

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
 Data File : VW020436.D
 Acq On : 11 Oct 2021 11:47
 Operator : SY/VA
 Sample : M4070-15
 Misc : 6.19g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EW7D8

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\SFAMWLM100821SMA.M
 Title : SFAM01.0

Signal : TIC: VW020436.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.349	67	73	83	rVB	82531	133140	14.04%	1.369%
2	2.495	93	97	103	rBV	3320	6537	0.69%	0.067%
3	2.818	146	150	154	rVB2	1301	2177	0.23%	0.022%
4	2.885	154	161	171	rBV	34827	74041	7.81%	0.761%
5	4.019	336	347	358	rBV	112098	321179	33.88%	3.302%
6	4.129	358	365	377	rVB	16850	44349	4.68%	0.456%
7	4.385	399	407	419	rVV3	2081	6817	0.72%	0.070%
8	4.915	483	494	508	rBV2	20974	69021	7.28%	0.710%
9	5.787	631	637	639	rBV3	963	1576	0.17%	0.016%
10	5.952	655	664	668	rBV3	872	2024	0.21%	0.021%
11	7.080	841	849	858	rVV	21316	59372	6.26%	0.610%
12	7.177	860	865	878	rVV2	5584	16603	1.75%	0.171%
13	7.653	933	943	957	rBV	163833	395833	41.75%	4.070%
14	7.945	986	991	998	rVB6	742	1530	0.16%	0.016%
15	8.281	1035	1046	1060	rBV2	330986	948060	100.00%	9.747%
16	8.707	1110	1116	1121	rBV3	442	902	0.10%	0.009%
17	8.848	1130	1139	1149	rBV	312400	606563	63.98%	6.236%
18	9.085	1172	1178	1182	rVB3	1161	2187	0.23%	0.022%
19	9.274	1201	1209	1218	rBV	223953	460056	48.53%	4.730%
20	9.512	1245	1248	1253	rVB2	348	549	0.06%	0.006%
21	9.610	1258	1264	1269	rVB3	469	875	0.09%	0.009%
22	9.665	1269	1273	1281	rVB3	369	848	0.09%	0.009%
23	9.750	1281	1287	1292	rBV4	786	1293	0.14%	0.013%
24	9.854	1299	1304	1307	rVB2	421	560	0.06%	0.006%
25	10.036	1326	1334	1344	rBV	110380	196908	20.77%	2.024%
26	10.122	1347	1348	1355	rVB3	445	599	0.06%	0.006%
27	10.329	1372	1382	1388	rBV	447595	800040	84.39%	8.225%
28	10.390	1388	1392	1398	rVB	8844	15256	1.61%	0.157%
29	10.506	1404	1411	1416	rBV8	1720	4275	0.45%	0.044%
30	10.579	1416	1423	1432	rVV	71046	119872	12.64%	1.232%
31	10.664	1432	1437	1440	rVV3	2743	4371	0.46%	0.045%
32	10.707	1440	1444	1447	rVV2	3045	4826	0.51%	0.050%
33	10.750	1447	1451	1458	rVB5	2406	5256	0.55%	0.054%
34	10.866	1463	1470	1474	rVV	19471	34934	3.68%	0.359%
35	10.927	1474	1480	1494	rVV	84676	142041	14.98%	1.460%
36	11.067	1498	1503	1508	rVV2	3675	6632	0.70%	0.068%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
 Title : SFAM01.0

37	11.146	1512	1516	1518	rVV3	1742	2788	0.29%	0.029%
38	11.176	1518	1521	1526	rVV3	3282	5693	0.60%	0.059%
39	11.231	1526	1530	1535	rVV3	1772	3069	0.32%	0.032%
40	11.280	1535	1538	1543	rVV5	440	833	0.09%	0.009%
41	11.341	1543	1548	1553	rVB6	520	875	0.09%	0.009%
42	11.634	1588	1596	1604	rBV	429212	710614	74.95%	7.306%
43	11.725	1607	1611	1617	rVB3	4192	7120	0.75%	0.073%
44	11.847	1625	1631	1641	rBV	4815	11495	1.21%	0.118%
45	12.036	1656	1662	1663	rBV2	603	631	0.07%	0.006%
46	12.164	1674	1683	1692	rVB6	2874	6907	0.73%	0.071%
47	12.341	1707	1712	1717	rVV4	4507	7253	0.77%	0.075%
48	12.414	1722	1724	1727	rVB3	639	709	0.07%	0.007%
49	12.463	1727	1732	1738	rVB2	5654	8950	0.94%	0.092%
50	12.518	1738	1741	1745	rBV3	435	621	0.07%	0.006%
51	12.603	1749	1755	1763	rVB3	3750	7215	0.76%	0.074%
52	12.695	1763	1770	1777	rBV	134747	221391	23.35%	2.276%
53	12.786	1777	1785	1786	rBV7	2132	4276	0.45%	0.044%
54	12.865	1793	1798	1801	rBV6	781	1351	0.14%	0.014%
55	12.896	1801	1803	1806	rVV3	1440	2347	0.25%	0.024%
56	12.932	1806	1809	1817	rVB7	2925	5698	0.60%	0.059%
57	13.103	1829	1837	1839	rBV	4876	7736	0.82%	0.080%
58	13.133	1839	1842	1846	rVB2	5189	7336	0.77%	0.075%
59	13.249	1857	1861	1866	rBV2	5679	8667	0.91%	0.089%
60	13.383	1874	1883	1889	rBV4	7803	17671	1.86%	0.182%
61	13.499	1897	1902	1905	rBV	4571	6394	0.67%	0.066%
62	13.560	1905	1912	1922	rVV	396117	638453	67.34%	6.564%
63	13.652	1923	1927	1933	rVB2	7282	11498	1.21%	0.118%
64	13.719	1933	1938	1940	rBV	13469	20614	2.17%	0.212%
65	13.755	1940	1944	1949	rVV	46481	74107	7.82%	0.762%
66	13.798	1949	1951	1954	rVV	6592	8930	0.94%	0.092%
67	13.853	1954	1960	1971	rVB	341188	593024	62.55%	6.097%
68	13.950	1971	1976	1980	rVB	6246	9641	1.02%	0.099%
69	14.011	1980	1986	1988	rBV	29680	47235	4.98%	0.486%
70	14.042	1988	1991	1995	rVV	25655	39113	4.13%	0.402%
71	14.097	1995	2000	2005	rVB	84561	128028	13.50%	1.316%
72	14.158	2005	2010	2013	rBV	22020	35926	3.79%	0.369%
73	14.206	2013	2018	2024	rVB2	117684	189857	20.03%	1.952%
74	14.341	2033	2040	2044	rBV3	28987	50820	5.36%	0.522%
75	14.420	2044	2053	2055	rBV	96120	154611	16.31%	1.590%
76	14.456	2055	2059	2064	rVB	152419	219074	23.11%	2.252%

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 Title : SFAM01.0

77	14.499	2064	2066	2069	rVB	9091	8387	0.88%	0.086%
78	14.536	2069	2072	2075	rBV	6247	8201	0.87%	0.084%
79	14.566	2075	2077	2081	rVB	11541	13963	1.47%	0.144%
80	14.639	2082	2089	2092	rBV2	14056	25027	2.64%	0.257%
81	14.688	2092	2097	2101	rBV3	33237	61150	6.45%	0.629%
82	14.725	2101	2103	2108	rVB	23327	33622	3.55%	0.346%
83	14.804	2108	2116	2122	rBV2	171684	328272	34.63%	3.375%
84	14.859	2122	2125	2132	rVB4	16206	34597	3.65%	0.356%
85	14.956	2136	2141	2148	rBV2	23740	43381	4.58%	0.446%
86	15.060	2149	2158	2162	rBV2	79966	178252	18.80%	1.833%
87	15.176	2170	2177	2186	rVB2	131003	297178	31.35%	3.055%
88	15.267	2188	2192	2196	rBV2	6029	9210	0.97%	0.095%
89	15.322	2196	2201	2205	rBV3	9725	17539	1.85%	0.180%
90	15.432	2211	2219	2220	rBV3	13655	27151	2.86%	0.279%
91	15.468	2221	2225	2229	rVV3	18174	33358	3.52%	0.343%
92	15.651	2248	2255	2261	rBV	65125	125796	13.27%	1.293%
93	15.731	2264	2268	2272	rVV4	7570	13944	1.47%	0.143%
94	15.822	2275	2283	2292	rVV2	55904	145326	15.33%	1.494%
95	15.944	2298	2303	2308	rBV	37536	64394	6.79%	0.662%
96	16.023	2308	2316	2322	rVV3	40671	104087	10.98%	1.070%
97	16.176	2332	2341	2353	rBV3	46380	124500	13.13%	1.280%
98	16.377	2363	2374	2378	rBV4	19912	55506	5.85%	0.571%
99	16.432	2379	2383	2387	rBV2	8894	15498	1.63%	0.159%
100	16.639	2408	2417	2429	rBV3	71483	188795	19.91%	1.941%

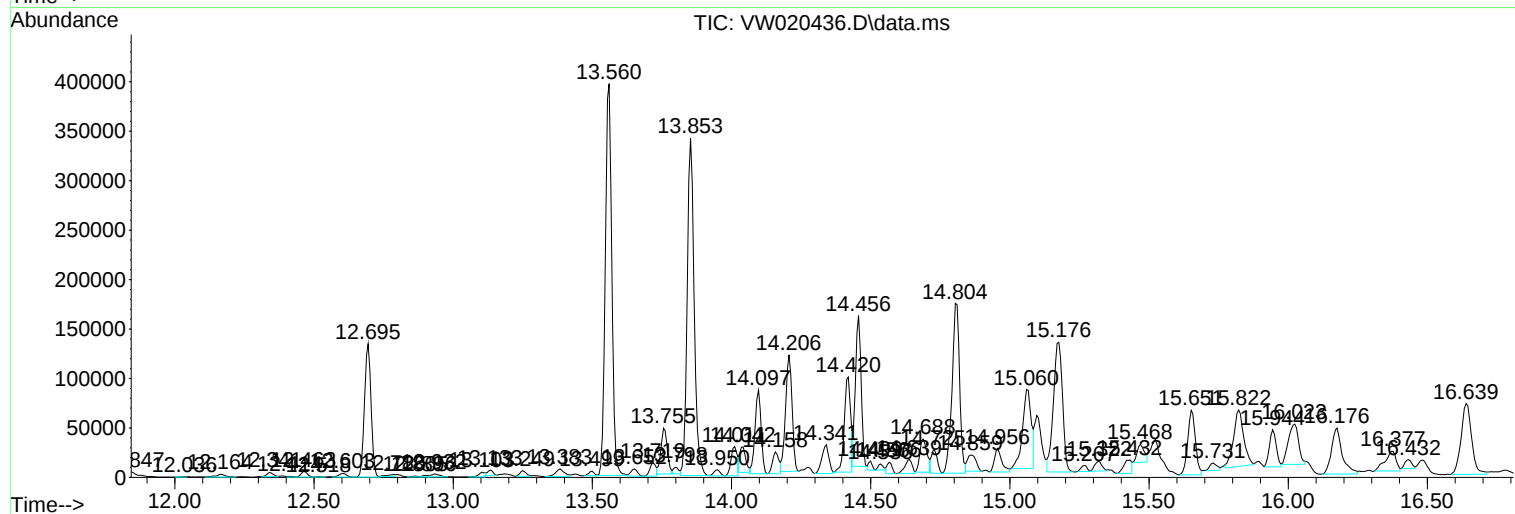
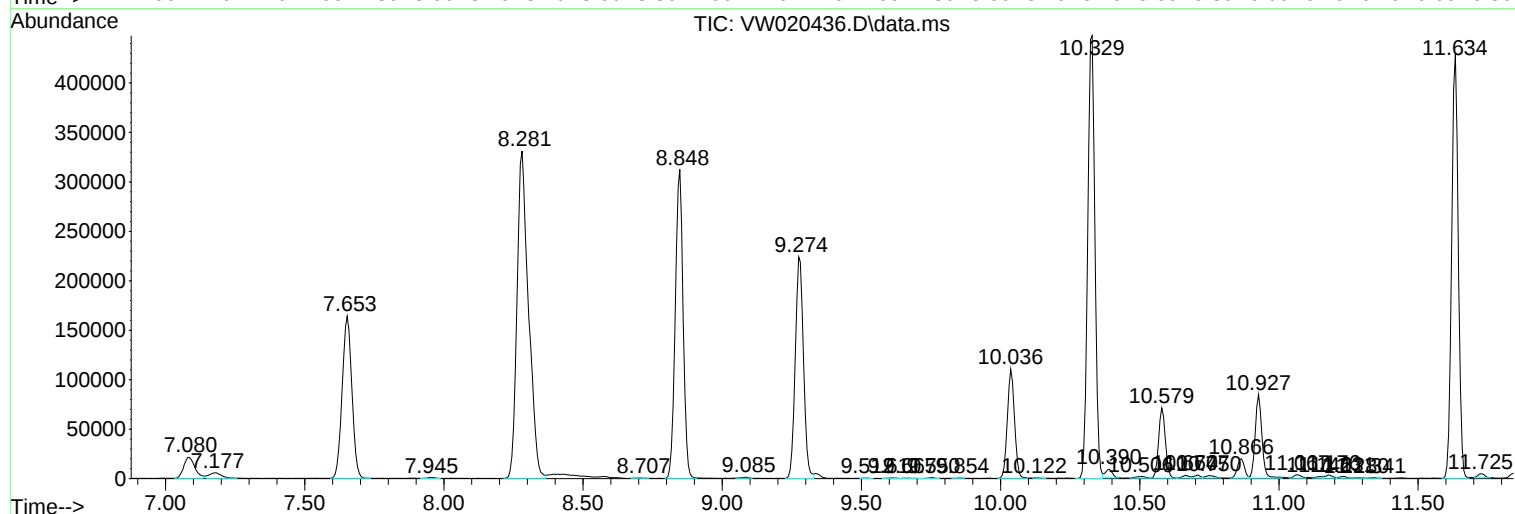
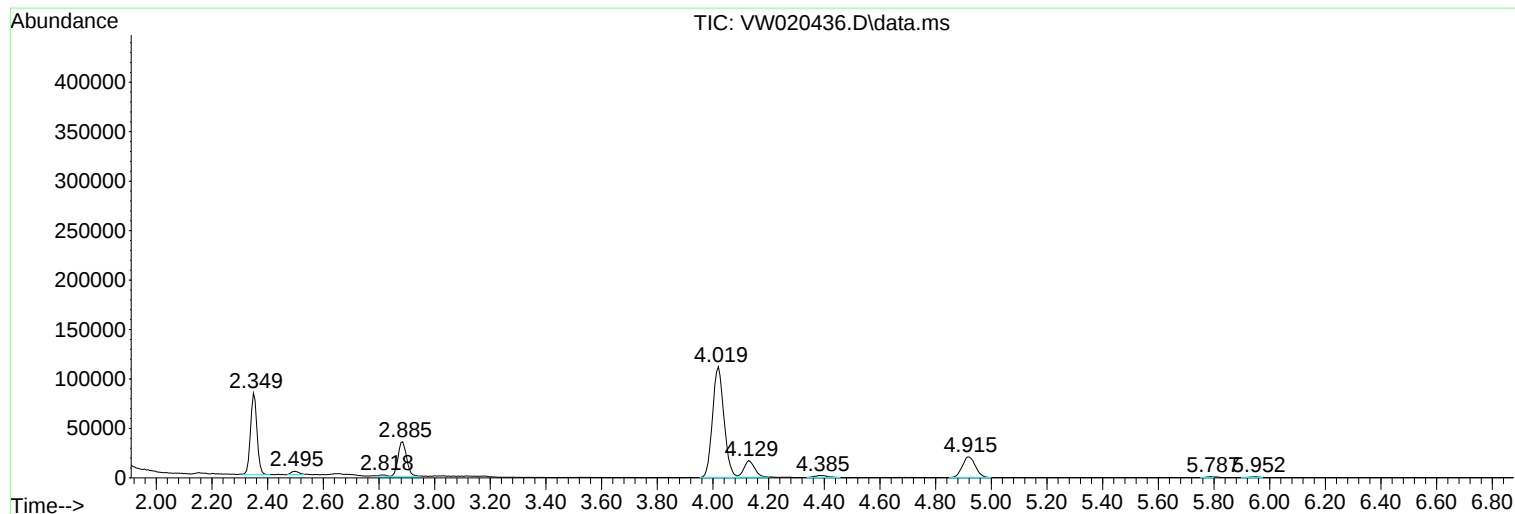
Sum of corrected areas: 9726807

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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



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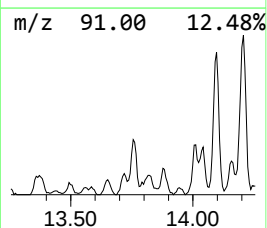
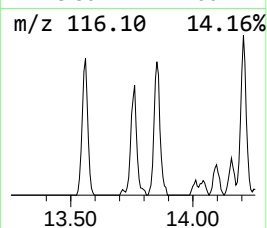
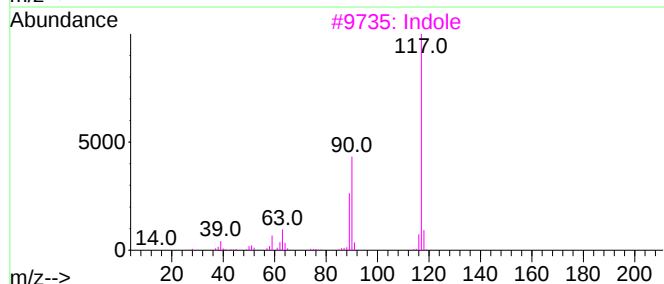
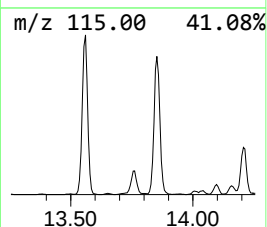
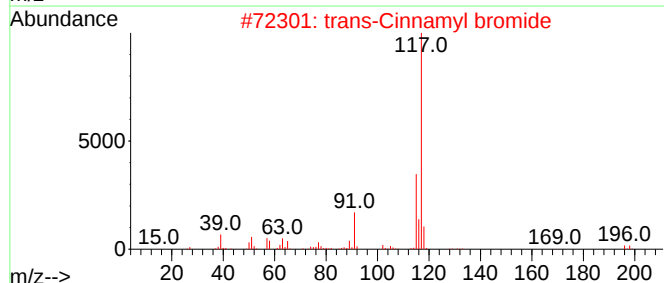
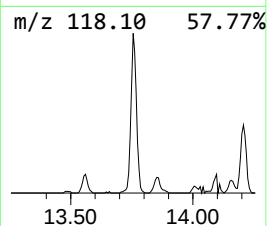
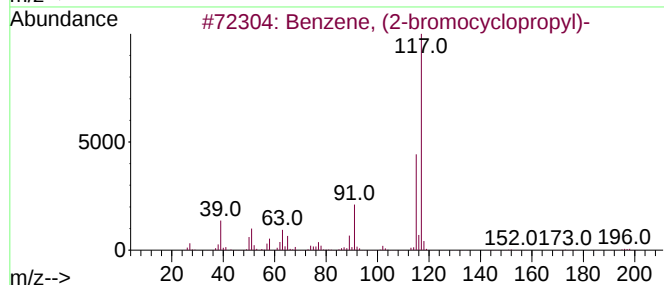
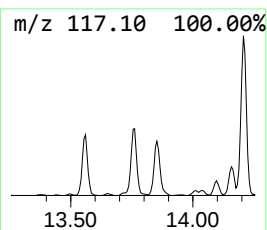
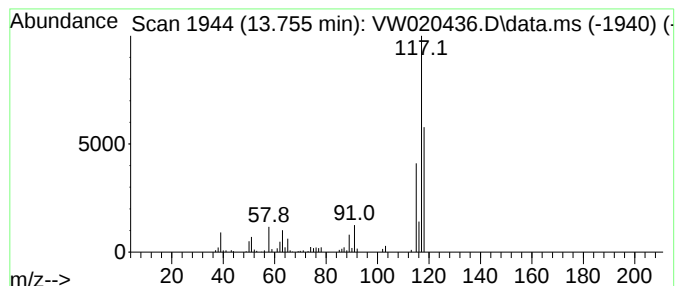
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Benzene, (2-bromocyclopropyl)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.755	2.90 ug/L	74107	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	59
2			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	50
3			Indole	117	C8H7N	000120-72-9	43
4			Benzene, (isocyanomethyl)-	117	C8H7N	010340-91-7	43
5			3-Bromo-1-phenyl-1-propene	196	C9H9Br	004392-24-9	42



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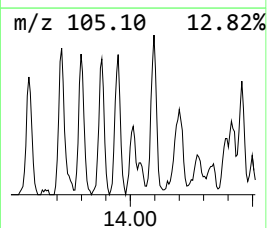
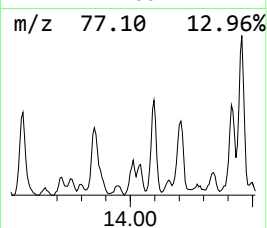
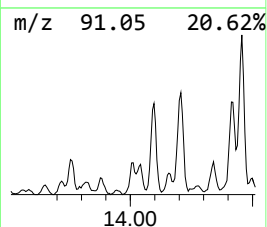
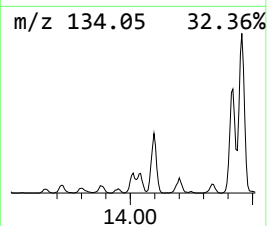
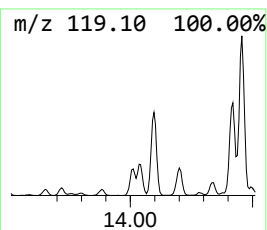
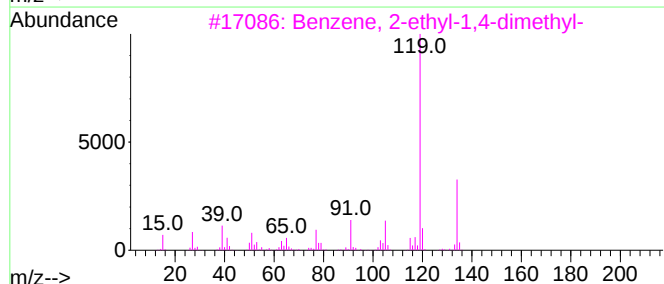
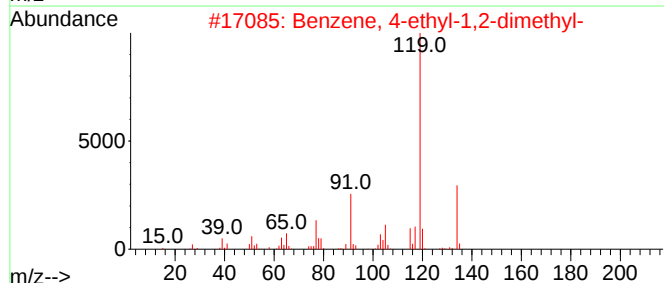
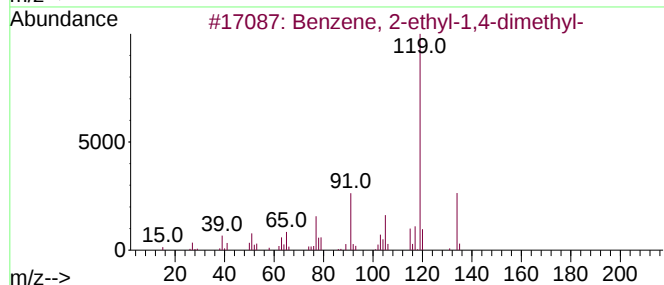
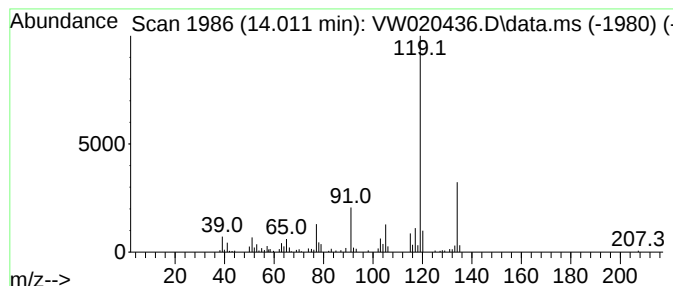
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.011	1.85 ug/L	47235	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



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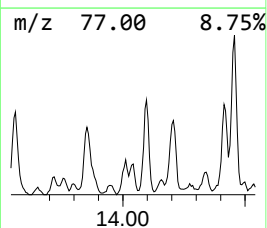
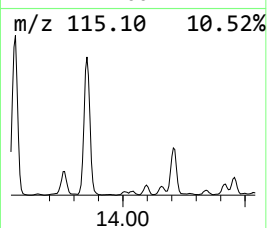
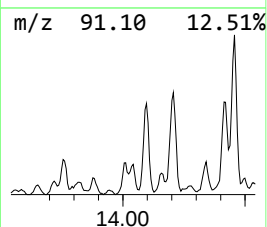
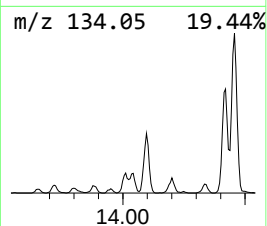
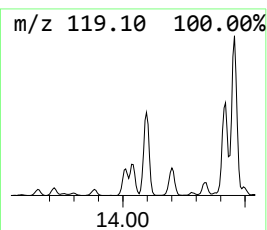
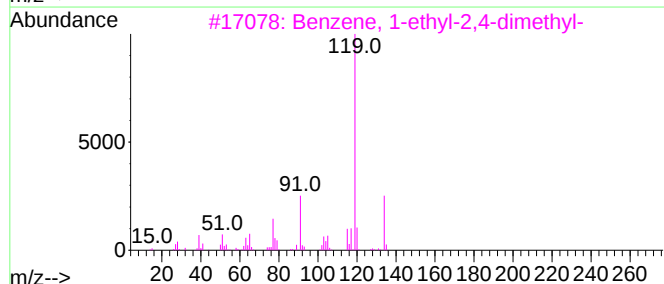
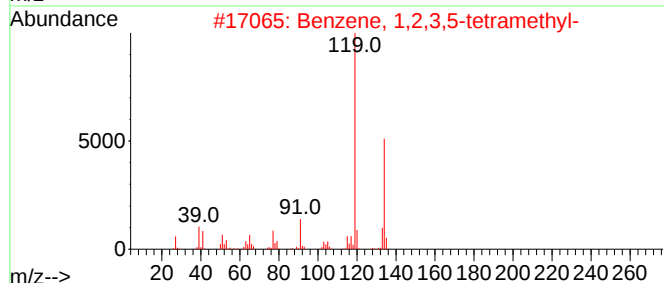
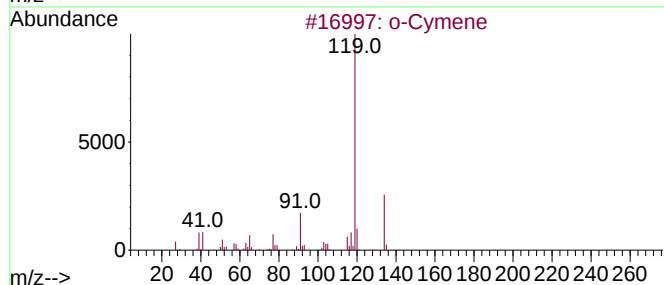
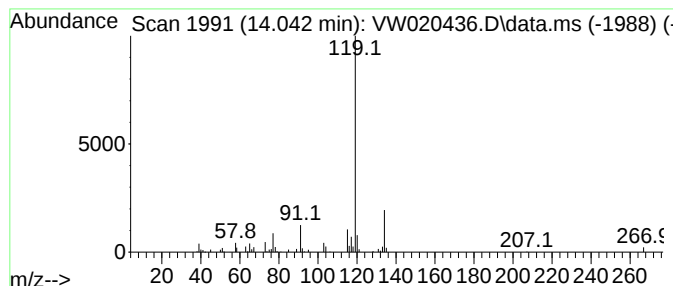
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Quant Title : SFAM01.0

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

Peak Number 5 o-Cymene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.042	1.53 ug/L	39113	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	93
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

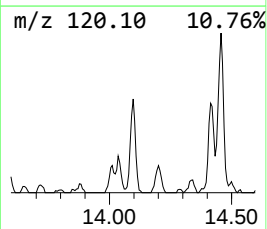
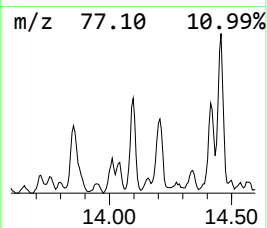
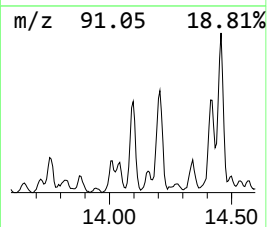
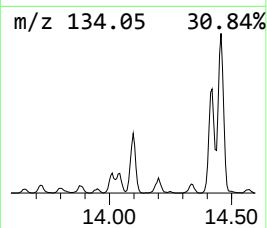
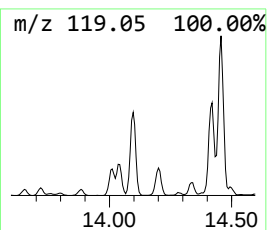
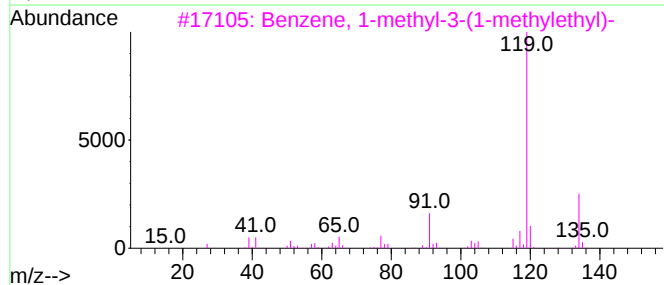
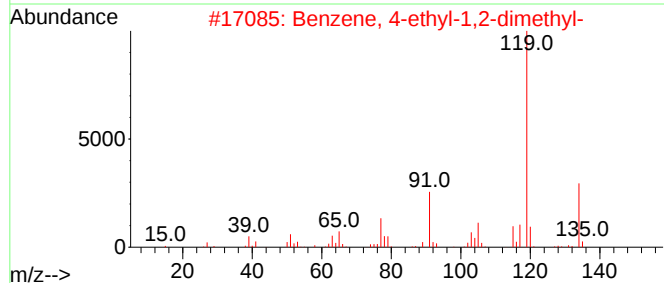
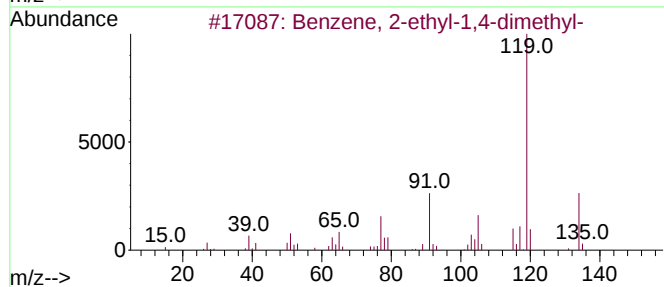
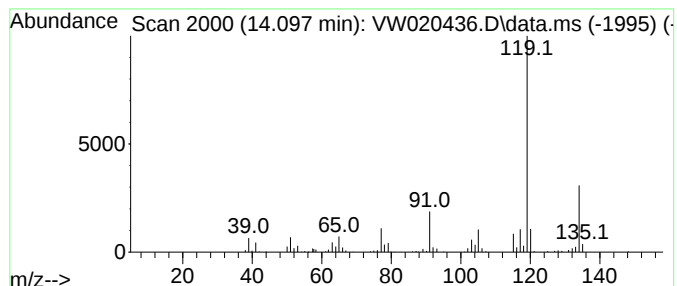
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Quant Title : SFAM01.0

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.097	5.01 ug/L	128028	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

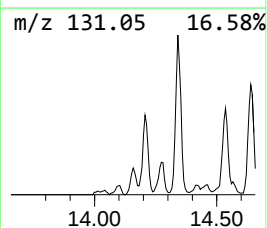
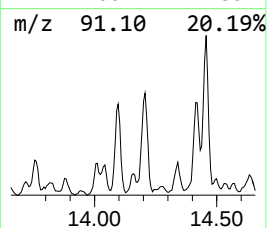
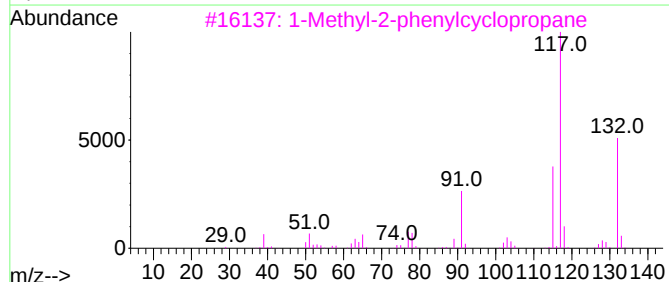
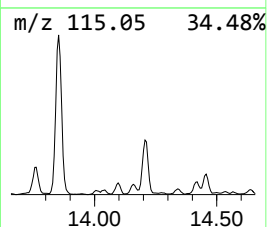
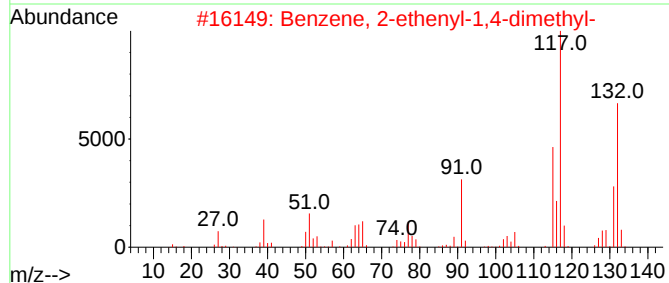
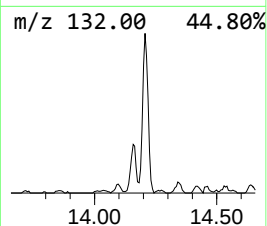
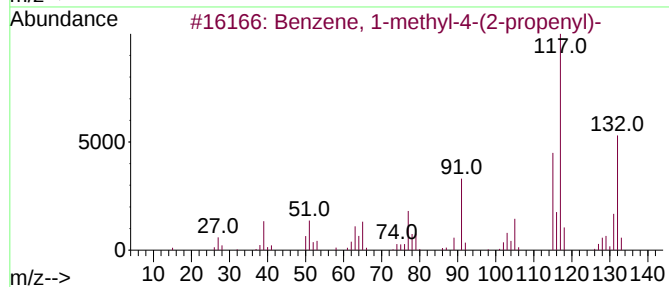
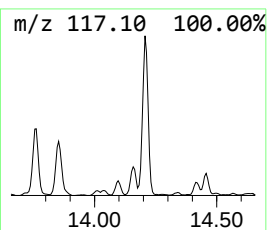
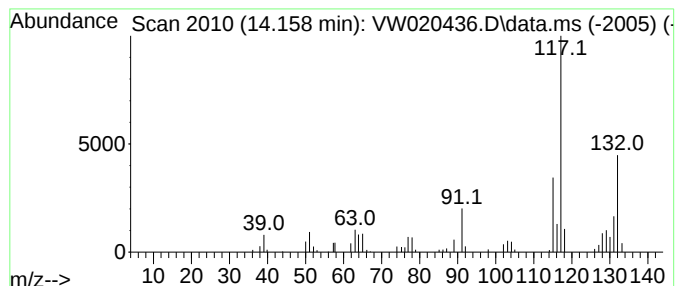
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Quant Title : SFAM01.0

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

Peak Number 7 Benzene, 1-methyl-4-(2-prop... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.158	1.41 ug/L	35926	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	96
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	93
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

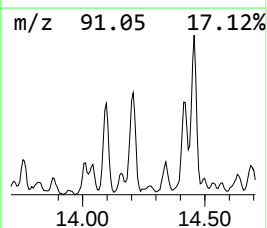
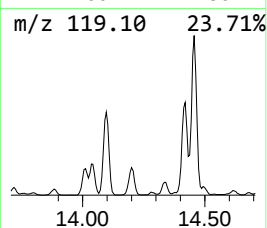
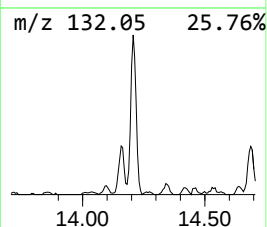
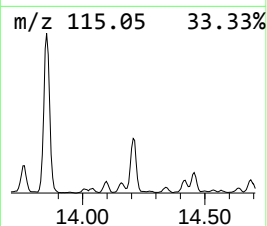
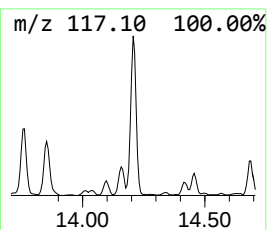
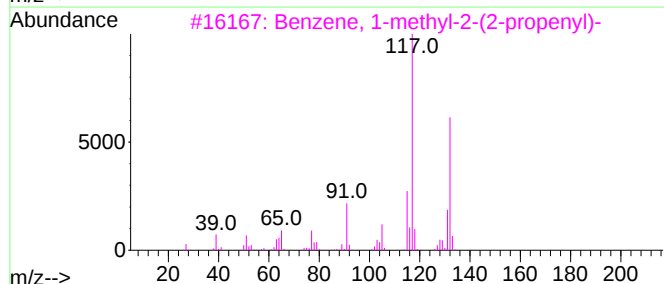
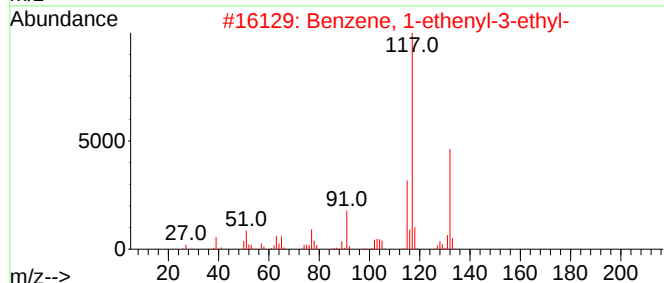
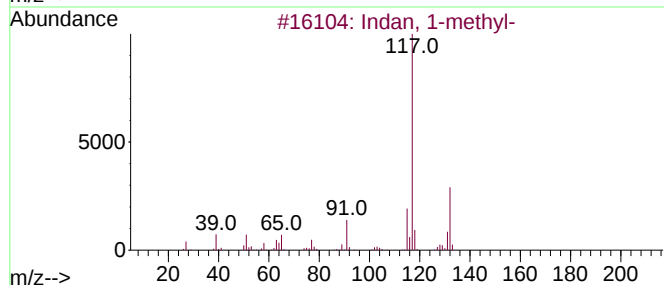
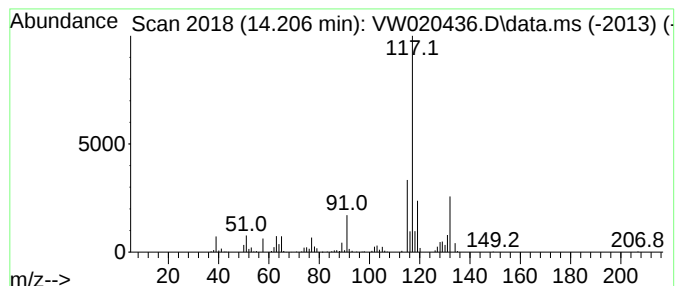
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Quant Title : SFAM01.0

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

Peak Number 8 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.206	7.43 ug/L	189857	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	76
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	72
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

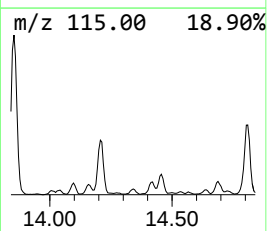
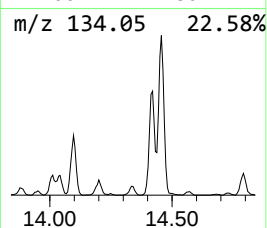
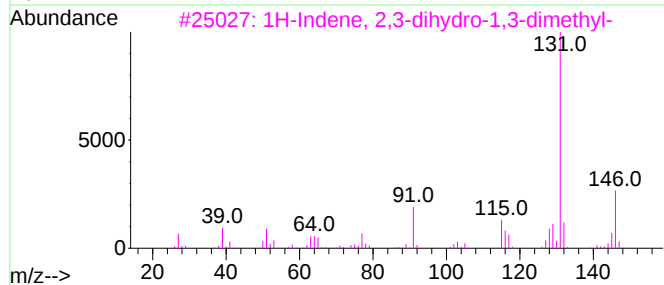
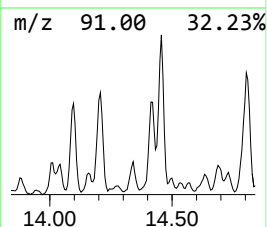
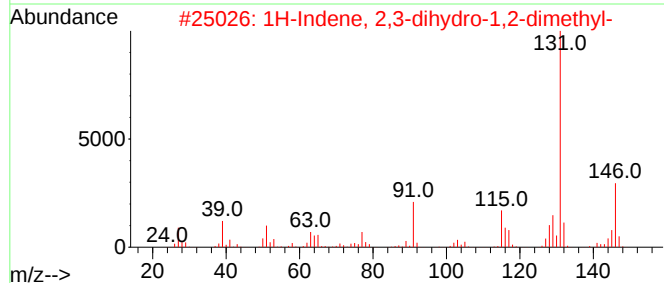
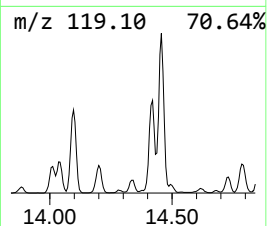
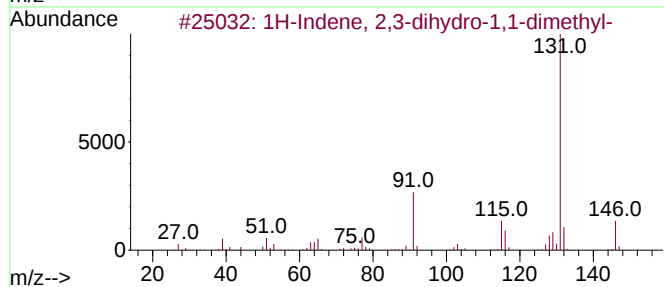
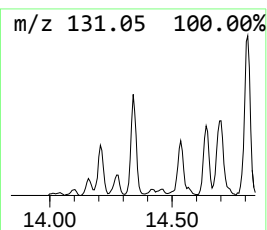
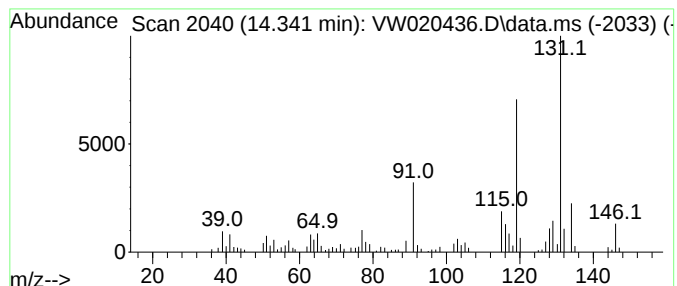
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Quant Title : SFAM01.0

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.341	1.99 ug/L	50820	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	92
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	86
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	86
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	86
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

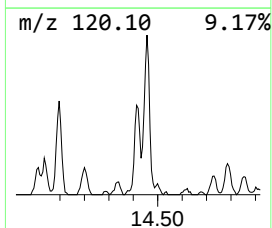
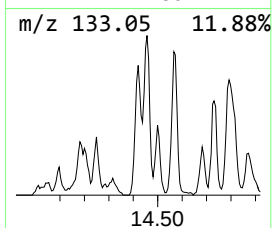
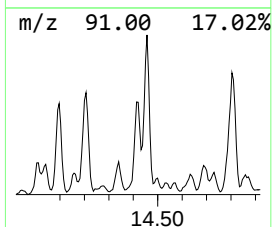
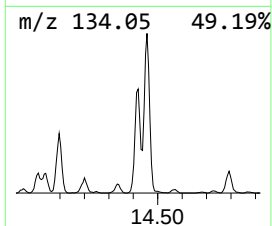
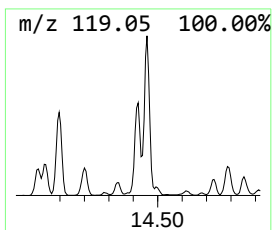
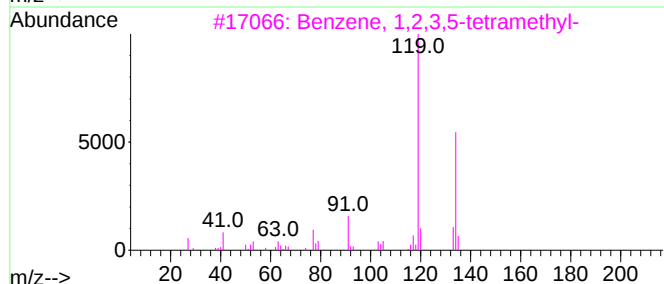
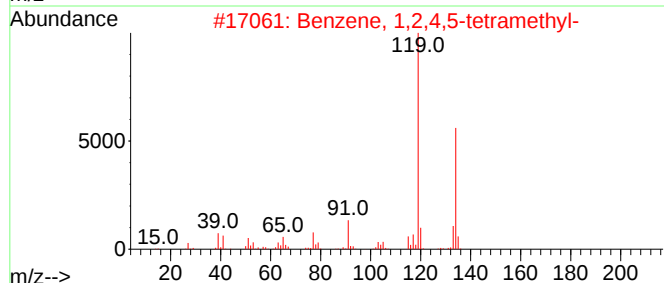
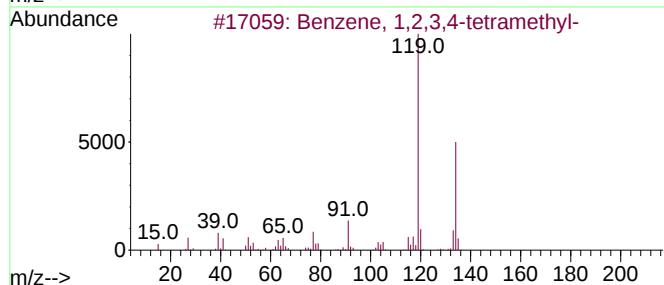
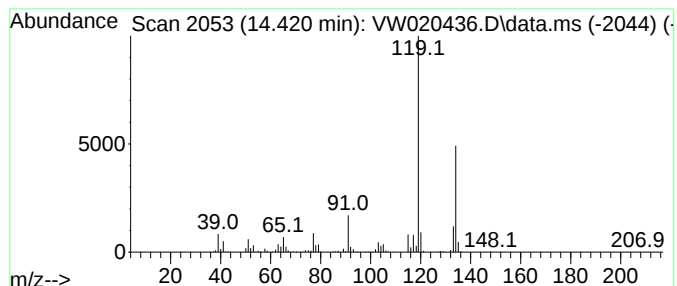
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Quant Title : SFAM01.0

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

Peak Number 10 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.420	6.05 ug/L	154611	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

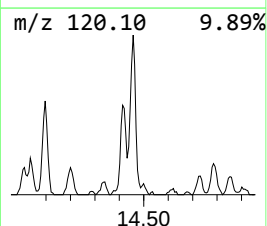
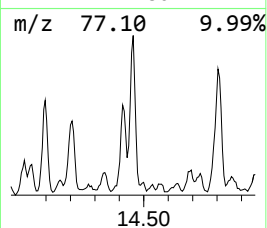
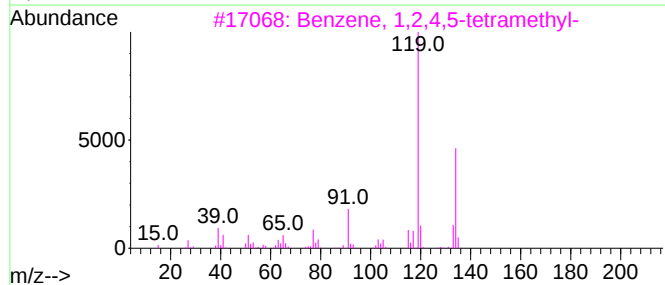
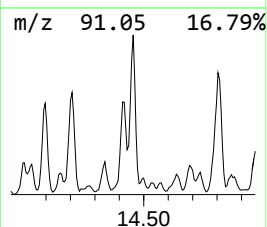
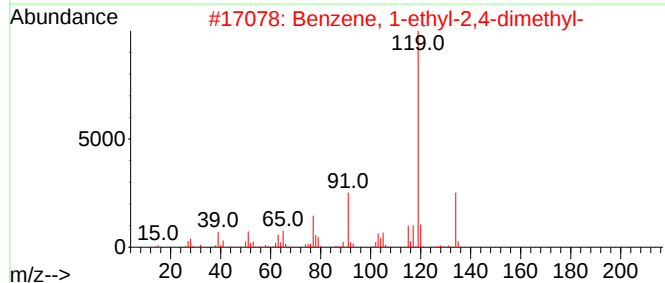
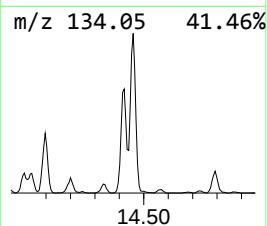
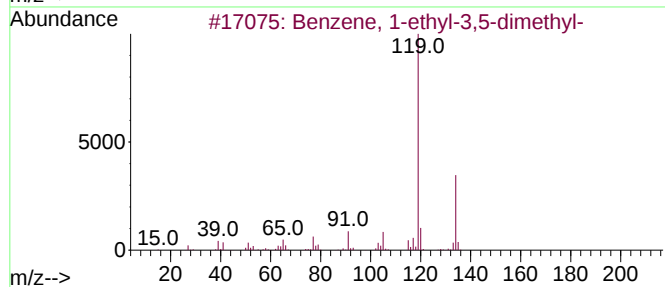
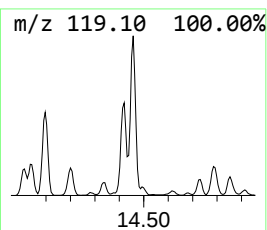
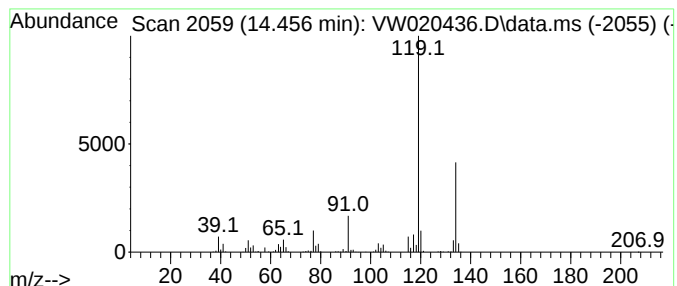
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Quant Title : SFAM01.0

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

Peak Number 11 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.456	8.58 ug/L	219074	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

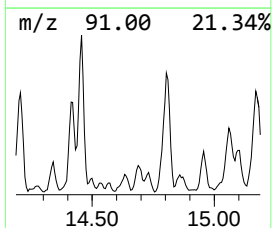
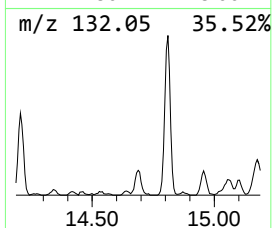
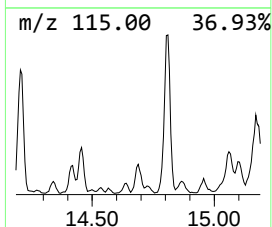
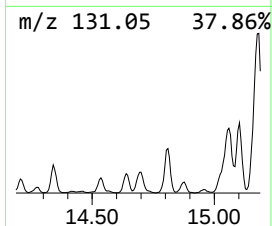
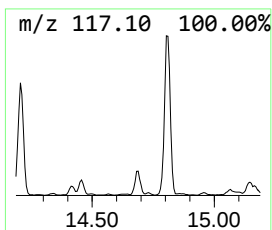
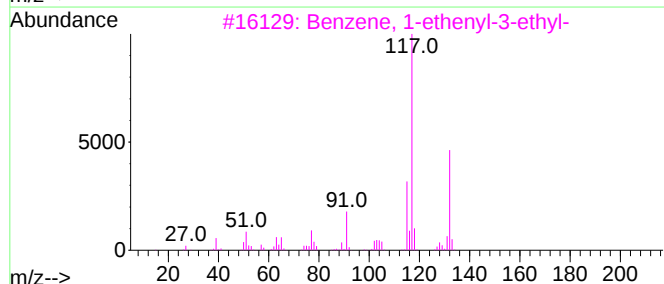
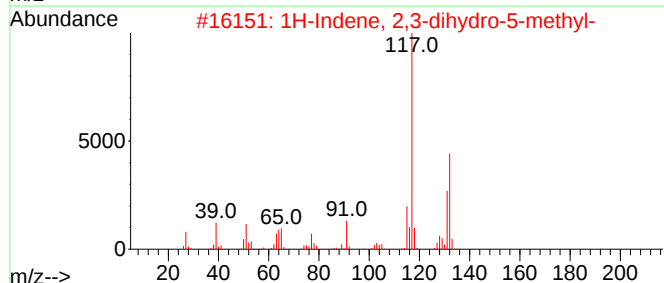
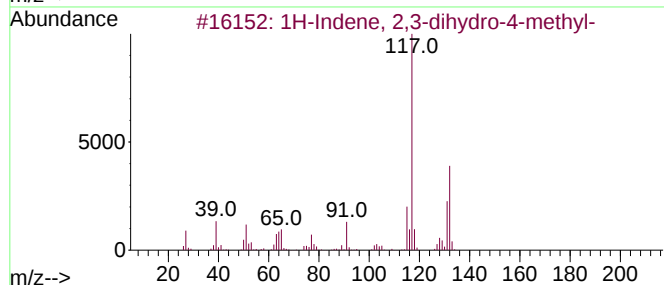
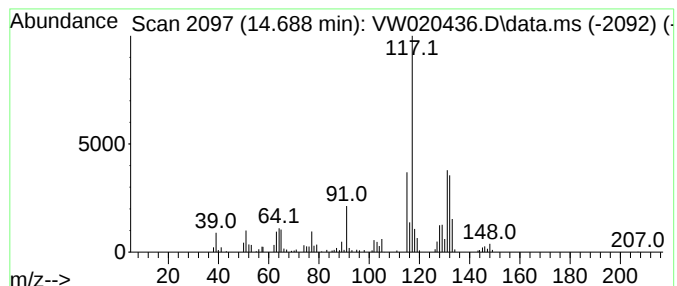
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Quant Title : SFAM01.0

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

Peak Number 12 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.688	2.39 ug/L	61150	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	80
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	80
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	74
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

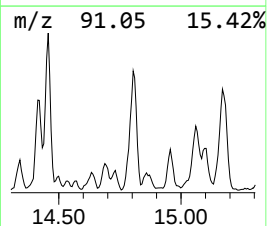
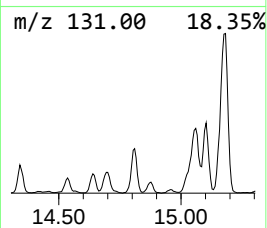
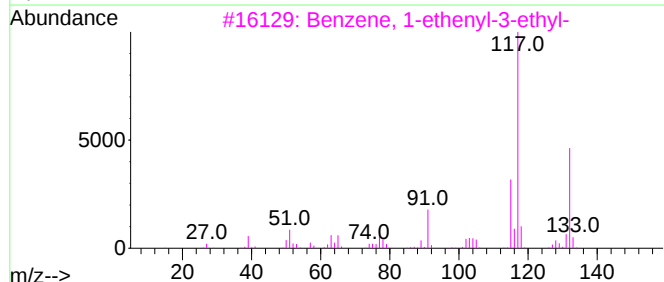
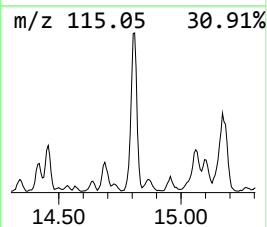
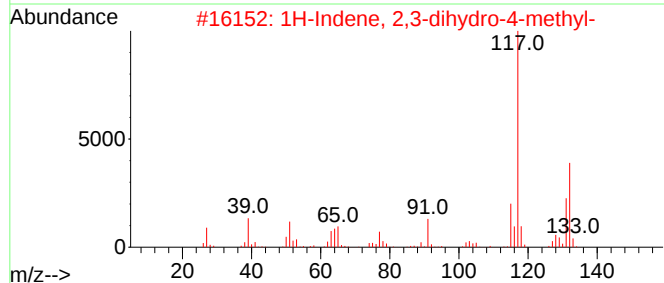
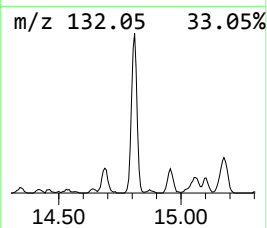
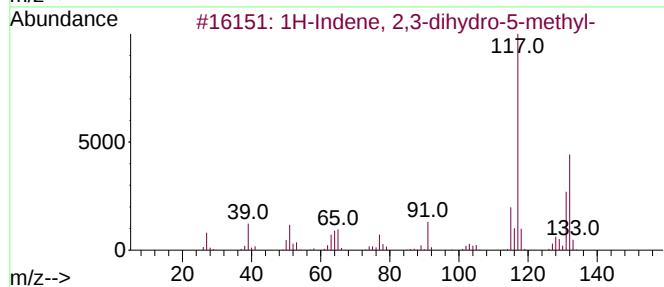
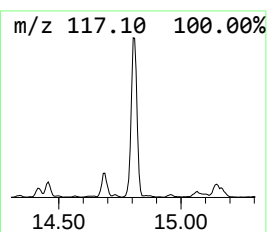
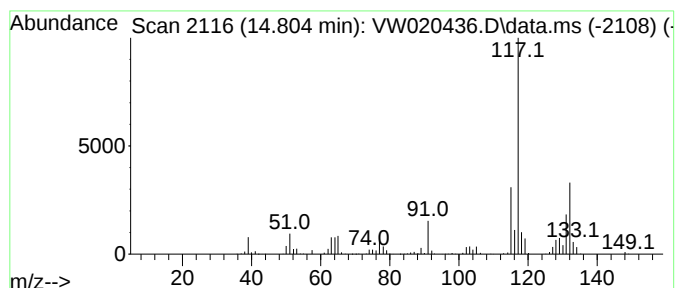
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Quant Title : SFAM01.0

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

Peak Number 14 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.804	12.85 ug/L	328272	1,4-Dichlorobenzene-d4			13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

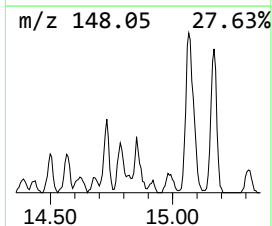
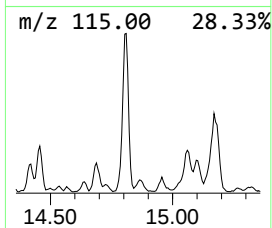
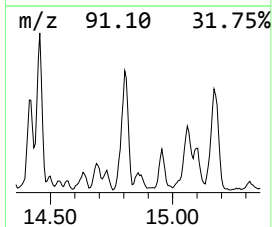
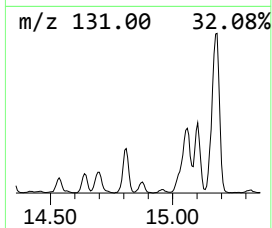
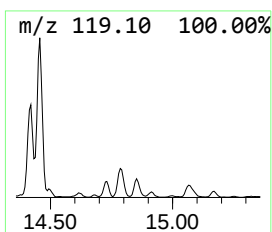
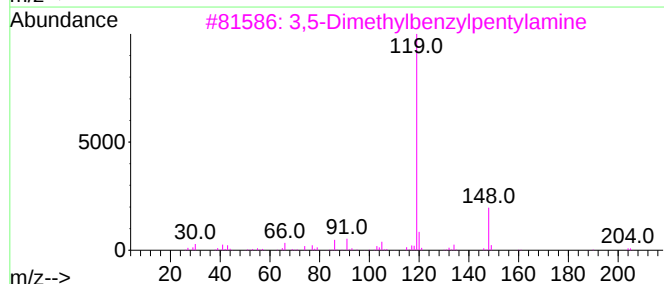
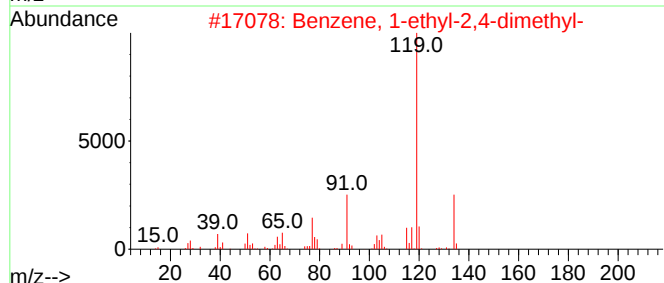
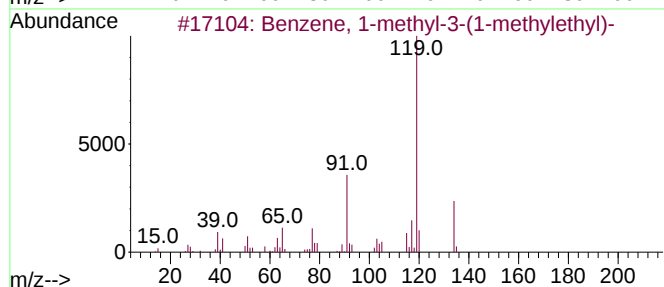
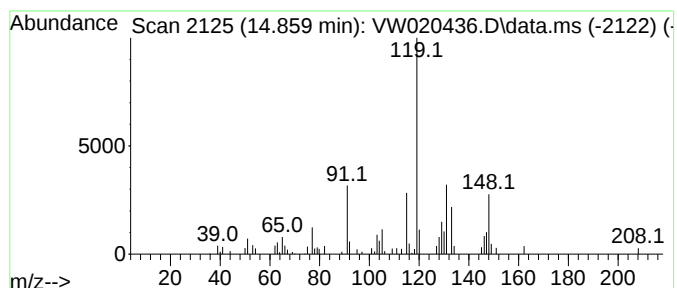
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Quant Title : SFAM01.0

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

Peak Number 15 Benzene, 1-methyl-3-(1-meth... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.859	1.35 ug/L	34597	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	50
3			3,5-Dimethylbenzylpentylamine	205	C14H23N	1000491-60-8	43
4			p-Cymene	134	C10H14	000099-87-6	43
5			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

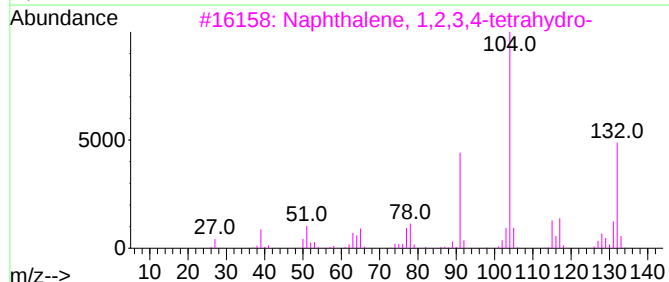
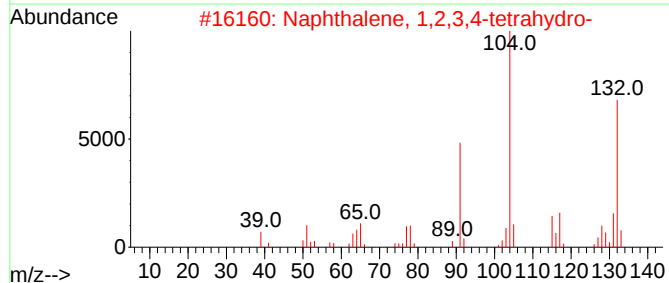
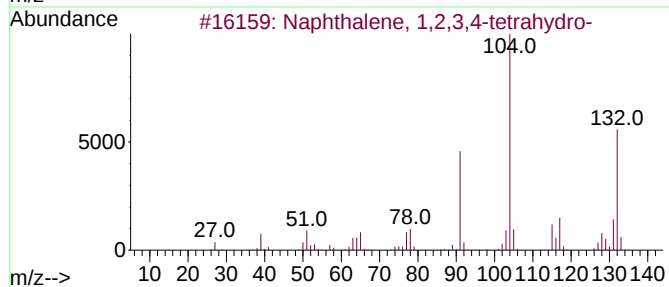
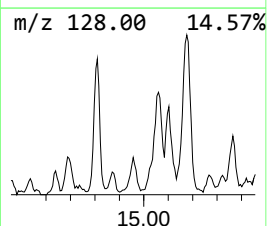
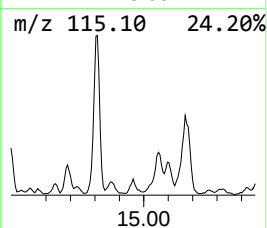
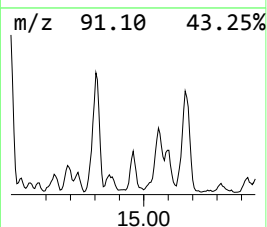
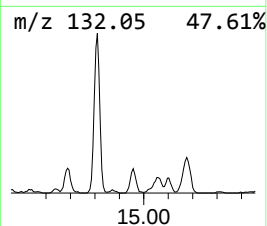
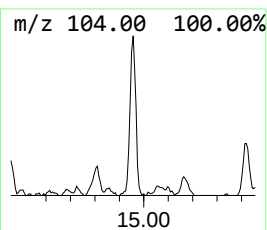
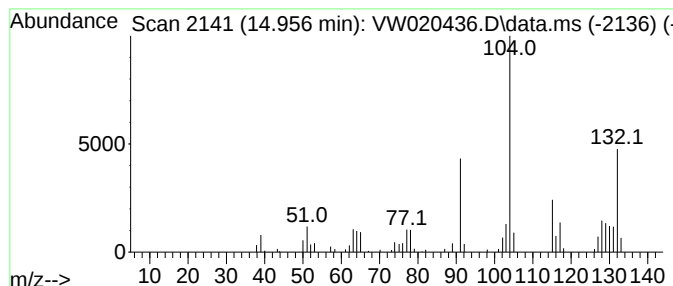
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Quant Title : SFAM01.0

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

Peak Number 16 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.956	1.70 ug/L	43381	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

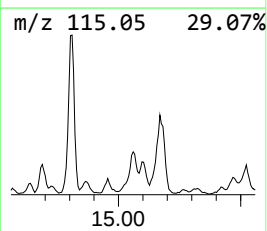
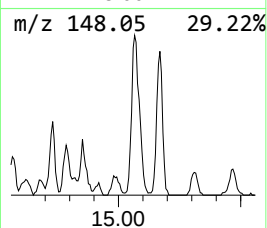
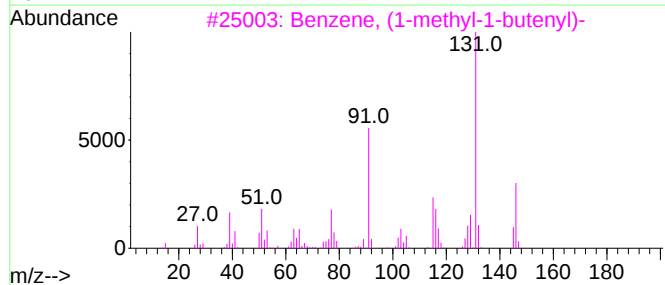
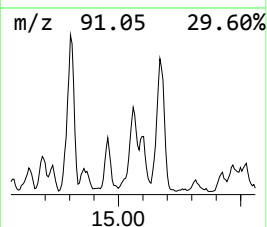
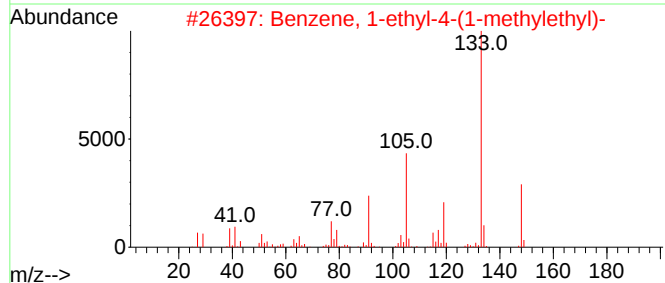
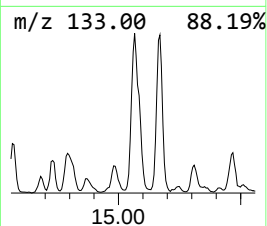
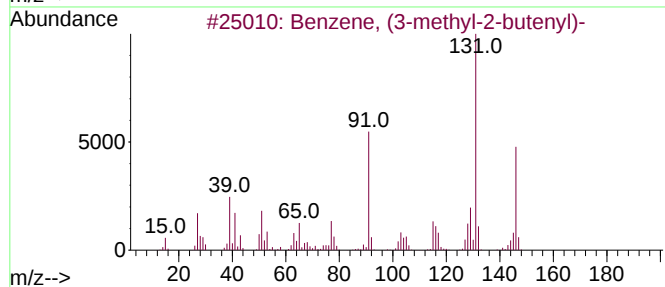
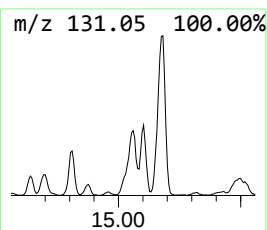
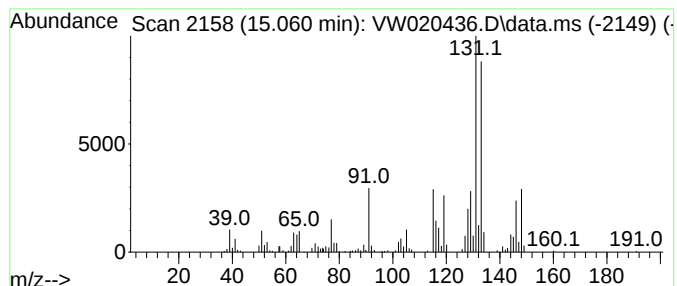
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Quant Title : SFAM01.0

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

Peak Number 17 Benzene, (3-methyl-2-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.060	6.98 ug/L	178252	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	50
2			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	50
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	46
4			Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	46
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

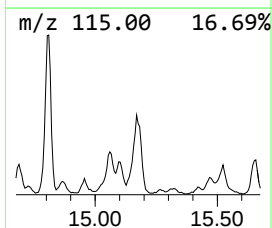
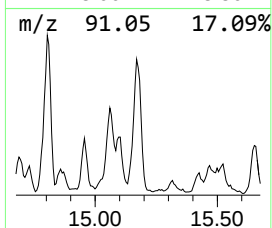
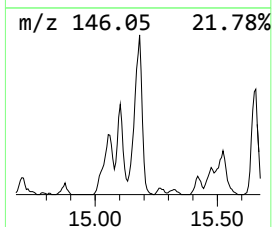
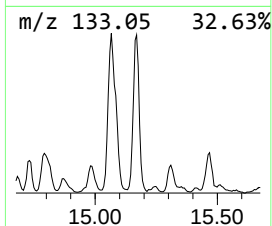
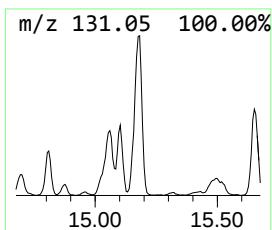
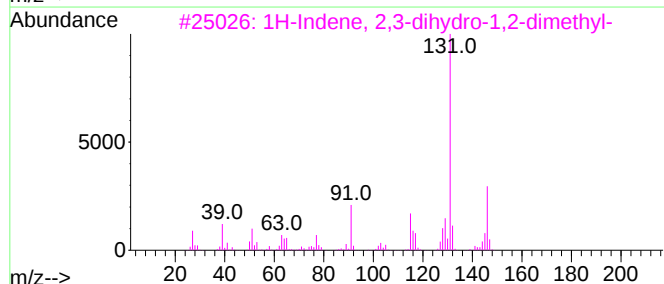
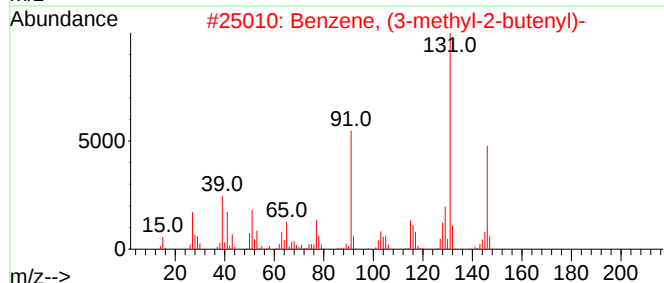
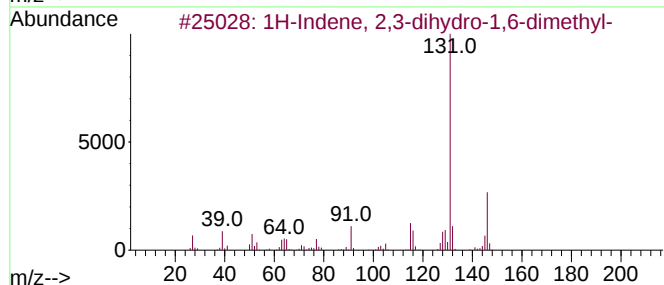
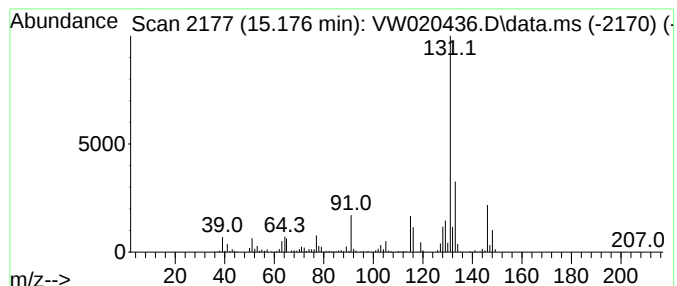
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Quant Title : SFAM01.0

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

Peak Number 18 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.176	11.64 ug/L	297178	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	81
3			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

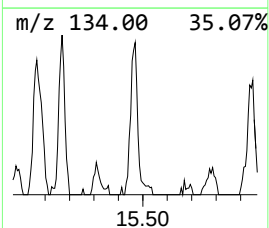
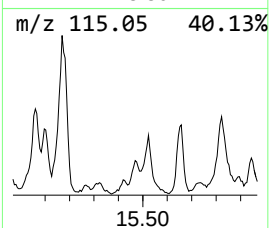
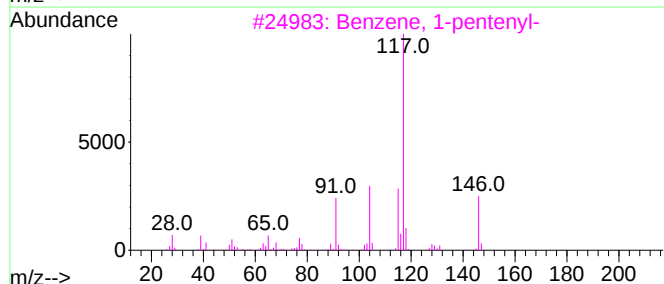
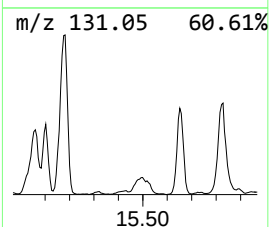
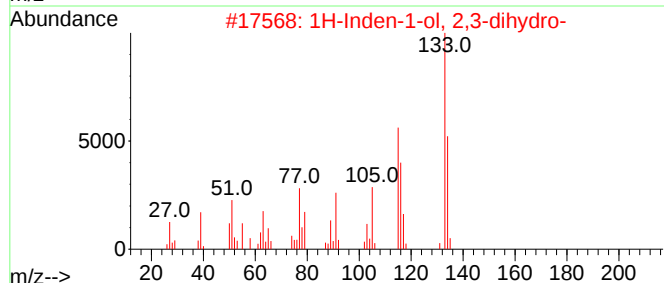
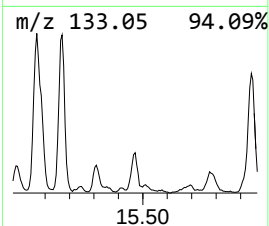
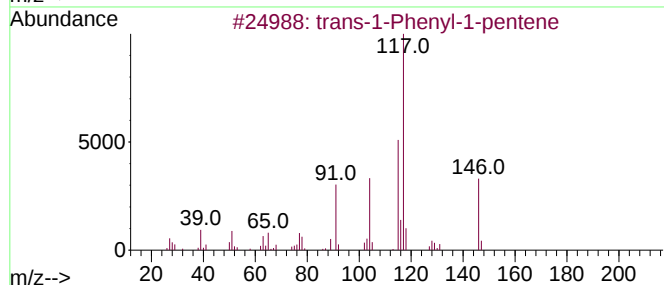
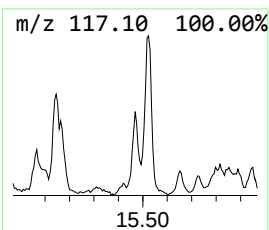
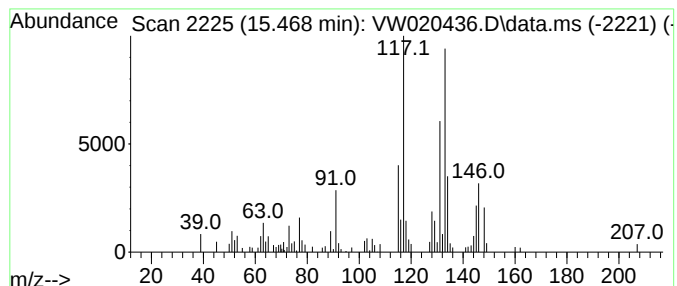
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Quant Title : SFAM01.0

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

Peak Number 19 unknown-01 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.468	1.31 ug/L	33358	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-1-Phenyl-1-pentene	146	C11H14	016002-93-0	42
2			1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	41
3			Benzene, 1-pentenyl-	146	C11H14	000826-18-6	38
4			Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	38
5			3-Buten-1-ol, 4-phenyl-	148	C10H12O	000937-58-6	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

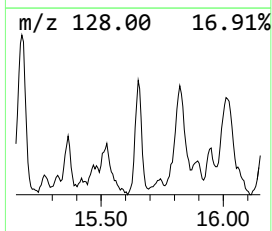
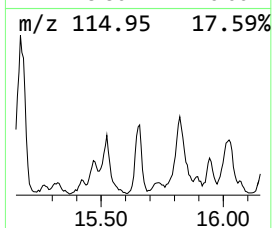
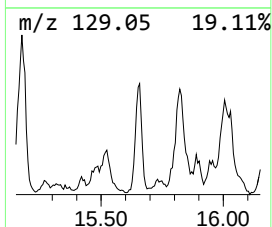
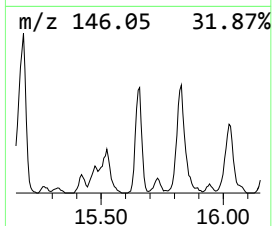
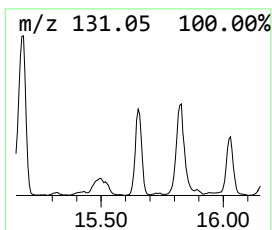
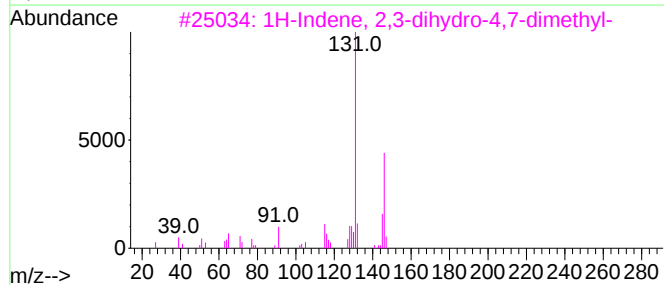
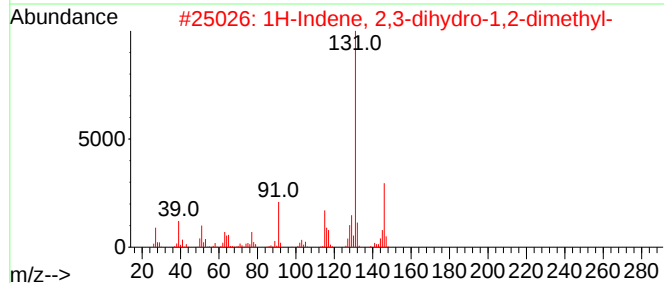
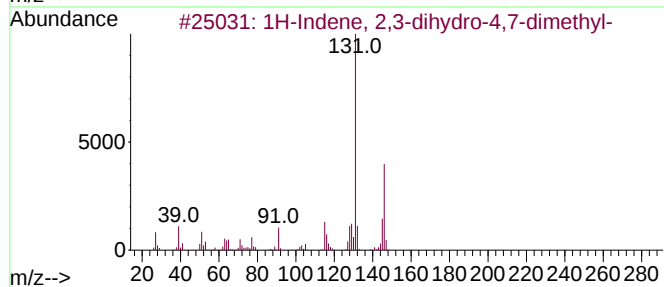
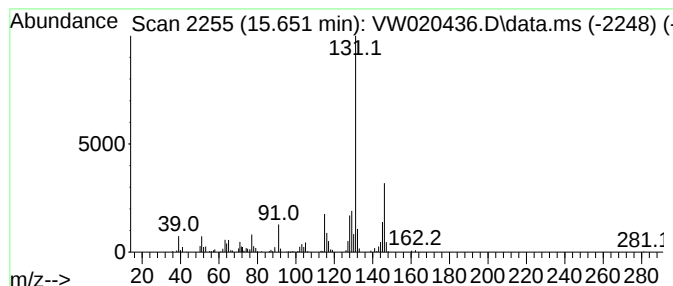
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Quant Title : SFAM01.0

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

Peak Number 20 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	4.93 ug/L	125796	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
4			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	91
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

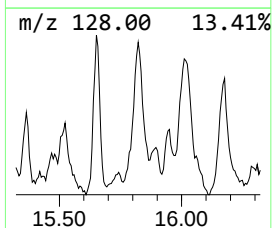
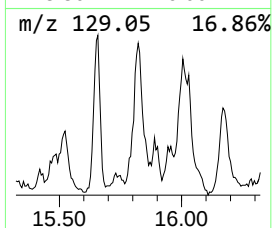
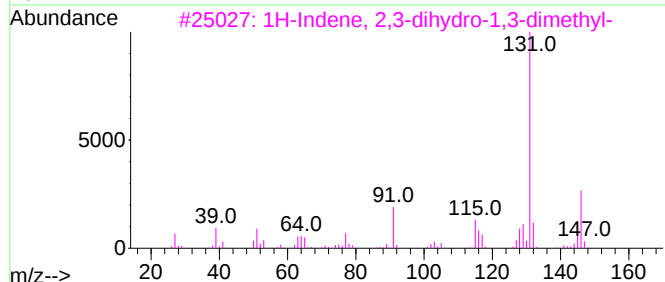
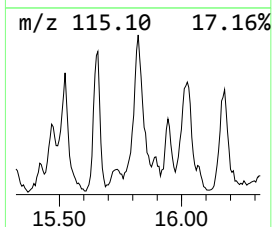
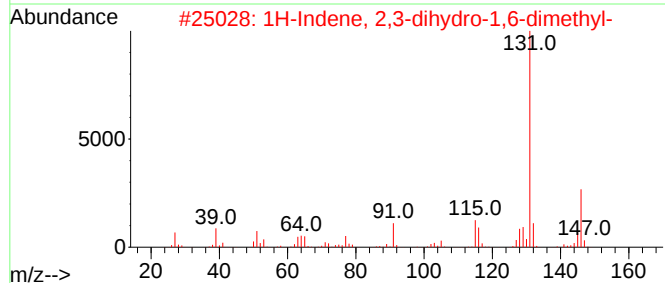
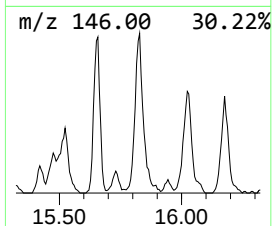
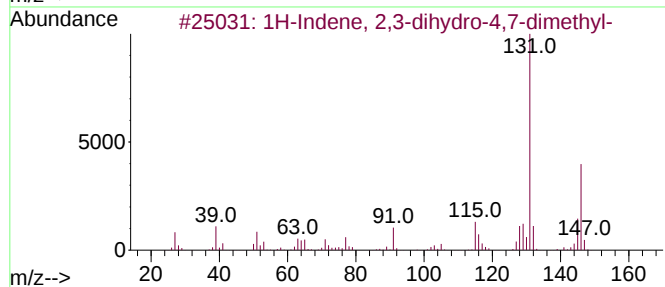
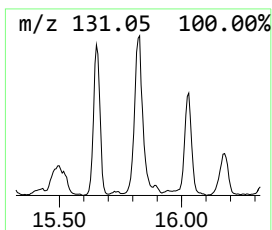
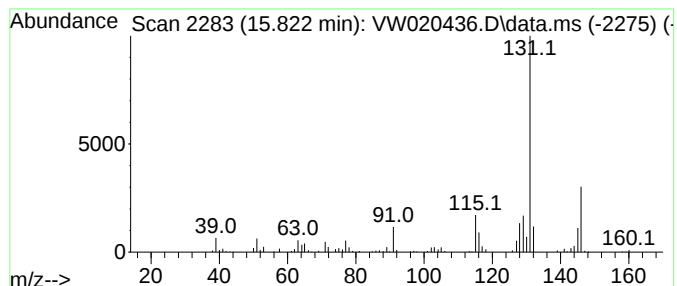
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Quant Title : SFAM01.0

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

Peak Number 21 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.822	5.69 ug/L	145326	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

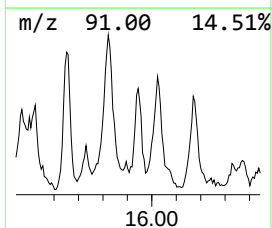
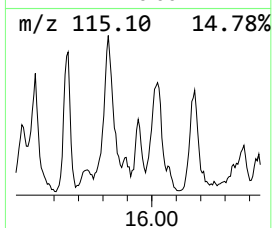
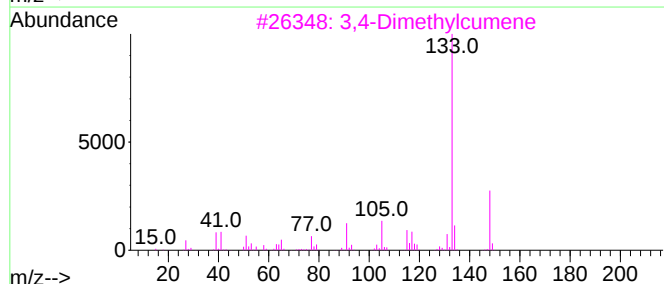
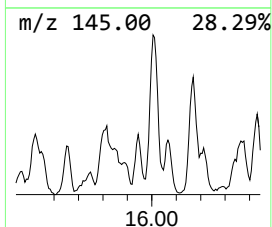
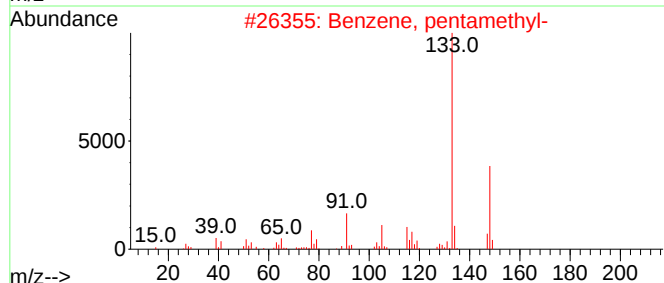
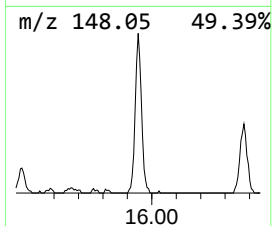
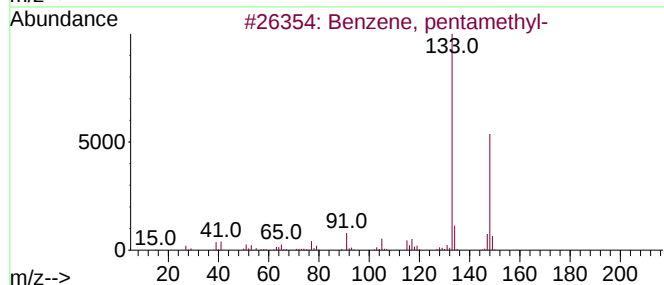
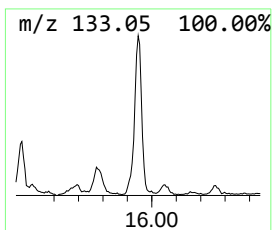
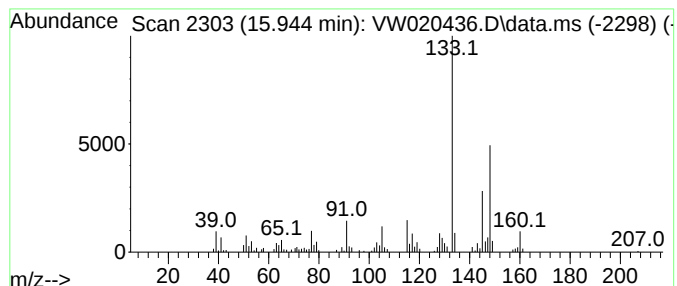
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Quant Title : SFAM01.0

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

Peak Number 22 Benzene, pentamethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.944	2.52 ug/L	64394	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	89
2			Benzene, pentamethyl-	148	C11H16	000700-12-9	76
3			3,4-Dimethylcumene	148	C11H16	1000370-34-1	70
4			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	62
5			4-tert-Butyltoluene	148	C11H16	000098-51-1	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

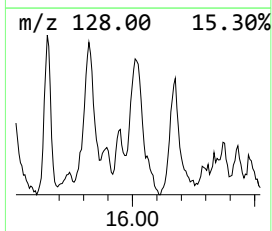
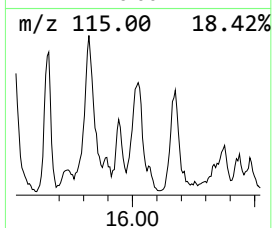
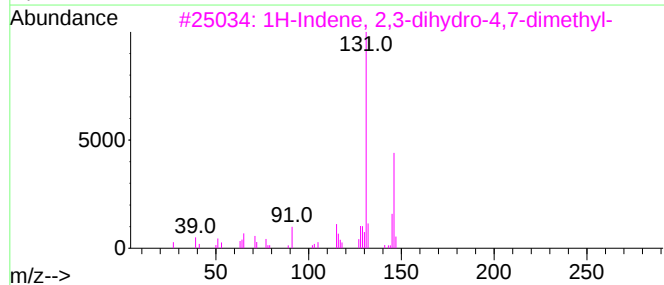
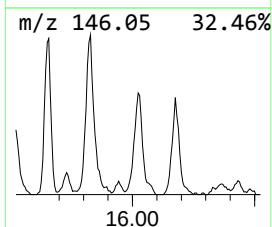
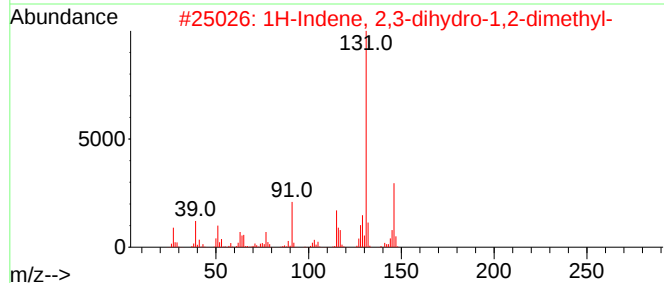
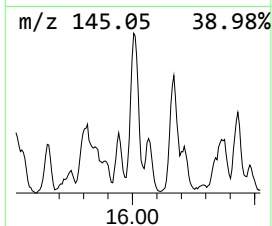
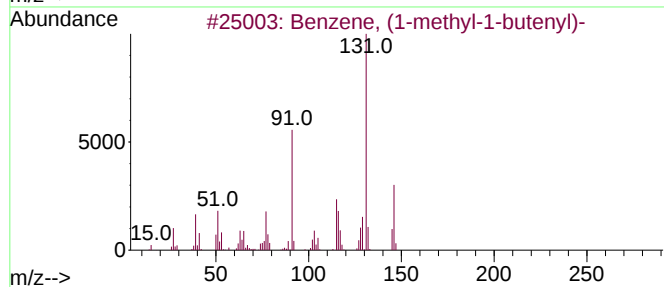
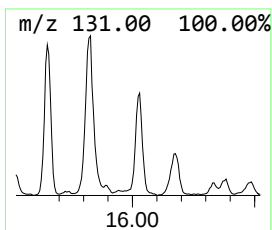
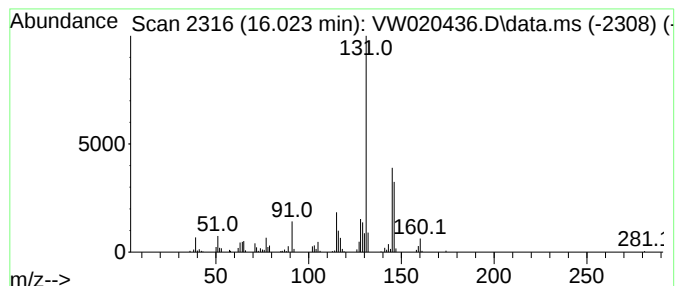
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Quant Title : SFAM01.0

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

Peak Number 23 Benzene, (1-methyl-1-butenyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.023	4.08 ug/L	104087	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	89
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

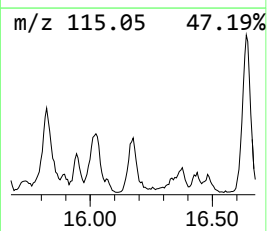
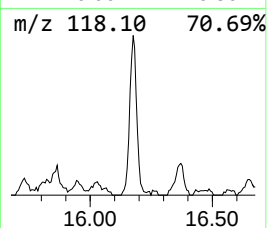
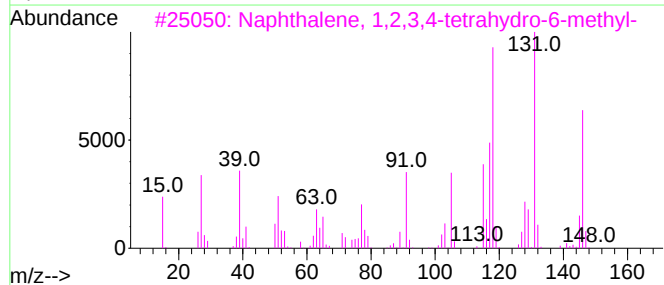
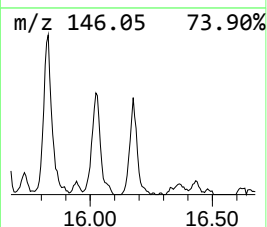
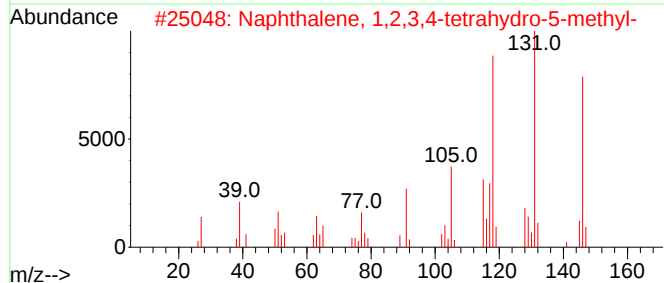
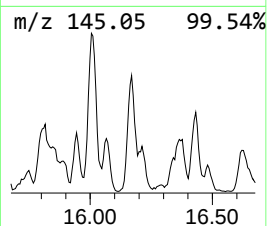
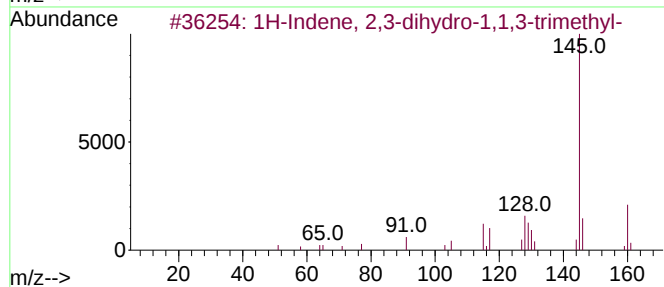
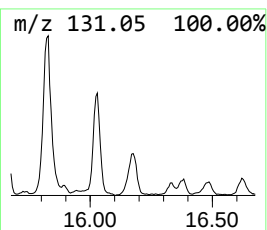
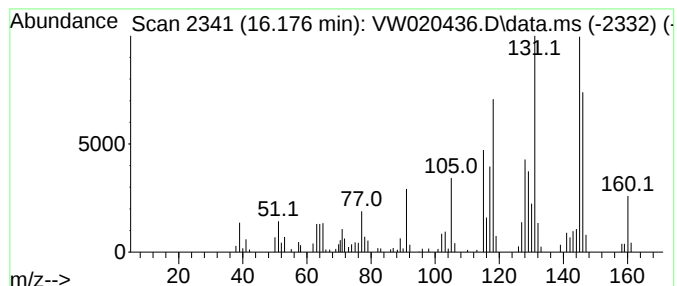
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Quant Title : SFAM01.0

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

Peak Number 24 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.176	4.88 ug/L	124500	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	60
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	60
4			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	60
5			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

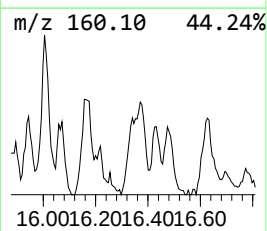
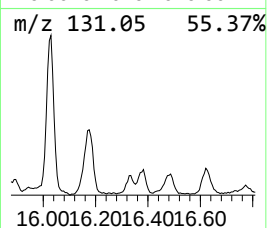
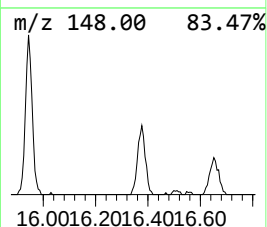
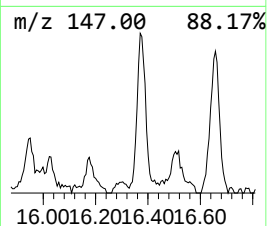
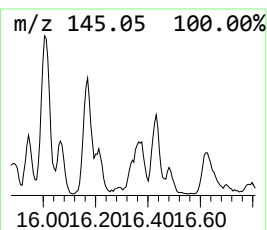
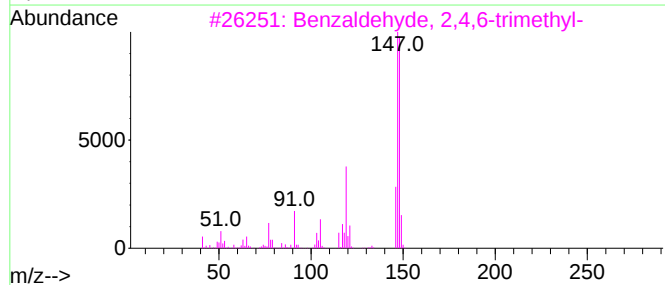
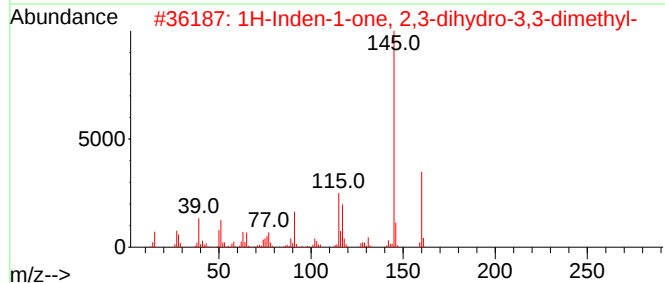
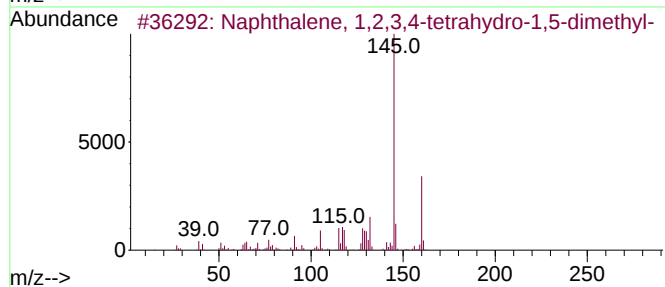
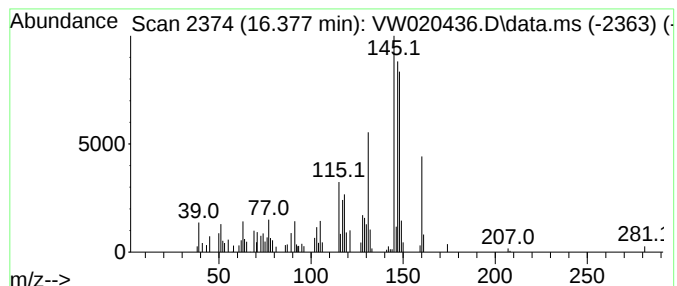
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Quant Title : SFAM01.0

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

Peak Number 25 unknown-02 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.377	2.17 ug/L	55506	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	46
2			1H-Inden-1-one, 2,3-dihydro-3,3-...	160	C11H12O	026465-81-6	42
3			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	41
4			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	41
5			Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	41



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
Data File : VW020436.D
Acq On : 11 Oct 2021 11:47
Operator : SY/VA
Sample : M4070-15
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW7D8

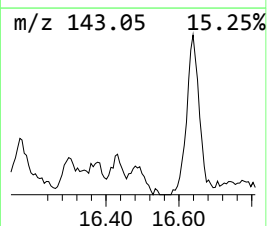
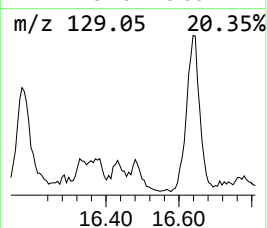
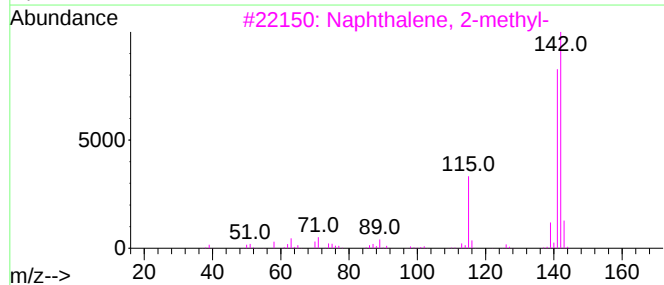
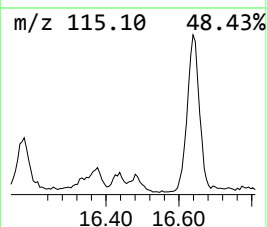
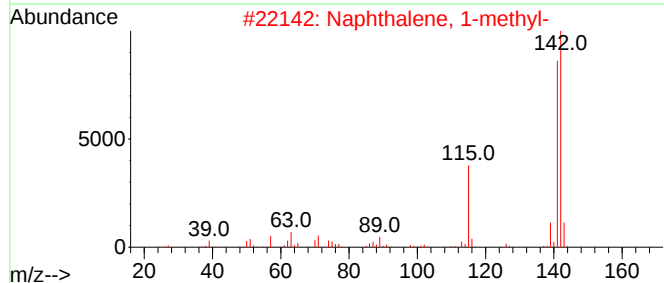
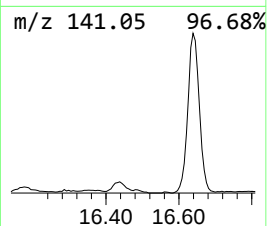
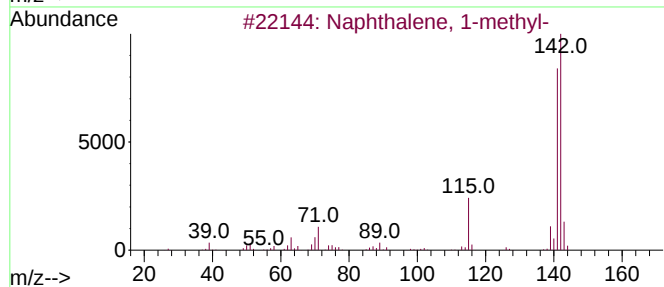
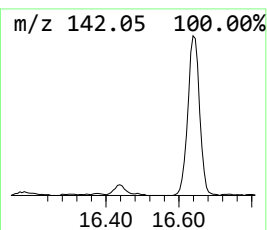
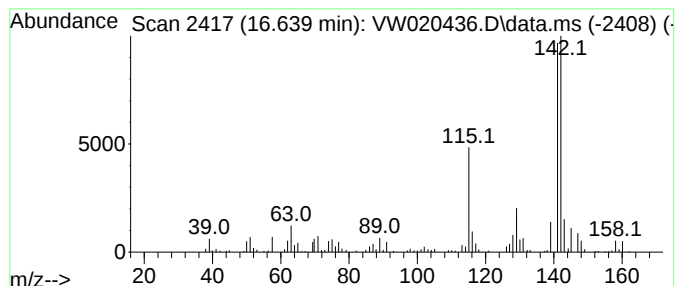
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Quant Title : SFAM01.0

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

Peak Number 26 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	7.39 ug/L	188795	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
4			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
 Data File : VW020436.D
 Acq On : 11 Oct 2021 11:47
 Operator : SY/VA
 Sample : M4070-15
 Misc : 6.19g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EW7D8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.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
Benzene, (2-bro...	13.755	2.9	ug/L	74107	3	13.560	638453	25.0
Benzene, 2-ethy...	14.011	1.9	ug/L	47235	3	13.560	638453	25.0
o-Cymene	14.042	1.5	ug/L	39113	3	13.560	638453	25.0
Benzene, 4-ethy...	14.097	5.0	ug/L	128028	3	13.560	638453	25.0
Benzene, 1-meth...	14.158	1.4	ug/L	35926	3	13.560	638453	25.0
Indan, 1-methyl-	14.206	7.4	ug/L	189857	3	13.560	638453	25.0
1H-Indene, 2,3-...	14.341	2.0	ug/L	50820	3	13.560	638453	25.0
Benzene, 1,2,3,...	14.420	6.0	ug/L	154611	3	13.560	638453	25.0
Benzene, 1-ethy...	14.456	8.6	ug/L	219074	3	13.560	638453	25.0
1H-Indene, 2,3-...	14.688	2.4	ug/L	61150	3	13.560	638453	25.0
1H-Indene, 2,3-...	14.804	12.9	ug/L	328272	3	13.560	638453	25.0
Benzene, 1-meth...	14.859	1.4	ug/L	34597	3	13.560	638453	25.0
Naphthalene, 1,...	14.956	1.7	ug/L	43381	3	13.560	638453	25.0
Benzene, (3-met...	15.060	7.0	ug/L	178252	3	13.560	638453	25.0
1H-Indene, 2,3-...	15.176	11.6	ug/L	297178	3	13.560	638453	25.0
unknown-01	15.468	1.3	ug/L	33358	3	13.560	638453	25.0
1H-Indene, 2,3-...	15.651	4.9	ug/L	125796	3	13.560	638453	25.0
1H-Indene, 2,3-...	15.822	5.7	ug/L	145326	3	13.560	638453	25.0
Benzene, pentam...	15.944	2.5	ug/L	64394	3	13.560	638453	25.0
Benzene, (1-met...	16.023	4.1	ug/L	104087	3	13.560	638453	25.0
1H-Indene, 2,3-...	16.176	4.9	ug/L	124500	3	13.560	638453	25.0
unknown-02	16.377	2.2	ug/L	55506	3	13.560	638453	25.0
Naphthalene, 1-...	16.639	7.4	ug/L	188795	3	13.560	638453	25.0