

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
 Data File : VW019447.D
 Acq On : 12 Jul 2021 22:37
 Operator : SY/VA
 Sample : M2973-04
 Misc : 3.90G/10ML/MSVOA_W/SOIL
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BGEB2

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\SFAMWLM070921SMA.M
 Title : SFAM01.0

Signal : TIC: VW019447.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.361	66	75	86	rVB	96002	185502	11.79%	1.430%
2	2.745	136	138	140	rVB3	752	362	0.02%	0.003%
3	2.898	155	163	178	rVB	70035	171554	10.90%	1.322%
4	3.398	238	245	253	rVB4	2216	5353	0.34%	0.041%
5	3.458	253	255	257	rVB	374	306	0.02%	0.002%
6	3.495	257	261	262	rBV2	326	314	0.02%	0.002%
7	3.550	266	270	271	rBV2	329	361	0.02%	0.003%
8	3.684	288	292	295	rBV3	460	795	0.05%	0.006%
9	3.733	298	300	306	rVB3	544	830	0.05%	0.006%
10	4.025	334	348	359	rBV	166488	509671	32.38%	3.928%
11	4.135	360	366	386	rVB	34891	105699	6.72%	0.815%
12	4.391	398	408	420	rBV	18762	53851	3.42%	0.415%
13	4.745	463	466	471	rBV	239	302	0.02%	0.002%
14	4.922	482	495	519	rVB	66083	230156	14.62%	1.774%
15	5.440	577	580	581	rVV2	258	311	0.02%	0.002%
16	5.793	635	638	640	rVV3	353	477	0.03%	0.004%
17	5.946	653	663	675	rBV	25208	76985	4.89%	0.593%
18	6.970	829	831	832	rBV	383	377	0.02%	0.003%
19	6.994	832	835	840	rVB4	817	1243	0.08%	0.010%
20	7.086	840	850	861	rBV	39811	129780	8.25%	1.000%
21	7.653	932	943	956	rBV	261497	653287	41.50%	5.035%
22	7.756	958	960	962	rBV2	270	363	0.02%	0.003%
23	7.958	989	993	996	rVB4	573	1008	0.06%	0.008%
24	8.281	1035	1046	1063	rBV2	535716	1573999	100.00%	12.131%
25	8.573	1091	1094	1100	rVB3	617	929	0.06%	0.007%
26	8.634	1100	1104	1107	rVB2	288	507	0.03%	0.004%
27	8.701	1107	1115	1122	rBV4	955	2279	0.14%	0.018%
28	8.842	1130	1138	1153	rBV	568801	1136836	72.23%	8.761%
29	9.079	1170	1177	1184	rVB4	2364	4840	0.31%	0.037%
30	9.274	1199	1209	1227	rBV	361583	733599	46.61%	5.654%
31	9.445	1233	1237	1239	rVV2	558	834	0.05%	0.006%
32	9.665	1268	1273	1281	rVB2	2338	5105	0.32%	0.039%
33	9.756	1287	1288	1294	rVB3	334	391	0.02%	0.003%
34	9.909	1308	1313	1315	rBV	227	341	0.02%	0.003%
35	9.957	1317	1321	1326	rBV3	782	1228	0.08%	0.009%
36	10.037	1326	1334	1344	rVV	167865	298835	18.99%	2.303%

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37	10.122	1345	1348	1356	rVB5	1870	3926	0.25%	0.030%
38	10.323	1373	1381	1388	rBV	743991	1321367	83.95%	10.184%
39	10.390	1388	1392	1398	rVB	26703	46822	2.97%	0.361%
40	10.506	1404	1411	1416	rVB6	1755	3828	0.24%	0.030%
41	10.579	1416	1423	1431	rBV2	103201	179476	11.40%	1.383%
42	10.665	1431	1437	1440	rVV4	1813	3217	0.20%	0.025%
43	10.701	1440	1443	1448	rVB2	2622	4090	0.26%	0.032%
44	10.762	1448	1453	1460	rVB6	1243	2269	0.14%	0.017%
45	10.829	1460	1464	1465	rBV2	412	610	0.04%	0.005%
46	10.860	1467	1469	1473	rVB2	540	557	0.04%	0.004%
47	10.927	1473	1480	1496	rBV	157036	271832	17.27%	2.095%
48	11.067	1500	1503	1507	rBV5	598	952	0.06%	0.007%
49	11.177	1512	1521	1526	rBV4	3475	6744	0.43%	0.052%
50	11.225	1526	1529	1536	rVB3	2555	4483	0.28%	0.035%
51	11.286	1536	1539	1542	rBV2	733	873	0.06%	0.007%
52	11.341	1542	1548	1551	rVB5	1221	1962	0.12%	0.015%
53	11.439	1559	1564	1568	rVB3	871	1612	0.10%	0.012%
54	11.512	1573	1576	1578	rBV3	738	820	0.05%	0.006%
55	11.530	1578	1579	1582	rBV3	355	306	0.02%	0.002%
56	11.561	1582	1584	1588	rVB2	271	310	0.02%	0.002%
57	11.634	1588	1596	1605	rBV	763249	1284868	81.63%	9.902%
58	11.725	1607	1611	1617	rVB4	2730	4267	0.27%	0.033%
59	11.835	1622	1629	1632	rVV6	2446	5287	0.34%	0.041%
60	11.872	1633	1635	1647	rVB6	2863	7462	0.47%	0.058%
61	11.994	1647	1655	1659	rBV6	1154	3047	0.19%	0.023%
62	12.030	1659	1661	1665	rBV3	435	485	0.03%	0.004%
63	12.140	1672	1679	1682	rBV4	2533	5176	0.33%	0.040%
64	12.237	1691	1695	1701	rVB4	3012	5629	0.36%	0.043%
65	12.347	1701	1713	1714	rBV6	6147	13179	0.84%	0.102%
66	12.475	1728	1734	1739	rBB7	3202	5900	0.37%	0.045%
67	12.597	1748	1754	1761	rVB4	5297	12091	0.77%	0.093%
68	12.695	1761	1770	1777	rBV	232064	395752	25.14%	3.050%
69	12.780	1777	1784	1792	rVB8	4601	11577	0.74%	0.089%
70	12.865	1792	1798	1802	rBV7	2442	5785	0.37%	0.045%
71	12.938	1804	1810	1819	rBV5	10316	25177	1.60%	0.194%
72	13.103	1833	1837	1844	rVB	9345	15557	0.99%	0.120%
73	13.188	1848	1851	1858	rVB2	7838	13534	0.86%	0.104%
74	13.249	1858	1861	1865	rBV3	3523	6026	0.38%	0.046%
75	13.298	1865	1869	1873	rVV2	5411	8087	0.51%	0.062%
76	13.408	1883	1887	1891	rBV2	10157	13915	0.88%	0.107%

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77	13.554	1905	1911	1918	rBV	615214	993371	63.11%	7.656%
78	13.646	1924	1926	1932	rVB4	8443	11086	0.70%	0.085%
79	13.853	1953	1960	1971	rBV	483181	807967	51.33%	6.227%
80	13.944	1972	1975	1979	rVB	7769	11508	0.73%	0.089%
81	14.097	1991	2000	2006	rVB2	86590	156142	9.92%	1.203%
82	14.158	2007	2010	2013	rBV	4142	4663	0.30%	0.036%
83	14.207	2013	2018	2024	rBV	60156	97845	6.22%	0.754%
84	14.322	2033	2037	2043	rVB6	4960	10162	0.65%	0.078%
85	14.414	2043	2052	2062	rBV	96402	162515	10.32%	1.252%
86	14.499	2062	2066	2068	rBV3	5333	8292	0.53%	0.064%
87	14.566	2074	2077	2083	rVB	26786	37853	2.40%	0.292%
88	14.688	2091	2097	2101	rBV	107726	166737	10.59%	1.285%
89	14.810	2109	2117	2122	rBV	252426	418671	26.60%	3.227%
90	15.060	2153	2158	2163	rVB	57854	95303	6.05%	0.734%
91	15.176	2172	2177	2184	rVB2	31156	67399	4.28%	0.519%
92	15.280	2189	2194	2198	rBV	18417	34810	2.21%	0.268%
93	15.469	2221	2225	2229	rBV3	19039	38003	2.41%	0.293%
94	15.517	2231	2233	2239	rVB	29386	44958	2.86%	0.346%
95	15.651	2249	2255	2262	rVB	33819	62128	3.95%	0.479%
96	15.822	2276	2283	2293	rBV2	26776	66856	4.25%	0.515%
97	15.956	2299	2305	2306	rBV3	10066	18048	1.15%	0.139%
98	16.029	2312	2317	2323	rVB	22538	43372	2.76%	0.334%
99	16.267	2352	2356	2362	rVB6	5289	11689	0.74%	0.090%
100	16.371	2369	2373	2380	rVB4	9312	18640	1.18%	0.144%

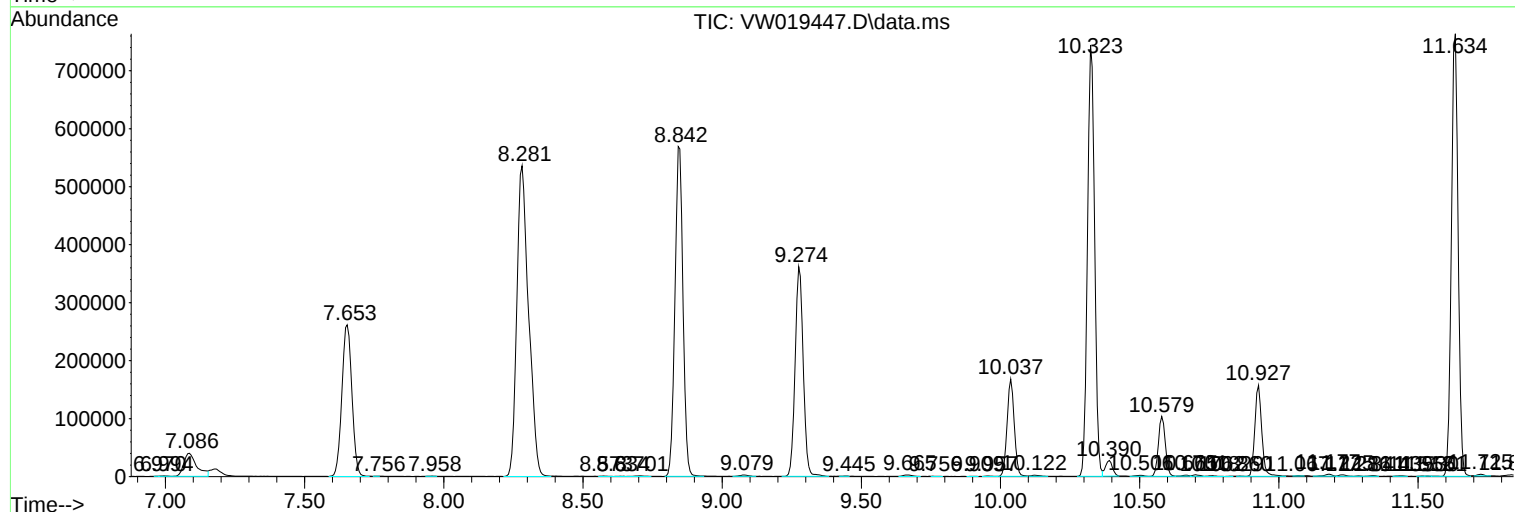
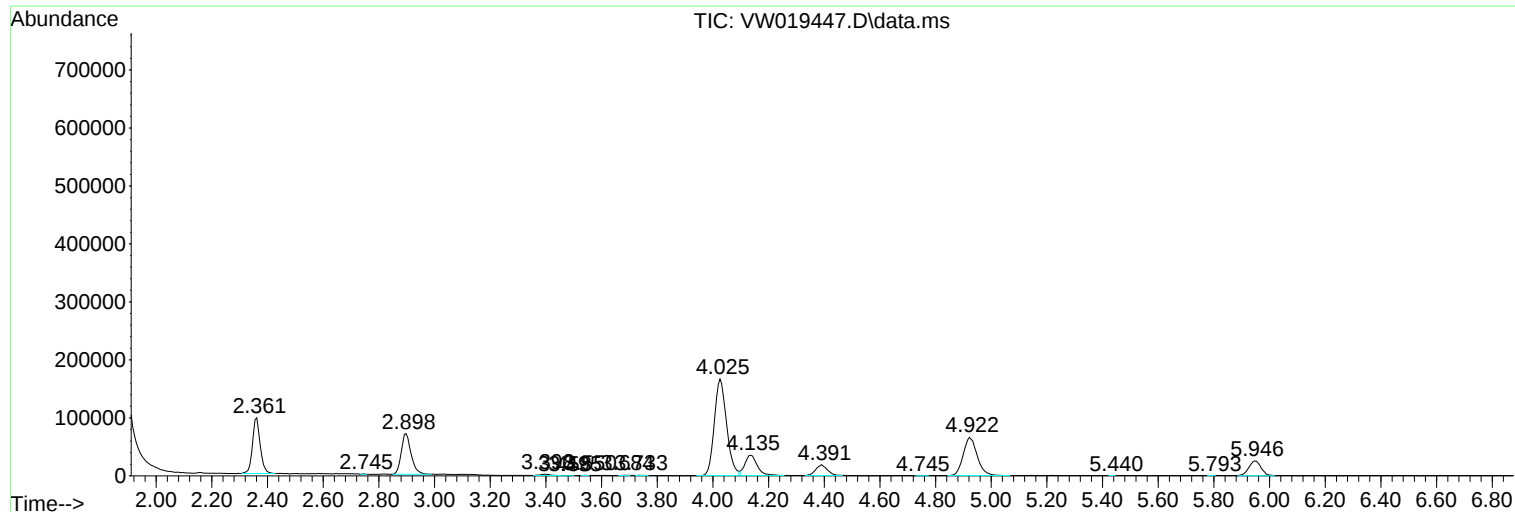
Sum of corrected areas: 12975485

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070921SMA.M
Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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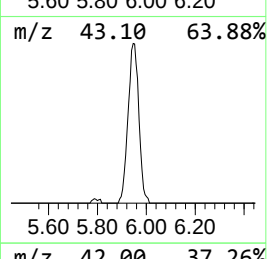
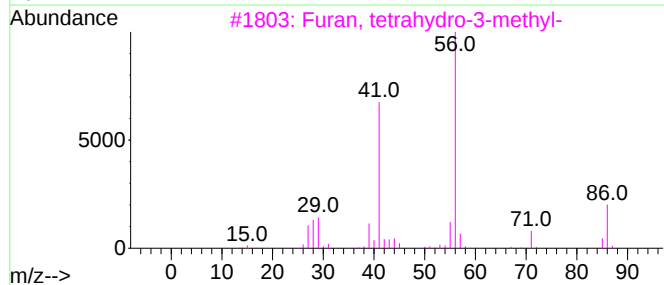
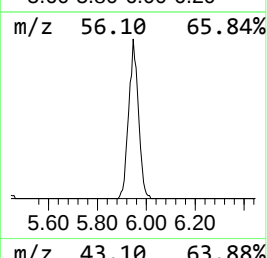
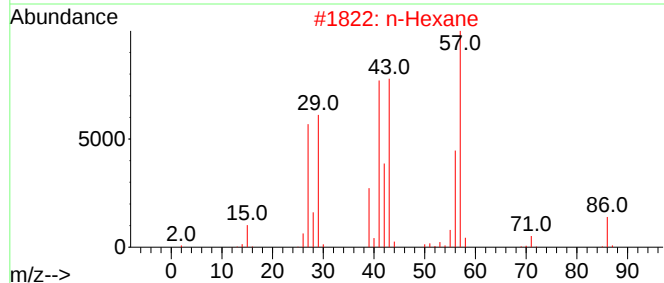
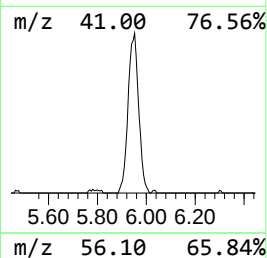
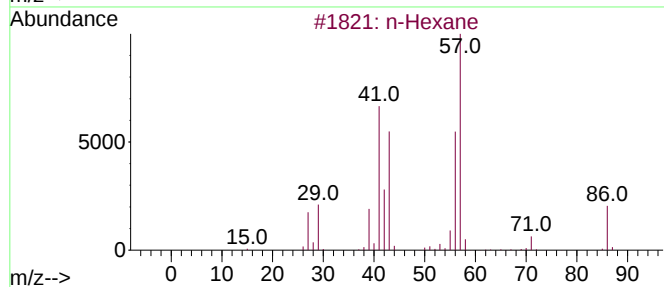
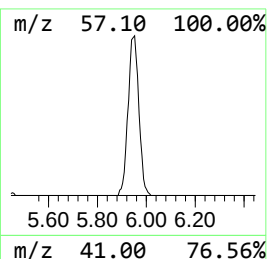
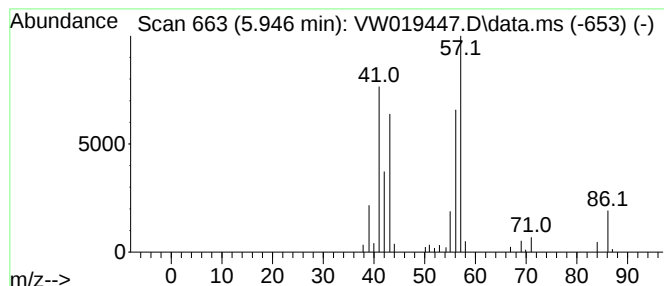
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070921SMA.M
Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.946	1.69 ug/L	76985	1,4-Difluorobenzene	8.848

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	90
2			n-Hexane	86	C6H14	000110-54-3	59
3			Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	50
4			Cyclopropane, propyl-	84	C6H12	002415-72-7	47
5			n-Hexane	86	C6H14	000110-54-3	45



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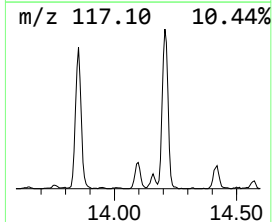
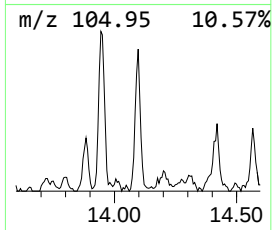
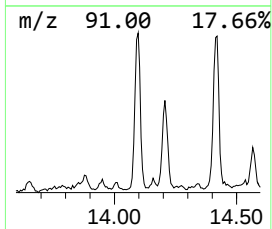
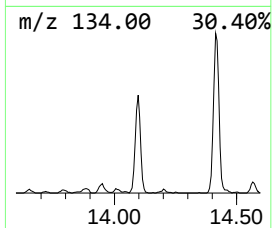
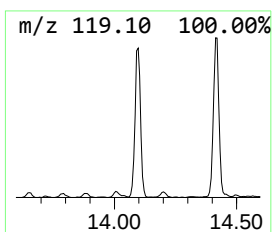
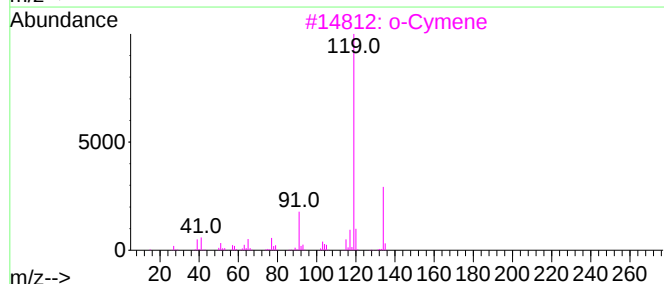
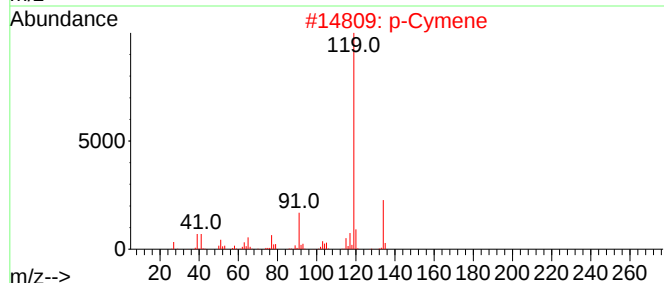
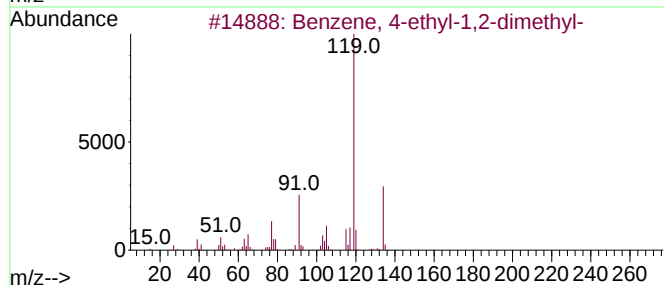
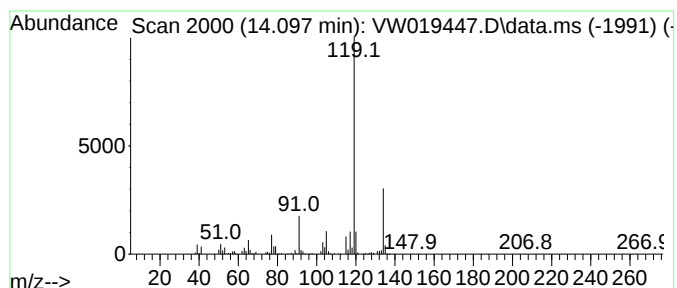
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.097	3.93 ug/L	156142	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		p-Cymene	134	C10H14	000099-87-6	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



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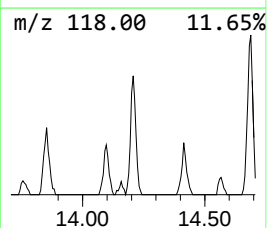
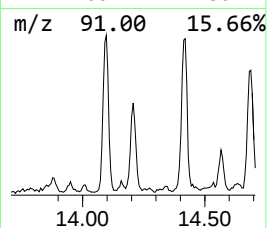
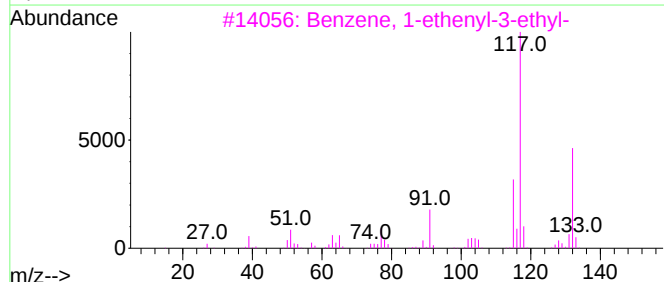
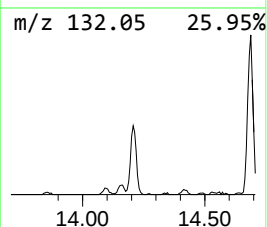
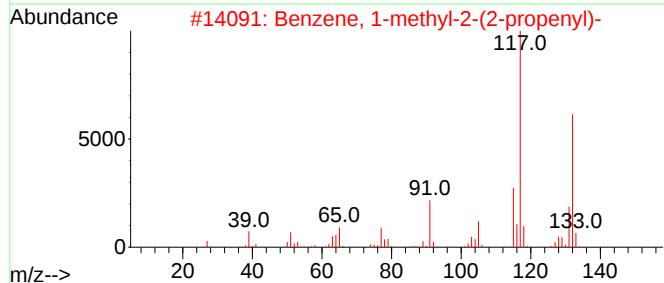
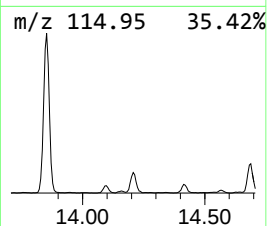
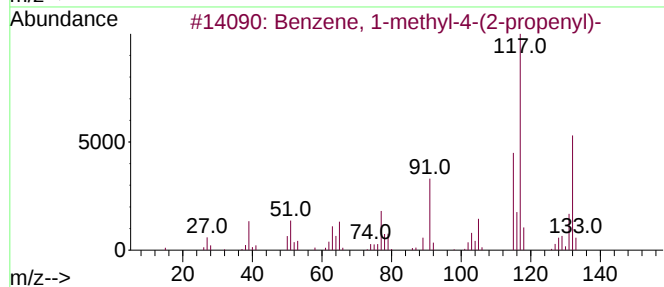
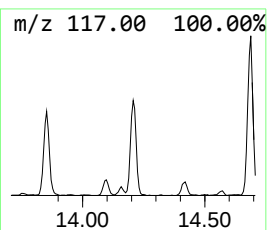
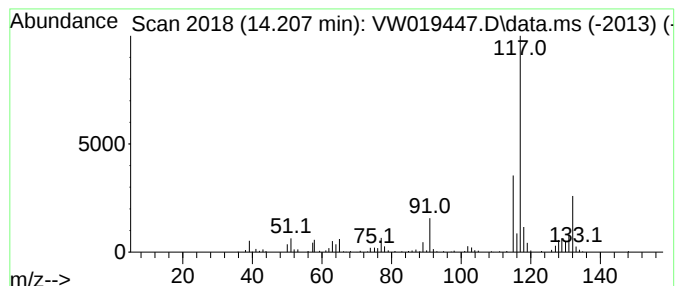
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Quant Title : SFAM01.0

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Peak Number 4 Benzene, 1-methyl-4-(2-prop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.207	2.46 ug/L	97845	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	87
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	86
5			Indan, 1-methyl-	132	C10H12	000767-58-8	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

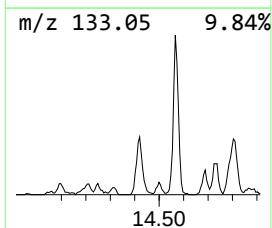
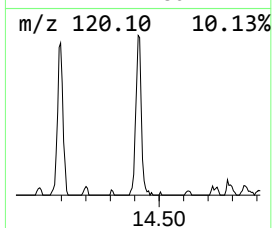
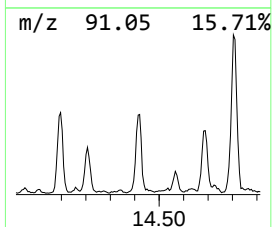
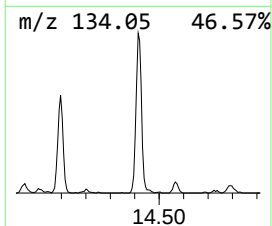
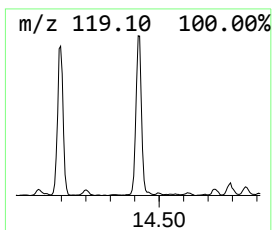
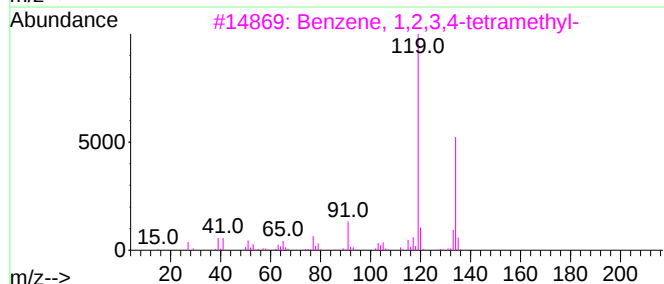
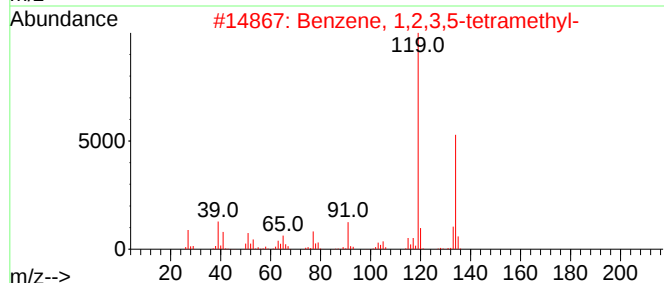
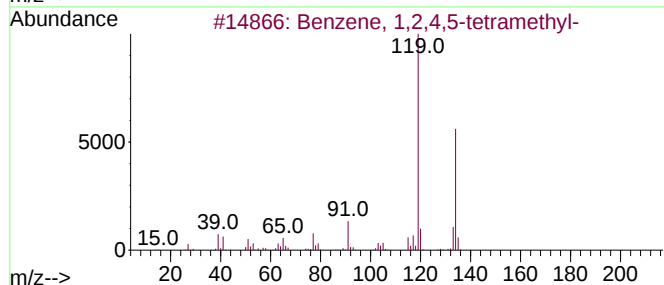
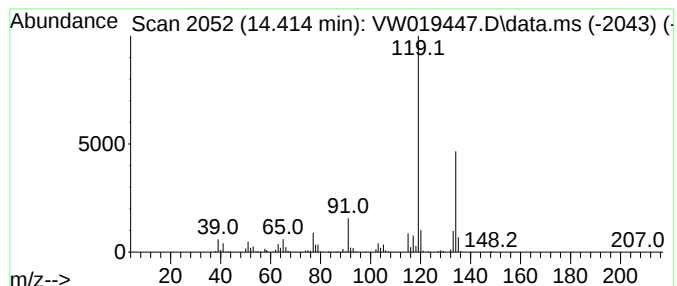
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.414	4.09 ug/L	162515	1,4-Dichlorobenzene-d4	13.560

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

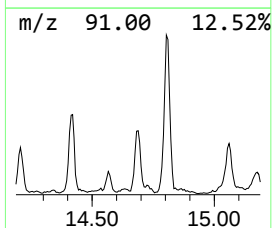
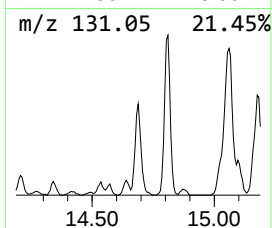
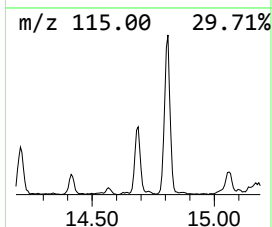
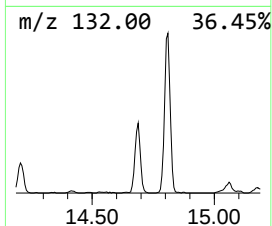
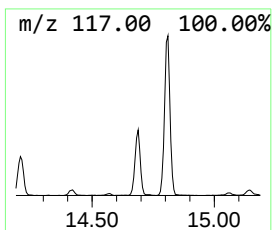
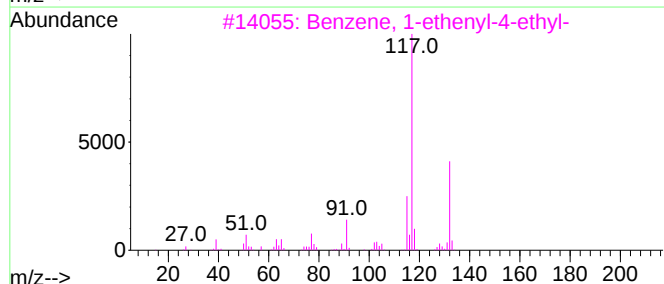
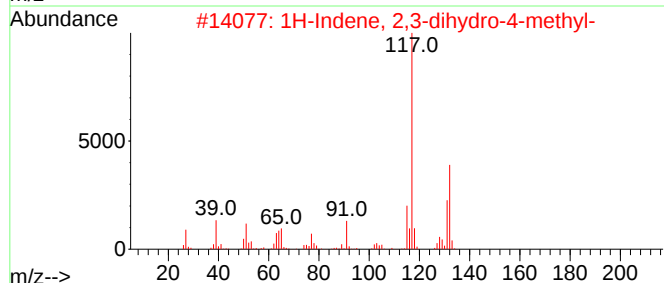
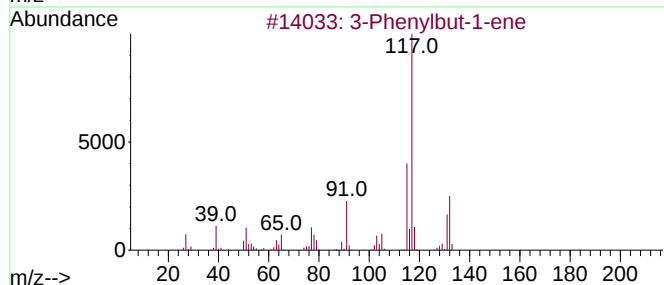
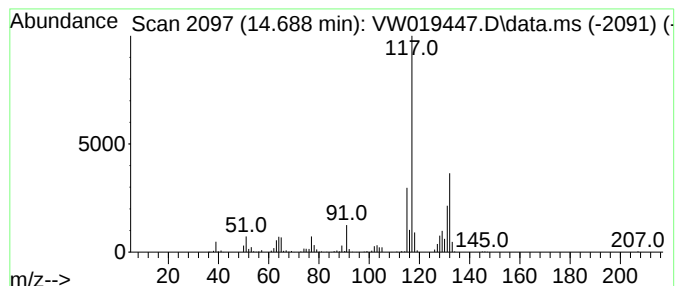
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 3-Phenylbut-1-ene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.688	4.20 ug/L	166737	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

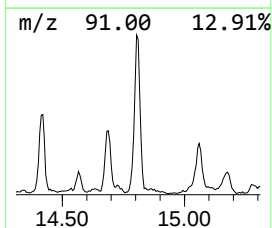
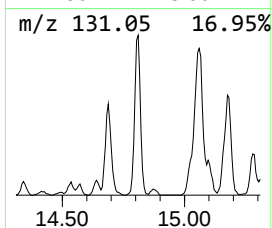
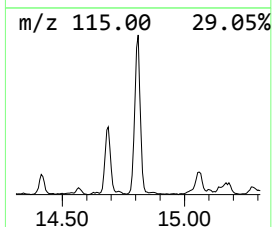
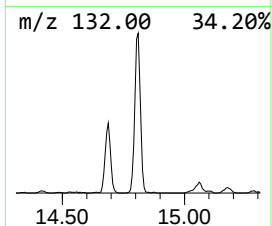
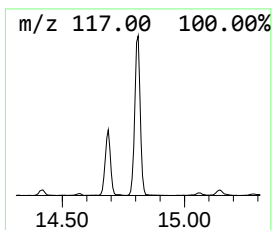
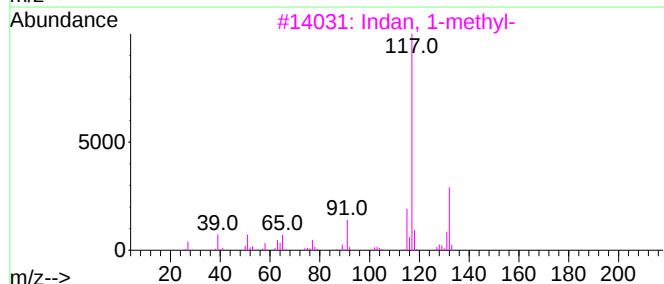
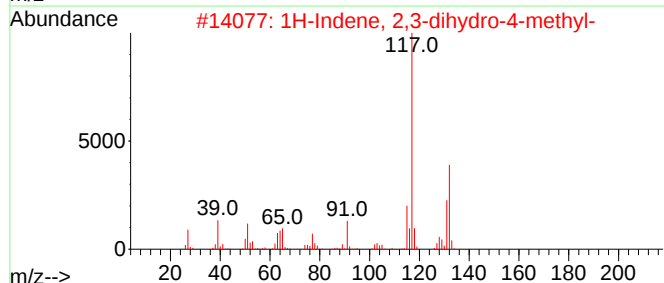
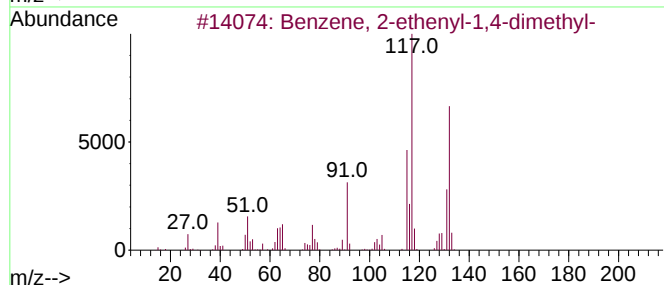
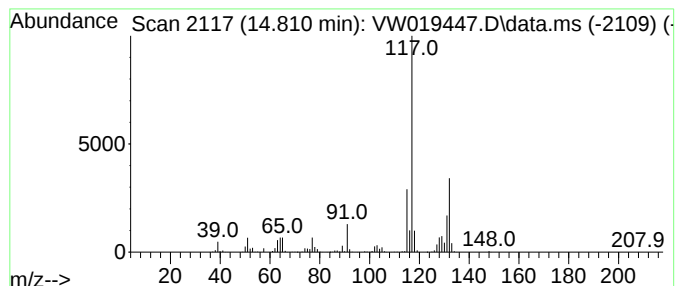
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.810	10.54 ug/L	418671	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

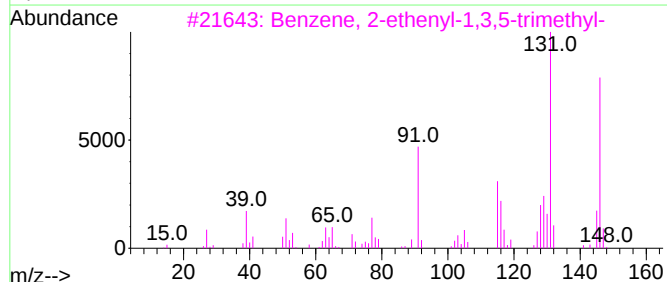
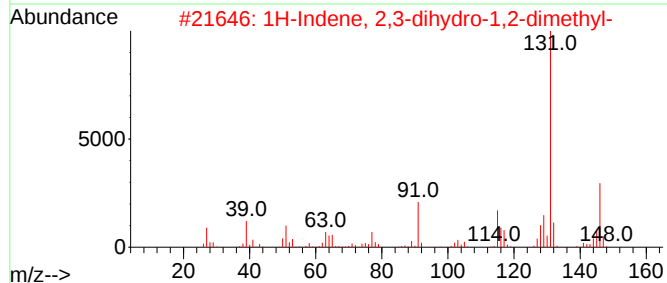
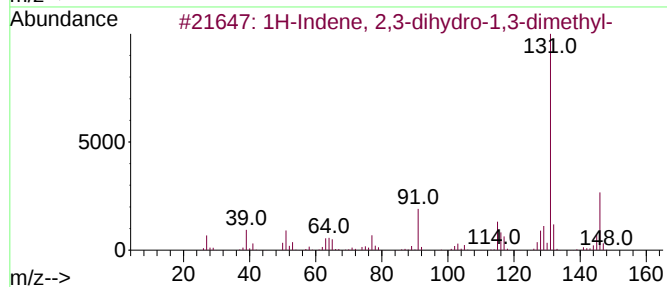
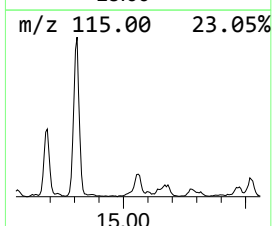
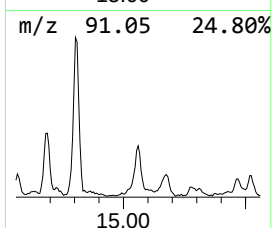
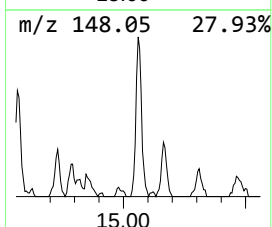
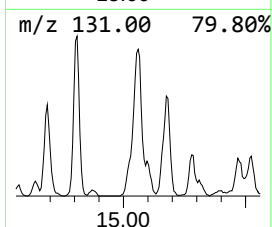
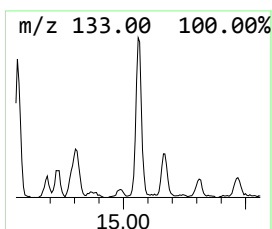
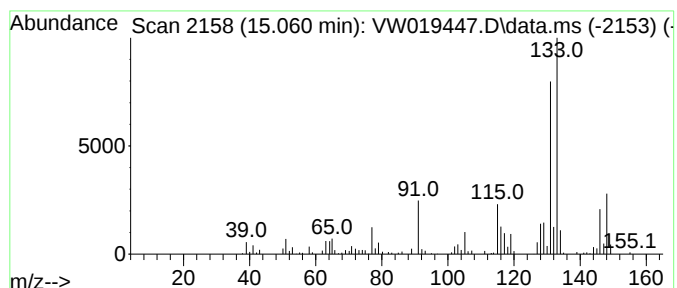
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.060	2.40 ug/L	95303	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	55
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	50
5			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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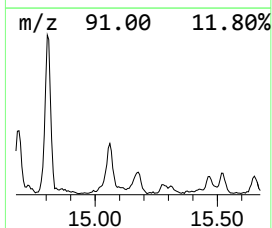
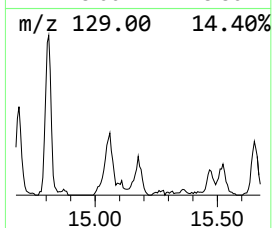
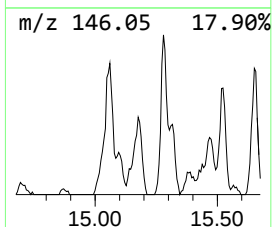
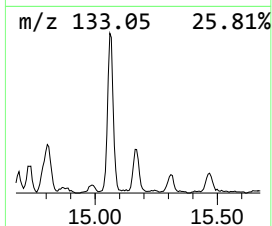
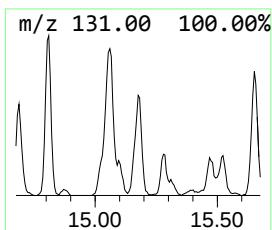
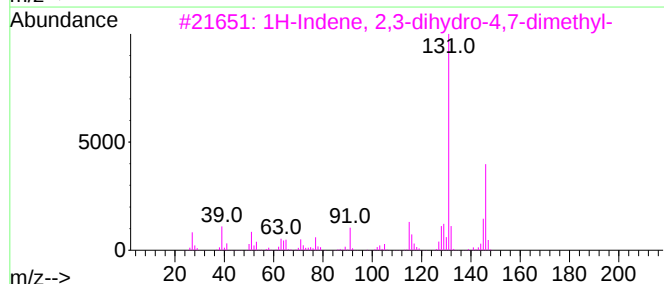
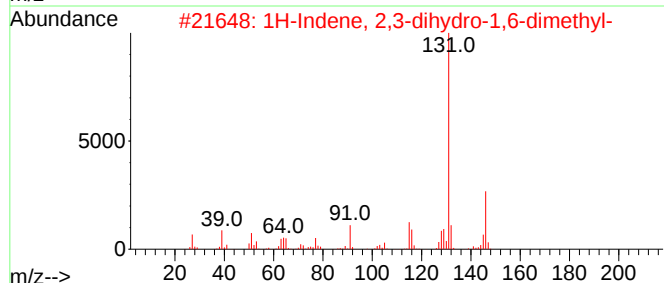
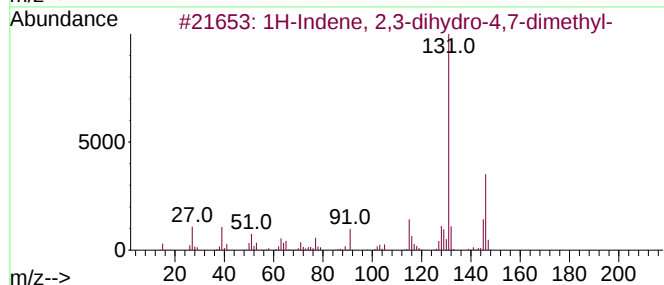
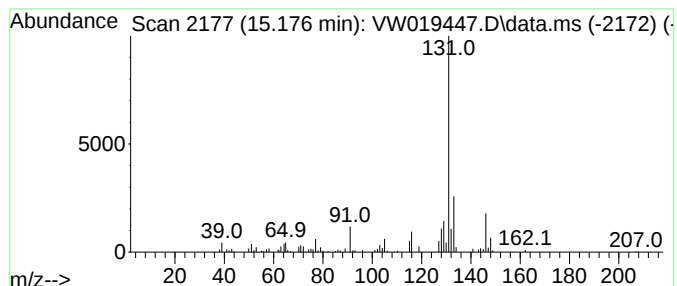
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.176	1.70 ug/L	67399	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	87
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
4			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

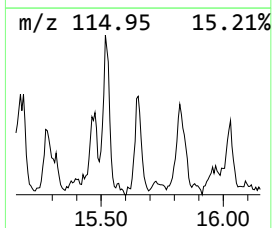
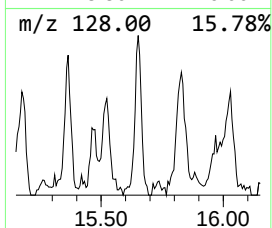
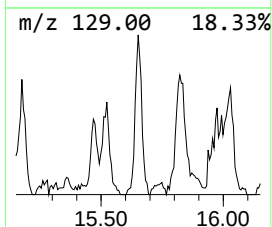
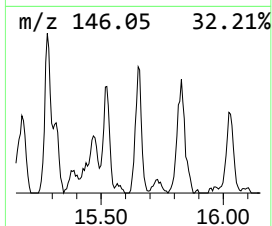
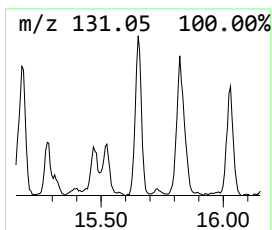
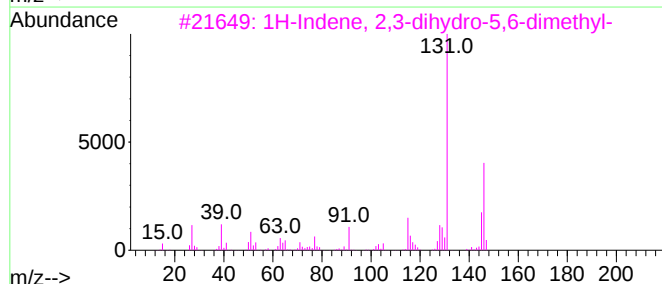
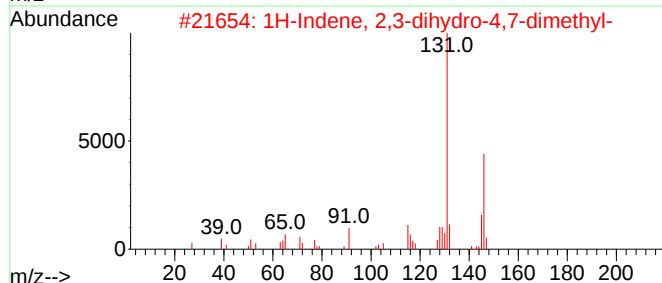
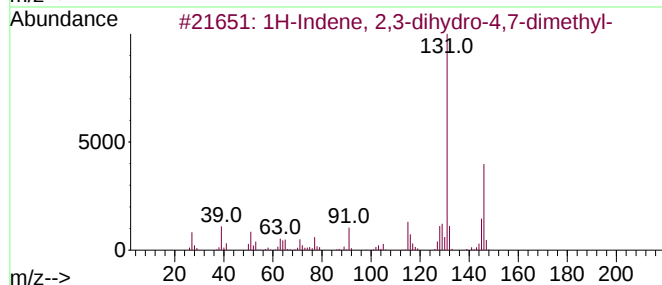
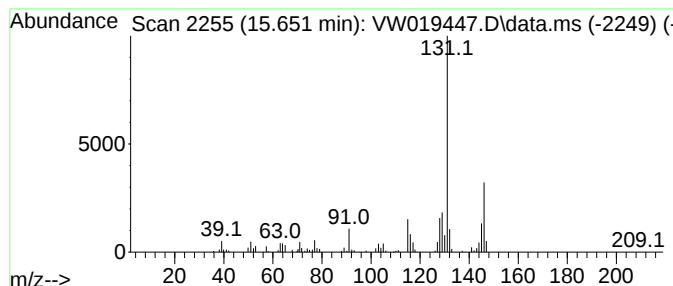
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 2,3-dihydro-5,6-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	1.56 ug/L	62128	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	95		
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95		
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93		
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91		
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

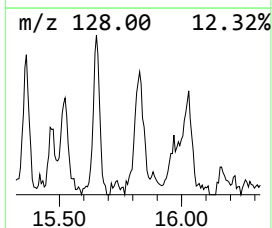
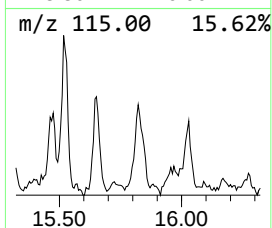
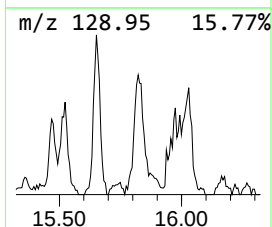
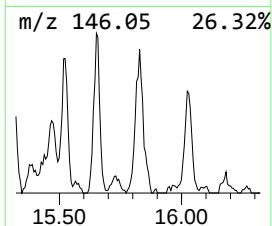
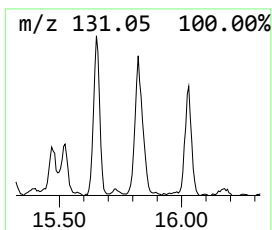
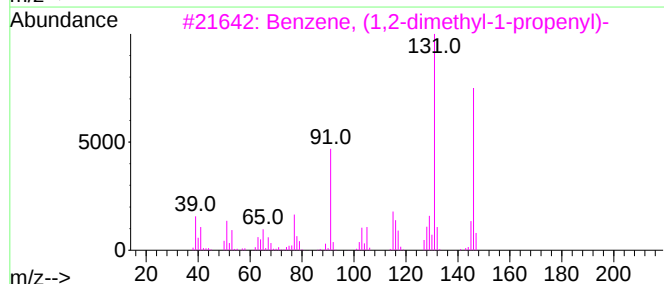
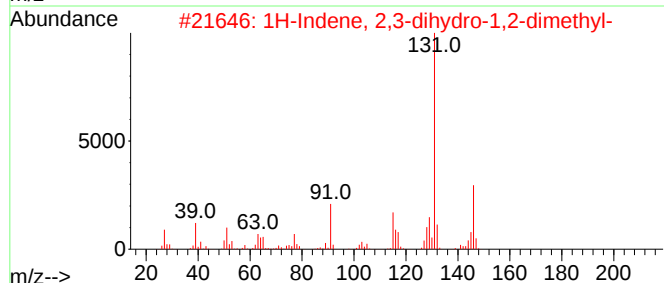
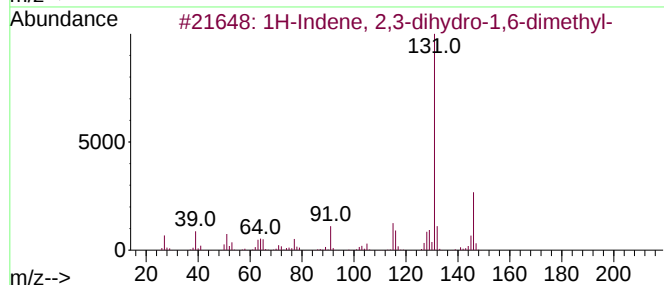
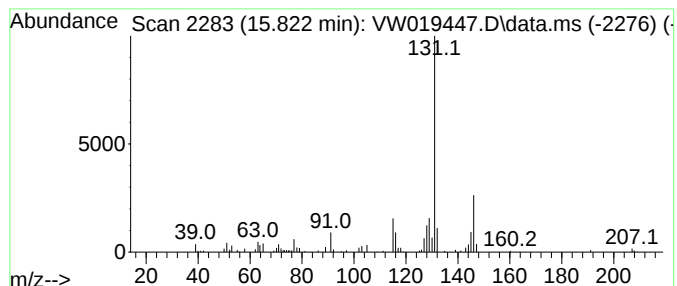
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Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.822	1.68 ug/L	66856	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	95
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071221\
Data File : VW019447.D
Acq On : 12 Jul 2021 22:37
Operator : SY/VA
Sample : M2973-04
Misc : 3.90G/10ML/MSVOA_W/SOIL
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGEB2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070921SMA.M
Quant Title : SFAM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.946	1.7	ug/L	76985	1	8.848	1136840	25.0
Benzene, 4-ethy...	14.097	3.9	ug/L	156142	3	13.560	993371	25.0
Benzene, 1-meth...	14.207	2.5	ug/L	97845	3	13.560	993371	25.0
Benzene, 1,2,4,...	14.414	4.1	ug/L	162515	3	13.560	993371	25.0
3-Phenylbut-1-ene	14.688	4.2	ug/L	166737	3	13.560	993371	25.0
Benzene, 2-ethe...	14.810	10.5	ug/L	418671	3	13.560	993371	25.0
1H-Indene, 2,3-...	15.060	2.4	ug/L	95303	3	13.560	993371	25.0
1H-Indene, 2,3-...	15.176	1.7	ug/L	67399	3	13.560	993371	25.0
1H-Indene, 2,3-...	15.651	1.6	ug/L	62128	3	13.560	993371	25.0
1H-Indene, 2,3-...	15.822	1.7	ug/L	66856	3	13.560	993371	25.0