

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071922\
 Data File : VW023876.D
 Acq On : 19 Jul 2022 11:51
 Operator : SY/VA
 Sample : N3709-06
 Misc : 6.00g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 H0B22

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\SFAMWLM071322SMA.M
 Title : SFAM01.0

Signal : TIC: VW023876.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.959	6	9	17	rVB2	32426	53931	5.38%	0.632%
2	2.178	42	45	48	rVB3	4366	5114	0.51%	0.060%
3	2.251	55	57	59	rBV3	3617	3240	0.32%	0.038%
4	2.349	65	73	81	rBV	101716	192787	19.25%	2.258%
5	2.635	116	120	125	rBV6	10104	23491	2.35%	0.275%
6	2.879	153	160	169	rVB	66681	135839	13.56%	1.591%
7	3.373	235	241	251	rVB3	12310	30865	3.08%	0.361%
8	4.013	335	346	356	rBV2	78948	242607	24.22%	2.841%
9	4.123	356	364	377	rVV	86682	235897	23.55%	2.763%
10	4.385	398	407	419	rVV3	10348	35382	3.53%	0.414%
11	4.915	481	494	503	rBV3	39405	146983	14.67%	1.721%
12	5.184	534	538	544	rVV8	2162	5084	0.51%	0.060%
13	5.409	572	575	576	rBV2	1757	1978	0.20%	0.023%
14	5.787	635	637	642	rVB6	1510	2109	0.21%	0.025%
15	5.940	652	662	673	rVB6	14977	46780	4.67%	0.548%
16	6.135	691	694	697	rVB4	1264	1746	0.17%	0.020%
17	6.561	761	764	767	rVV4	1353	1909	0.19%	0.022%
18	6.970	829	831	839	rVB9	2224	4886	0.49%	0.057%
19	7.080	839	849	857	rBV	37404	107224	10.70%	1.256%
20	7.171	858	864	876	rVB	30447	86311	8.62%	1.011%
21	7.500	916	918	921	rVV3	1478	2008	0.20%	0.024%
22	7.647	933	942	955	rVB	199148	490843	49.00%	5.749%
23	7.909	978	985	986	rBV7	2281	4149	0.41%	0.049%
24	7.945	988	991	999	rVB8	4068	8922	0.89%	0.104%
25	8.146	1018	1024	1025	rBV6	1749	2776	0.28%	0.033%
26	8.275	1036	1045	1059	rBV2	321379	1001648	100.00%	11.731%
27	8.561	1088	1092	1093	rBV4	2495	2737	0.27%	0.032%
28	8.695	1106	1114	1124	rBV4	13308	29548	2.95%	0.346%
29	8.841	1129	1138	1146	rBV	410184	798351	79.70%	9.350%
30	9.012	1163	1166	1169	rBV4	941	1640	0.16%	0.019%
31	9.073	1169	1176	1184	rVB3	13019	29757	2.97%	0.349%
32	9.140	1184	1187	1190	rVB4	1274	1743	0.17%	0.020%
33	9.274	1198	1209	1216	rBV	258925	538327	53.74%	6.305%
34	9.512	1245	1248	1255	rVB8	2990	5917	0.59%	0.069%
35	9.597	1255	1262	1267	rBV8	1918	4706	0.47%	0.055%
36	9.658	1269	1272	1277	rVB6	2339	3989	0.40%	0.047%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
 Title : SFAM01.0

37	9.738	1281	1285	1286	rBV4	1571	2008	0.20%	0.024%
38	9.957	1315	1321	1326	rVB7	3898	8460	0.84%	0.099%
39	10.036	1326	1334	1344	rBV	70849	138241	13.80%	1.619%
40	10.207	1357	1362	1367	rBV8	3337	6693	0.67%	0.078%
41	10.323	1372	1381	1388	rBV	441329	770928	76.97%	9.029%
42	10.500	1406	1410	1416	rVB3	8308	14204	1.42%	0.166%
43	10.573	1417	1422	1430	rVB2	49089	84727	8.46%	0.992%
44	10.658	1432	1436	1440	rBV6	2761	4879	0.49%	0.057%
45	10.762	1448	1453	1459	rVB8	4395	8313	0.83%	0.097%
46	10.841	1461	1466	1472	rVB7	3232	6846	0.68%	0.080%
47	10.926	1473	1480	1492	rBV	108546	205572	20.52%	2.408%
48	11.097	1506	1508	1512	rVV4	1927	2670	0.27%	0.031%
49	11.176	1512	1521	1525	rVV5	9061	18357	1.83%	0.215%
50	11.225	1525	1529	1535	rVB7	5456	10298	1.03%	0.121%
51	11.274	1535	1537	1542	rBV6	1881	2482	0.25%	0.029%
52	11.335	1542	1547	1552	rVB7	1950	4252	0.42%	0.050%
53	11.426	1554	1562	1567	rBV10	4901	12429	1.24%	0.146%
54	11.628	1588	1595	1605	rBV	480340	822922	82.16%	9.638%
55	11.725	1606	1611	1614	rVV7	4643	8730	0.87%	0.102%
56	11.841	1621	1630	1633	rBV2	13092	31903	3.19%	0.374%
57	11.975	1649	1652	1657	rBV7	1715	2679	0.27%	0.031%
58	12.030	1658	1661	1666	rBV6	2109	4230	0.42%	0.050%
59	12.134	1669	1678	1681	rBV6	12291	28053	2.80%	0.329%
60	12.225	1691	1693	1700	rVB8	5440	9907	0.99%	0.116%
61	12.304	1700	1706	1707	rBV6	5080	8287	0.83%	0.097%
62	12.341	1707	1712	1716	rBV4	13521	22575	2.25%	0.264%
63	12.463	1727	1732	1737	rVB	15823	28112	2.81%	0.329%
64	12.603	1751	1755	1762	rVB2	20769	35735	3.57%	0.419%
65	12.688	1762	1769	1776	rBV	220340	379474	37.88%	4.444%
66	12.938	1805	1810	1818	rVB7	15790	35554	3.55%	0.416%
67	13.042	1822	1827	1833	rBV2	41058	72148	7.20%	0.845%
68	13.182	1845	1850	1852	rBV4	15513	25717	2.57%	0.301%
69	13.255	1858	1862	1866	rBV6	4308	8006	0.80%	0.094%
70	13.316	1870	1872	1875	rBV4	2361	3715	0.37%	0.044%
71	13.371	1876	1881	1886	rVB4	16176	36470	3.64%	0.427%
72	13.554	1905	1911	1921	rBV2	295599	560313	55.94%	6.562%
73	13.646	1923	1926	1932	rVB4	7914	14758	1.47%	0.173%
74	13.749	1940	1943	1944	rBV3	2565	2452	0.24%	0.029%
75	13.798	1946	1951	1954	rVV	17936	30967	3.09%	0.363%
76	13.853	1954	1960	1971	rVV	231125	461286	46.05%	5.403%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
 Title : SFAM01.0

77	13.944	1971	1975	1983	rVB	28936	50228	5.01%	0.588%
78	14.024	1983	1988	1991	rBV5	5037	8896	0.89%	0.104%
79	14.060	1991	1994	1996	rVV	7712	10325	1.03%	0.121%
80	14.091	1996	1999	2003	rVB2	12743	19886	1.99%	0.233%
81	14.200	2011	2017	2023	rBV2	30810	56506	5.64%	0.662%
82	14.267	2026	2028	2032	rVB5	6069	7837	0.78%	0.092%
83	14.328	2033	2038	2044	rBV5	12892	27003	2.70%	0.316%
84	14.383	2044	2047	2051	rVV5	4504	6258	0.62%	0.073%
85	14.493	2062	2065	2068	rBV4	3843	5194	0.52%	0.061%
86	14.615	2083	2085	2086	rBV2	3133	2639	0.26%	0.031%
87	14.682	2092	2096	2101	rBV4	15425	26744	2.67%	0.313%
88	14.786	2109	2113	2121	rVB2	11848	23358	2.33%	0.274%
89	14.853	2122	2124	2128	rVV4	1710	1820	0.18%	0.021%
90	14.907	2131	2133	2138	rVV6	1561	1809	0.18%	0.021%
91	15.170	2173	2176	2183	rVB6	5048	9718	0.97%	0.114%
92	15.359	2202	2207	2209	rBV6	2420	3701	0.37%	0.043%
93	15.401	2209	2214	2220	rVV4	10137	16200	1.62%	0.190%
94	15.474	2222	2226	2232	rVB9	3077	6978	0.70%	0.082%
95	15.791	2274	2278	2283	rVB6	6125	8277	0.83%	0.097%
96	15.962	2298	2306	2311	rVB10	3415	7317	0.73%	0.086%
97	16.017	2311	2315	2318	rBV5	1820	3327	0.33%	0.039%
98	16.169	2335	2340	2343	rVB6	1815	3084	0.31%	0.036%
99	16.224	2347	2349	2352	rVB4	1910	2048	0.20%	0.024%
100	16.419	2375	2381	2383	rBV5	1604	2438	0.24%	0.029%

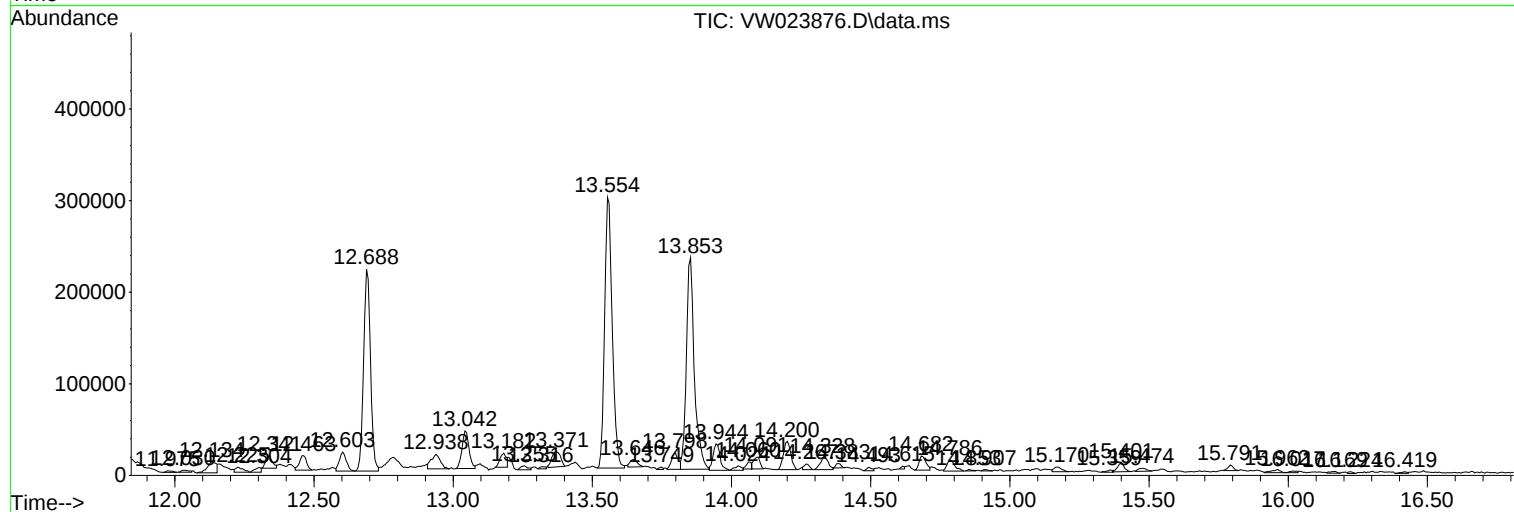
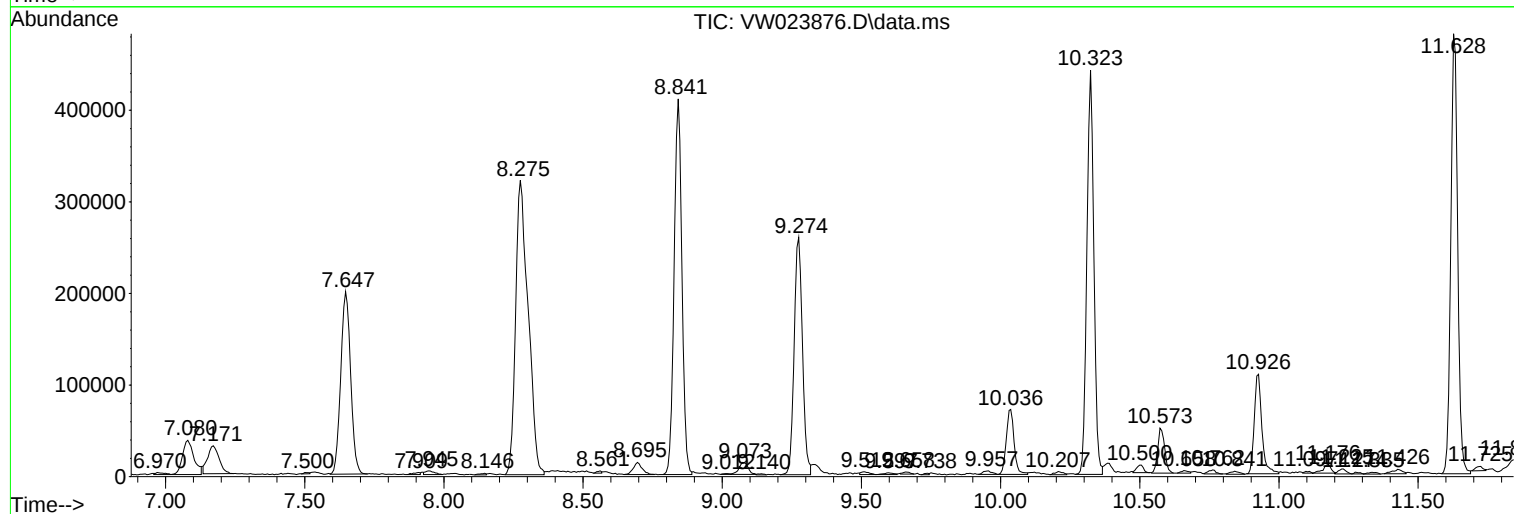
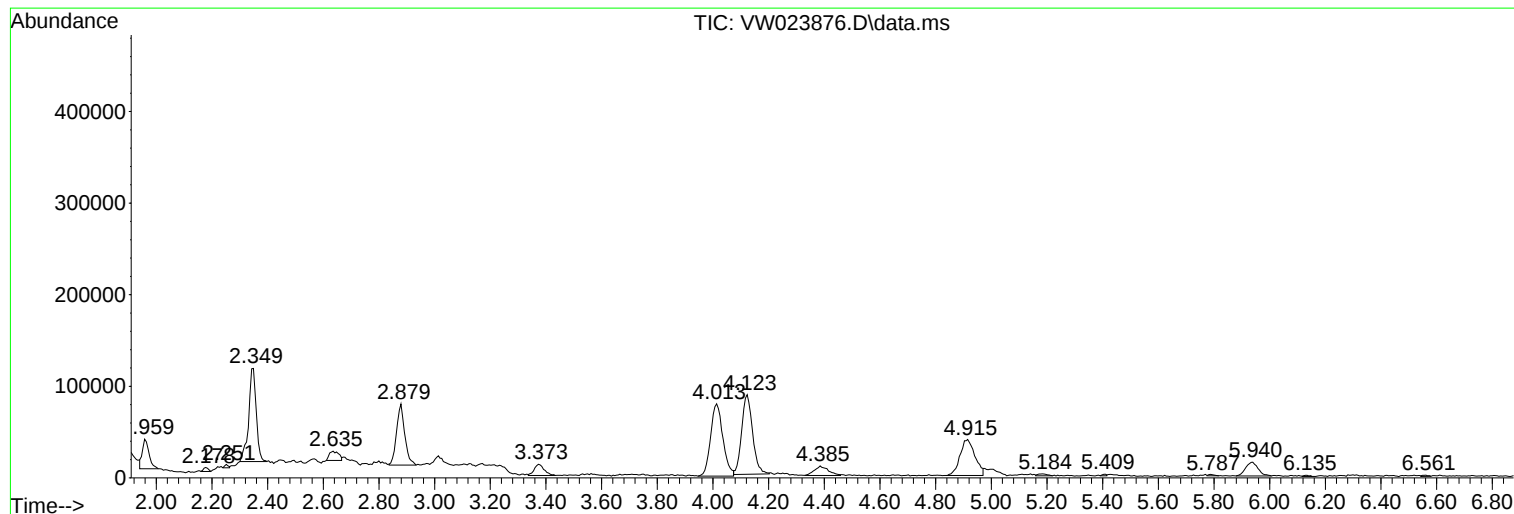
Sum of corrected areas: 8538167

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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



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Instrument :
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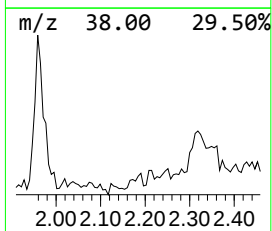
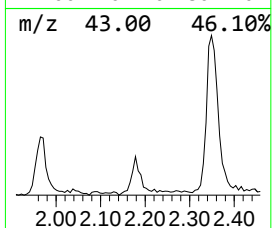
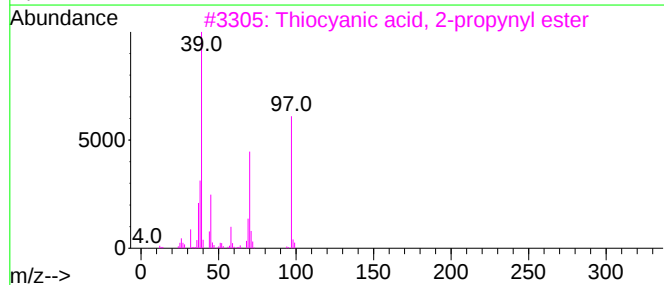
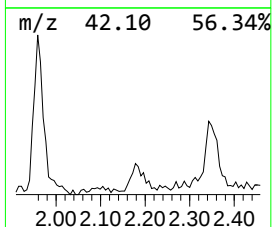
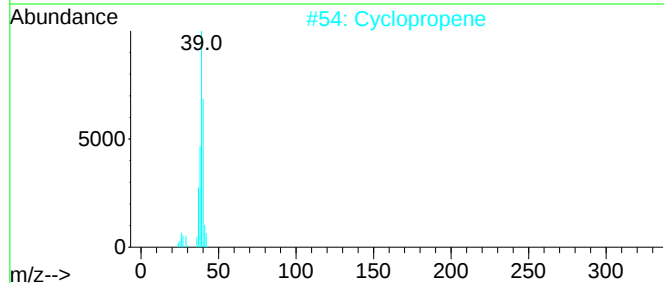
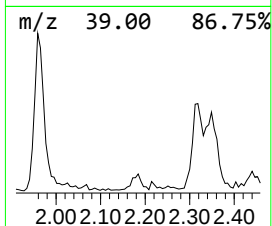
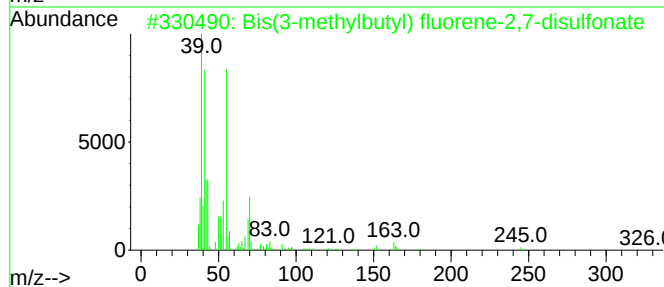
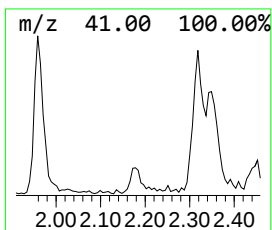
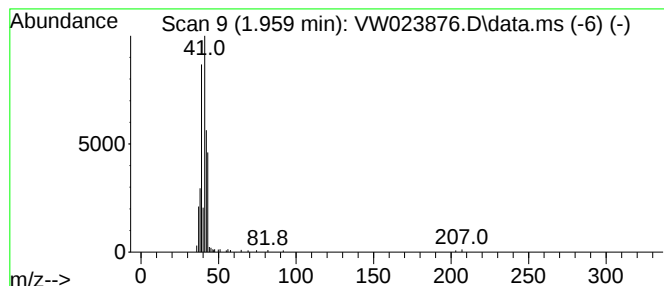
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.959	1.69 ug/L	53931	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(3-methylbutyl) fluorene-2,7-...	466	C23H30O6S2	253664-95-8	40
2			Cyclopropene	40	C3H4	002781-85-3	40
3			Thiocyanic acid, 2-propynyl ester	97	C4H3NS	024309-48-6	40
4			2-Butynal	68	C4H4O	001119-19-3	40
5			2-Butenal, (E)-	70	C4H6O	000123-73-9	40



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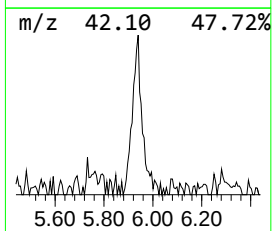
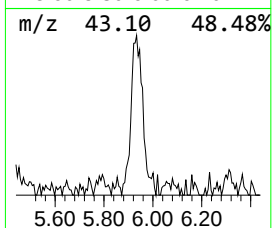
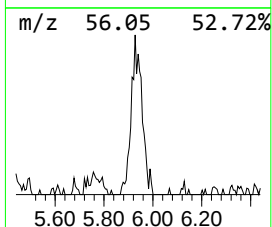
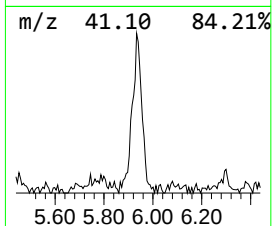
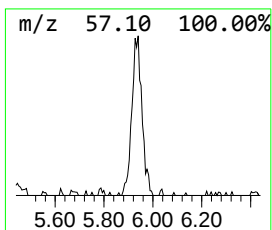
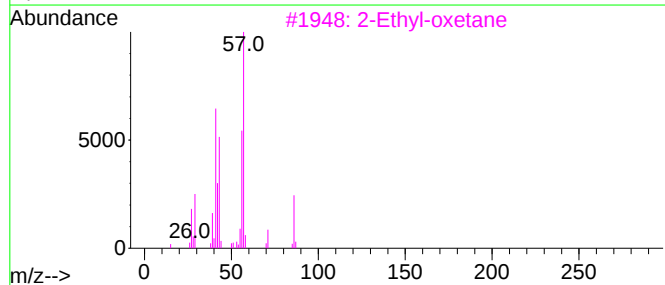
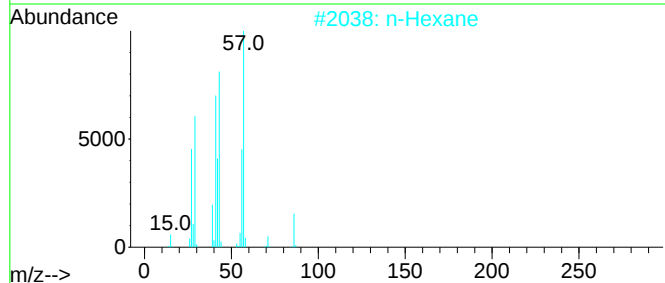
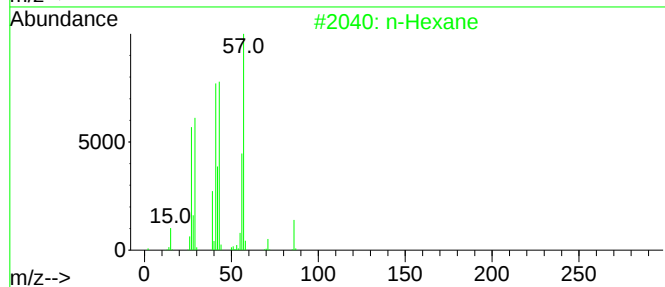
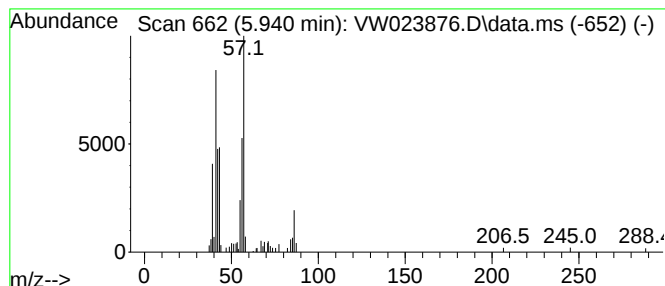
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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.
5.940	1.46 ug/L	46780	1,4-Difluorobenzene	8.841

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexane	86	C6H14	000110-54-3	52
2		n-Hexane	86	C6H14	000110-54-3	50
3		2-Ethyl-oxetane	86	C5H10O	1010386-40-2	49
4		n-Hexane	86	C6H14	000110-54-3	42
5		1-Hexene	84	C6H12	000592-41-6	42



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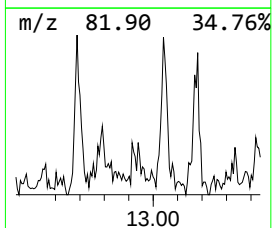
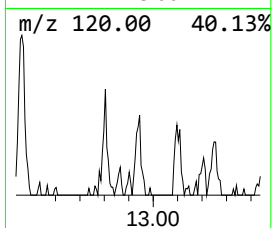
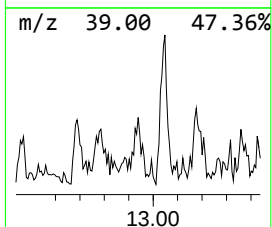
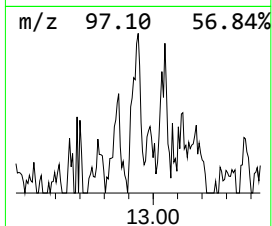
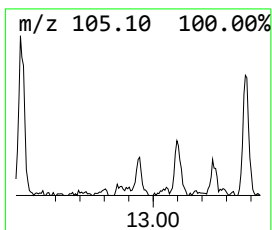
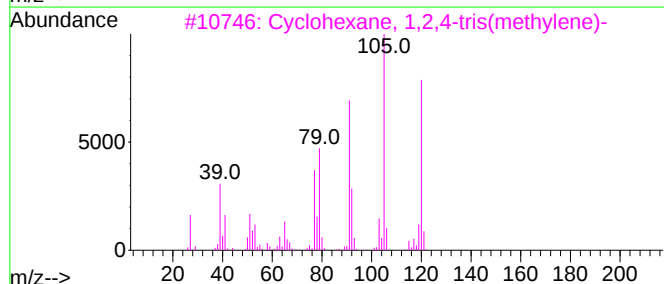
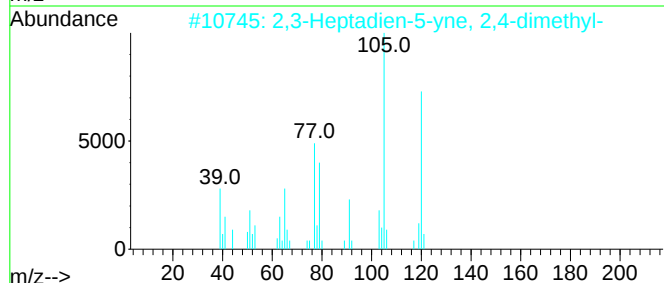
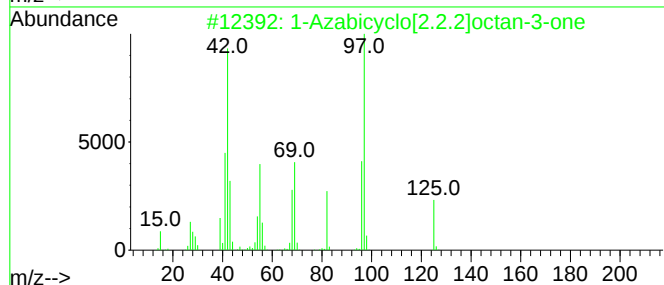
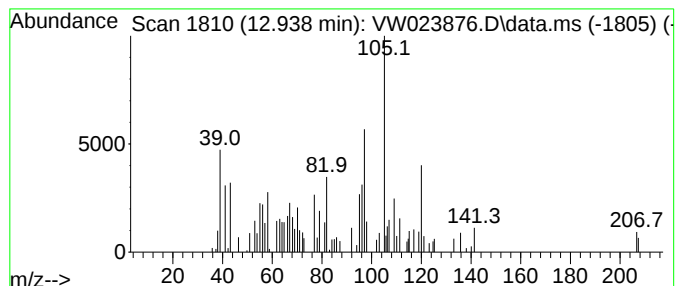
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.938	1.59 ug/L	35554	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Azabicyclo[2.2.2]octan-3-one	125	C7H11NO	003731-38-2	14
2			2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	11
3			Cyclohexane, 1,2,4-tris(methylene)-	120	C9H12	014296-81-2	11
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	10
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071922\
Data File : VW023876.D
Acq On : 19 Jul 2022 11:51
Operator : SY/VA
Sample : N3709-06
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
H0B22

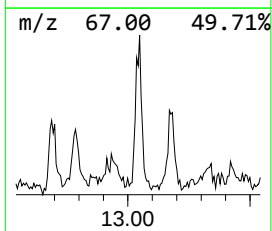
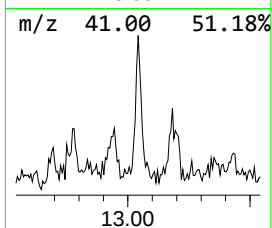
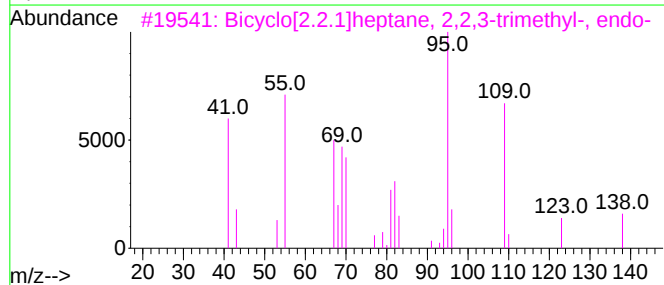
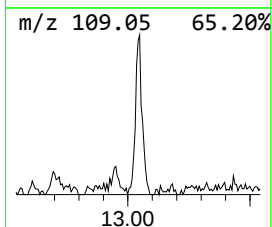
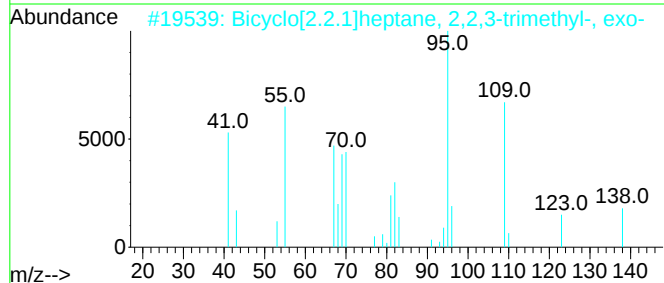
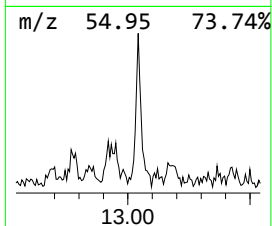
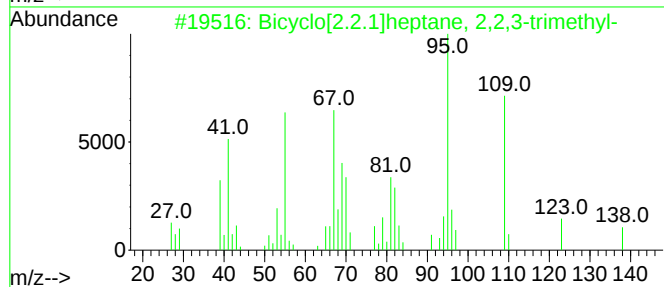
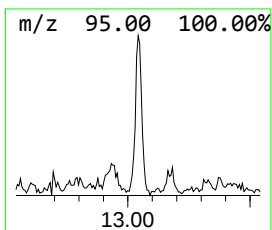
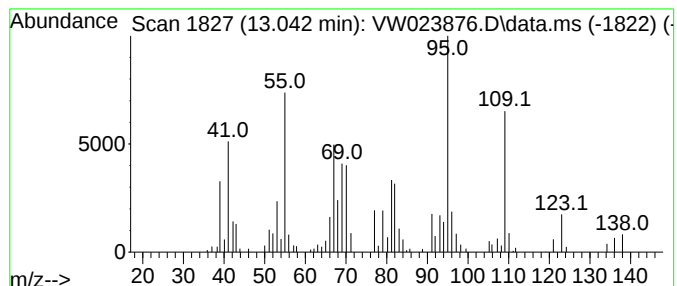
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Quant Title : SFAM01.0

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

Peak Number 6 Bicyclo[2.2.1]heptane, 2,2,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.042	3.22 ug/L	72148	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	95
2			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-41-8	80
3			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	80
4			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	53
5			2-Nonen-4-yn-1-ol, (Z)-	138	C9H14O	134225-90-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071922\
Data File : VW023876.D
Acq On : 19 Jul 2022 11:51
Operator : SY/VA
Sample : N3709-06
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
H0B22

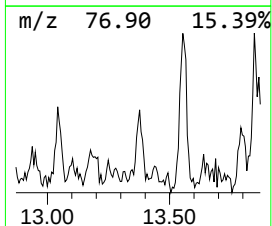
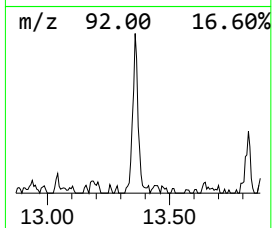
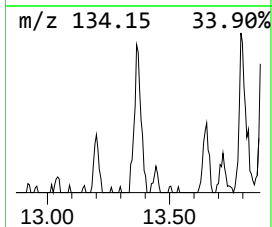
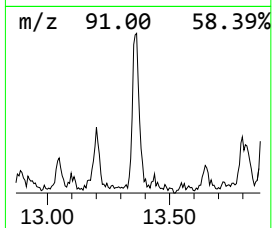
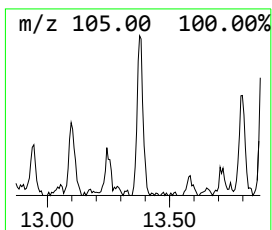
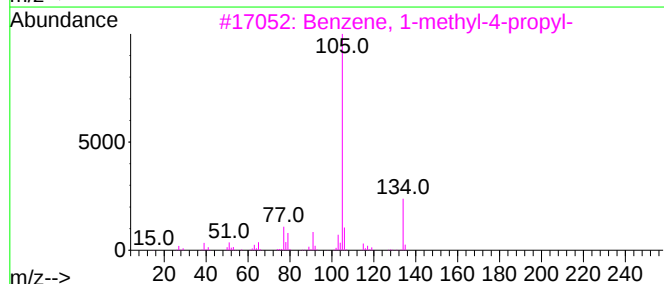
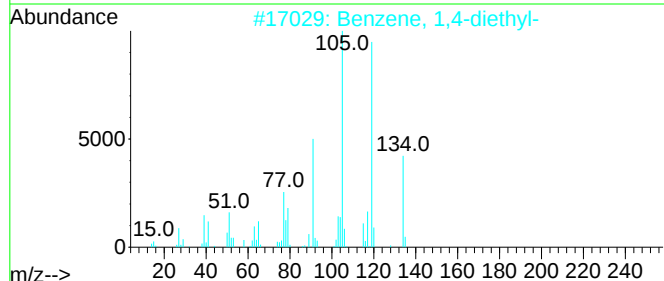
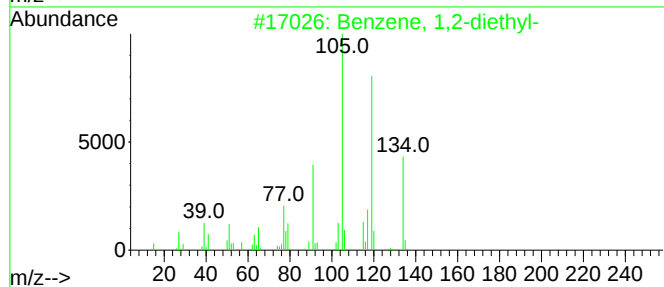
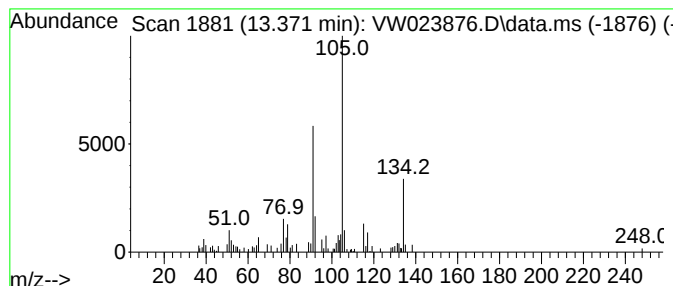
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Quant Title : SFAM01.0

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

Peak Number 7 Benzene, 1,2-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.371	1.63 ug/L	36470	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	64
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	62
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	58
4			2-Tolyloxirane	134	C9H10O	002783-26-8	58
5			p-Mentha-1,5,8-triene	134	C10H14	021195-59-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071922\
Data File : VW023876.D
Acq On : 19 Jul 2022 11:51
Operator : SY/VA
Sample : N3709-06
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
H0B22

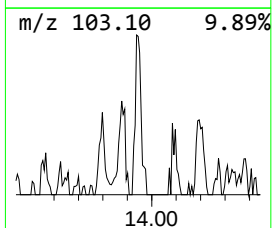
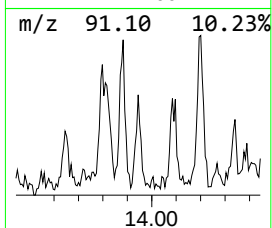
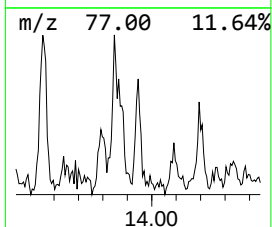
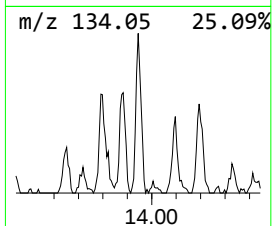
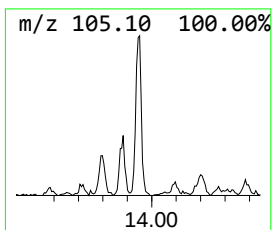
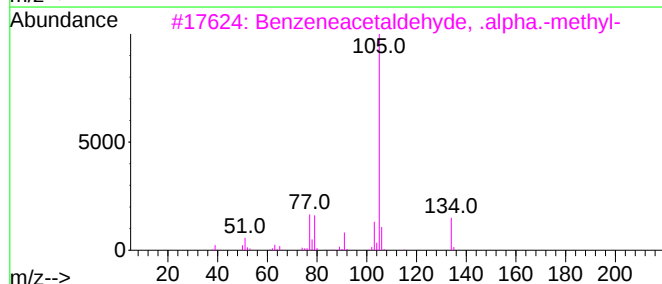
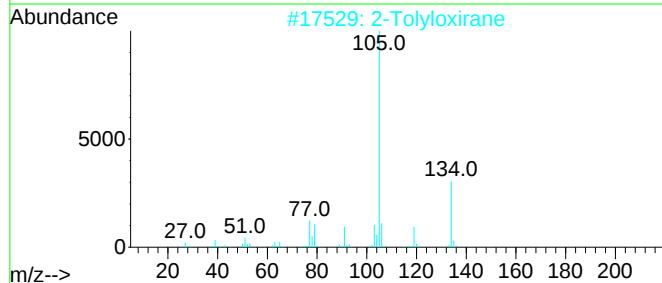
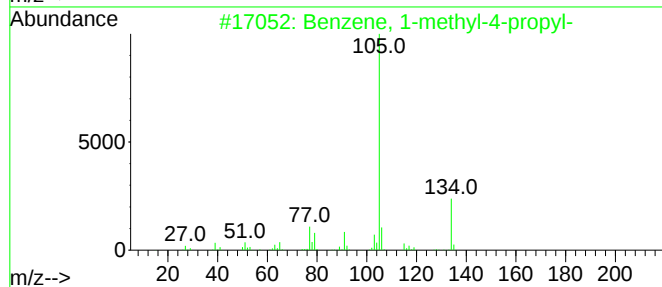
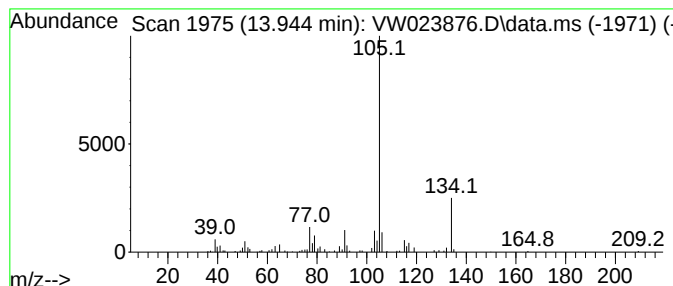
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Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-methyl-4-propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.944	2.24 ug/L	50228	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
2			2-Tolyloxirane	134	C9H10O	002783-26-8	87
3			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	87
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	80
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071922\
Data File : VW023876.D
Acq On : 19 Jul 2022 11:51
Operator : SY/VA
Sample : N3709-06
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
H0B22

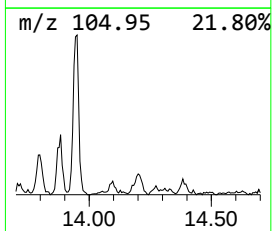
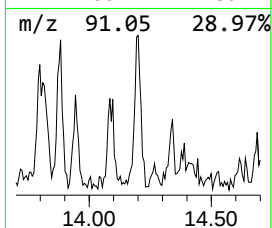
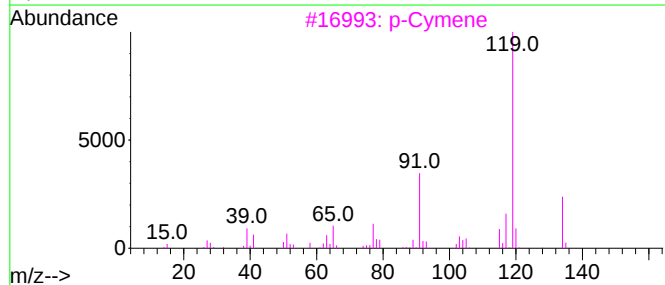
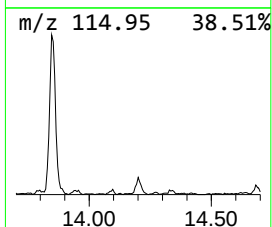
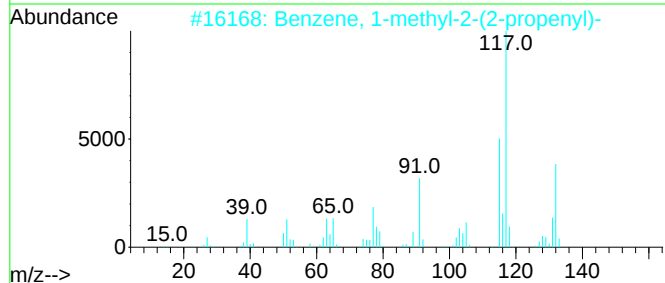
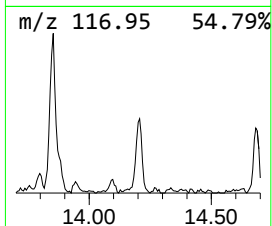
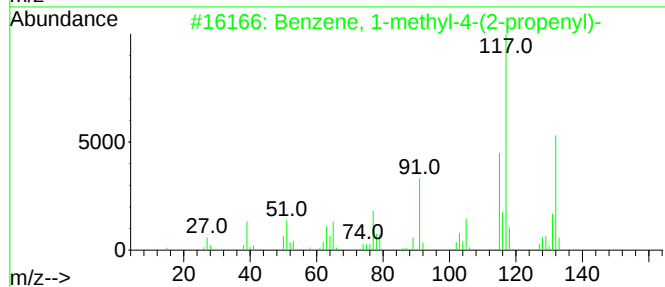
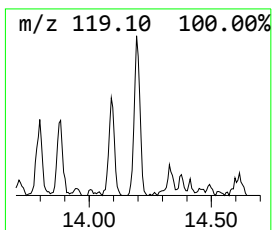
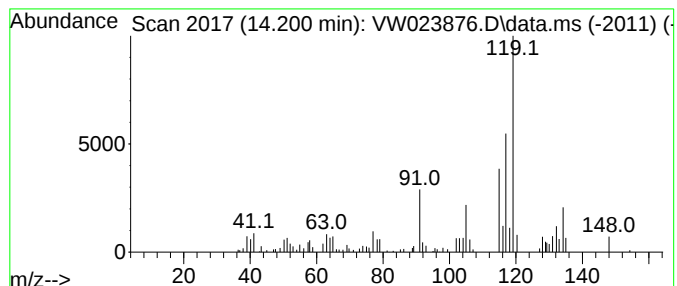
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Quant Title : SFAM01.0

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

Peak Number 9 Benzene, 1-methyl-4-(2-prop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.200	2.52 ug/L	56506	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	50
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	46
3			p-Cymene	134	C10H14	000099-87-6	43
4			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	42
5			2,4-Dimethylstyrene	132	C10H12	002234-20-0	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071922\
Data File : VW023876.D
Acq On : 19 Jul 2022 11:51
Operator : SY/VA
Sample : N3709-06
Misc : 6.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
H0B22

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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					#	RT	Resp	Conc
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(DEL) Alkane: S...	5.940	1.5	ug/L	46780	1	8.841	798351	25.0
unknown-02	12.938	1.6	ug/L	35554	3	13.554	560313	25.0
Bicyclo[2.2.1]h...	13.042	3.2	ug/L	72148	3	13.554	560313	25.0
Benzene, 1,2-di...	13.371	1.6	ug/L	36470	3	13.554	560313	25.0
Benzene, 1-meth...	13.944	2.2	ug/L	50228	3	13.554	560313	25.0
Benzene, 1-meth...	14.200	2.5	ug/L	56506	3	13.554	560313	25.0