

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060421\
 Data File : VW019135.D
 Acq On : 04 Jun 2021 17:53
 Operator : SY/VA
 Sample : M2618-08
 Misc : 4.36G/10ML/MSVOA_W/SOIL
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BG5S6

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\SFAMWLM060221SMA.M
 Title : SFAM01.0

Signal : TIC: VW019135.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	65	75	86	rBV	185057	371079	15.26%	1.926%
2	2.501	93	98	108	rBV	26454	68813	2.83%	0.357%
3	2.898	154	163	176	rBV	125600	313160	12.88%	1.626%
4	3.404	239	246	257	rVB5	2140	6382	0.26%	0.033%
5	4.025	336	348	358	rBV	270817	853076	35.09%	4.429%
6	4.135	358	366	385	rVB	242821	712577	29.31%	3.699%
7	4.446	403	417	435	rVV3	11856	60580	2.49%	0.314%
8	4.922	482	495	515	rBV2	32058	122462	5.04%	0.636%
9	5.196	539	540	555	rVB5	3290	7980	0.33%	0.041%
10	5.793	625	638	652	rBV2	9033	32286	1.33%	0.168%
11	5.946	654	663	674	rVB3	5077	15551	0.64%	0.081%
12	6.110	682	690	699	rBV6	2961	10273	0.42%	0.053%
13	6.915	811	822	829	rBV2	12388	39601	1.63%	0.206%
14	7.086	841	850	854	rBV2	51507	133936	5.51%	0.695%
15	7.171	855	864	879	rVB2	126852	430298	17.70%	2.234%
16	7.537	916	924	932	rBV8	2356	7455	0.31%	0.039%
17	7.653	932	943	957	rVV	419546	1026512	42.22%	5.329%
18	7.958	983	993	999	rVB9	5098	13386	0.55%	0.069%
19	8.159	1019	1026	1033	rVB5	3014	7390	0.30%	0.038%
20	8.281	1033	1046	1062	rBV2	834350	2431288	100.00%	12.622%
21	8.476	1070	1078	1087	rVV4	6102	23202	0.95%	0.120%
22	8.555	1087	1091	1099	rVB7	4063	10268	0.42%	0.053%
23	8.707	1107	1116	1130	rBV	46290	106862	4.40%	0.555%
24	8.842	1130	1138	1153	rBV	774390	1597238	65.70%	8.292%
25	9.079	1168	1177	1183	rBV8	4626	13975	0.57%	0.073%
26	9.201	1192	1197	1201	rBV3	3860	6733	0.28%	0.035%
27	9.274	1201	1209	1226	rVV2	590450	1245838	51.24%	6.468%
28	9.433	1227	1235	1245	rVB3	15140	37857	1.56%	0.197%
29	9.671	1268	1274	1284	rBV5	9758	24238	1.00%	0.126%
30	9.762	1284	1289	1298	rVB3	7110	14166	0.58%	0.074%
31	9.890	1306	1310	1316	rVB2	8156	15949	0.66%	0.083%
32	10.036	1326	1334	1341	rBV	296111	515403	21.20%	2.676%
33	10.213	1355	1363	1372	rBV	40609	79593	3.27%	0.413%
34	10.323	1372	1381	1389	rVV	1048117	1883893	77.49%	9.780%
35	10.390	1389	1392	1395	rVV	22615	37410	1.54%	0.194%
36	10.421	1395	1397	1404	rVB	16541	26272	1.08%	0.136%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM060221SMA.M
 Title : SFAM01.0

37	10.500	1404	1410	1416	rVB7	10342	23604	0.97%	0.123%
38	10.579	1416	1423	1436	rBV	177425	312250	12.84%	1.621%
39	10.707	1437	1444	1452	rBV	69198	126208	5.19%	0.655%
40	10.829	1459	1464	1467	rBV	23376	42373	1.74%	0.220%
41	10.859	1467	1469	1474	rVB2	21618	29159	1.20%	0.151%
42	10.927	1474	1480	1485	rBV	213688	363932	14.97%	1.889%
43	10.969	1485	1487	1495	rVB	32940	48569	2.00%	0.252%
44	11.067	1498	1503	1508	rBV3	31700	56699	2.33%	0.294%
45	11.177	1518	1521	1528	rVB7	3518	5931	0.24%	0.031%
46	11.286	1534	1539	1544	rVB	8303	14756	0.61%	0.077%
47	11.341	1544	1548	1554	rVB2	5943	9502	0.39%	0.049%
48	11.439	1555	1564	1574	rBV3	43568	90843	3.74%	0.472%
49	11.530	1574	1579	1582	rVV3	8887	17463	0.72%	0.091%
50	11.567	1583	1585	1589	rVV2	10000	16562	0.68%	0.086%
51	11.634	1589	1596	1604	rVV	877874	1472866	60.58%	7.646%
52	11.731	1604	1612	1618	rVV3	28172	57506	2.37%	0.299%
53	11.872	1619	1635	1642	rVB2	51593	144398	5.94%	0.750%
54	11.945	1642	1647	1650	rBV3	4583	7503	0.31%	0.039%
55	11.993	1650	1655	1662	rVB2	43741	77493	3.19%	0.402%
56	12.060	1662	1666	1670	rBV4	6556	12223	0.50%	0.063%
57	12.164	1678	1683	1690	rVB2	28498	52690	2.17%	0.274%
58	12.237	1690	1695	1703	rBV	17345	27566	1.13%	0.143%
59	12.347	1703	1713	1722	rBV9	13761	47518	1.95%	0.247%
60	12.463	1727	1732	1737	rVB3	9603	16773	0.69%	0.087%
61	12.530	1737	1743	1746	rBV5	9972	18394	0.76%	0.095%
62	12.603	1750	1755	1762	rBV	86992	144243	5.93%	0.749%
63	12.695	1764	1770	1776	rBV	301641	490403	20.17%	2.546%
64	12.755	1776	1780	1783	rBV2	13096	19886	0.82%	0.103%
65	12.804	1786	1788	1794	rVB4	10724	15862	0.65%	0.082%
66	12.865	1794	1798	1799	rBV	11482	11609	0.48%	0.060%
67	12.914	1800	1806	1822	rVB3	178865	447844	18.42%	2.325%
68	13.054	1822	1829	1834	rBV6	21361	62251	2.56%	0.323%
69	13.194	1847	1852	1855	rBV3	8812	16259	0.67%	0.084%
70	13.249	1856	1861	1864	rVV	35641	58360	2.40%	0.303%
71	13.292	1864	1868	1874	rVB3	40527	71707	2.95%	0.372%
72	13.402	1875	1886	1890	rBV8	28293	74705	3.07%	0.388%
73	13.444	1890	1893	1895	rVV3	5840	6151	0.25%	0.032%
74	13.554	1905	1911	1924	rBV	512041	912364	37.53%	4.736%
75	13.646	1924	1926	1933	rVB7	12763	21659	0.89%	0.112%
76	13.713	1933	1937	1940	rBV4	12957	20872	0.86%	0.108%

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 Title : SFAM01.0

77	13.761	1940	1945	1948	rBV5	22206	37153	1.53%	0.193%
78	13.792	1948	1950	1954	rVV2	12461	17595	0.72%	0.091%
79	13.853	1954	1960	1973	rVV	462395	820411	33.74%	4.259%
80	13.975	1976	1980	1983	rVV3	14930	24799	1.02%	0.129%
81	14.017	1983	1987	1991	rVV7	13566	27325	1.12%	0.142%
82	14.066	1991	1995	2004	rVB3	36748	92175	3.79%	0.479%
83	14.145	2004	2008	2013	rVB7	9115	16228	0.67%	0.084%
84	14.206	2013	2018	2023	rBV5	12017	28078	1.15%	0.146%
85	14.328	2032	2038	2043	rBV5	17866	32414	1.33%	0.168%
86	14.420	2048	2053	2057	rBV3	25398	42991	1.77%	0.223%
87	14.463	2057	2060	2063	rVB4	4258	6181	0.25%	0.032%
88	14.560	2074	2076	2081	rBV5	4146	6041	0.25%	0.031%
89	14.621	2081	2086	2093	rVB2	43039	75930	3.12%	0.394%
90	14.694	2093	2098	2100	rBV4	10402	18624	0.77%	0.097%
91	14.731	2100	2104	2109	rVB6	11519	20844	0.86%	0.108%
92	14.804	2109	2116	2123	rBV3	17046	44136	1.82%	0.229%
93	15.066	2154	2159	2164	rVV3	11359	22058	0.91%	0.115%
94	15.176	2172	2177	2183	rVB4	18681	35596	1.46%	0.185%
95	15.365	2204	2208	2213	rVV2	8993	12981	0.53%	0.067%
96	15.432	2214	2219	2225	rVV2	38603	63383	2.61%	0.329%
97	15.481	2225	2227	2234	rVB8	5478	9345	0.38%	0.049%
98	15.810	2276	2281	2286	rVV8	7204	15042	0.62%	0.078%
99	16.432	2379	2383	2391	rVB4	7594	14705	0.60%	0.076%
100	16.633	2412	2416	2420	rBV6	4358	8505	0.35%	0.044%

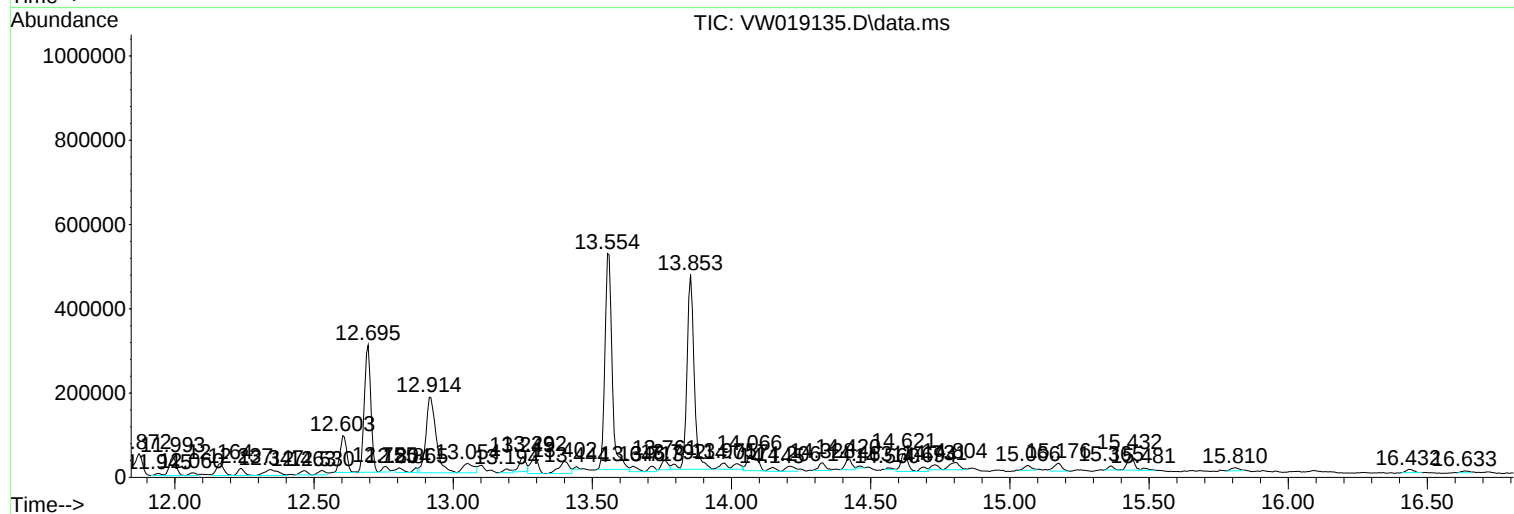
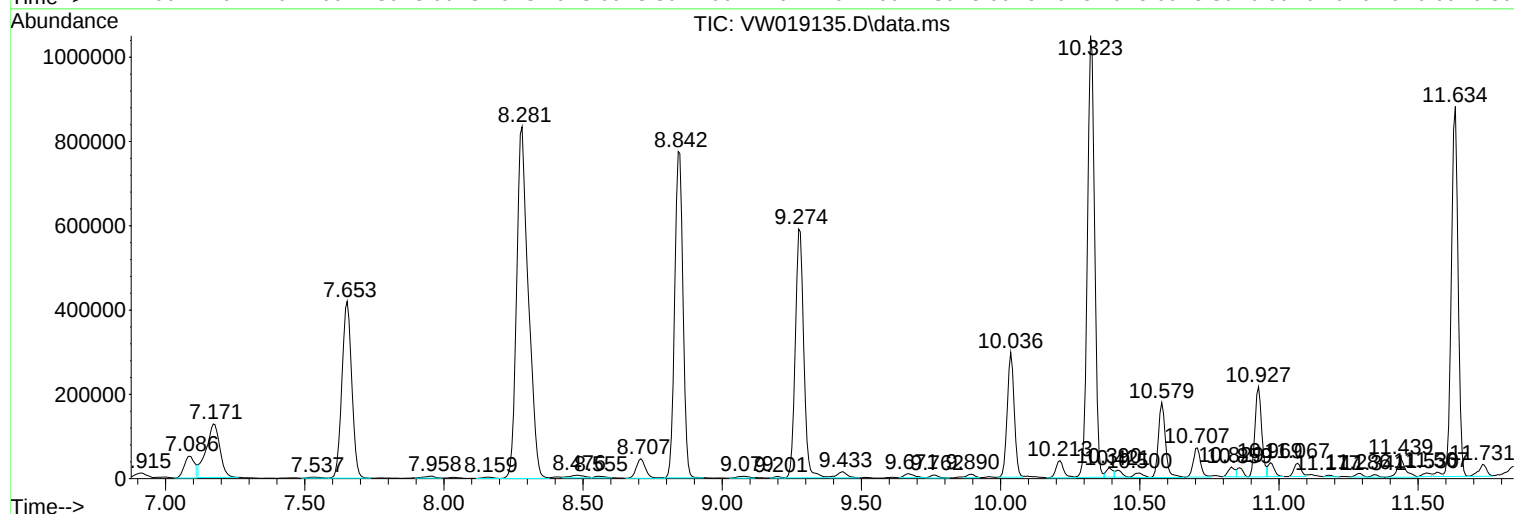
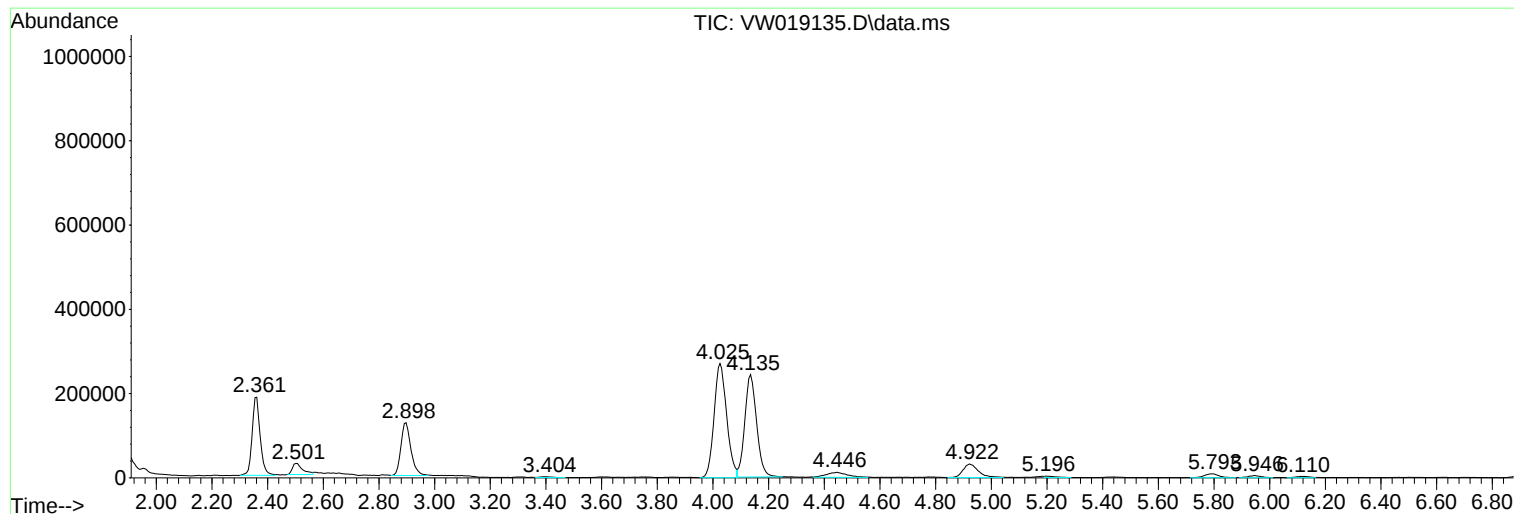
Sum of corrected areas: 19262478

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM060221SMA.M
Quant Title : SFAM01.0

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



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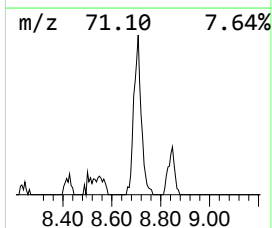
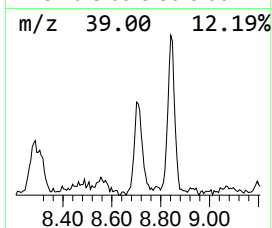
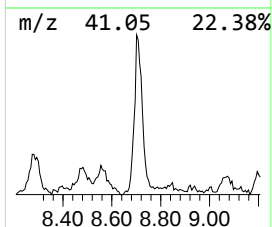
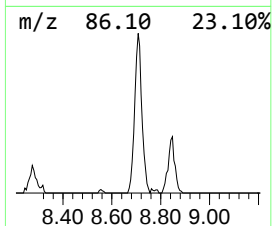
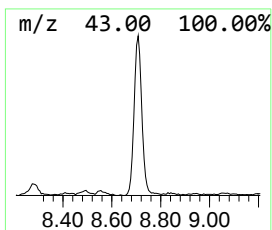
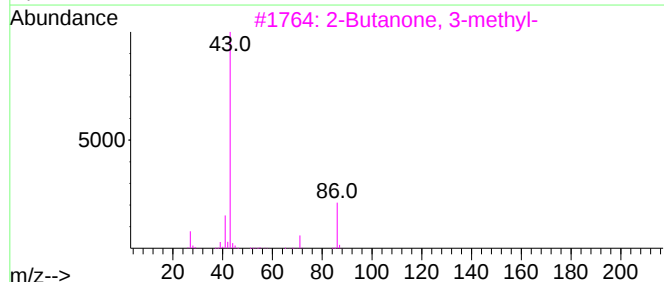
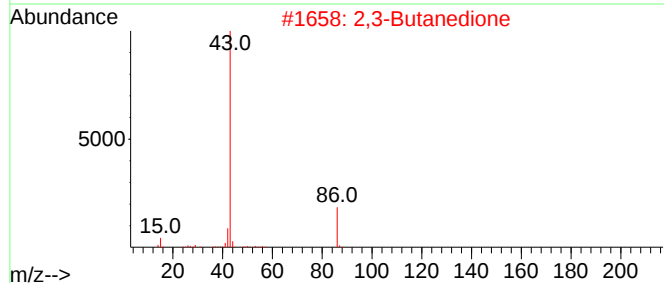
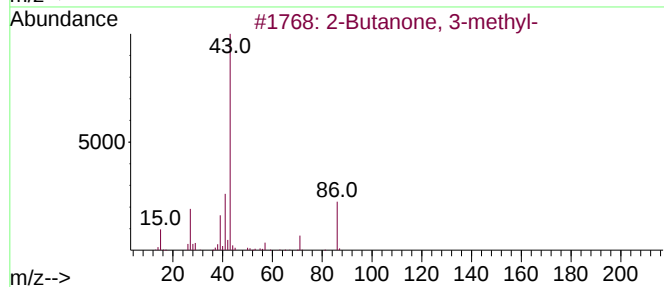
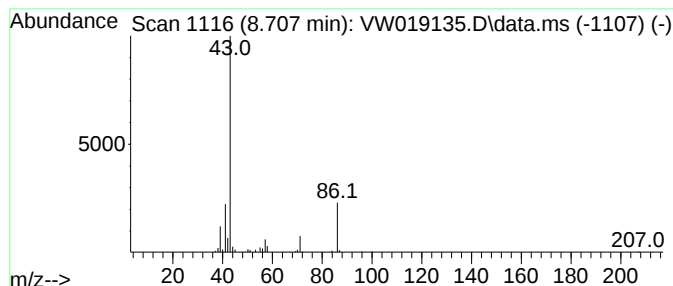
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM060221SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 2-Butanone, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.707	1.67 ug/L	106862	1,4-Difluorobenzene	8.848

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	72
2	2,3-Butanedione			86	C4H6O2	000431-03-8	64
3	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	53
4	2-Pentanone			86	C5H10O	000107-87-9	53
5	2-Pentanone			86	C5H10O	000107-87-9	45



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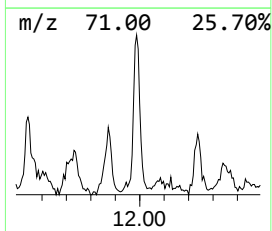
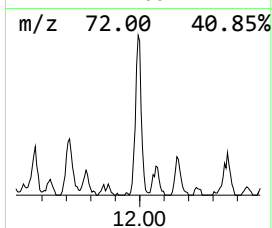
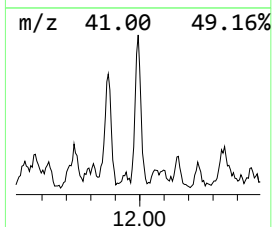
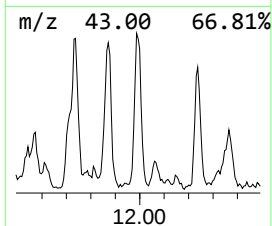
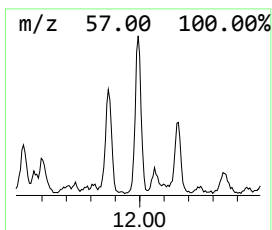
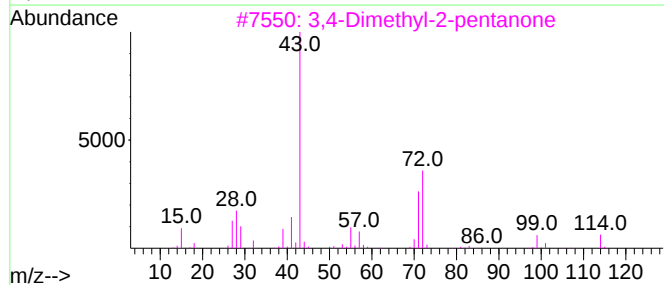
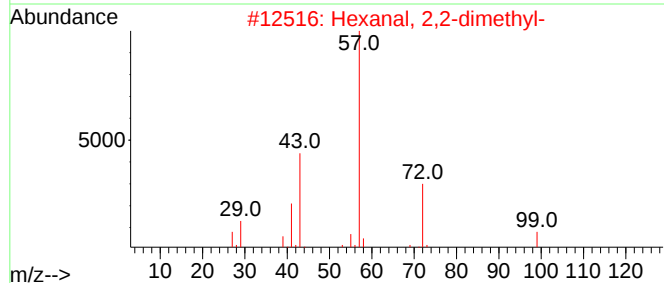
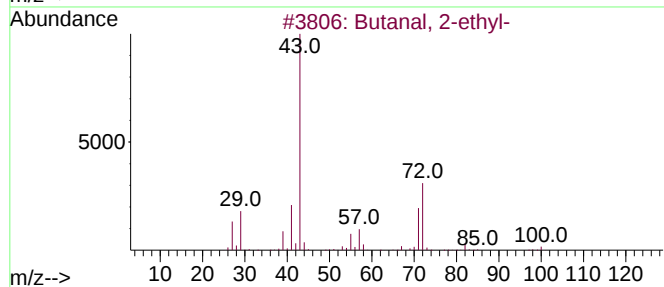
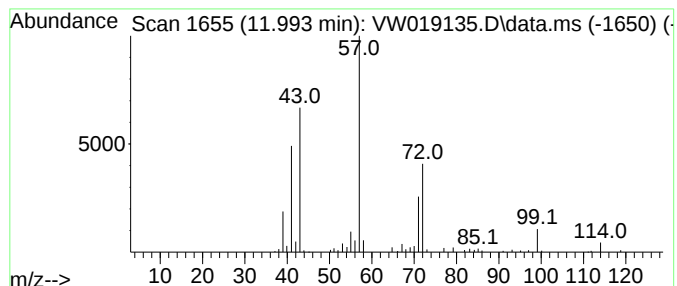
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Quant Title : SFAM01.0

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

Peak Number 5 Butanal, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.993	1.32 ug/L	77493	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanal, 2-ethyl-	100	C6H12O	000097-96-1	53
2			Hexanal, 2,2-dimethyl-	128	C8H16O	000996-12-3	40
3			3,4-Dimethyl-2-pentanone	114	C7H14O	1000202-23-1	38
4			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	38
5			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	35



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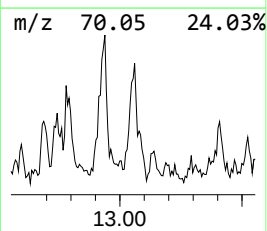
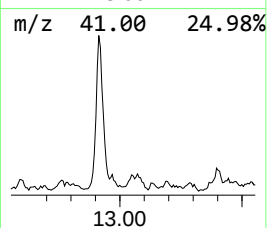
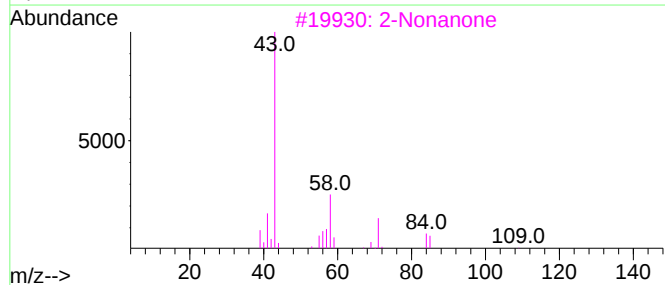
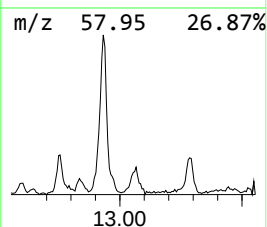
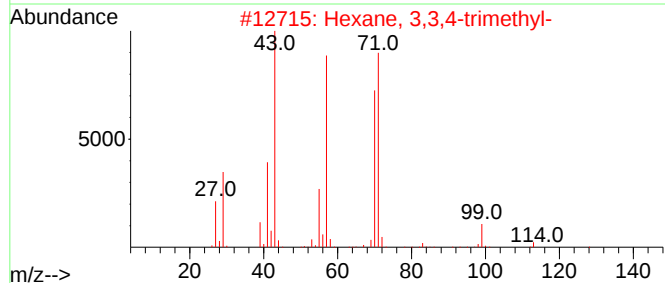
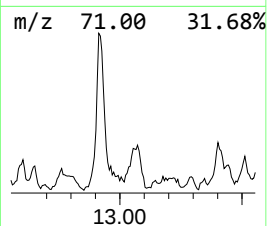
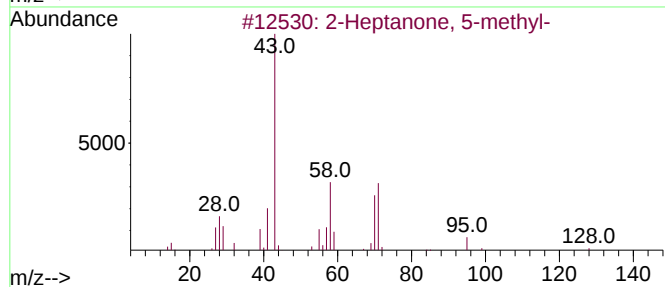
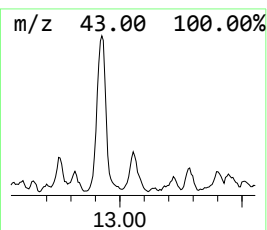
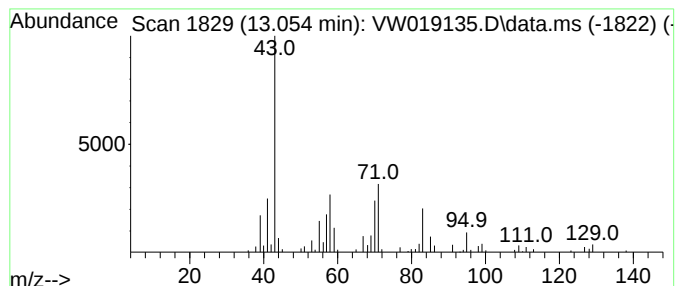
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Quant Title : SFAM01.0

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

Peak Number 7 2-Heptanone, 5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.054	1.71 ug/L	62251	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptanone, 5-methyl-	128	C8H16O	018217-12-4	53
2			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	22
3			2-Nonanone	142	C9H18O	000821-55-6	16
4			Pentane, 3-ethyl-	100	C7H16	000617-78-7	14
5			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	14



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060421\
Data File : VW019135.D
Acq On : 04 Jun 2021 17:53
Operator : SY/VA
Sample : M2618-08
Misc : 4.36G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BG5S6

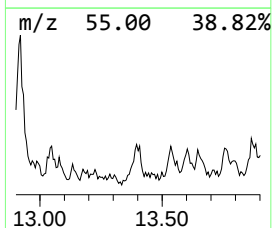
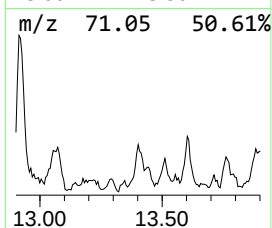
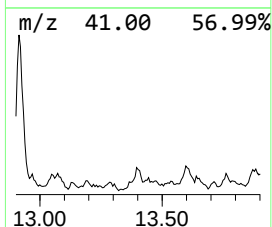
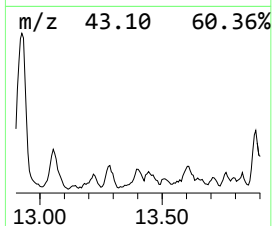
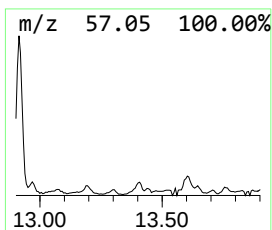
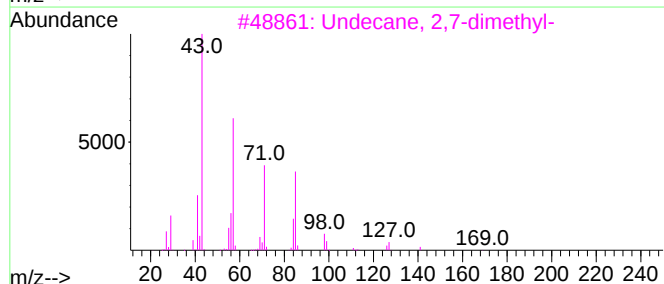
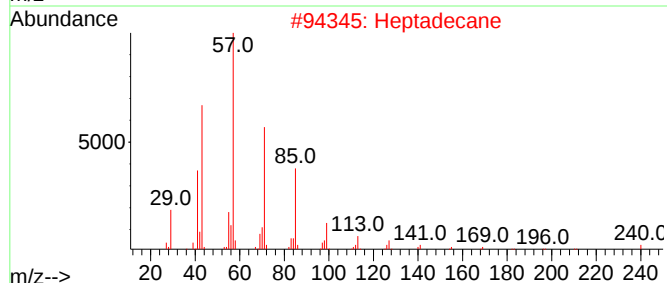
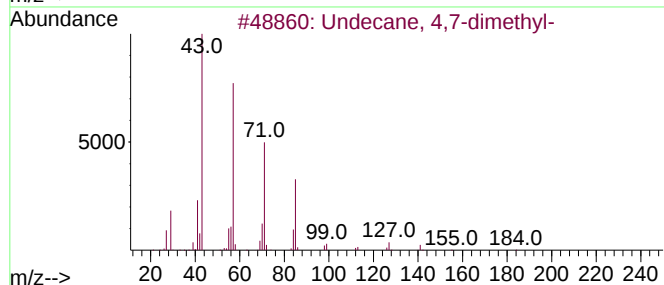
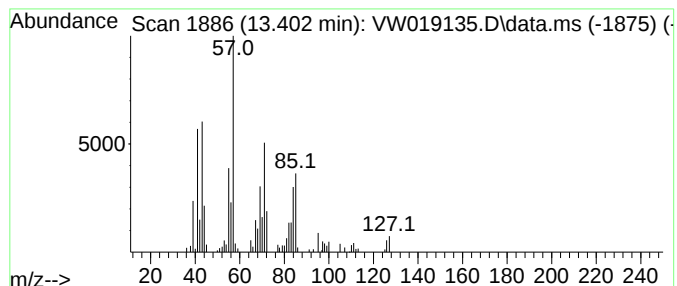
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Quant Title : SFAM01.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.402	2.05 ug/L	74705	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	50
2		Heptadecane	240	C17H36	000629-78-7	47
3		Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	47
4		Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	47
5		Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060421\
Data File : VW019135.D
Acq On : 04 Jun 2021 17:53
Operator : SY/VA
Sample : M2618-08
Misc : 4.36G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BG5S6

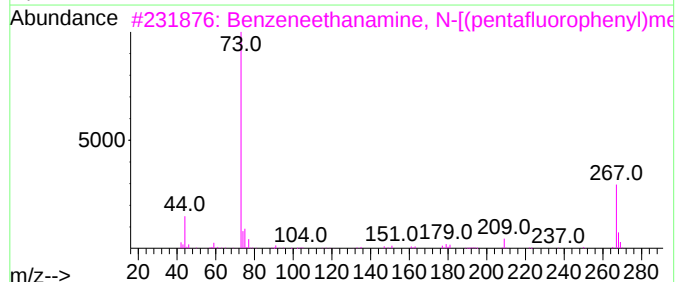
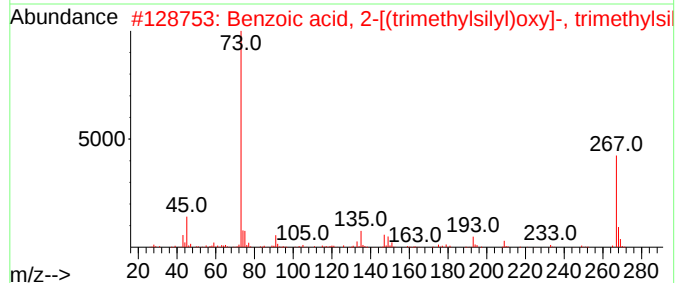
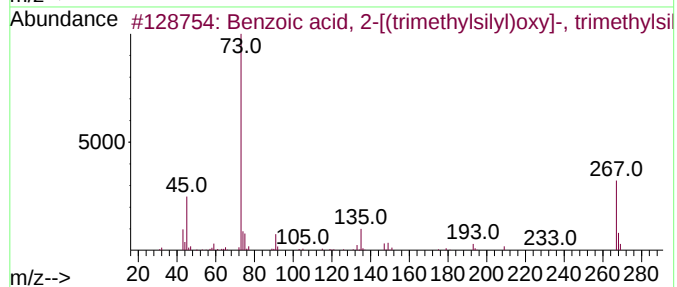
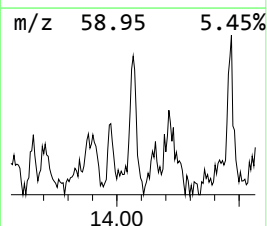
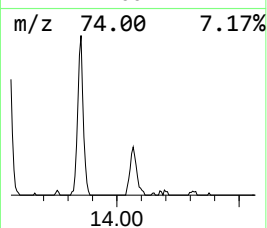
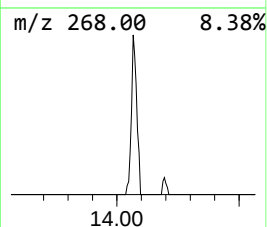
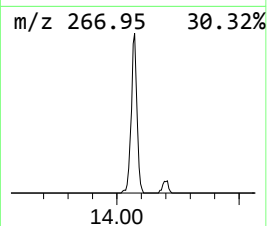
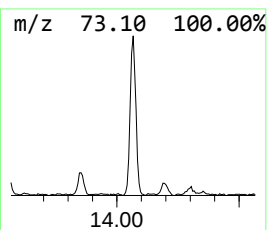
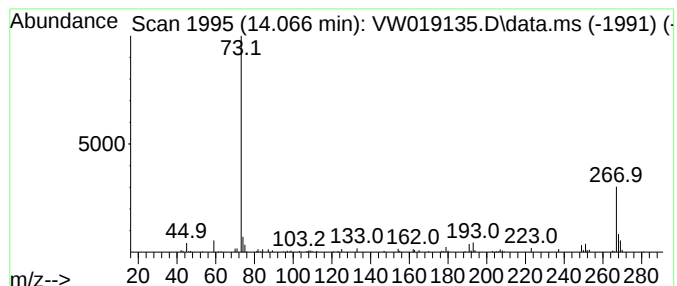
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Quant Title : SFAM01.0

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

Peak Number 9 Benzoic acid, 2-[(trimethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.066	2.53 ug/L	92175	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2-[(trimethylsilyl...]	282	C13H22O3Si2	003789-85-3	53
2			Benzoic acid, 2-[(trimethylsilyl...]	282	C13H22O3Si2	003789-85-3	45
3			Benzenethanamine, N-[(pentaflu...]	475	C21H26F5NO2Si2	055429-85-1	40
4			Benzoic acid, 2-[(trimethylsilyl...]	282	C13H22O3Si2	003789-85-3	38
5			Benzoic acid, 2-[(trimethylsilyl...]	282	C13H22O3Si2	003789-85-3	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060421\
Data File : VW019135.D
Acq On : 04 Jun 2021 17:53
Operator : SY/VA
Sample : M2618-08
Misc : 4.36G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BG5S6

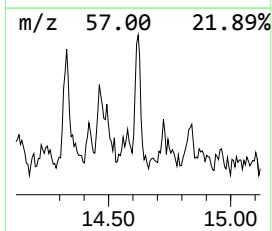
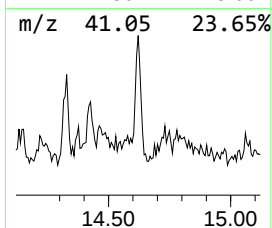
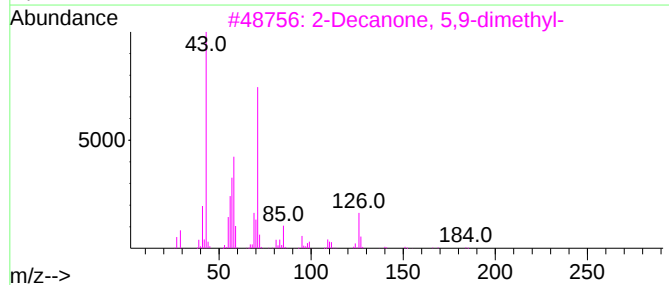
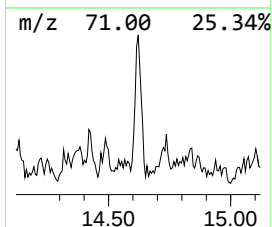
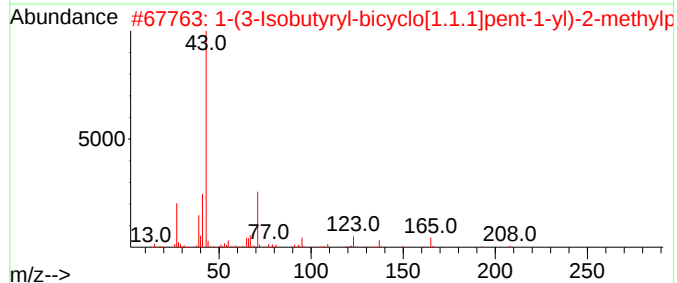
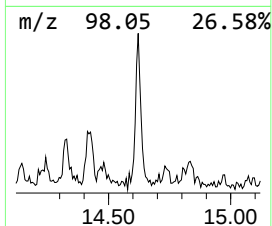
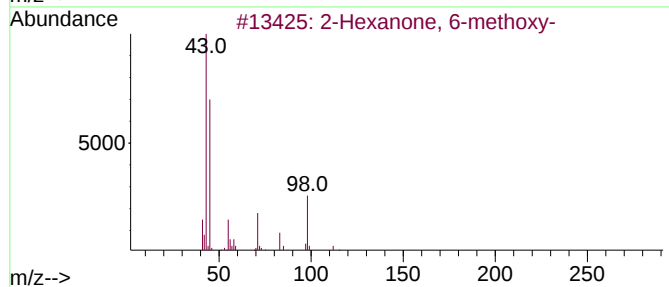
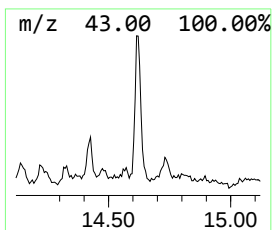
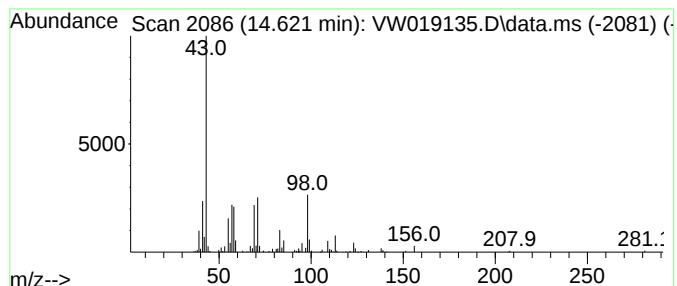
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.621	2.08 ug/L	75930	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone, 6-methoxy-		130	C7H14O2		029006-00-6	38
2	1-(3-Isobutyryl-bicyclo[1.1.1]pe...		208	C13H20O2		1000287-95-8	25
3	2-Decanone, 5,9-dimethyl-		184	C12H24O		033933-82-3	25
4	9-Octadecanone		268	C18H36O		018394-00-8	25
5	4-Decanone		156	C10H20O		000624-16-8	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060421\
Data File : VW019135.D
Acq On : 04 Jun 2021 17:53
Operator : SY/VA
Sample : M2618-08
Misc : 4.36G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BG5S6

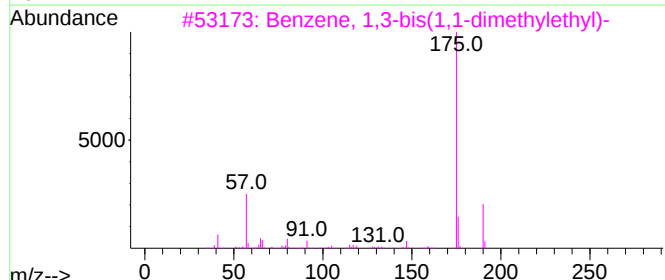
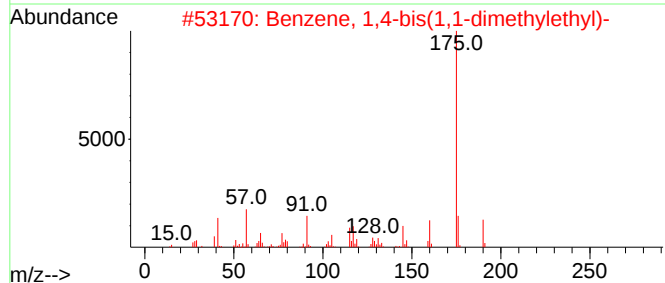
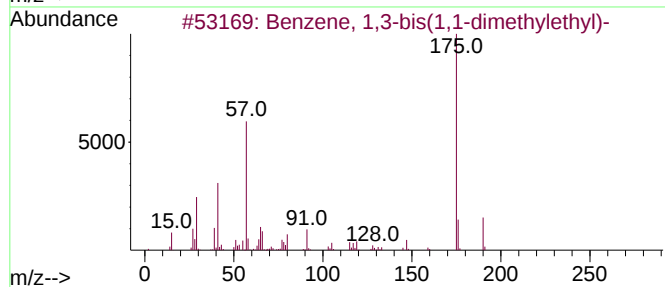
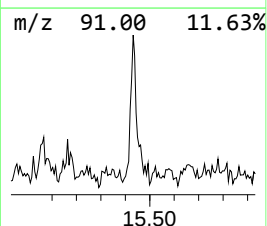
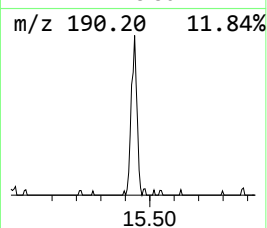
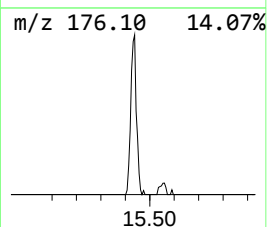
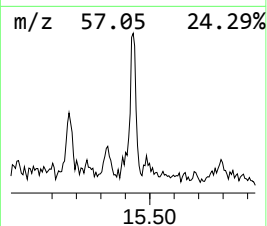
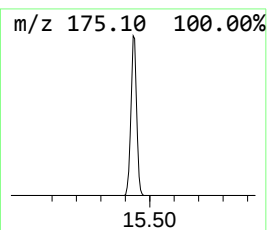
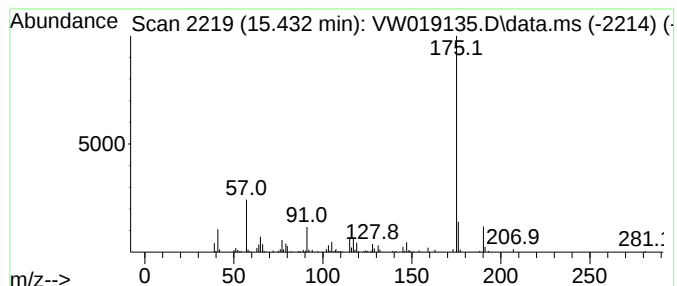
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Quant Title : SFAM01.0

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

Peak Number 11 Benzene, 1,3-bis(1,1-dimeth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.432	1.74 ug/L	63383	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	86
2			Benzene, 1,4-bis(1,1-dimethyleth...	190	C14H22	001012-72-2	86
3			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	86
4			Benzene, 1,4-bis(1,1-dimethyleth...	190	C14H22	001012-72-2	72
5			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060421\
Data File : VW019135.D
Acq On : 04 Jun 2021 17:53
Operator : SY/VA
Sample : M2618-08
Misc : 4.36G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BG5S6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM060221SMA.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
2-Butanone, 3-m...	8.707	1.7	ug/L	106862	1	8.848	1597240	25.0
Butanal, 2-ethyl-	11.993	1.3	ug/L	77493	2	11.634	1472870	25.0
2-Heptanone, 5-...	13.054	1.7	ug/L	62251	3	13.560	912364	25.0
(DEL) Alkane: S...	13.402	2.0	ug/L	74705	3	13.560	912364	25.0
Benzoic acid, 2...	14.066	2.5	ug/L	92175	3	13.560	912364	25.0
unknown-01	14.621	2.1	ug/L	75930	3	13.560	912364	25.0
Benzene, 1,3-bi...	15.432	1.7	ug/L	63383	3	13.560	912364	25.0