

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
 Data File : VW021128.D
 Acq On : 04 Dec 2021 19:16
 Operator : SY/VA
 Sample : M4885-14
 Misc : 3.93g/10.0mL/MSVOA_W/SOIL
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EW9M1

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\SFAMWLM120321SMA.M
 Title : SFAM01.0

Signal : TIC: VW021128.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.355	68	74	84	rVB	47179	76255	0.60%	0.318%
2	2.885	154	161	173	rVB	25083	54941	0.43%	0.229%
3	3.391	239	244	249	rVB3	695	1422	0.01%	0.006%
4	3.904	326	328	331	rBV	122	155	0.00%	0.001%
5	4.019	337	347	359	rBV	46913	133344	1.05%	0.557%
6	4.129	360	365	373	rVB	1267	2734	0.02%	0.011%
7	4.251	382	385	386	rBV	182	141	0.00%	0.001%
8	4.385	396	407	422	rVB	11153	33329	0.26%	0.139%
9	4.635	446	448	452	rBV	84	151	0.00%	0.001%
10	4.916	484	494	511	rVV5	5374	19376	0.15%	0.081%
11	5.105	523	525	527	rBV2	235	305	0.00%	0.001%
12	5.355	565	566	570	rVB	146	136	0.00%	0.001%
13	5.434	570	579	587	rBV3	2348	5970	0.05%	0.025%
14	5.940	656	662	671	rVB5	1182	3144	0.02%	0.013%
15	6.147	692	696	699	rBV	113	179	0.00%	0.001%
16	6.184	699	702	705	rVB	147	164	0.00%	0.001%
17	6.452	743	746	747	rVB	145	132	0.00%	0.001%
18	6.738	790	793	796	rBB	141	181	0.00%	0.001%
19	6.982	826	833	841	rBV3	2334	5533	0.04%	0.023%
20	7.080	841	849	854	rBV	9279	25832	0.20%	0.108%
21	7.653	933	943	955	rBV	62767	157621	1.24%	0.658%
22	7.952	984	992	999	rBV5	1982	5225	0.04%	0.022%
23	8.031	1002	1005	1009	rVB3	509	709	0.01%	0.003%
24	8.153	1020	1025	1027	rBV2	602	648	0.01%	0.003%
25	8.275	1032	1045	1060	rBV3	111338	351281	2.76%	1.466%
26	8.506	1079	1083	1089	rVB4	934	1654	0.01%	0.007%
27	8.580	1089	1095	1100	rVB5	1380	2924	0.02%	0.012%
28	8.695	1111	1114	1117	rVB	385	425	0.00%	0.002%
29	8.842	1131	1138	1147	rVV	59866	116685	0.92%	0.487%
30	9.092	1168	1179	1198	rBV	6794350	12719743	100.00%	53.097%
31	9.275	1201	1209	1215	rBV	73169	145735	1.15%	0.608%
32	9.433	1231	1235	1236	rVV2	94	132	0.00%	0.001%
33	9.518	1246	1249	1253	rBB4	493	786	0.01%	0.003%
34	9.616	1260	1265	1268	rBB2	372	662	0.01%	0.003%
35	9.695	1276	1278	1280	rBV	135	133	0.00%	0.001%
36	9.750	1283	1287	1292	rVB3	652	966	0.01%	0.004%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.M
 Title : SFAM01.0

37	9.890	1307	1310	1316	rBV4	611	888	0.01%	0.004%
38	9.957	1316	1321	1327	rVB3	495	812	0.01%	0.003%
39	10.037	1327	1334	1342	rBV	27695	49148	0.39%	0.205%
40	10.207	1358	1362	1367	rBV3	806	1155	0.01%	0.005%
41	10.250	1367	1369	1371	rVB	341	217	0.00%	0.001%
42	10.323	1372	1381	1387	rBV	98488	171550	1.35%	0.716%
43	10.390	1387	1392	1399	rVB	19265	33928	0.27%	0.142%
44	10.500	1403	1410	1414	rBV5	863	1428	0.01%	0.006%
45	10.579	1417	1423	1429	rVB	12656	20448	0.16%	0.085%
46	10.671	1433	1438	1439	rBV3	998	1717	0.01%	0.007%
47	10.701	1439	1443	1450	rVB2	7422	12887	0.10%	0.054%
48	10.786	1451	1457	1461	rBV2	4607	7246	0.06%	0.030%
49	10.866	1461	1470	1477	rBV	5727227	9516172	74.81%	39.724%
50	11.067	1497	1503	1510	rVB	3442	5374	0.04%	0.022%
51	11.177	1515	1521	1525	rVV3	1263	2046	0.02%	0.009%
52	11.232	1525	1530	1534	rVB2	1419	2301	0.02%	0.010%
53	11.299	1540	1541	1544	rVB2	169	132	0.00%	0.001%
54	11.341	1544	1548	1552	rBB2	326	574	0.00%	0.002%
55	11.439	1558	1564	1571	rBB	7677	12442	0.10%	0.052%
56	11.628	1589	1595	1602	rBV	35434	60614	0.48%	0.253%
57	11.731	1606	1612	1622	rBB2	3971	6030	0.05%	0.025%
58	11.835	1623	1629	1637	rBB2	5815	9962	0.08%	0.042%
59	11.902	1637	1640	1641	rBV	129	129	0.00%	0.001%
60	12.164	1677	1683	1690	rBB2	5518	8618	0.07%	0.036%
61	12.292	1701	1704	1708	rBB	133	216	0.00%	0.001%
62	12.335	1709	1711	1712	rBV	419	333	0.00%	0.001%
63	12.463	1728	1732	1735	rBV2	1043	1411	0.01%	0.006%
64	12.603	1749	1755	1763	rBB2	23501	37417	0.29%	0.156%
65	12.689	1763	1769	1777	rBV	30946	51764	0.41%	0.216%
66	12.768	1780	1782	1784	rVV	181	195	0.00%	0.001%
67	12.804	1784	1788	1792	rVB	637	985	0.01%	0.004%
68	12.865	1793	1798	1801	rBV2	501	606	0.00%	0.003%
69	12.926	1806	1808	1809	rBV	639	456	0.00%	0.002%
70	13.006	1818	1821	1825	rBB	137	146	0.00%	0.001%
71	13.103	1834	1837	1841	rVV2	365	561	0.00%	0.002%
72	13.195	1845	1852	1855	rBB3	575	1006	0.01%	0.004%
73	13.249	1855	1861	1864	rBV2	977	1605	0.01%	0.007%
74	13.292	1864	1868	1874	rVB2	3127	4700	0.04%	0.020%
75	13.408	1881	1887	1890	rBB2	763	1187	0.01%	0.005%
76	13.469	1894	1897	1898	rBB2	209	144	0.00%	0.001%

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 Title : SFAM01.0

77	13.505	1899	1903	1906	rBV2	382	469	0.00%	0.002%
78	13.554	1906	1911	1916	rBV	8933	14086	0.11%	0.059%
79	13.646	1925	1926	1929	rBV	307	286	0.00%	0.001%
80	13.762	1941	1945	1949	rBB3	506	830	0.01%	0.003%
81	13.798	1950	1951	1954	rBV	110	130	0.00%	0.001%
82	13.847	1954	1959	1965	rVV	10844	18798	0.15%	0.078%
83	13.902	1965	1968	1972	rVV	343	527	0.00%	0.002%
84	13.944	1972	1975	1978	rVV3	288	396	0.00%	0.002%
85	14.066	1989	1995	2002	rBB2	6846	11288	0.09%	0.047%
86	14.585	2079	2080	2085	rBB	128	161	0.00%	0.001%
87	14.719	2096	2102	2106	rBV3	1193	1826	0.01%	0.008%
88	14.822	2117	2119	2123	rVB	143	188	0.00%	0.001%
89	14.956	2139	2141	2144	rVB	139	145	0.00%	0.001%
90	15.042	2153	2155	2157	rBV	136	133	0.00%	0.001%
91	15.237	2183	2187	2189	rBV2	142	233	0.00%	0.001%
92	15.298	2194	2197	2200	rBV2	230	331	0.00%	0.001%
93	15.359	2203	2207	2213	rBB3	430	929	0.01%	0.004%
94	15.432	2213	2219	2223	rBV	2829	4524	0.04%	0.019%
95	15.487	2223	2228	2235	rVB3	1235	2155	0.02%	0.009%
96	16.121	2328	2332	2333	rBV	164	142	0.00%	0.001%
97	16.139	2333	2335	2338	rVB	145	138	0.00%	0.001%
98	16.212	2345	2347	2348	rBV	163	130	0.00%	0.001%
99	16.432	2381	2383	2386	rVB	351	374	0.00%	0.002%
100	16.639	2412	2417	2420	rVV2	227	358	0.00%	0.001%

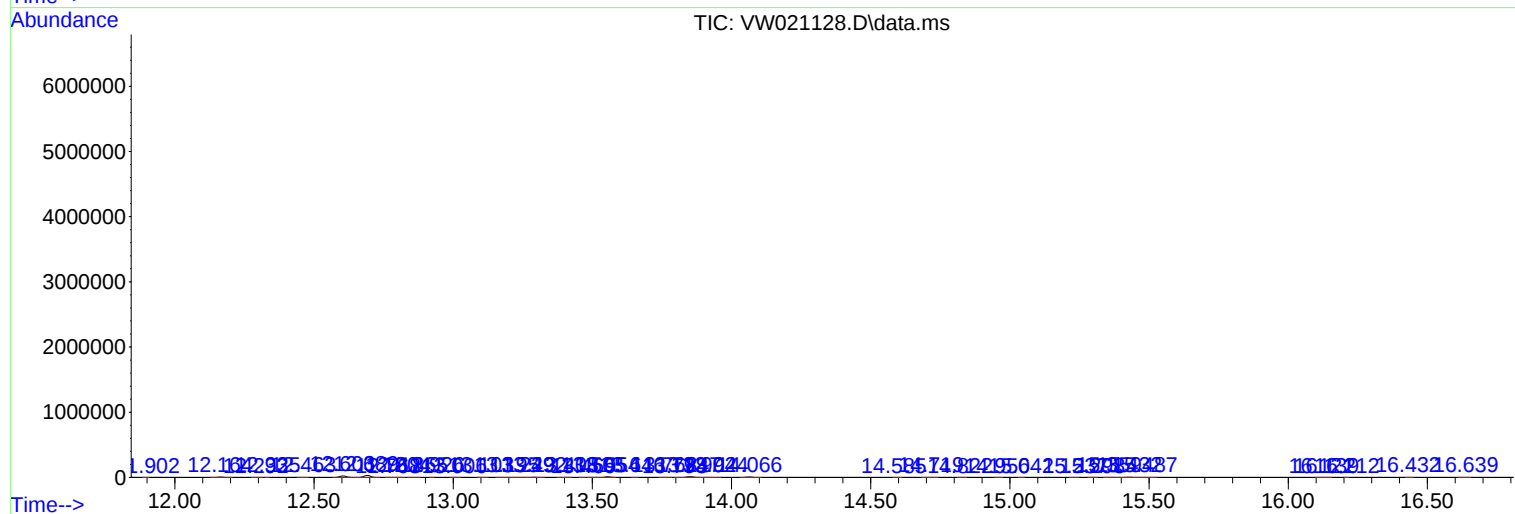
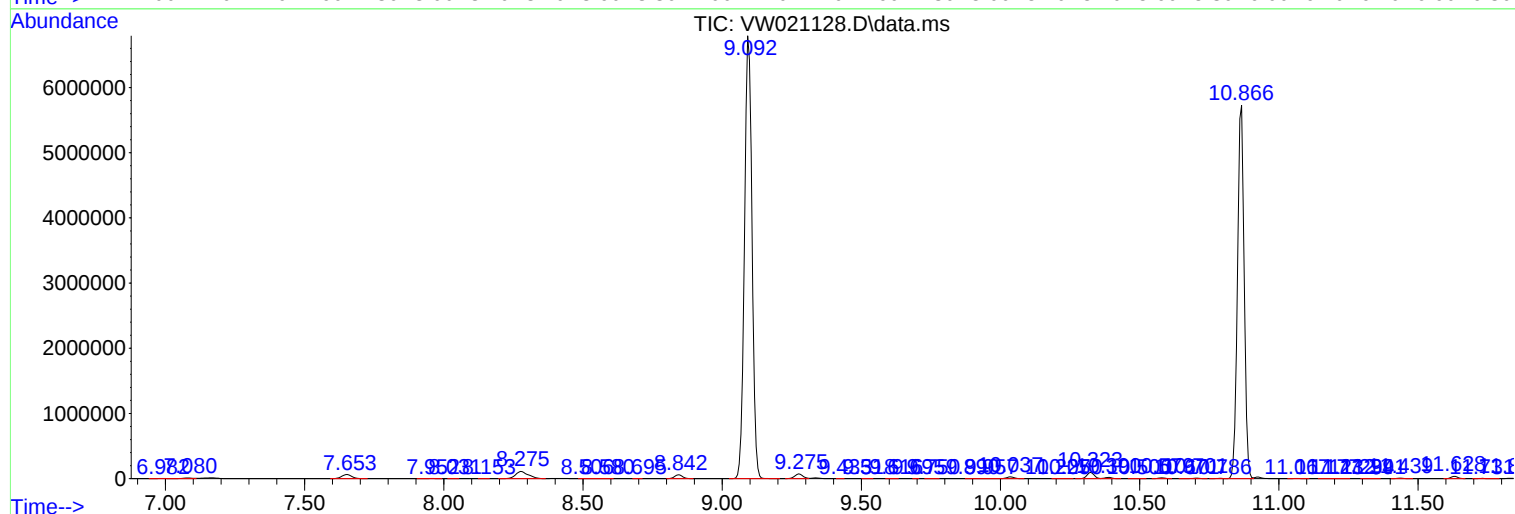
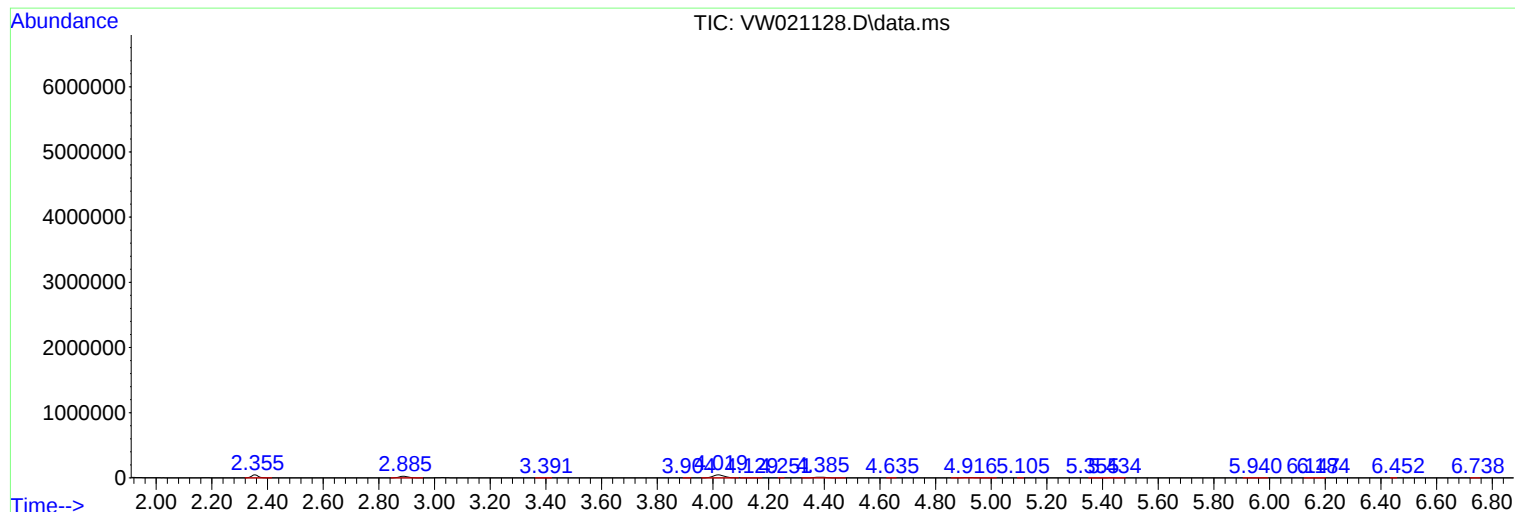
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Instrument :
MSVOA_W
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.M
Quant Title : SFAM01.0

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



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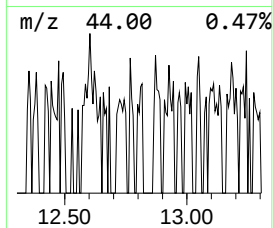
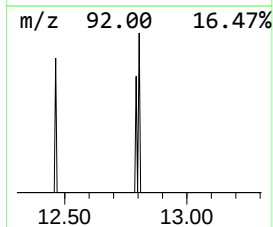
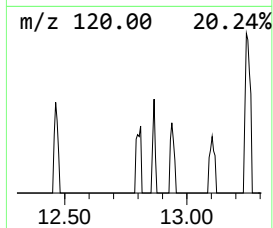
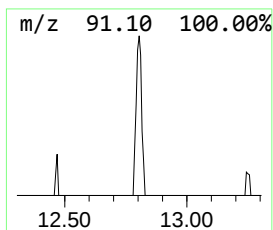
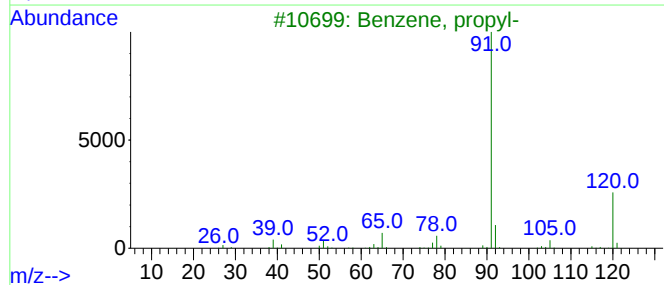
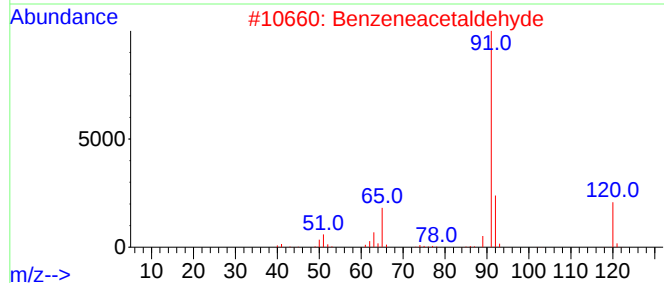
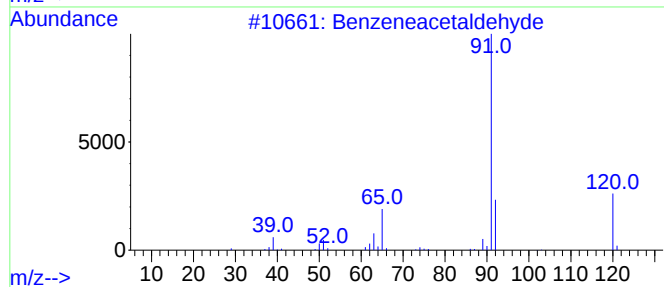
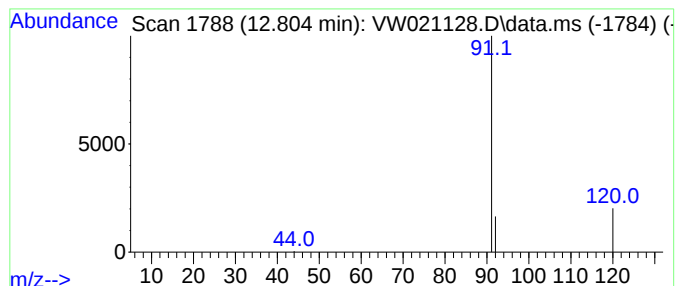
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.M
Quant Title : SFAM01.0

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

Peak Number 6 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.804	1.75 ug/L	985	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetaldehyde	120	C8H8O	000122-78-1	9
2			Benzeneacetaldehyde	120	C8H8O	000122-78-1	9
3			Benzene, propyl-	120	C9H12	000103-65-1	7
4			1,3,5-Cycloheptatriene, 7-ethyl-	120	C9H12	017634-51-4	7
5			Benzeneacetaldehyde	120	C8H8O	000122-78-1	4



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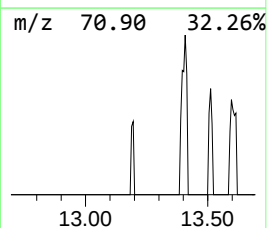
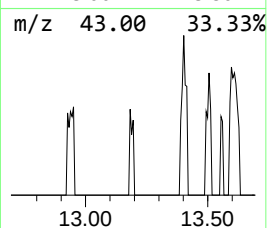
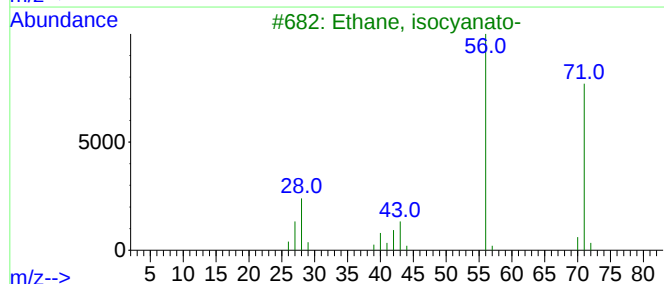
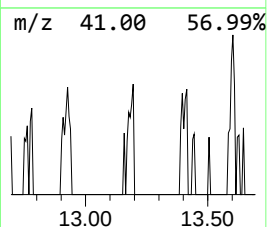
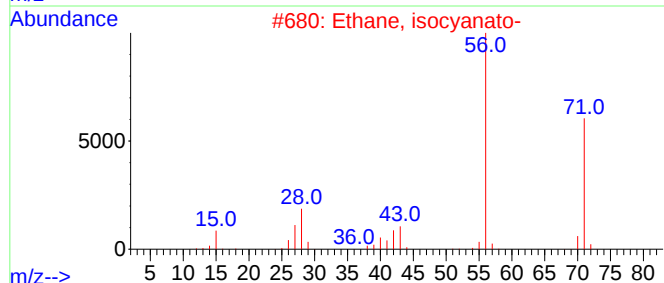
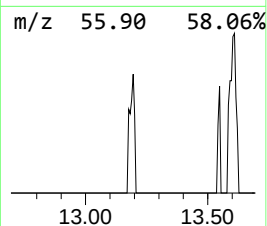
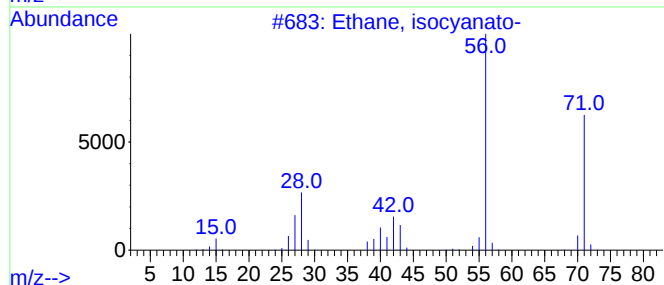
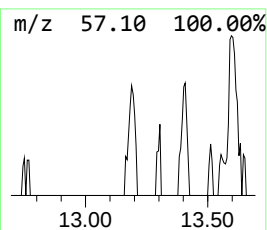
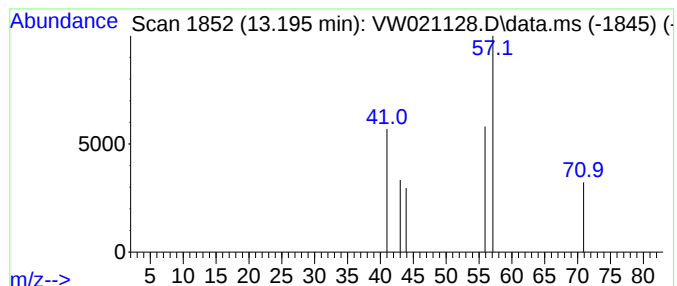
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.M
Quant Title : SFAM01.0

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

Peak Number 7 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.195	1.79 ug/L	1006	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, isocyanato-	71	C3H5NO	000109-90-0	4
2			Ethane, isocyanato-	71	C3H5NO	000109-90-0	4
3			Ethane, isocyanato-	71	C3H5NO	000109-90-0	4
4			Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	4
5			Azetidine	57	C3H7N	000503-29-7	4



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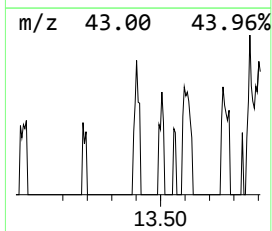
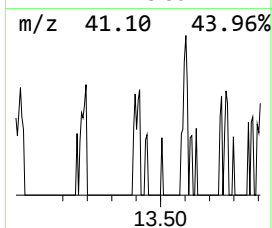
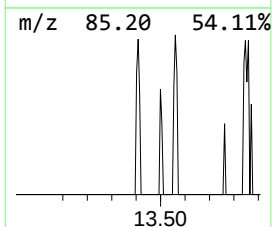
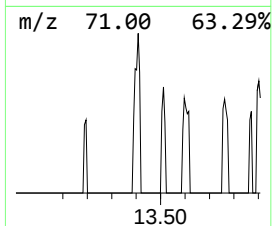
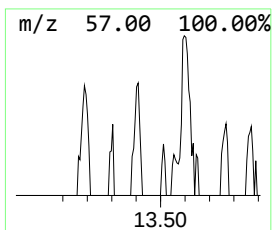
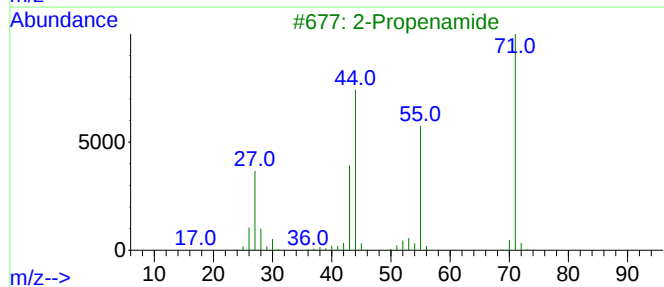
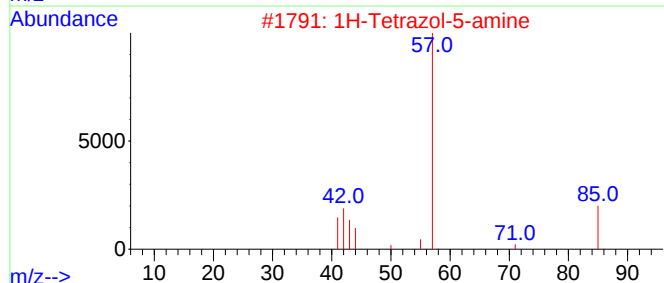
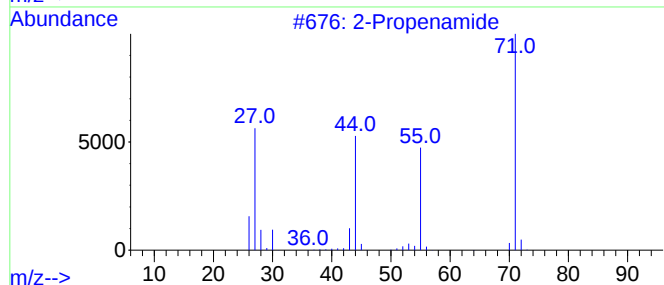
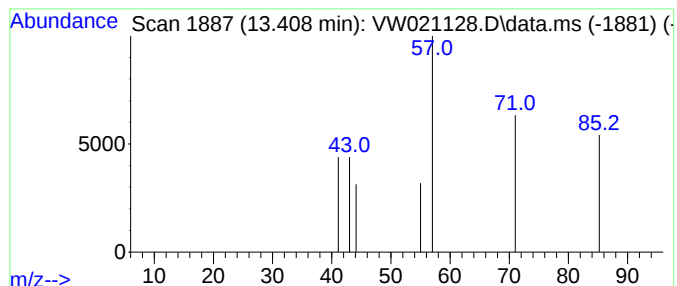
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.M
Quant Title : SFAM01.0

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

Peak Number 8 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.408	2.11 ug/L	1187	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenamide			71	C3H5NO	000079-06-1	9
2	1H-Tetrazol-5-amine			85	CH3N5	004418-61-5	7
3	2-Propenamide			71	C3H5NO	000079-06-1	7
4	Azetidine, 2-methyl-			71	C4H9N	019812-49-8	5
5	Azetidine, 1-methyl-			71	C4H9N	004923-79-9	5



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
Data File : VW021128.D
Acq On : 04 Dec 2021 19:16
Operator : SY/VA
Sample : M4885-14
Misc : 3.93g/10.0mL/MSVOA_W/SOIL
ALS Vial : 18 Sample Multiplier: 1

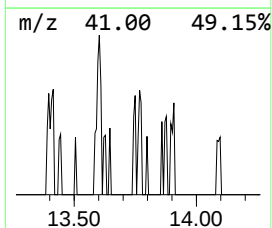
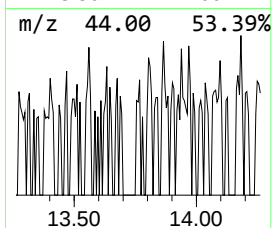
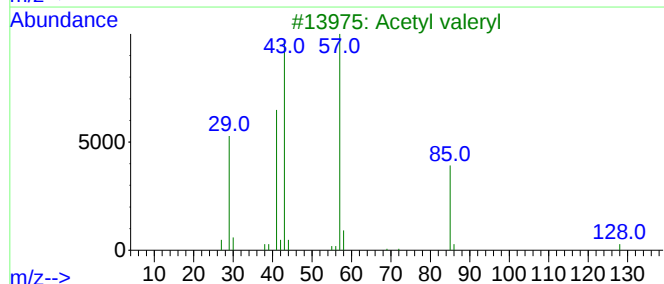
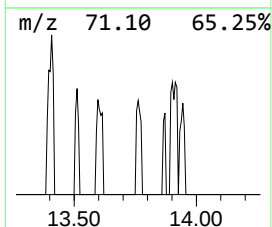
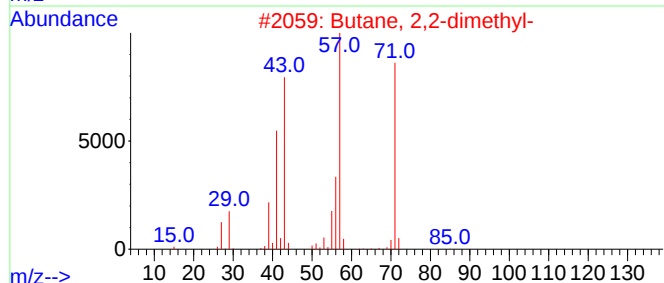
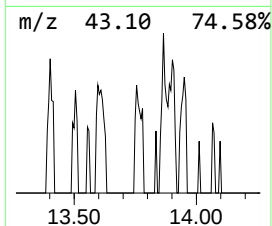
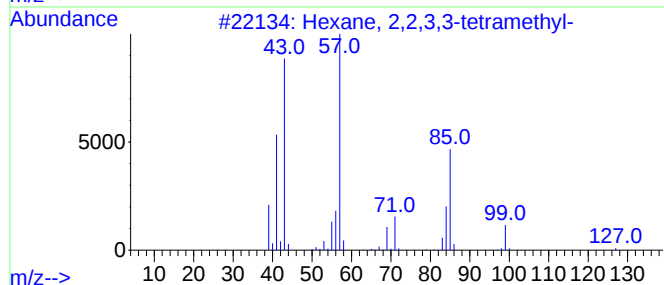
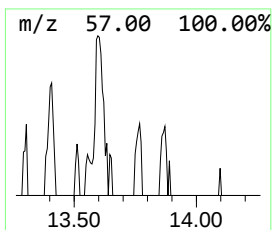
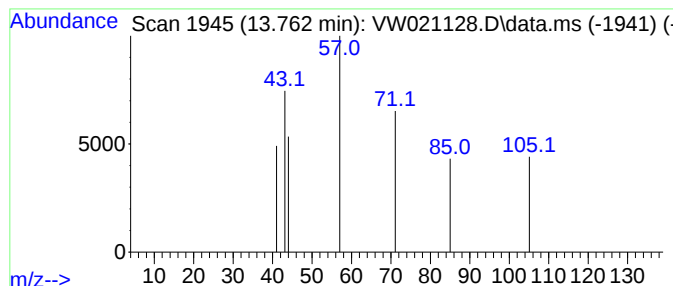
Instrument :
MSVOA_W
ClientSampleId :
EW9M1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.M
Quant Title : SFAM01.0

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.762	1.47 ug/L	830	1,4-Dichlorobenzene-d4		13.560	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexane, 2,2,3,3-tetramethyl-		142	C10H22	013475-81-5	9
2	Butane, 2,2-dimethyl-		86	C6H14	000075-83-2	9
3	Acetyl valeryl		128	C7H12O2	000096-04-8	9
4	Butane, 2,2-dimethyl-		86	C6H14	000075-83-2	9
5	Hexane, 3,3-dimethyl-		114	C8H18	000563-16-6	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
Data File : VW021128.D
Acq On : 04 Dec 2021 19:16
Operator : SY/VA
Sample : M4885-14
Misc : 3.93g/10.0mL/MSVOA_W/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW9M1

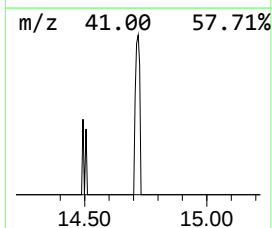
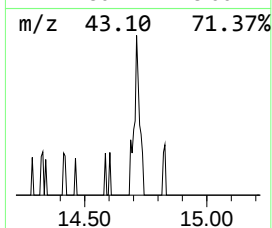
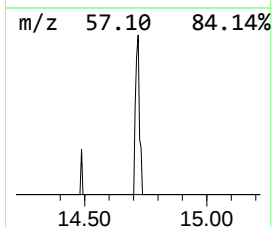
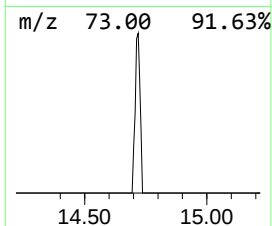
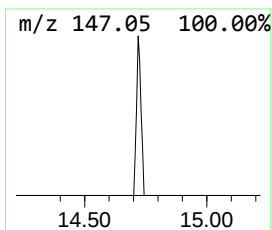
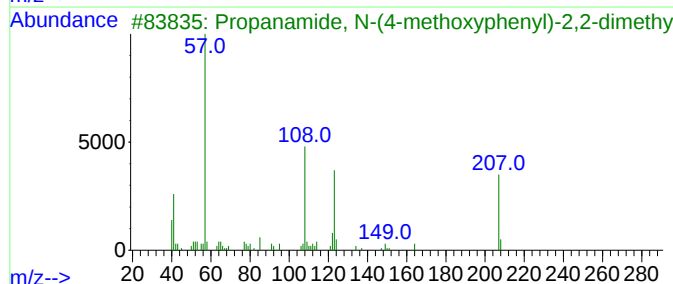
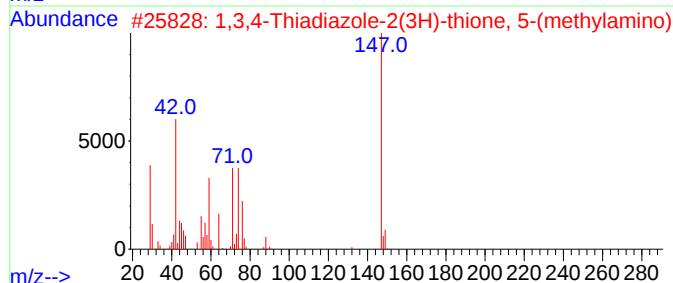
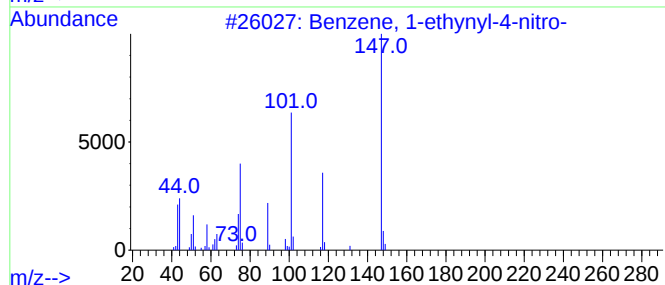
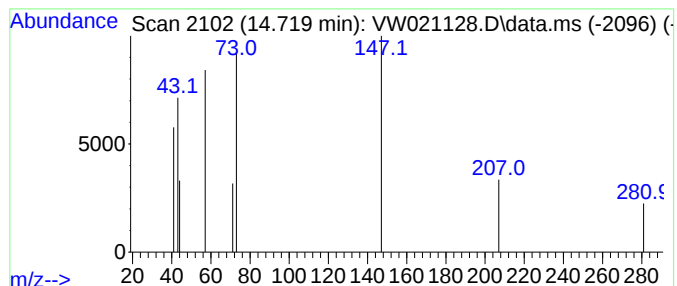
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Quant Title : SFAM01.0

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

Peak Number 11 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.719	3.24 ug/L	1826	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethynyl-4-nitro-	147	C8H5NO2	000937-31-5	4
2			1,3,4-Thiadiazole-2(3H)-thione, ...	147	C3H5N3S2	027386-01-2	4
3			Propanamide, N-(4-methoxyphenyl)...	207	C12H17NO2	056619-94-4	4
4			Propanamide, N-(3-methoxyphenyl)...	207	C12H17NO2	056619-93-3	4
5			o-Anisidine, N-trimethylacetyl-	207	C12H17NO2	033768-49-9	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
Data File : VW021128.D
Acq On : 04 Dec 2021 19:16
Operator : SY/VA
Sample : M4885-14
Misc : 3.93g/10.0mL/MSVOA_W/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW9M1

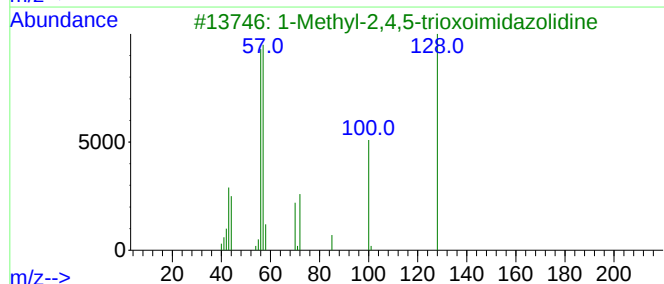
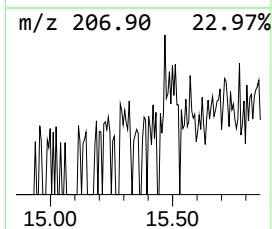
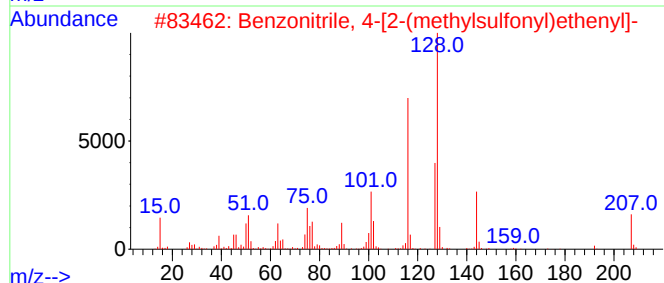
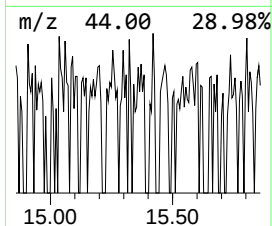
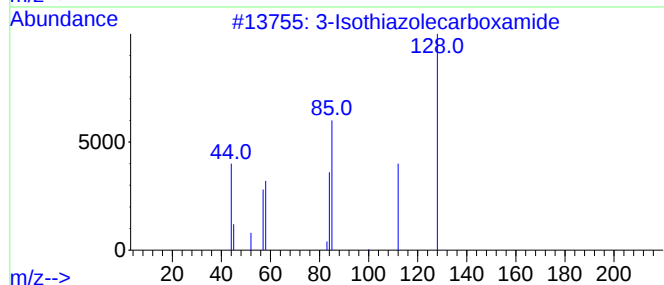
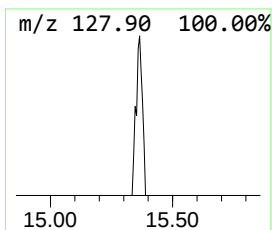
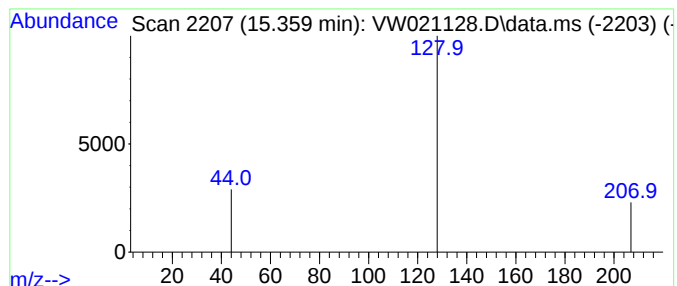
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Quant Title : SFAM01.0

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

Peak Number 12 unknown-05 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.359	1.65 ug/L	929	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Isothiazolecarboxamide	128	C4H4N2OS	024342-43-6	5
2			Benzonitrile, 4-[2-(methylsulfonyl)ethenyl]-	207	C10H9NO2S	077355-31-8	4
3			1-Methyl-2,4,5-trioximidazolidine	128	C4H4N2O3	003659-97-0	4
4			2,4,5-Trihydroxypyrimidine	128	C4H4N2O3	000496-76-4	4
5			5-Hydroxyuracil	128	C4H4N2O3	1000489-22-6	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
Data File : VW021128.D
Acq On : 04 Dec 2021 19:16
Operator : SY/VA
Sample : M4885-14
Misc : 3.93g/10.0mL/MSVOA_W/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW9M1

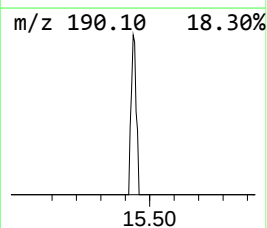
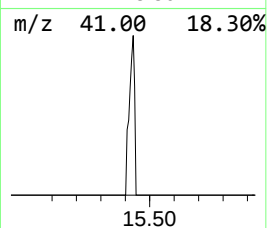
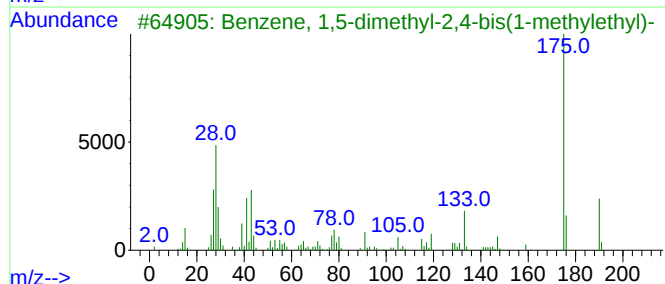
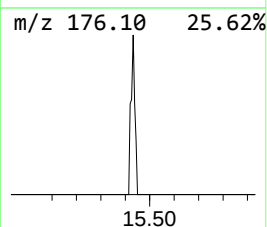
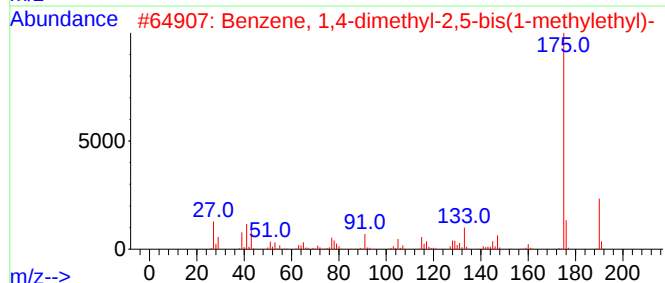
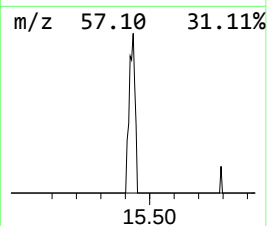
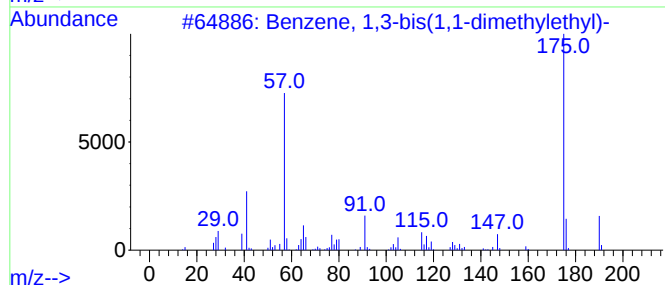
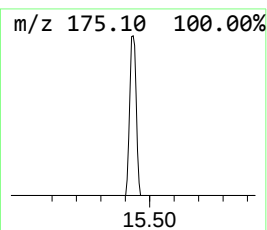
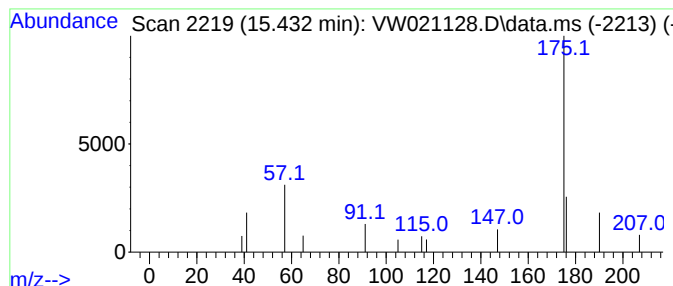
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Quant Title : SFAM01.0

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

Peak Number 13 Benzene, 1,3-bis(1,1-dimeth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.432	8.03 ug/L	4524	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	72
2			Benzene, 1,4-dimethyl-2,5-bis(1-...	190	C14H22	010375-96-9	50
3			Benzene, 1,5-dimethyl-2,4-bis(1-...	190	C14H22	005186-68-5	50
4			Benzenepropanal, 4-(1,1-dimethyl...	190	C13H18O	018127-01-0	50
5			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
Data File : VW021128.D
Acq On : 04 Dec 2021 19:16
Operator : SY/VA
Sample : M4885-14
Misc : 3.93g/10.0mL/MSVOA_W/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW9M1

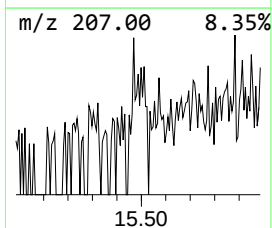
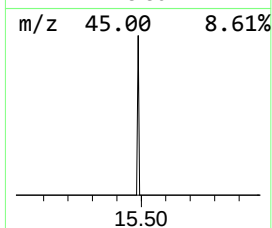
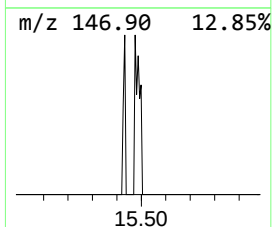
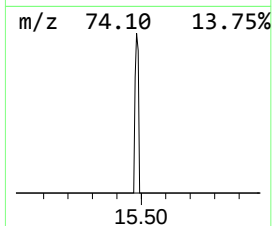
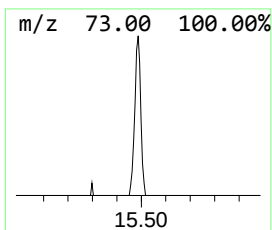
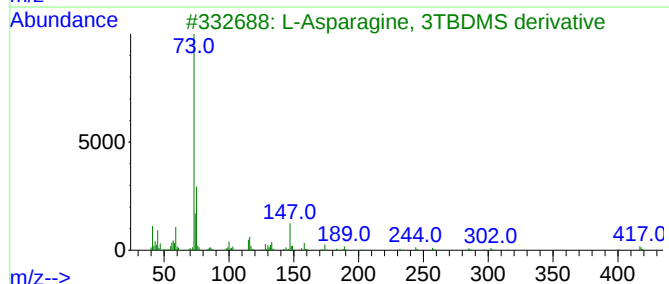
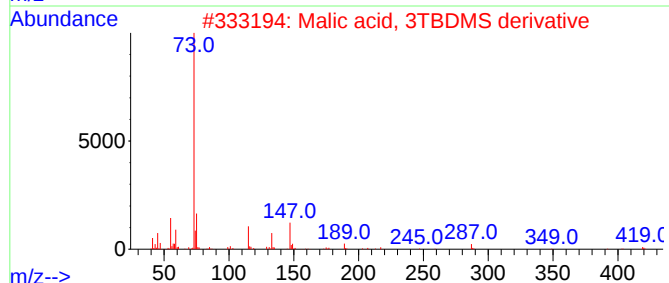
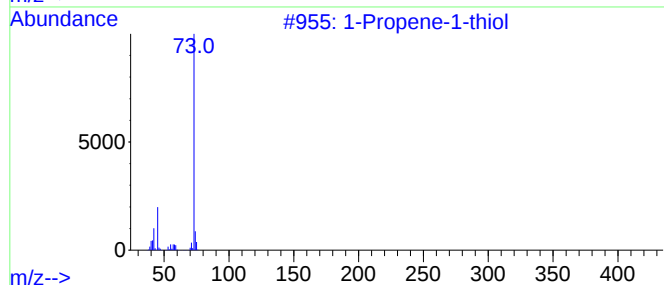
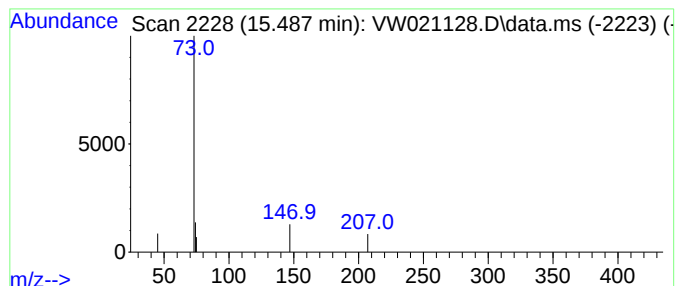
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Quant Title : SFAM01.0

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

Peak Number 14 unknown-06 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.487	3.82 ug/L	2155	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Propene-1-thiol	74	C3H6S	000925-89-3	4
2		Malic acid, 3TBDMS derivative	476	C22H48O5Si3	099461-86-6	4
3		L-Asparagine, 3TBDMS derivative	474	C22H50N2O3Si3	096381-41-8	4
4		Cyclopropane, 2-methylene-1-pent...	196	C12H24Si	167300-47-2	4
5		4-Chloro-3-nitro-2H-pyrazole	147	C3H2ClN3O2	1000411-34-5	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120321\
Data File : VW021128.D
Acq On : 04 Dec 2021 19:16
Operator : SY/VA
Sample : M4885-14
Misc : 3.93g/10.0mL/MSVOA_W/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EW9M1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120321SMA.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
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unknown-02	13.195	1.8	ug/L	1006	3	13.560	14086	25.0
unknown-03	13.408	2.1	ug/L	1187	3	13.560	14086	25.0
(DEL) Alkane: S...	13.762	1.5	ug/L	830	3	13.560	14086	25.0
unknown-04	14.719	3.2	ug/L	1826	3	13.560	14086	25.0
unknown-05	15.359	1.6	ug/L	929	3	13.560	14086	25.0
Benzene, 1,3-bi...	15.432	8.0	ug/L	4524	3	13.560	14086	25.0
unknown-06	15.487	3.8	ug/L	2155	3	13.560	14086	25.0