

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
 Data File : VW023766.D
 Acq On : 15 Jul 2022 23:54
 Operator : SY/VA
 Sample : N3700-06RE
 Misc : 6.41g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 F5C13RE

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Signal : TIC: VW023766.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.239	47	55	57	rBV6	11151	24837	37.27%	4.694%
2	3.788	308	309	313	rBV4	1372	1203	1.81%	0.227%
3	3.879	322	324	326	rBV3	1178	1002	1.50%	0.189%
4	3.946	334	335	339	rBV4	1140	1040	1.56%	0.197%
5	4.007	339	345	346	rBV5	5552	8217	12.33%	1.553%
6	4.269	386	388	391	rBV3	1285	1697	2.55%	0.321%
7	4.659	450	452	454	rBV3	1225	1415	2.12%	0.267%
8	4.757	467	468	472	rBV3	1374	1126	1.69%	0.213%
9	4.806	475	476	479	rVB3	1778	1199	1.80%	0.227%
10	4.848	481	483	486	rBV2	1390	965	1.45%	0.182%
11	4.885	486	489	490	rBV2	1377	1790	2.69%	0.338%
12	5.562	599	600	602	rBV2	1162	910	1.37%	0.172%
13	5.610	606	608	609	rBV2	1389	1068	1.60%	0.202%
14	5.653	613	615	619	rBV5	1349	1642	2.46%	0.310%
15	6.110	688	690	691	rBV2	1688	1472	2.21%	0.278%
16	6.251	710	713	714	rBV3	1027	937	1.41%	0.177%
17	6.397	735	737	741	rBV5	1238	2326	3.49%	0.440%
18	6.805	800	804	806	rBV5	896	1403	2.11%	0.265%
19	7.086	844	850	852	rBV6	3223	5590	8.39%	1.057%
20	7.360	893	895	898	rVB4	1625	1791	2.69%	0.339%
21	7.391	898	900	903	rBV4	830	1210	1.82%	0.229%
22	7.488	912	916	918	rBV4	891	1202	1.80%	0.227%
23	7.592	931	933	934	rBV2	1065	905	1.36%	0.171%
24	7.647	937	942	949	rVB3	11827	26624	39.95%	5.032%
25	7.805	966	968	971	rBV3	1067	1219	1.83%	0.230%
26	7.939	985	990	993	rBV7	928	1701	2.55%	0.322%
27	8.141	1021	1023	1024	rBV	1081	921	1.38%	0.174%
28	8.281	1036	1046	1056	rBV5	18275	66648	100.00%	12.597%
29	8.378	1060	1062	1063	rBV2	1787	1301	1.95%	0.246%
30	8.842	1133	1138	1148	rVB3	16662	35941	53.93%	6.793%
31	9.268	1198	1208	1218	rBV3	14264	37074	55.63%	7.007%
32	9.402	1227	1230	1231	rBV3	1012	1025	1.54%	0.194%
33	9.470	1240	1241	1244	rVB3	1227	1056	1.58%	0.200%
34	9.500	1244	1246	1249	rBV3	1421	1701	2.55%	0.322%
35	9.561	1252	1256	1257	rBV4	1352	1753	2.63%	0.331%
36	9.750	1282	1287	1290	rBV6	2054	2075	3.11%	0.392%

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Integrator: RTE

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Peak Location: TOP

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

37	9.805	1293	1296	1299	rVV5	1028	1079	1.62%	0.204%
38	10.037	1330	1334	1342	rVB4	5423	10787	16.19%	2.039%
39	10.323	1374	1381	1387	rBV2	18834	37482	56.24%	7.084%
40	10.463	1403	1404	1406	rBV2	1213	1068	1.60%	0.202%
41	10.512	1411	1412	1413	rBV	1559	844	1.27%	0.160%
42	10.530	1413	1415	1417	rVB3	1401	1046	1.57%	0.198%
43	10.622	1428	1430	1433	rVB4	1422	1436	2.15%	0.271%
44	10.652	1433	1435	1438	rVB4	1257	1105	1.66%	0.209%
45	10.707	1438	1444	1449	rBV7	2309	4906	7.36%	0.927%
46	10.811	1457	1461	1465	rBV7	1081	1919	2.88%	0.363%
47	10.884	1471	1473	1475	rBV3	1099	1291	1.94%	0.244%
48	10.927	1475	1480	1486	rBV2	7235	12491	18.74%	2.361%
49	11.024	1494	1496	1498	rBV2	1097	920	1.38%	0.174%
50	11.128	1512	1513	1518	rBV4	951	1231	1.85%	0.233%
51	11.170	1518	1520	1522	rVV2	1168	841	1.26%	0.159%
52	11.402	1555	1558	1560	rBV4	880	905	1.36%	0.171%
53	11.439	1560	1564	1569	rBV7	1430	3138	4.71%	0.593%
54	11.481	1569	1571	1573	rBV3	1072	922	1.38%	0.174%
55	11.542	1579	1581	1585	rVB5	1428	1841	2.76%	0.348%
56	11.634	1590	1596	1604	rVB	26286	48946	73.44%	9.251%
57	11.817	1624	1626	1628	rBV3	1053	920	1.38%	0.174%
58	11.920	1641	1643	1645	rVV3	910	925	1.39%	0.175%
59	11.969	1649	1651	1653	rVB3	1217	1003	1.50%	0.190%
60	11.993	1653	1655	1657	rBV3	1421	1144	1.72%	0.216%
61	12.085	1668	1670	1673	rBV4	1009	1030	1.55%	0.195%
62	12.250	1694	1697	1700	rBV5	989	1182	1.77%	0.223%
63	12.335	1709	1711	1714	rBV4	1041	881	1.32%	0.167%
64	12.365	1714	1716	1719	rBV4	852	908	1.36%	0.172%
65	12.603	1750	1755	1758	rBV5	2751	4207	6.31%	0.795%
66	12.634	1758	1760	1763	rVB3	1376	1096	1.64%	0.207%
67	12.688	1763	1769	1776	rBV	16527	30046	45.08%	5.679%
68	12.743	1776	1778	1781	rBV4	1339	1404	2.11%	0.265%
69	12.792	1784	1786	1787	rVB2	1635	1006	1.51%	0.190%
70	12.920	1805	1807	1811	rBV5	1060	1048	1.57%	0.198%
71	12.999	1818	1820	1823	rVB4	1264	1485	2.23%	0.281%
72	13.298	1867	1869	1873	rVB4	2074	1693	2.54%	0.320%
73	13.341	1873	1876	1878	rBV4	1283	1366	2.05%	0.258%
74	13.512	1901	1904	1906	rBV4	892	1009	1.51%	0.191%
75	13.554	1906	1911	1919	rVV2	19066	34224	51.35%	6.469%
76	13.749	1942	1943	1947	rVV4	1743	1745	2.62%	0.330%

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

77	13.786	1947	1949	1954	rVB6	1050	1351	2.03%	0.255%
78	13.847	1954	1959	1966	rBV2	18831	32494	48.75%	6.142%
79	14.011	1984	1986	1987	rVV2	1348	845	1.27%	0.160%
80	14.060	1991	1994	1995	rBV2	1804	2108	3.16%	0.398%
81	14.103	1998	2001	2003	rVB4	1023	1209	1.81%	0.229%
82	14.133	2003	2006	2010	rBV5	1108	1452	2.18%	0.274%
83	14.170	2010	2012	2016	rVB5	1040	1089	1.63%	0.206%
84	14.213	2016	2019	2021	rVB4	954	1108	1.66%	0.209%
85	14.322	2034	2037	2039	rBV4	1129	1297	1.95%	0.245%
86	14.456	2057	2059	2062	rBV3	683	879	1.32%	0.166%
87	14.536	2070	2072	2075	rBV3	1208	1418	2.13%	0.268%
88	14.566	2075	2077	2079	rVV3	1272	1508	2.26%	0.285%
89	14.621	2084	2086	2087	rVV2	1171	1112	1.67%	0.210%
90	14.767	2108	2110	2114	rVB5	1019	883	1.32%	0.167%
91	14.877	2127	2128	2131	rVB2	1451	919	1.38%	0.174%
92	14.950	2137	2140	2143	rBV5	1151	1870	2.81%	0.353%
93	15.529	2232	2235	2237	rBV4	1325	1554	2.33%	0.294%
94	15.633	2250	2252	2255	rBV3	1427	1165	1.75%	0.220%
95	15.694	2260	2262	2265	rVB4	1426	1418	2.13%	0.268%
96	15.731	2265	2268	2269	rBV2	1347	1181	1.77%	0.223%
97	16.426	2380	2382	2387	rBV5	785	1091	1.64%	0.206%
98	16.462	2387	2388	2391	rBV3	960	1149	1.72%	0.217%
99	16.541	2397	2401	2406	rBV8	1348	2571	3.86%	0.486%
100	16.633	2415	2416	2419	rVV3	943	899	1.35%	0.170%

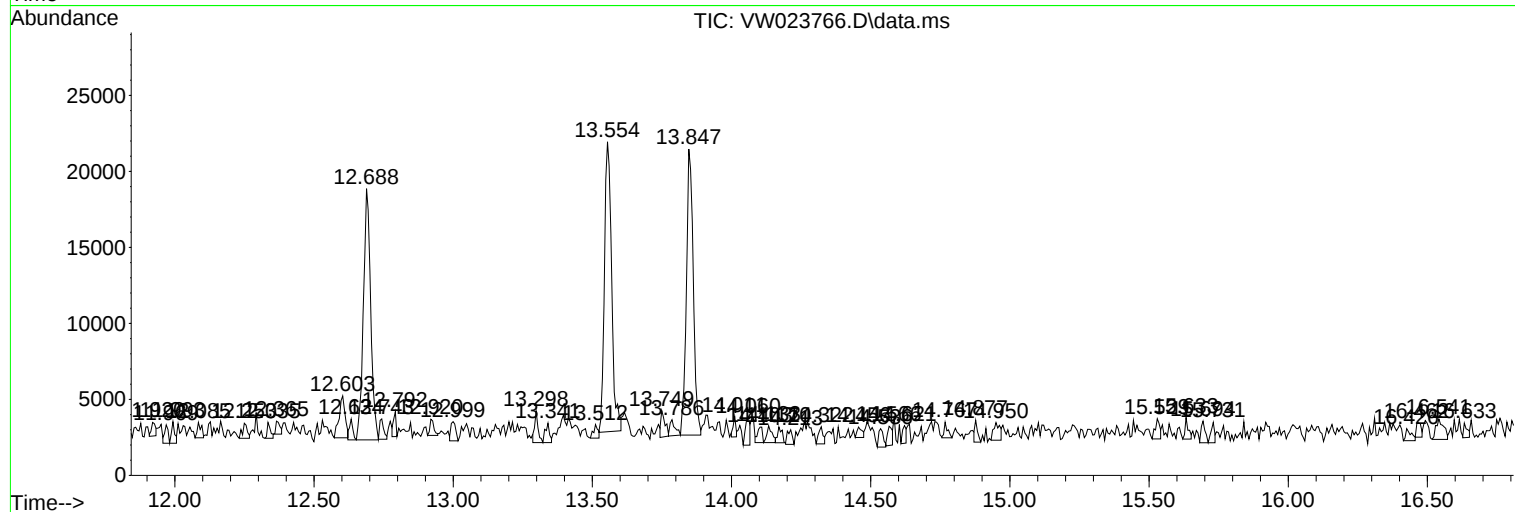
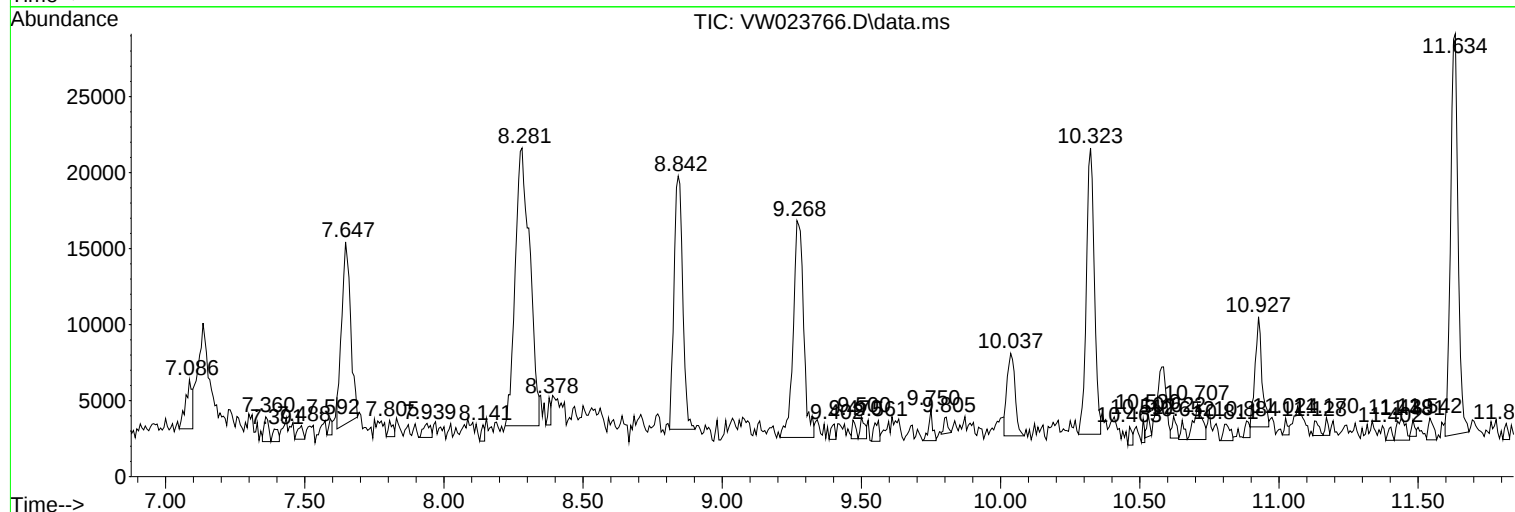
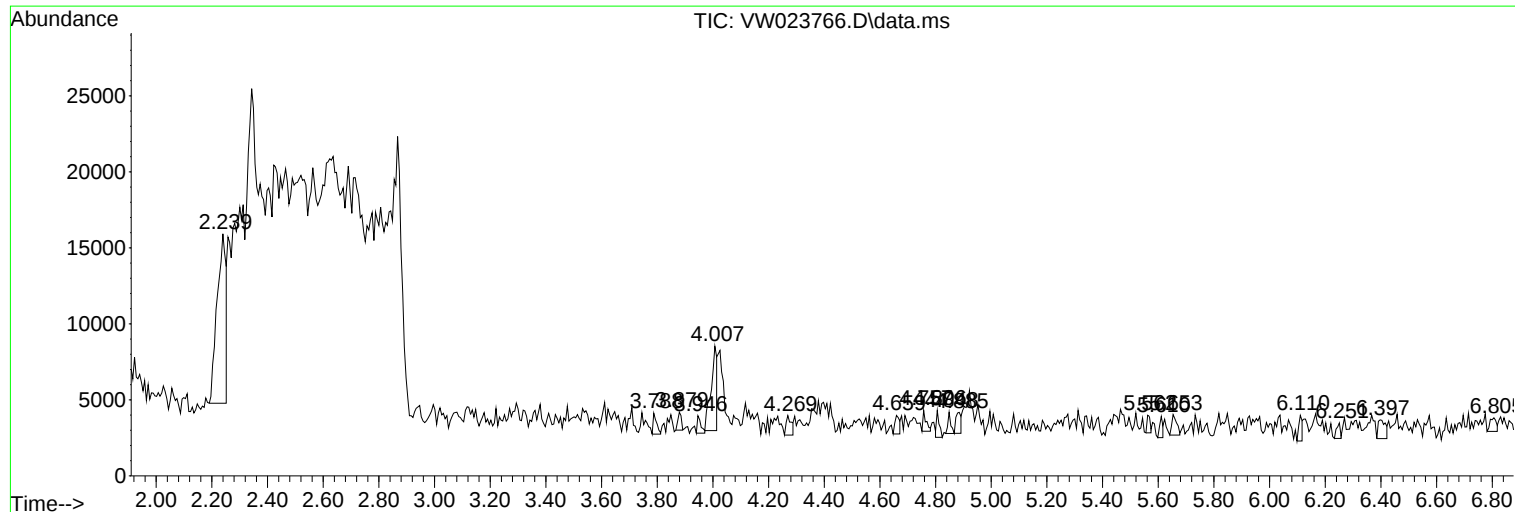
Sum of corrected areas: 529076

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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

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

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



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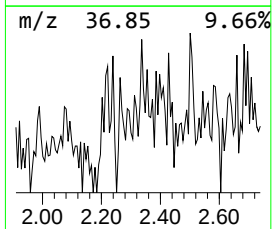
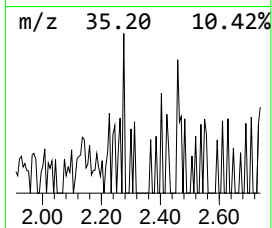
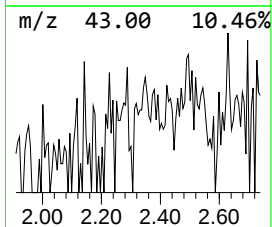
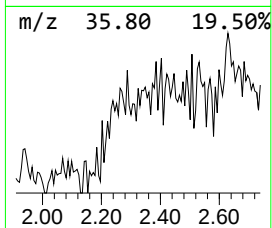
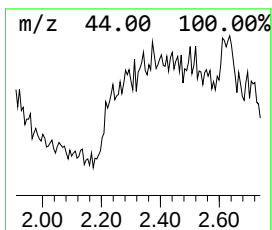
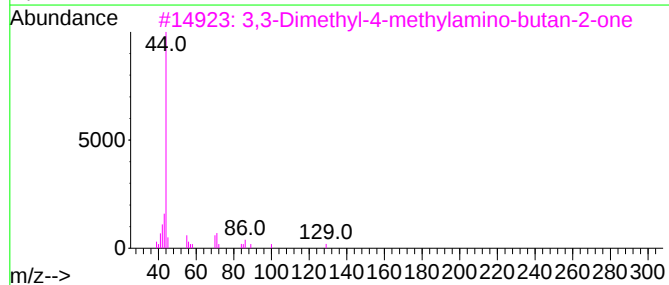
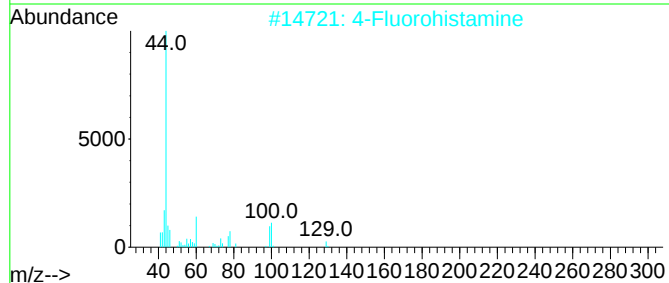
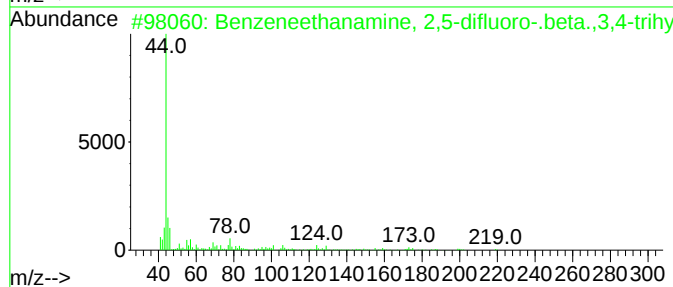
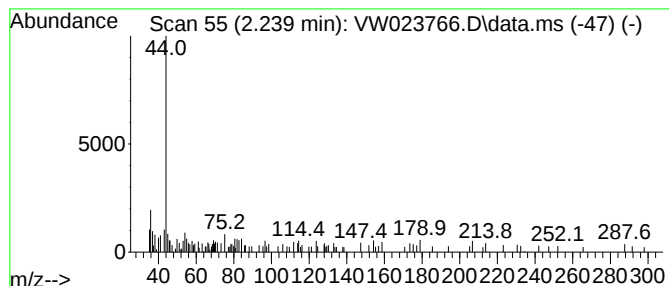
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 Benzeneethanamine, 2,5-difl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.239	17.28 ug/L	24837	1,4-Difluorobenzene	8.842

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzeneethanamine, 2,5-difluoro-...	219	C9H11F2NO3	152434-78-1	53
2		4-Fluorohistamine	129	C5H8FN3	049872-60-8	53
3		3,3-Dimethyl-4-methylamino-butan...	129	C7H15NO	123528-99-4	47
4		2-Butenediamide, (E)-	114	C4H6N2O2	000627-64-5	42
5		3-Methyl-3,5--(cyanoethyl)tetrah...	236	C12H16N2OS	1000140-32-3	38



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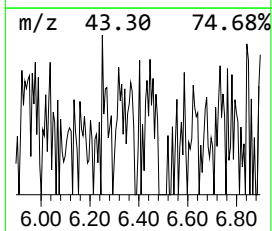
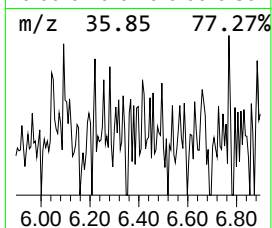
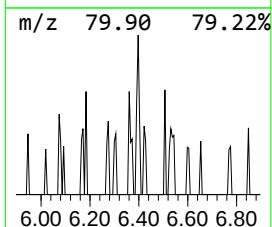
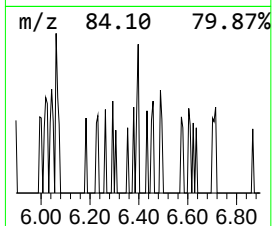
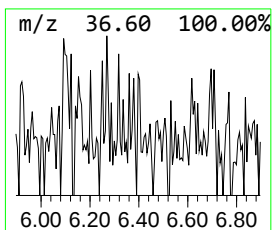
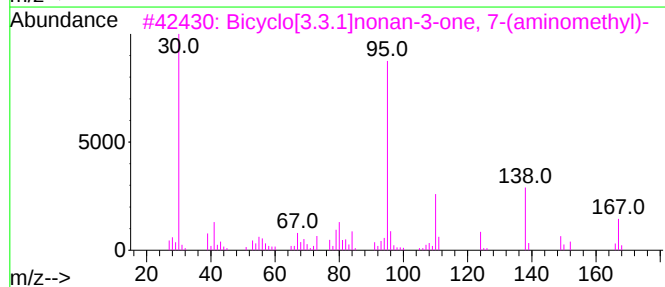
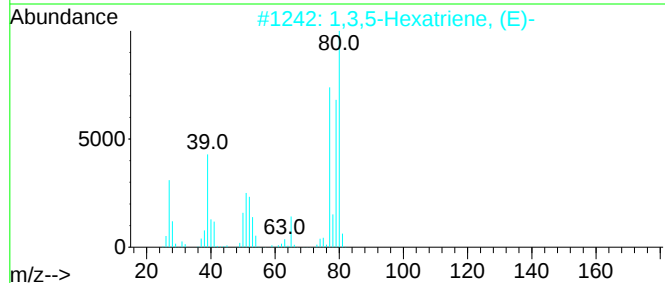
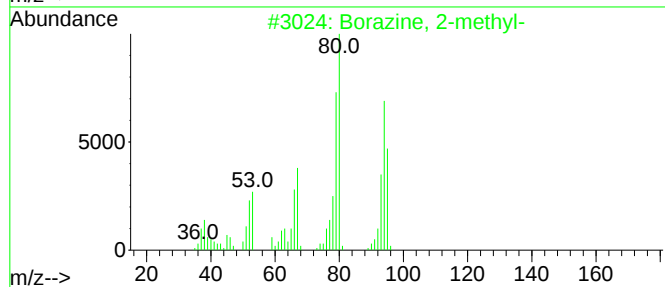
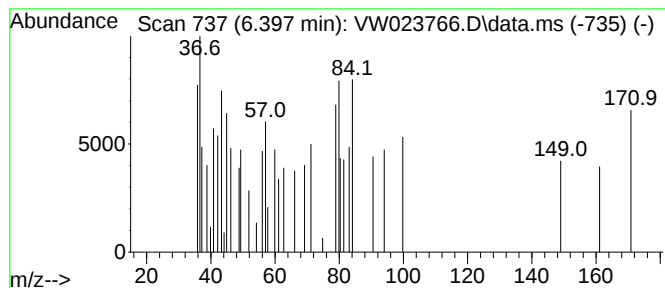
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.397	1.62 ug/L	2326	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Borazine, 2-methyl-	95	CH8B3N3	021127-95-7	12
2			1,3,5-Hexatriene, (E)-	80	C6H8	000821-07-8	10
3			Bicyclo[3.3.1]nonan-3-one, 7-(am...	167	C10H17NO	056701-31-6	10
4			4,7-Methano-1H-indene, octahydro...	148	C11H16	068284-23-1	10
5			s-Triazaborane	81	B3H6N3	006569-51-3	10



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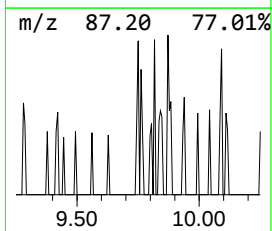
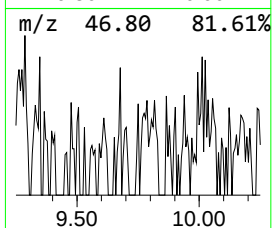
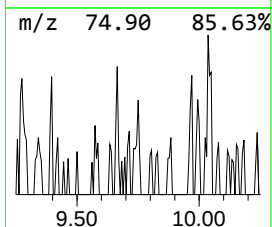
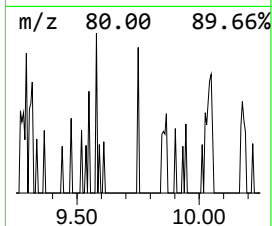
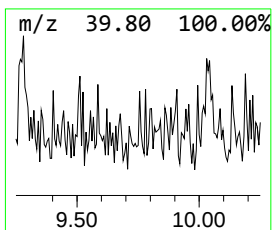
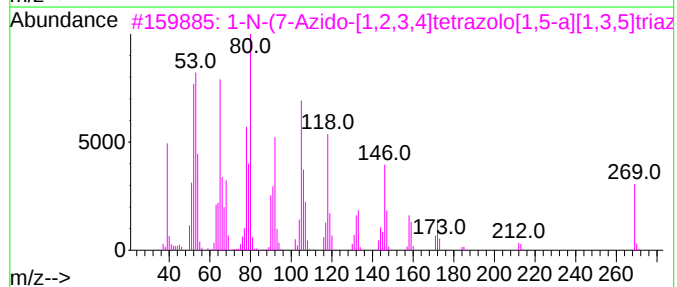
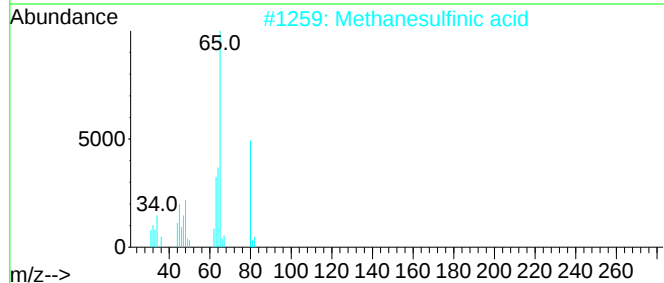
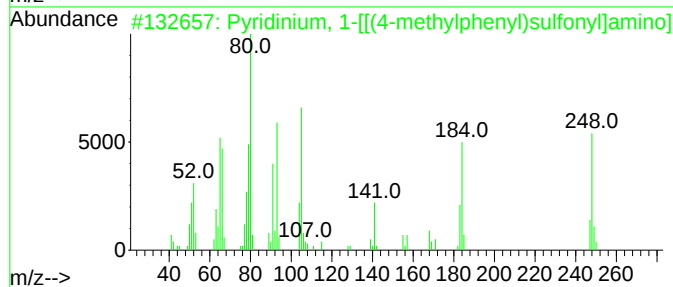
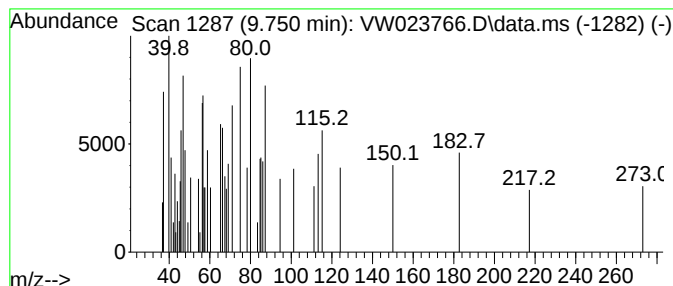
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Quant Title : SFAM01.0

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

Peak Number 3 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.750	1.44 ug/L	2075	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridinium, 1-[[4-methylphenyl]...	248	C12H12N2O2S	040949-56-2	33
2			Methanesulfinic acid	80	CH4O2S	017696-73-0	16
3			1-N-(7-Azido-[1,2,3,4]tetrazolo[...	269	C9H7N11	1000388-12-8	16
4			Bicyclo[2.2.2]octanone	124	C8H12O	002716-23-6	14
5			Cycloocta-2,4-dien-1-ol	124	C8H12O	1010164-53-3	11



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

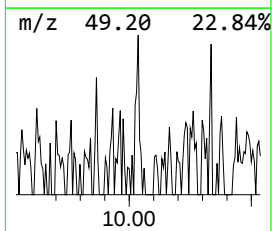
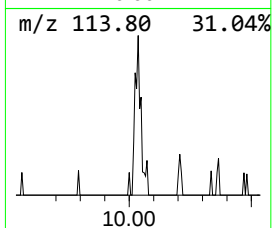
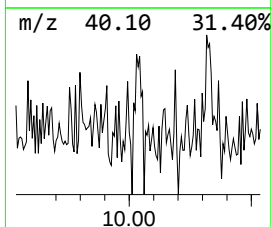
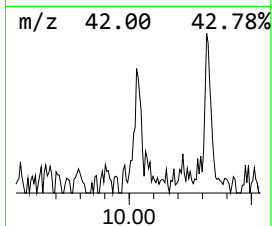
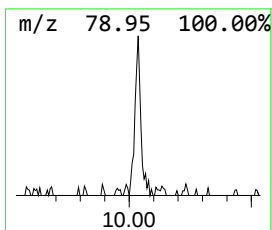
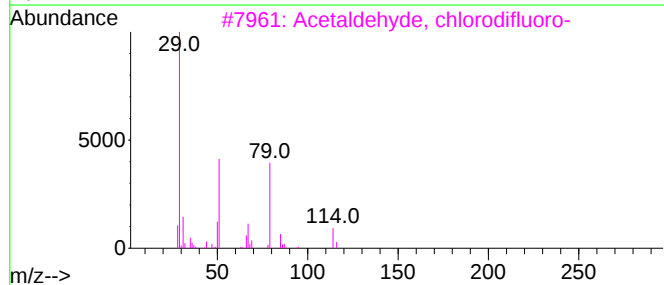
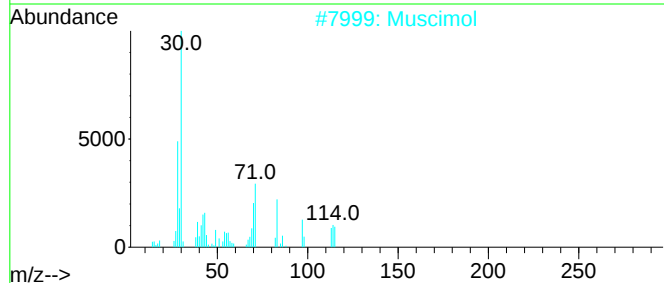
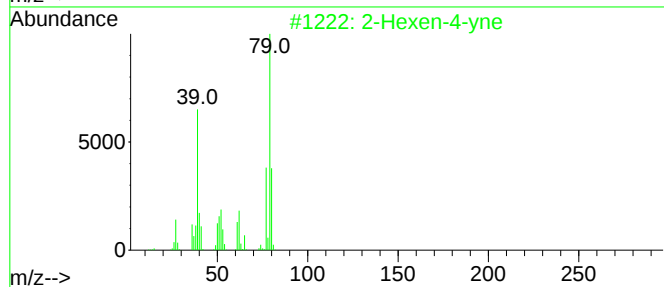
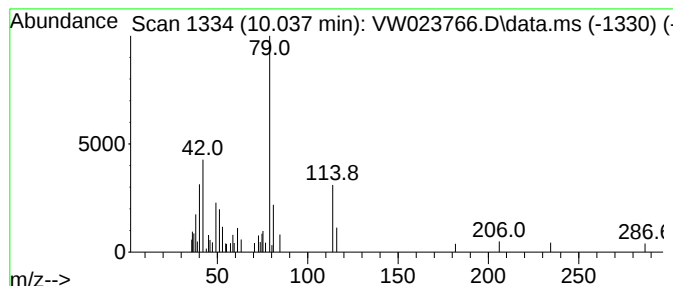
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Quant Title : SFAM01.0

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

Peak Number 4 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.037	7.50 ug/L	10787	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexen-4-yne			80	C6H8	014092-20-7	10
2	Muscimol			114	C4H6N2O2	002763-96-4	9
3	Acetaldehyde, chlorodifluoro-			114	C2HClF2O	000811-96-1	9
4	Thiophosgene			114	CCl2S	000463-71-8	9
5	3-Cyclohexene-1-methanol			112	C7H12O	001679-51-2	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

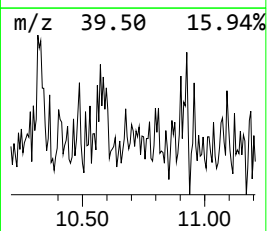
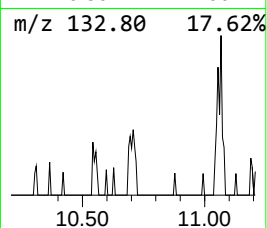
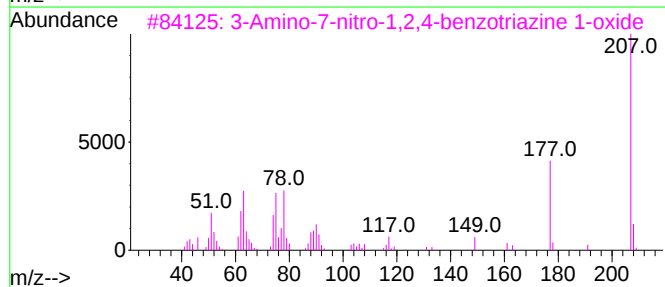
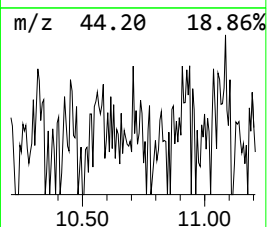
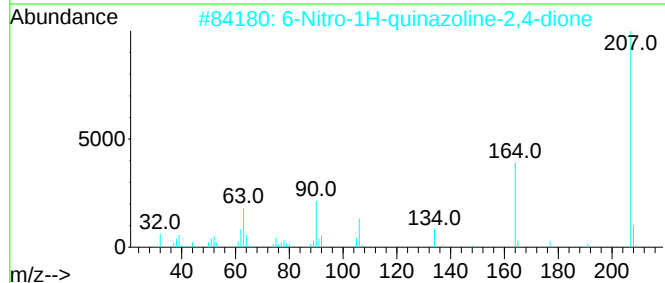
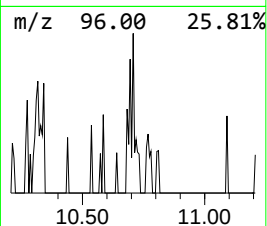
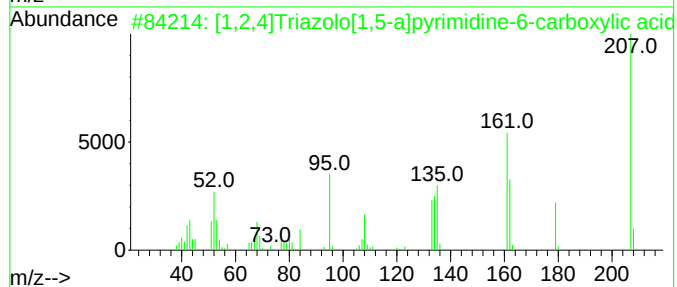
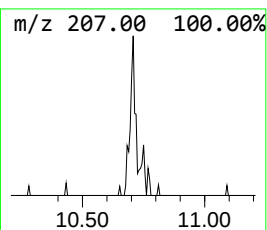
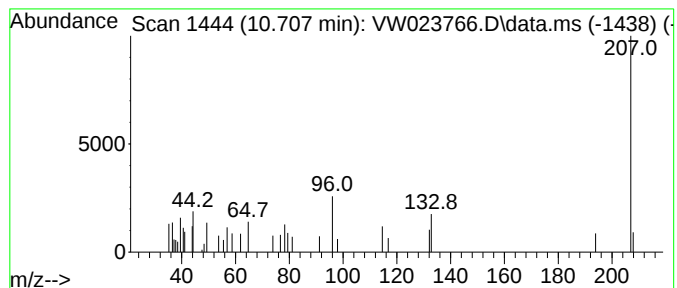
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.707	2.51 ug/L	4906	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			[1,2,4]Triazolo[1,5-a]pyrimidine...	207	C8H9N5O2	1000351-62-2	38
2			6-Nitro-1H-quinazoline-2,4-dione	207	C8H5N3O4	1000318-16-1	32
3			3-Amino-7-nitro-1,2,4-benzotriaz...	207	C7H5N5O3	1010256-54-1	25
4			1-(2,4-difluorophenyl)-1H-pyrrol...	207	C11H7F2NO	125126-74-1	9
5			2-Amino-6-methoxy-4-(2H-1,2,3,4-...	207	C8H9N5O2	1010387-42-7	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

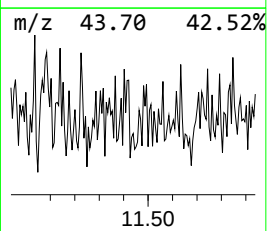
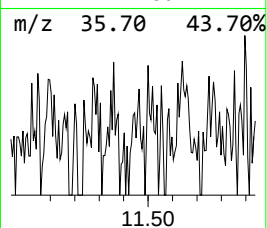
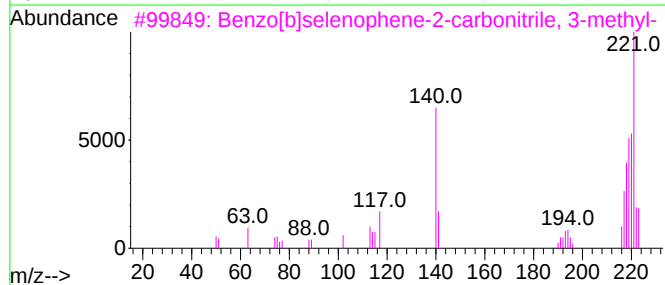
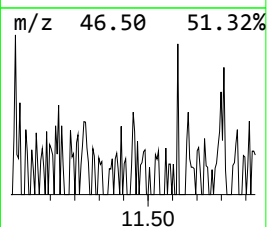
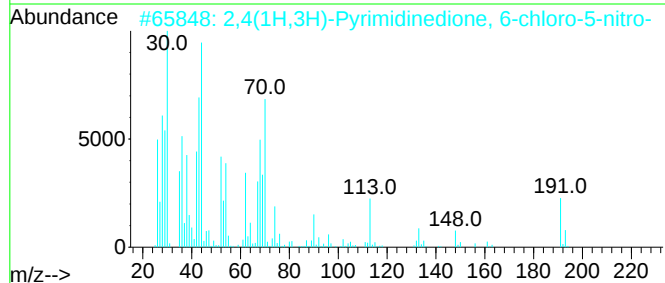
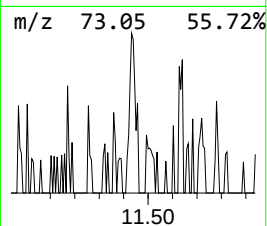
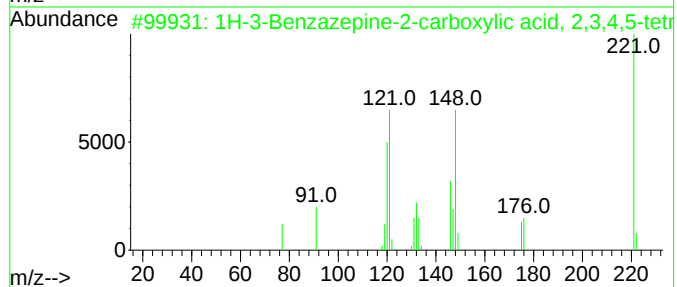
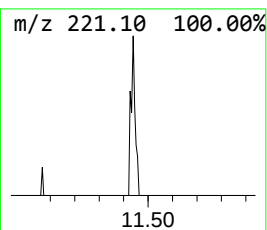
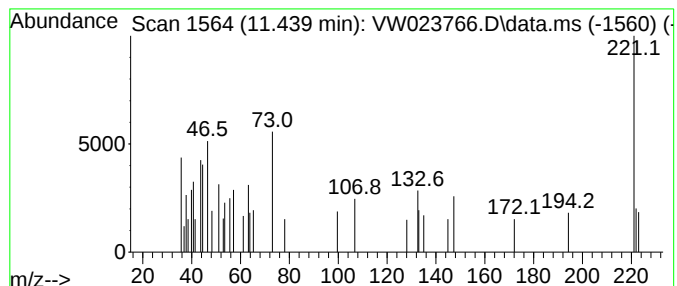
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Quant Title : SFAM01.0

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

Peak Number 6 1H-3-Benzazepine-2-carboxyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.439	1.60 ug/L	3138	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-3-Benzazepine-2-carboxylic ac...	221	C11H11NO4	017639-47-3	53		
2	2,4(1H,3H)-Pyrimidinedione, 6-ch...	191	C4H2ClN3O4	006630-30-4	25		
3	Benzo[b]selenophene-2-carbonitri...	221	C10H7NSe	037007-55-9	16		
4	4-Quinolinol, 2-phenyl-	221	C15H11NO	001144-20-3	10		
5	2-Iodo-3-pyridinol	221	C5H4INO	040263-57-8	9		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

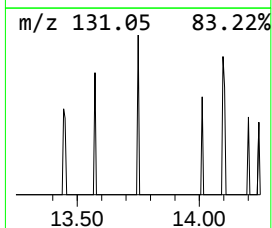
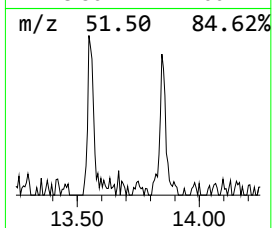
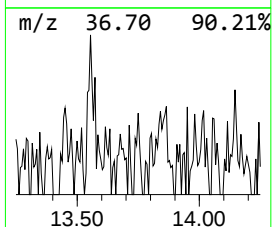
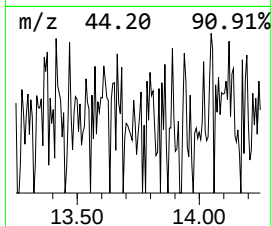
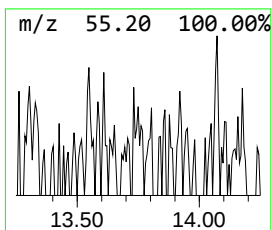
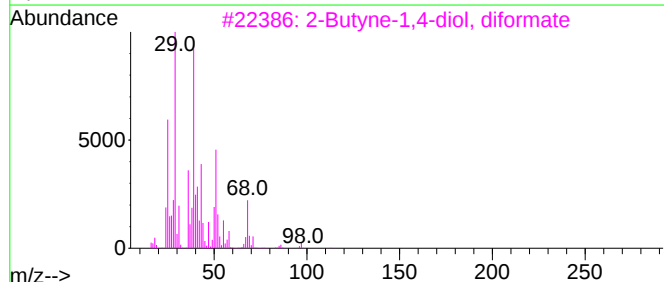
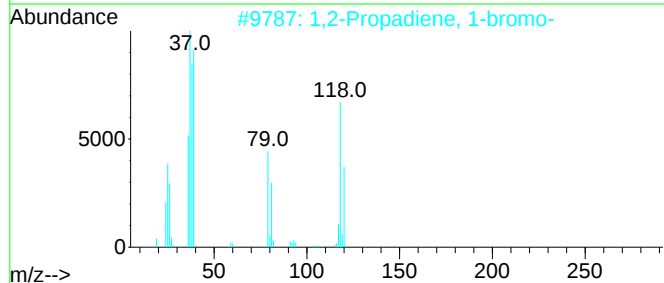
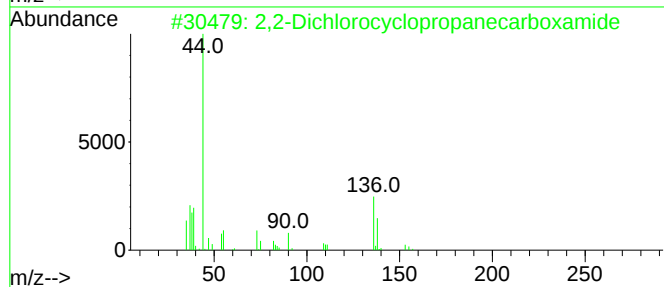
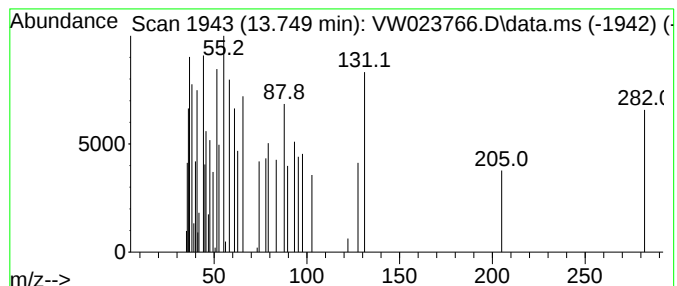
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-05 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.749	1.27 ug/L	1745	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-Dichlorocyclopropanecarboxamide	153	C4H5Cl2NO	075885-60-8	17
2			1,2-Propadiene, 1-bromo-	118	C3H3Br	010024-18-7	10
3			2-Butyne-1,4-diol, diformate	142	C6H6O4	036677-73-3	9
4			Ethyl 2-(2-chloroacetamido)-3,3,...	263	C7H9ClF3NO4	1000224-74-2	9
5			5-Amino-1-(2-methoxyphenyl)-1H-1...	233	C10H11N5O2	1000387-16-8	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

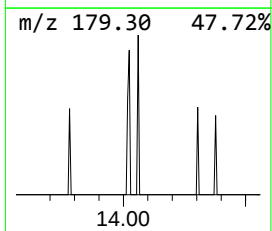
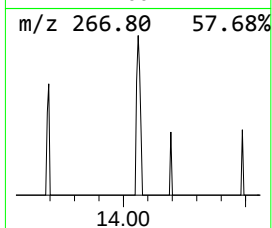
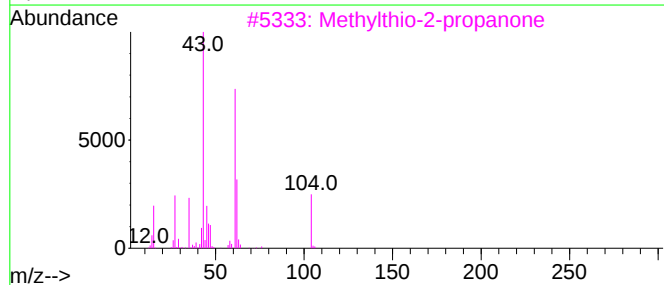
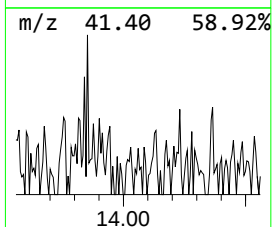
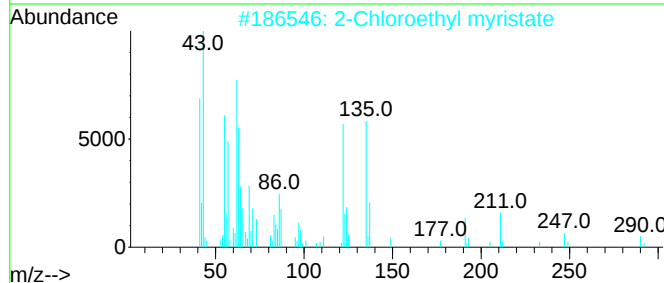
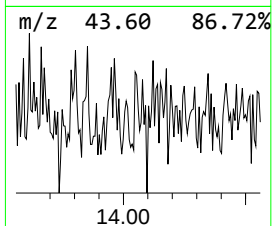
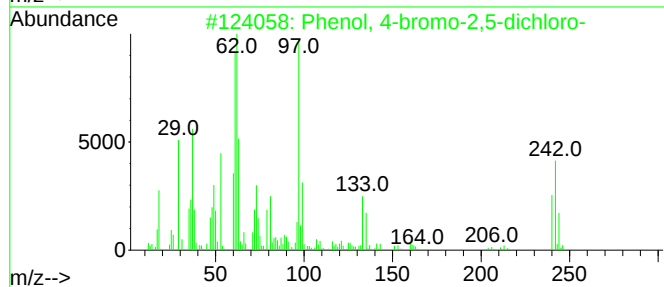
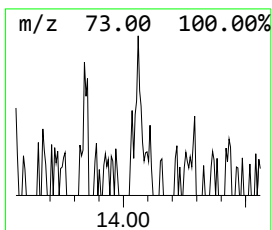
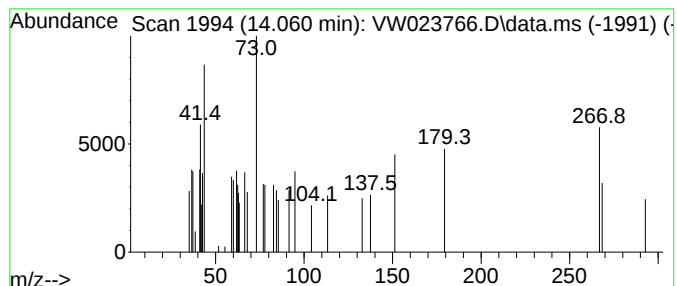
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.060	1.54 ug/L	2108	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-bromo-2,5-dichloro-	240	C6H3BrCl2O	001940-42-7	37
2			2-Chloroethyl myristate	290	C16H31ClO2	051479-36-8	25
3			Methylthio-2-propanone	104	C4H8OS	014109-72-9	9
4			Dimethylfluoroamine	63	C2H6FN	014722-43-1	9
5			Carbamic acid, acetyl-, ethyl ester	131	C5H9NO3	002597-54-8	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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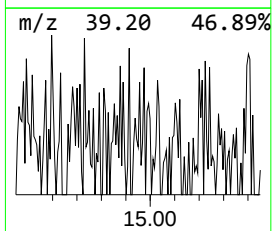
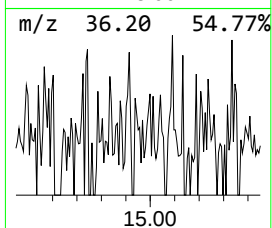
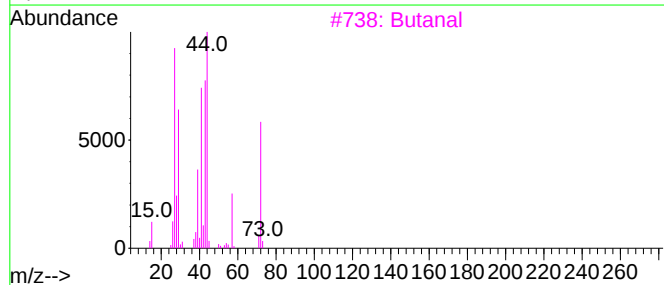
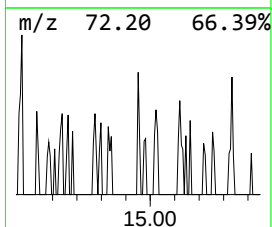
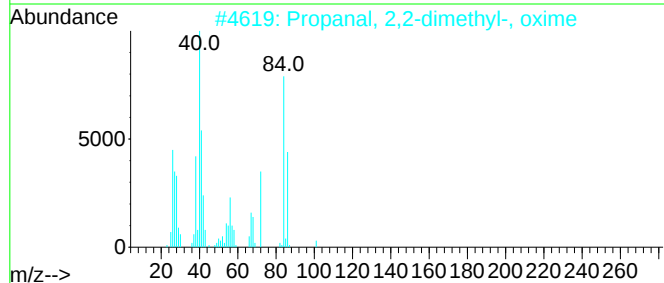
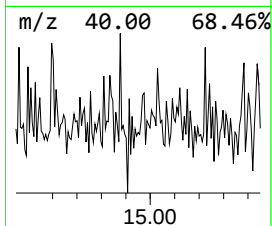
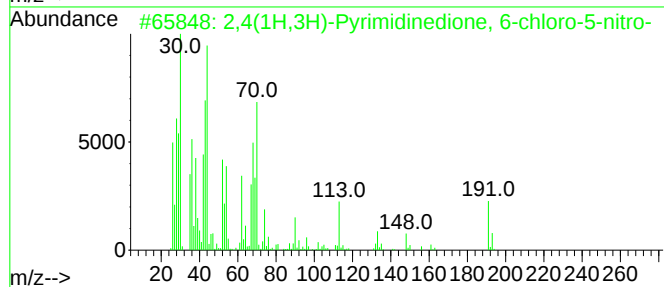
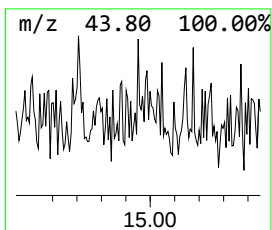
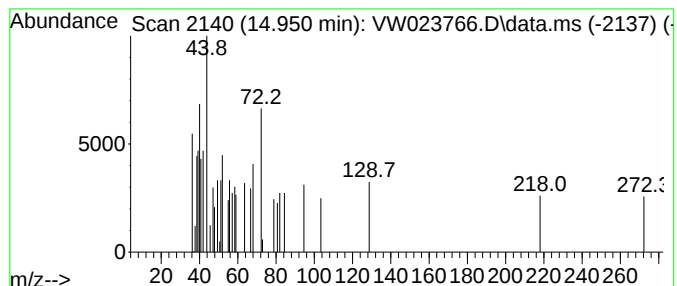
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-07 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.950	1.37 ug/L	1870	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4(1H,3H)-Pyrimidinedione, 6-ch...	191	C4H2ClN3O4	006630-30-4	38
2			Propanal, 2,2-dimethyl-, oxime	101	C5H11NO	000637-91-2	14
3			Butanal	72	C4H8O	000123-72-8	11
4			Butanal	72	C4H8O	000123-72-8	11
5			Butanal	72	C4H8O	000123-72-8	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
Data File : VW023766.D
Acq On : 15 Jul 2022 23:54
Operator : SY/VA
Sample : N3700-06RE
Misc : 6.41g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5C13RE

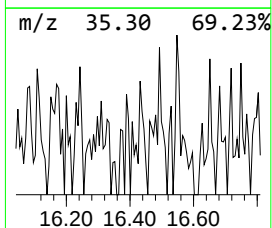
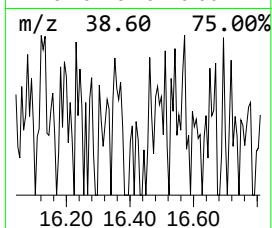
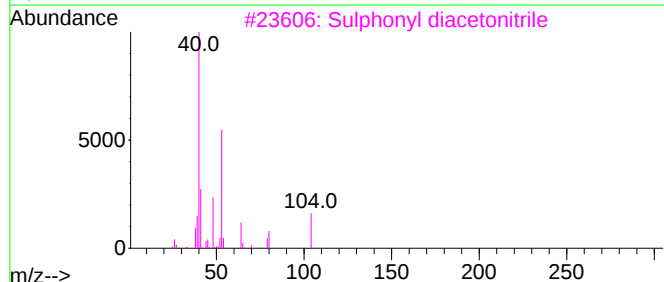
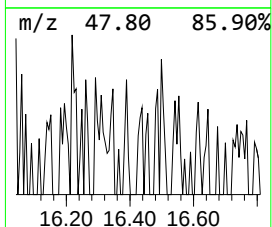
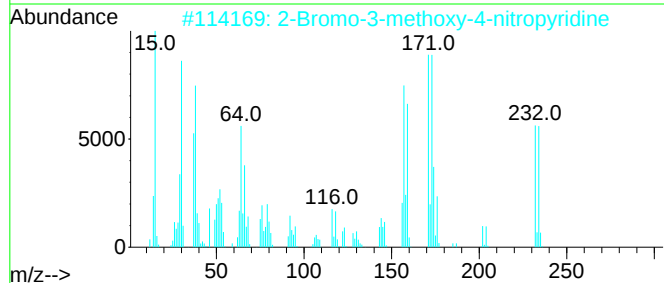
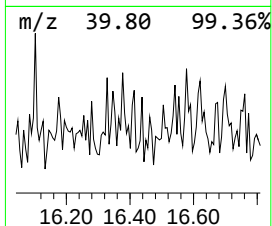
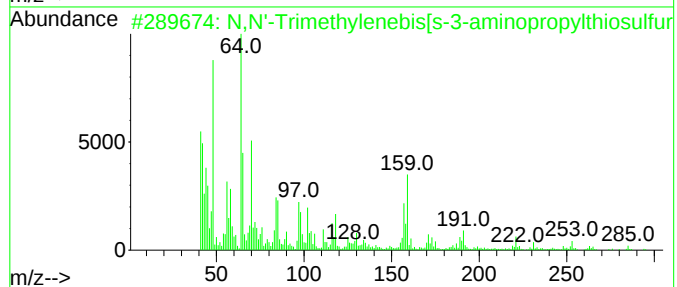
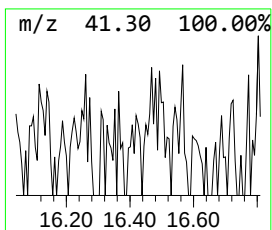
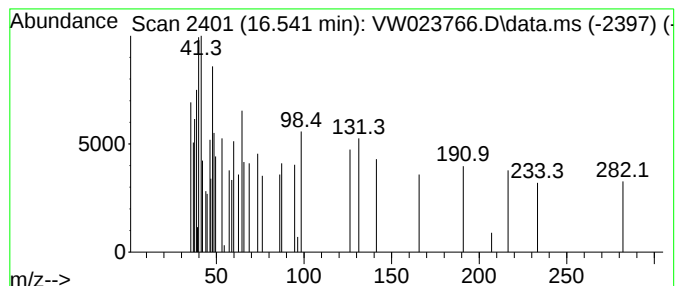
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Quant Title : SFAM01.0

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

Peak Number 11 unknown-08 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.541	1.88 ug/L	2571	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N,N'-Trimethylenebis[s-3-aminopr...	382	C9H22N2O6S4	035871-60-4	9
2			2-Bromo-3-methoxy-4-nitropyridine	232	C6H5BrN2O3	1000444-83-2	9
3			Sulphonyl diacetonitrile	144	C4H4N2O2S	1000342-33-0	9
4			Methyl 3-(5-bromo-1H-1,2,4-triaz...	231	C6H6BrN3O2	1000443-27-3	9
5			6-Hydroxy-5,8-dioxo-carbostyryl-...	264	C10H8N4O3S	1000213-46-3	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071522\
 Data File : VW023766.D
 Acq On : 15 Jul 2022 23:54
 Operator : SY/VA
 Sample : N3700-06RE
 Misc : 6.41g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 F5C13RE

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanami...	2.239	17.3	ug/L	24837	1	8.842	35941	25.0
unknown-01	6.397	1.6	ug/L	2326	1	8.842	35941	25.0
unknown-02	9.750	1.4	ug/L	2075	1	8.842	35941	25.0
unknown-03	10.037	7.5	ug/L	10787	1	8.842	35941	25.0
unknown-04	10.707	2.5	ug/L	4906	2	11.634	48946	25.0
1H-3-Benzazepin...	11.439	1.6	ug/L	3138	2	11.634	48946	25.0
unknown-05	13.749	1.3	ug/L	1745	3	13.560	34224	25.0
unknown-06	14.060	1.5	ug/L	2108	3	13.560	34224	25.0
unknown-07	14.950	1.4	ug/L	1870	3	13.560	34224	25.0
unknown-08	16.541	1.9	ug/L	2571	3	13.560	34224	25.0