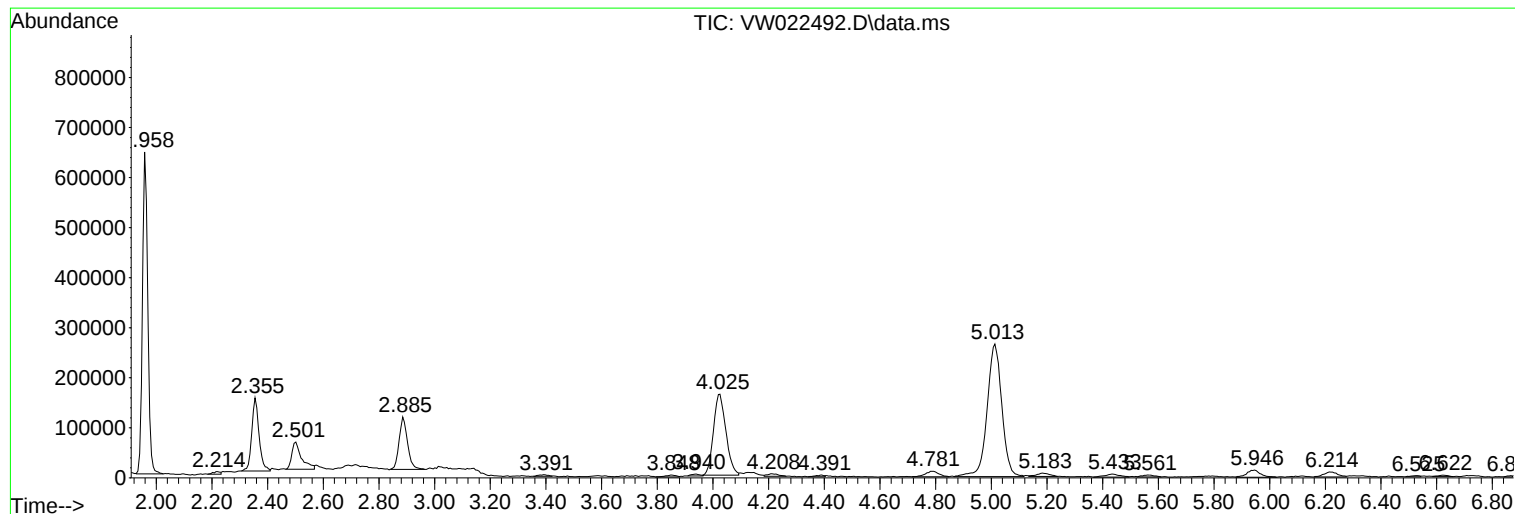
Sum of corrected areas: 19547662

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Instrument :
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ETNR3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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



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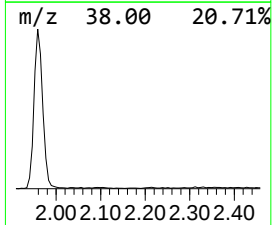
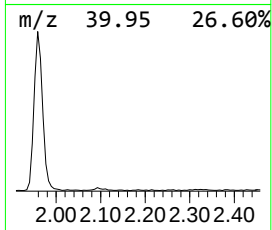
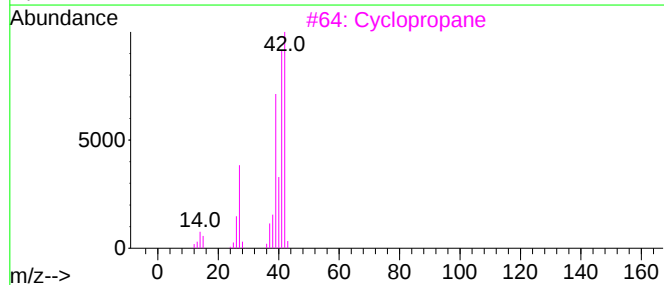
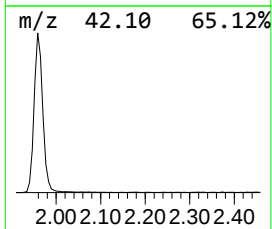
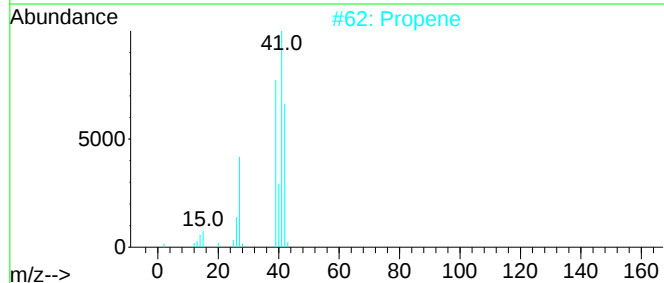
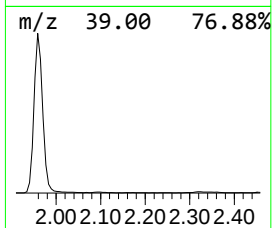
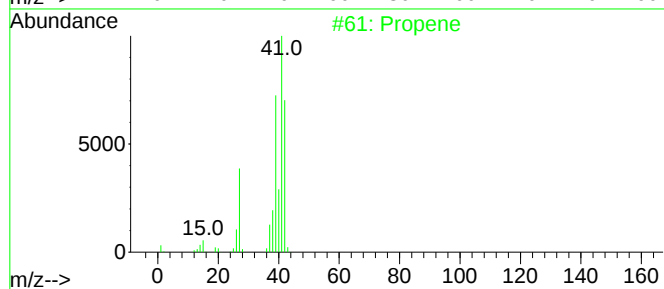
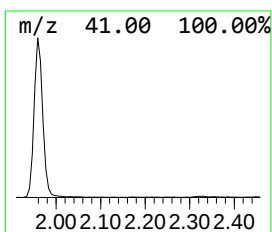
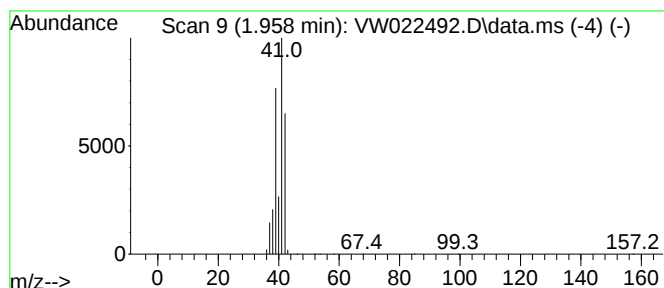
Instrument :
MSVOA_W
ClientSampleId :
ETNR3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.958	19.23 ug/L	860491	1,4-Difluorobenzene	8.841	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Propene	42	C3H6	000115-07-1	59
3	Cyclopropane	42	C3H6	000075-19-4	10
4	Cyclopropene	40	C3H4	002781-85-3	10
5	Cyclopropane	42	C3H6	000075-19-4	7



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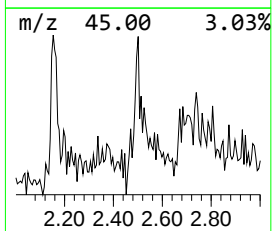
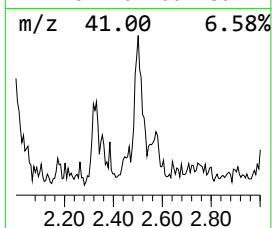
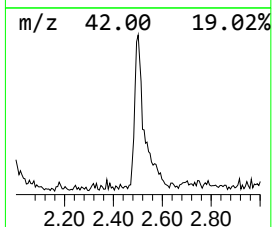
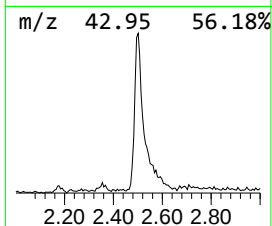
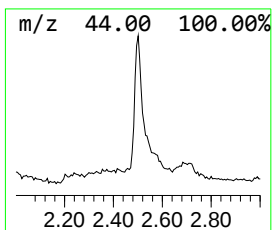
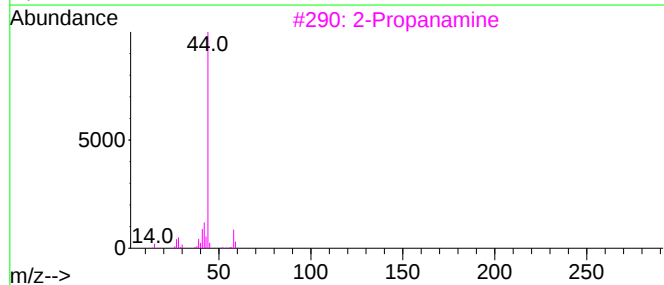
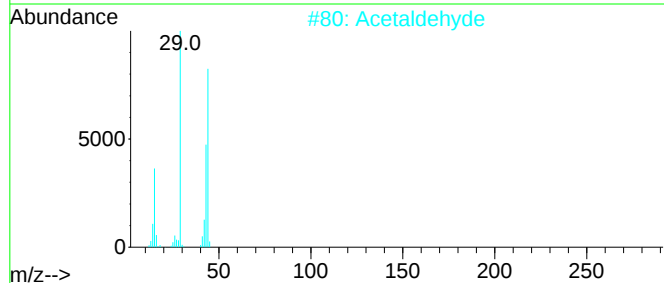
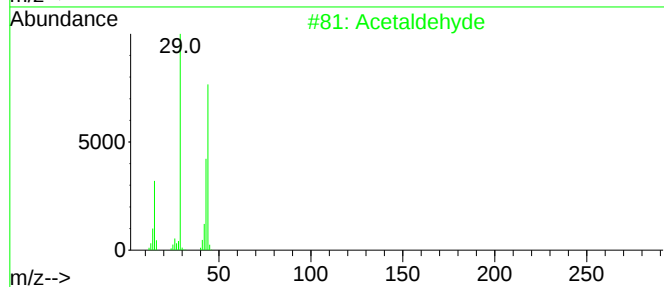
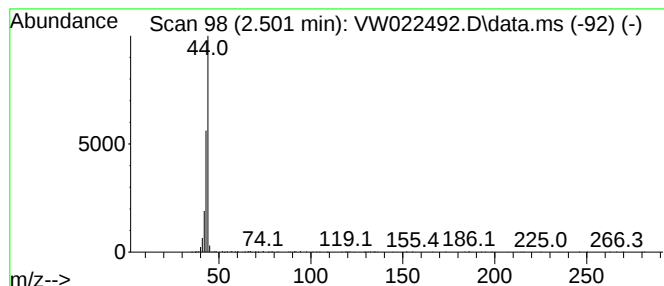
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 Acetaldehyde Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.501	2.96 ug/L	132546	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetaldehyde	44	C2H4O	000075-07-0	56
2			Acetaldehyde	44	C2H4O	000075-07-0	56
3			2-Propanamine	59	C3H9N	000075-31-0	9
4			1,2-Propanediamine	74	C3H10N2	000078-90-0	7
5			Acetic acid, [(aminocarbonyl)ami...	132	C3H4N2O4	000585-05-7	7



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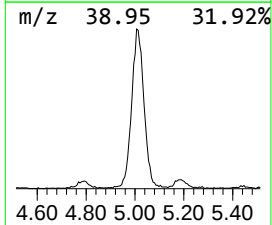
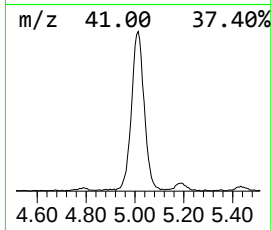
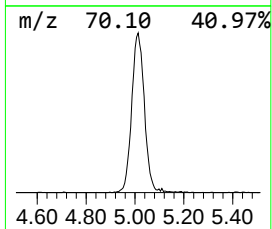
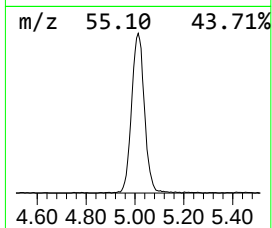
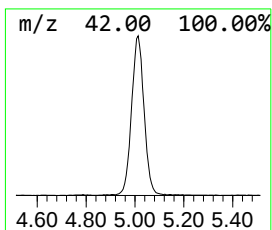
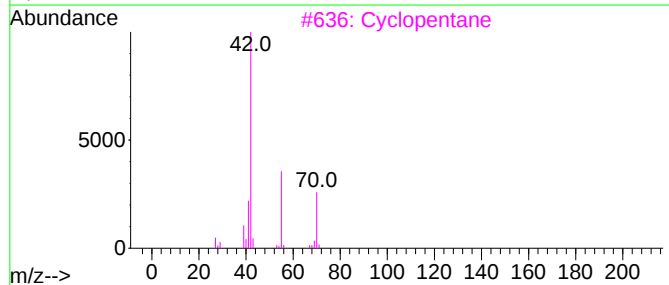
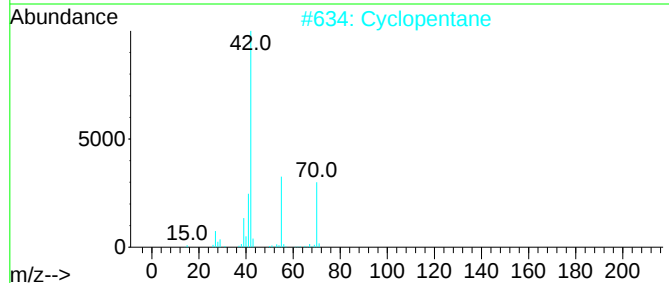
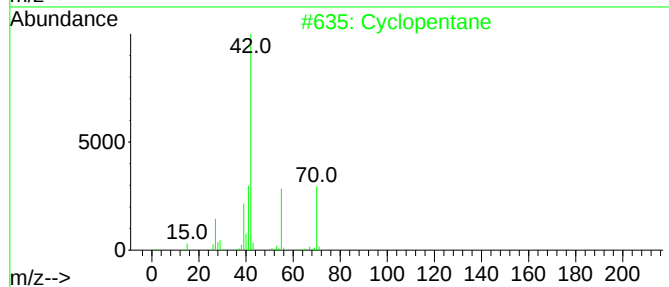
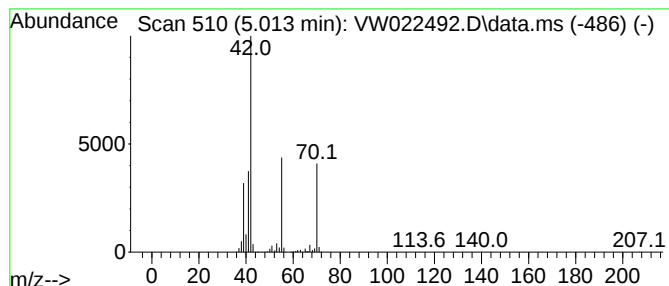
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Quant Title : SFAM01.0

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

Peak Number 3 (DEL) Alkane: Cyclic5.013 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.013	21.69 ug/L	970610	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane	70	C5H10	000287-92-3	90
2			Cyclopentane	70	C5H10	000287-92-3	86
3			Cyclopentane	70	C5H10	000287-92-3	72
4			Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
5			1-Pentene	70	C5H10	000109-67-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR3

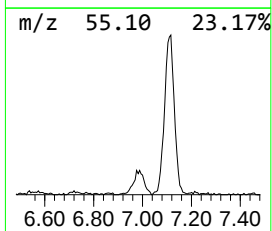
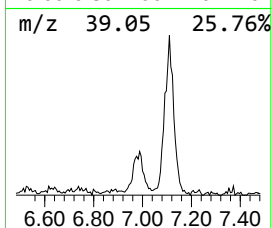
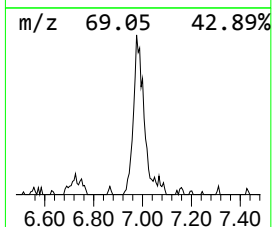
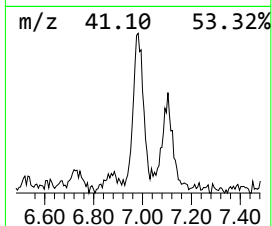
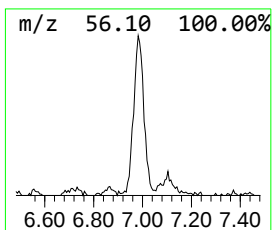
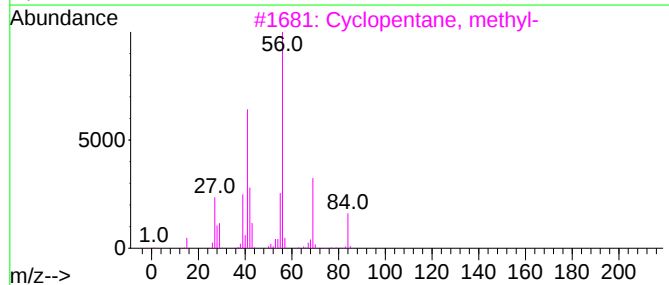
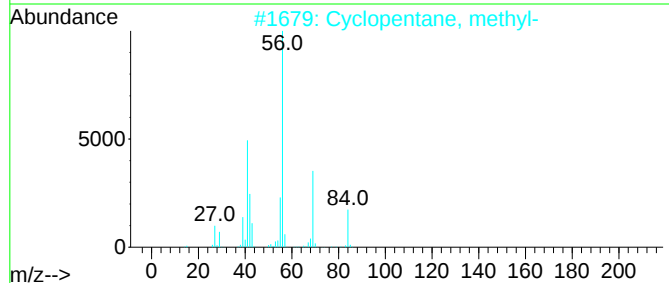
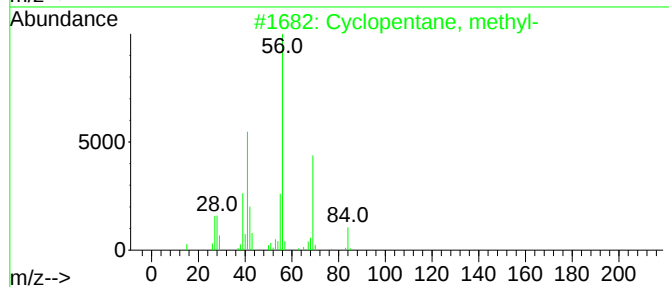
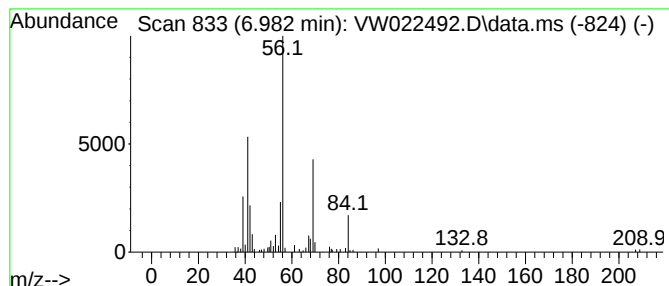
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Quant Title : SFAM01.0

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

Peak Number 4 (DEL) Alkane: Cyclic6.982 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.982	1.61 ug/L	72054	1,4-Difluorobenzene	8.841

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	80
3			Cyclopentane, methyl-	84	C6H12	000096-37-7	80
4			Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR3

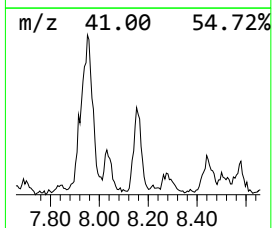
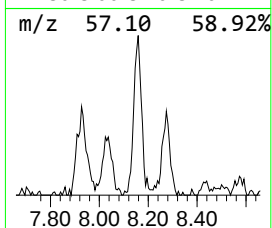
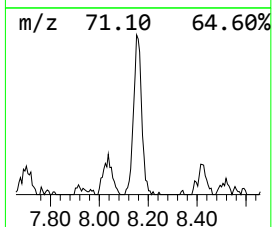
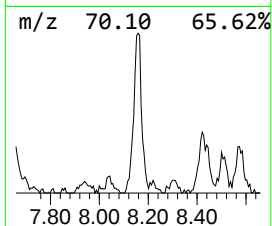
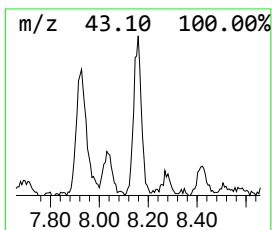
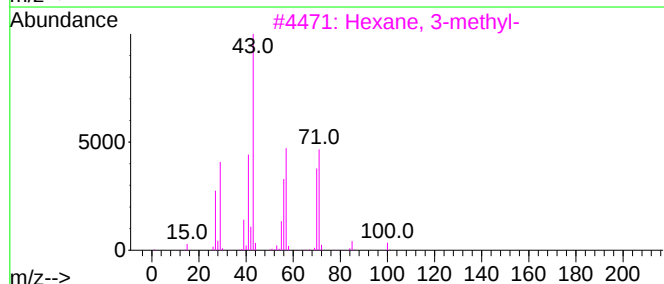
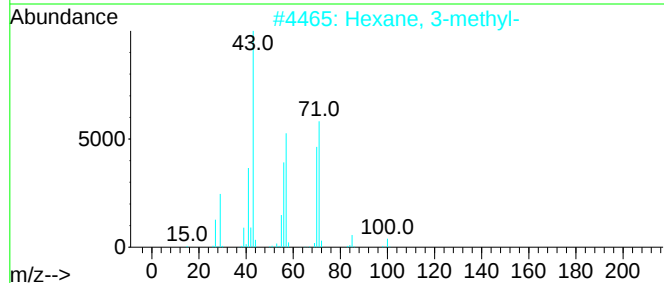
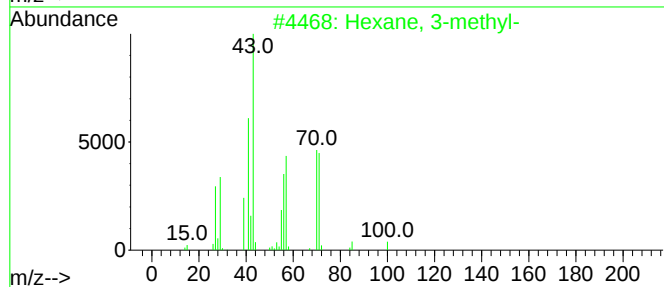
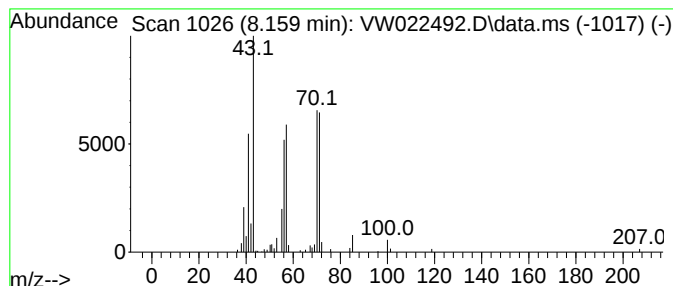
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Quant Title : SFAM01.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.159	1.35 ug/L	60208	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
3			Hexane, 3-methyl-	100	C7H16	000589-34-4	87
4			Heptane	100	C7H16	000142-82-5	58
5			Heptane, 4-methyl-	114	C8H18	000589-53-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR3

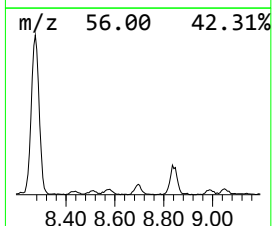
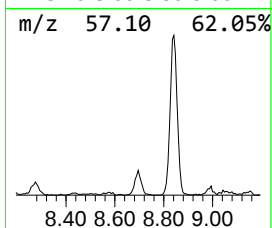
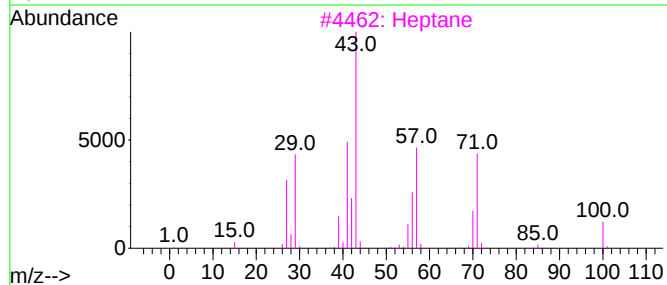
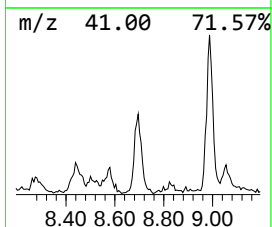
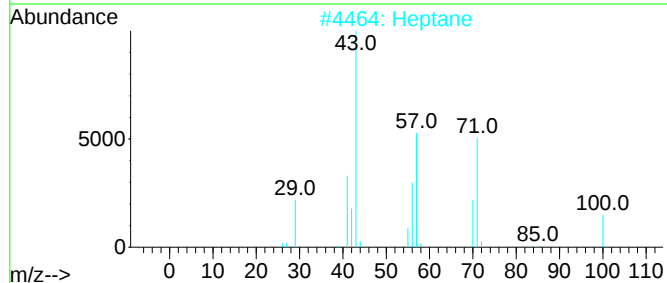
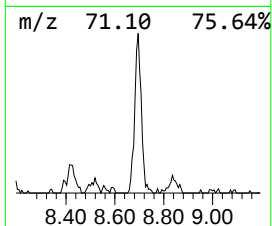
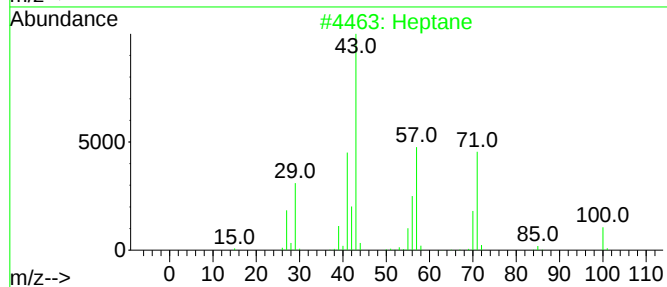
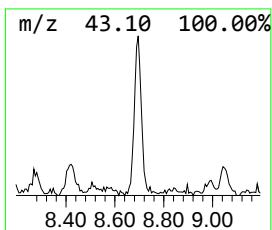
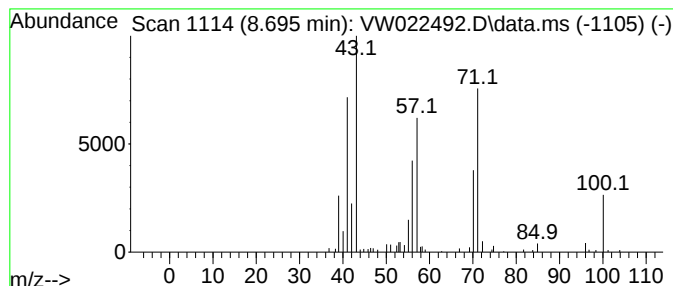
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Quant Title : SFAM01.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.695	1.30 ug/L	58230	1,4-Difluorobenzene	8.841

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	86
3		Heptane	100	C7H16	000142-82-5	64
4		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	47
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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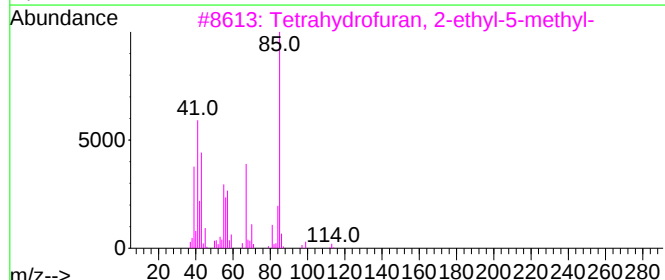
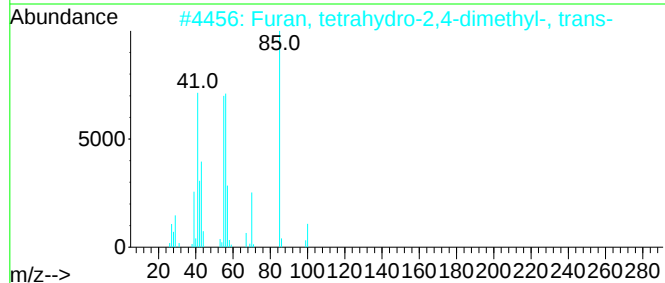
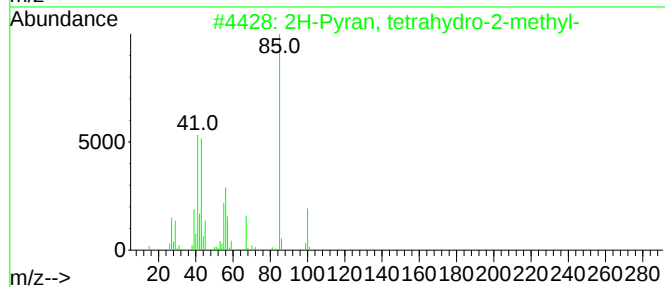
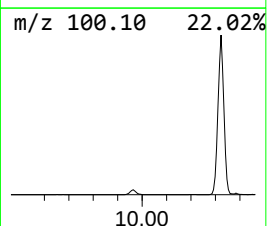
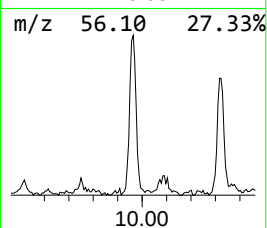
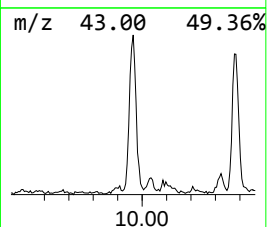
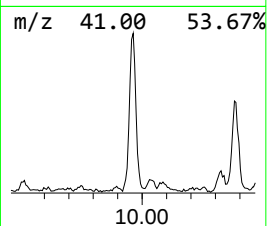
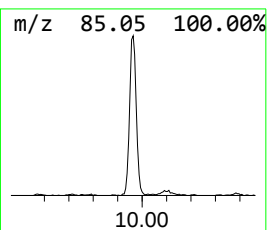
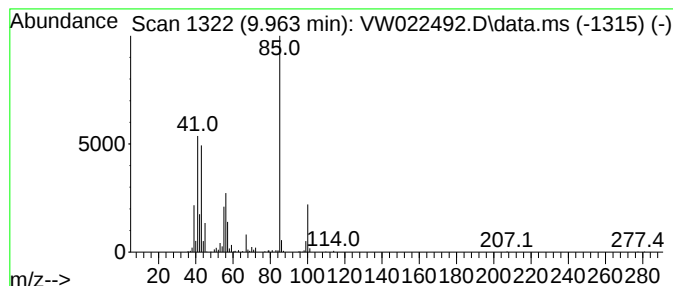
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Quant Title : SFAM01.0

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

Peak Number 7 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	4.52 ug/L	202136	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	94
2	Furan, tetrahydro-2,4-dimethyl-,...			100	C6H12O	039168-02-0	78
3	Tetrahydrofuran, 2-ethyl-5-methyl-			114	C7H14O	000931-39-5	64
4	1-Methoxy-3-methyl-2-butene			100	C6H12O	022093-99-8	58
5	1,3-Butadiene, 1-(methylthio)-			100	C5H8S	010574-97-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR3

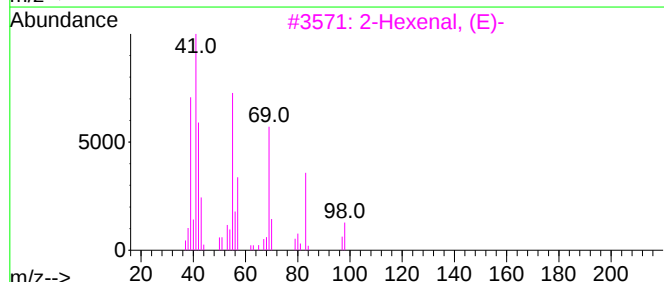
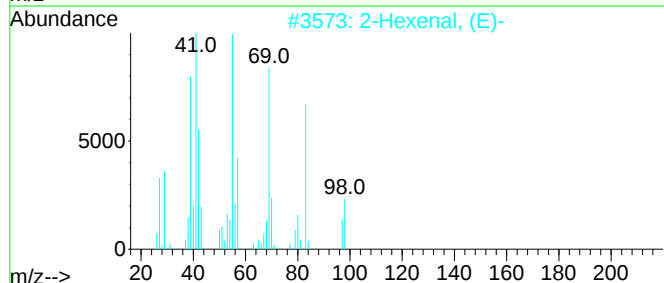
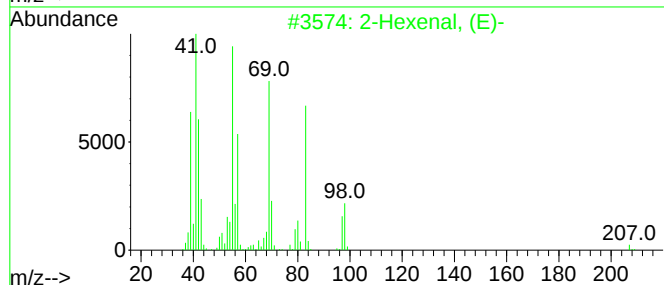
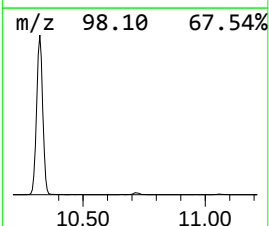
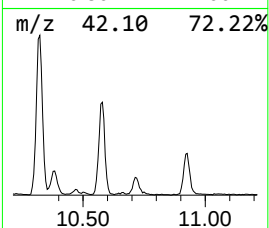
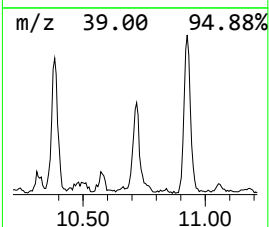
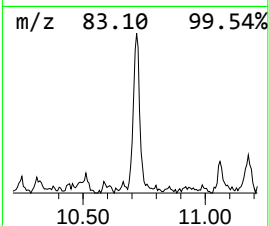
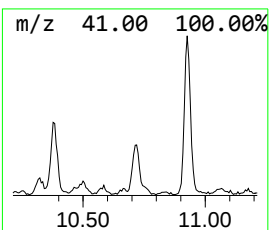
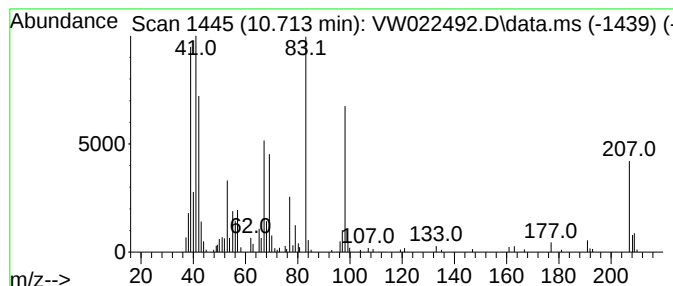
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Quant Title : SFAM01.0

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

Peak Number 9 2-Hexenal, (E)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.713	1.88 ug/L	128680	Chlorobenzene-d5	11.627

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexenal, (E)-			98	C6H10O	006728-26-3	89
2	2-Hexenal, (E)-			98	C6H10O	006728-26-3	76
3	2-Hexenal, (E)-			98	C6H10O	006728-26-3	68
4	2-Hexenal, (E)-			98	C6H10O	006728-26-3	64
5	2-Pentenal, 2-methyl-			98	C6H10O	000623-36-9	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR3

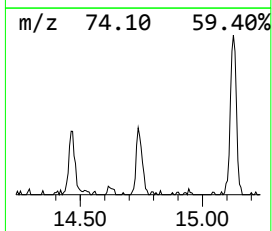
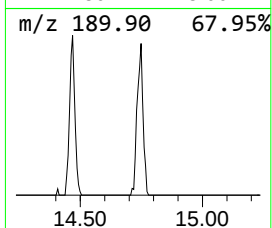
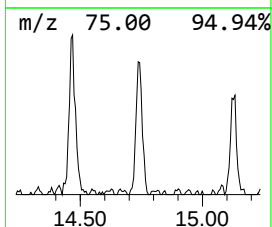
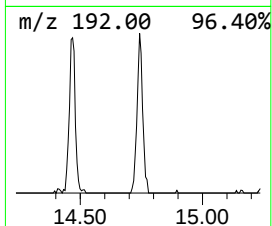
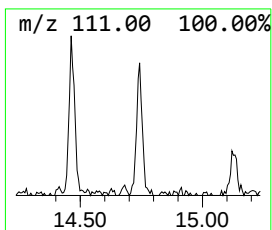
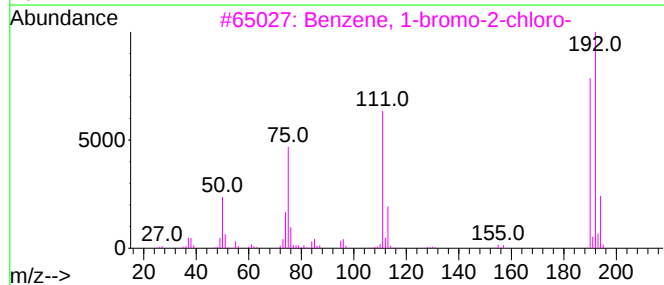
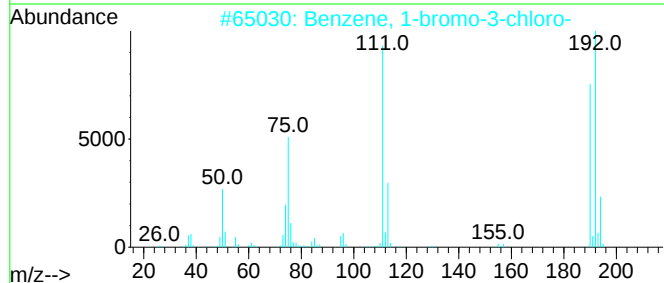
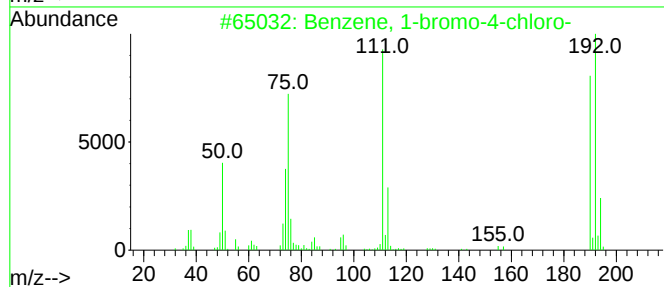
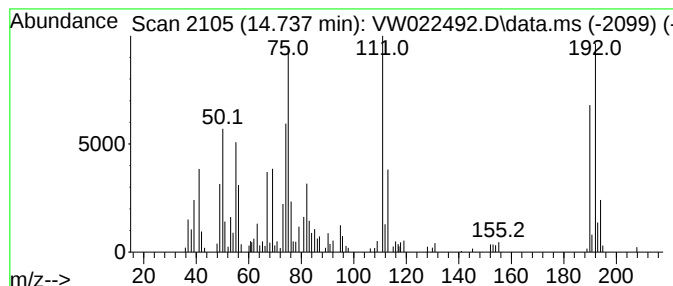
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Quant Title : SFAM01.0

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

Peak Number 10 Benzene, 1-bromo-4-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.737	1.30 ug/L	75523	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
2			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	93
5			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022492.D
Acq On : 11 Apr 2022 18:57
Operator : SY/VA
Sample : N2316-13
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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...	1.958	19.2	ug/L	860491	1	8.841	1118630	25.0
Acetaldehyde	2.501	3.0	ug/L	132546	1	8.841	1118630	25.0
(DEL) Alkane: C...	5.013	21.7	ug/L	970610	1	8.841	1118630	25.0
(DEL) Alkane: C...	6.982	1.6	ug/L	72054	1	8.841	1118630	25.0
(DEL) Alkane: S...	8.159	1.4	ug/L	60208	1	8.841	1118630	25.0
(DEL) Alkane: S...	8.695	1.3	ug/L	58230	1	8.841	1118630	25.0
2H-Pyran, tetra...	9.963	4.5	ug/L	202136	1	8.841	1118630	25.0
2-Hexenal, (E)-	10.713	1.9	ug/L	128680	2	11.627	1712940	25.0
Benzene, 1-brom...	14.737	1.3	ug/L	75523	3	13.554	1457070	25.0