

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
 Data File : VW026917.D
 Acq On : 26 Aug 2023 01:03
 Operator : SY/MD
 Sample : 04175-05
 Misc : 5.76g/10mL/MSVOA_W/SOIL
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EZYX0

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\SFAMWLM081023SMA.M
 Title : SFAM01.0

Signal : TIC: VW026917.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.113	35	37	38	rBV2	1656	1353	2.65%	0.326%
2	2.241	52	58	60	rBV6	6341	14900	29.15%	3.590%
3	2.527	103	105	107	rBV3	2314	2234	4.37%	0.538%
4	2.649	123	125	127	rBV3	1863	1680	3.29%	0.405%
5	2.887	161	164	165	rBV3	2748	3885	7.60%	0.936%
6	3.881	324	327	330	rBV5	2603	3820	7.47%	0.920%
7	3.960	338	340	343	rBV4	1995	1932	3.78%	0.465%
8	4.119	364	366	369	rVB4	2085	1780	3.48%	0.429%
9	4.558	435	438	440	rBV4	1767	2036	3.98%	0.491%
10	4.783	472	475	479	rBV6	1654	2709	5.30%	0.653%
11	4.984	506	508	510	rBV3	1633	1547	3.03%	0.373%
12	5.124	527	531	533	rBV5	2027	2516	4.92%	0.606%
13	5.185	536	541	542	rBV5	1859	2595	5.08%	0.625%
14	5.265	552	554	555	rBV2	2023	1620	3.17%	0.390%
15	5.411	574	578	581	rVB6	2094	3334	6.52%	0.803%
16	5.478	587	589	593	rBV5	1728	2407	4.71%	0.580%
17	5.515	593	595	598	rVV4	1353	1378	2.70%	0.332%
18	5.606	608	610	611	rBV2	1831	1244	2.43%	0.300%
19	5.691	622	624	627	rBV4	1350	1620	3.17%	0.390%
20	5.856	648	651	653	rBV4	2012	1849	3.62%	0.445%
21	6.082	686	688	691	rBV4	1755	1994	3.90%	0.480%
22	6.283	718	721	722	rBV3	1530	1779	3.48%	0.429%
23	6.332	727	729	731	rVV3	1693	1841	3.60%	0.444%
24	6.490	753	755	757	rBV3	1246	1364	2.67%	0.329%
25	6.716	789	792	794	rBV4	1585	1963	3.84%	0.473%
26	6.862	814	816	818	rBV3	1212	1253	2.45%	0.302%
27	6.929	825	827	830	rVB4	1478	1716	3.36%	0.413%
28	7.069	847	850	851	rBV3	1944	1806	3.53%	0.435%
29	7.197	869	871	874	rVB4	1819	1687	3.30%	0.406%
30	7.228	874	876	879	rBV3	1301	1687	3.30%	0.406%
31	7.496	919	920	922	rBV2	1854	1618	3.17%	0.390%
32	7.575	928	933	934	rBV5	2071	1752	3.43%	0.422%
33	7.648	937	945	952	rBV4	8855	28784	56.31%	6.934%
34	7.746	959	961	964	rVV4	1323	1569	3.07%	0.378%
35	7.801	968	970	972	rBV3	1860	1407	2.75%	0.339%
36	7.844	972	977	980	rBV7	1911	3509	6.86%	0.845%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
 Title : SFAM01.0

37	7.959	993	996	1000	rVB6	1256	1823	3.57%	0.439%
38	8.039	1006	1009	1010	rBV3	2013	1884	3.69%	0.454%
39	8.081	1013	1016	1018	rBV4	1559	2210	4.32%	0.532%
40	8.154	1024	1028	1032	rBV7	2334	3781	7.40%	0.911%
41	8.197	1032	1035	1036	rBV3	1484	1839	3.60%	0.443%
42	8.276	1041	1048	1060	rVV2	14697	51117	100.00%	12.315%
43	8.508	1083	1086	1087	rBV3	1422	1861	3.64%	0.448%
44	8.593	1098	1100	1102	rBV3	1549	1804	3.53%	0.435%
45	8.697	1114	1117	1121	rBV6	1308	2408	4.71%	0.580%
46	8.764	1126	1128	1131	rBV4	1399	1488	2.91%	0.358%
47	8.843	1137	1141	1150	rVB	12803	25009	48.93%	6.025%
48	9.020	1168	1170	1174	rVB5	2133	2691	5.26%	0.648%
49	9.057	1174	1176	1177	rBV2	1608	1523	2.98%	0.367%
50	9.130	1186	1188	1189	rVB2	2225	1312	2.57%	0.316%
51	9.234	1202	1205	1206	rBV3	1484	1643	3.21%	0.396%
52	9.276	1206	1212	1216	rBV4	11049	22032	43.10%	5.308%
53	9.490	1244	1247	1249	rBV4	1211	1868	3.65%	0.450%
54	9.648	1270	1273	1275	rBV4	1470	2128	4.16%	0.513%
55	9.666	1275	1276	1279	rVB3	1826	1241	2.43%	0.299%
56	9.800	1297	1298	1300	rBV3	1530	1324	2.59%	0.319%
57	9.989	1326	1329	1330	rVB3	1198	1189	2.33%	0.286%
58	10.038	1332	1337	1339	rVV4	3757	5624	11.00%	1.355%
59	10.105	1347	1348	1350	rBV2	2274	1475	2.89%	0.355%
60	10.148	1353	1355	1358	rBV4	1533	1651	3.23%	0.398%
61	10.197	1362	1363	1367	rVB4	1799	1608	3.15%	0.387%
62	10.233	1367	1369	1373	rVB5	2129	3050	5.97%	0.735%
63	10.270	1373	1375	1376	rBV2	1656	1405	2.75%	0.338%
64	10.325	1379	1384	1391	rVB2	13987	23785	46.53%	5.730%
65	10.672	1436	1441	1442	rBV5	1167	1690	3.31%	0.407%
66	10.752	1453	1454	1459	rVB5	1108	1171	2.29%	0.282%
67	10.800	1459	1462	1465	rVB4	1613	2130	4.17%	0.513%
68	10.995	1492	1494	1496	rVB3	1692	1311	2.56%	0.316%
69	11.026	1496	1499	1500	rBV3	1443	1311	2.56%	0.316%
70	11.148	1517	1519	1521	rVV3	1411	1233	2.41%	0.297%
71	11.172	1521	1523	1526	rVB4	2007	1689	3.30%	0.407%
72	11.325	1545	1548	1550	rVB4	2359	1863	3.64%	0.449%
73	11.355	1552	1553	1557	rBV4	1914	2352	4.60%	0.567%
74	11.453	1568	1569	1572	rBV3	1314	1162	2.27%	0.280%
75	11.629	1592	1598	1605	rBV	14253	25711	50.30%	6.194%
76	11.709	1609	1611	1614	rBV4	1748	2012	3.94%	0.485%

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 Title : SFAM01.0

77	11.818	1627	1629	1631	rBV3	1029	1169	2.29%	0.282%
78	11.965	1650	1653	1654	rBV3	1776	1525	2.98%	0.367%
79	12.068	1666	1670	1673	rBV6	2101	3395	6.64%	0.818%
80	12.135	1680	1681	1683	rBV2	1637	1455	2.85%	0.351%
81	12.233	1695	1697	1699	rVB3	2122	1520	2.97%	0.366%
82	12.434	1727	1730	1731	rBV3	1457	1301	2.55%	0.313%
83	12.538	1744	1747	1748	rVB3	1924	1369	2.68%	0.330%
84	12.562	1748	1751	1753	rBV4	1972	2429	4.75%	0.585%
85	12.599	1755	1757	1759	rVV3	3438	3052	5.97%	0.735%
86	12.696	1768	1773	1779	rVB6	6199	12627	24.70%	3.042%
87	12.867	1798	1801	1802	rBV3	1859	1700	3.33%	0.410%
88	12.989	1819	1821	1823	rVB3	2097	1207	2.36%	0.291%
89	13.556	1909	1914	1919	rBV3	8798	16905	33.07%	4.073%
90	13.672	1932	1933	1938	rBV5	2176	3653	7.15%	0.880%
91	13.849	1957	1962	1971	rBV5	8663	17741	34.71%	4.274%
92	14.025	1987	1991	1995	rBV7	1703	3594	7.03%	0.866%
93	14.056	1995	1996	2000	rVB4	1944	1493	2.92%	0.360%
94	14.178	2014	2016	2019	rVB4	1709	1653	3.23%	0.398%
95	14.623	2087	2089	2091	rBV3	1670	1186	2.32%	0.286%
96	14.812	2119	2120	2122	rBV2	2351	1755	3.43%	0.423%
97	14.909	2134	2136	2137	rBV2	2217	1423	2.78%	0.343%
98	15.501	2231	2233	2237	rBV5	1872	3287	6.43%	0.792%
99	16.458	2388	2390	2392	rBV3	2142	1367	2.67%	0.329%
100	16.555	2404	2406	2409	rVB4	1881	1704	3.33%	0.411%

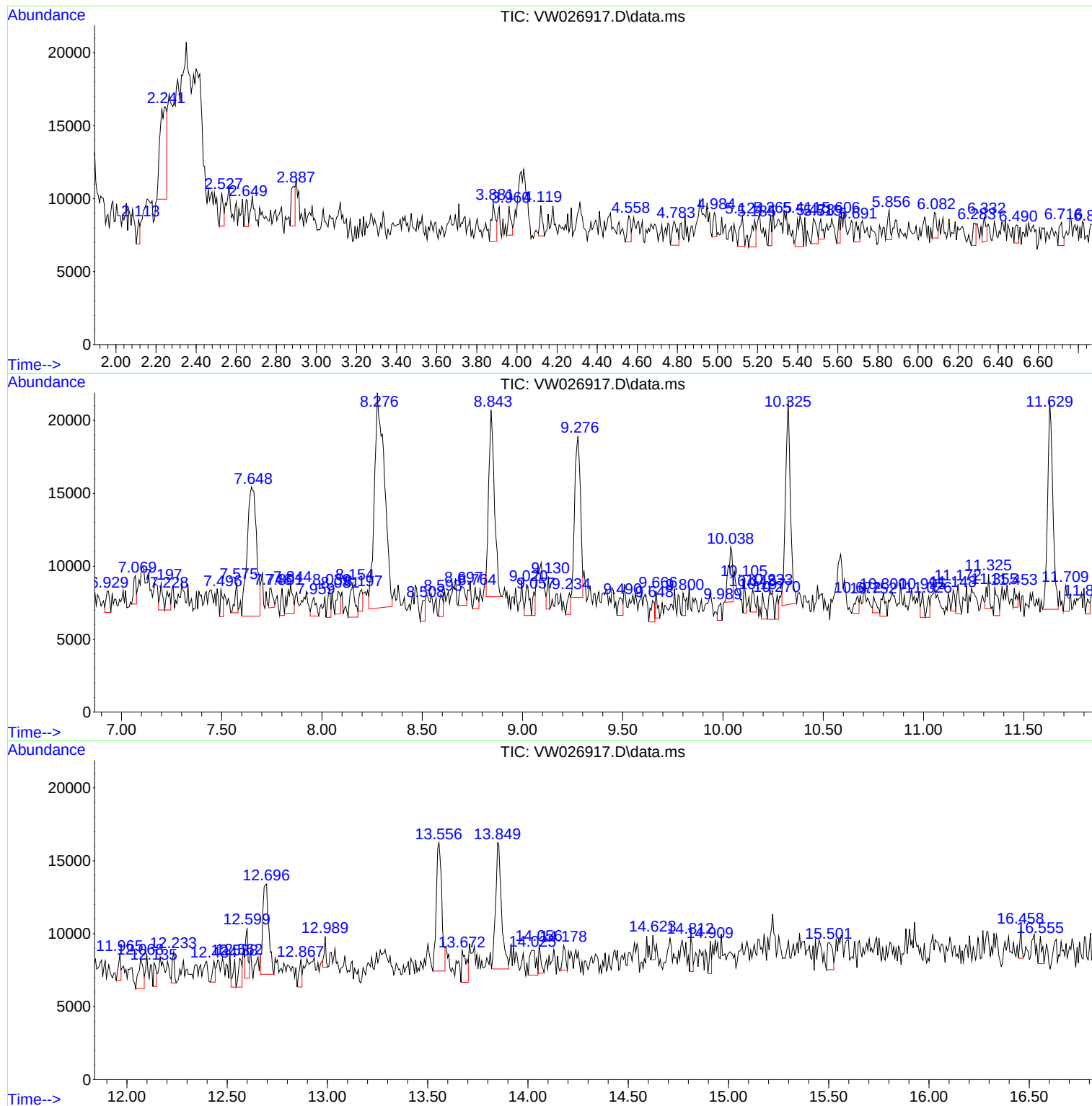
Sum of corrected areas: 415086

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Instrument :
MSVOA_W
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
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Instrument :
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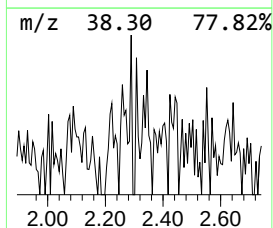
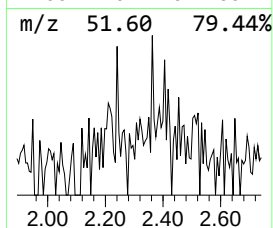
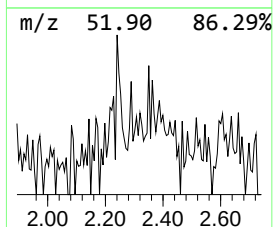
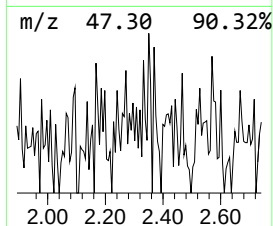
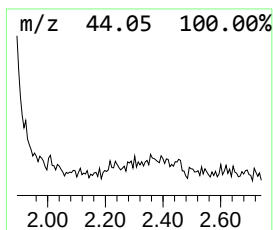
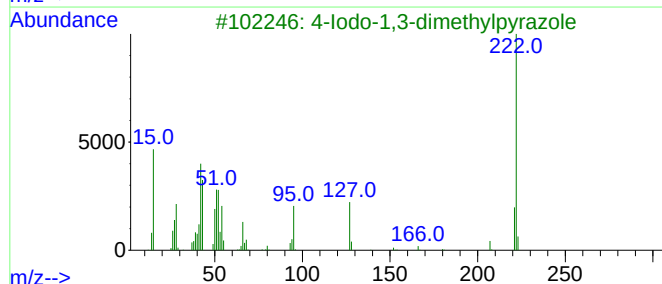
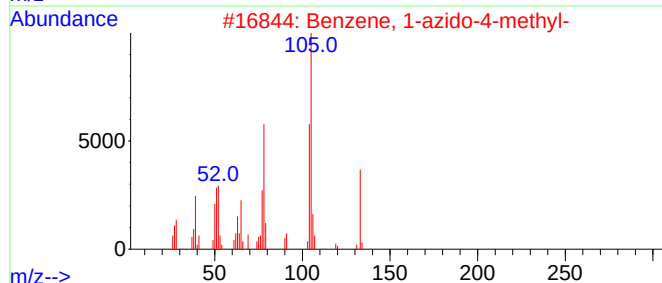
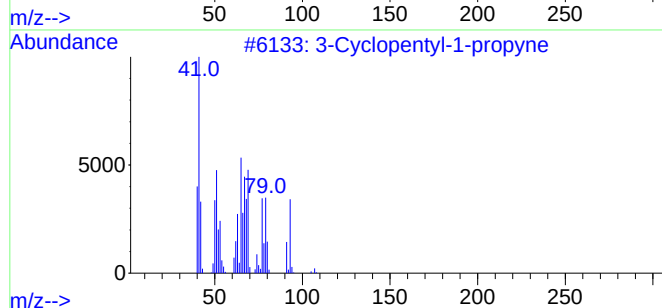
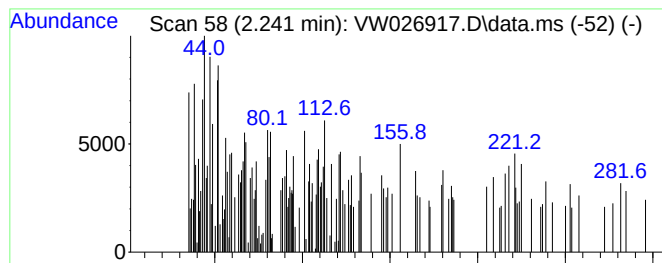
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.241	14.89 ug/L	14900	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Cyclopentyl-1-propyne	108	C8H12	116279-08-4	35
2			Benzene, 1-azido-4-methyl-	133	C7H7N3	002101-86-2	22
3			4-Iodo-1,3-dimethylpyrazole	222	C5H7IN2	006647-97-8	17
4			Quinoline, 2-methyl-4-[5-(2-meth...	291	C17H13N3O2	1000337-11-3	16
5			Biphenyl, 4-(furfurylidenamino)-...	292	C17H12N2O3	1000259-33-4	12



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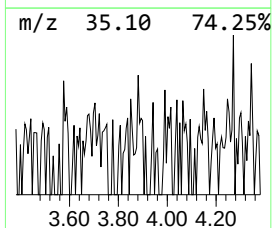
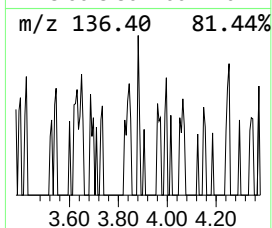
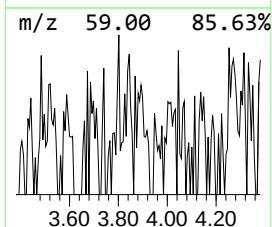
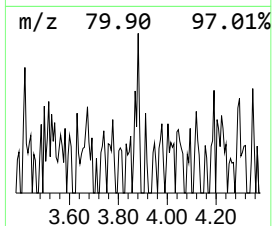
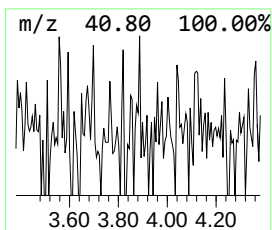
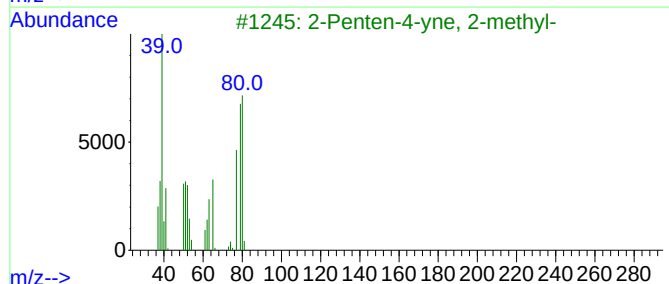
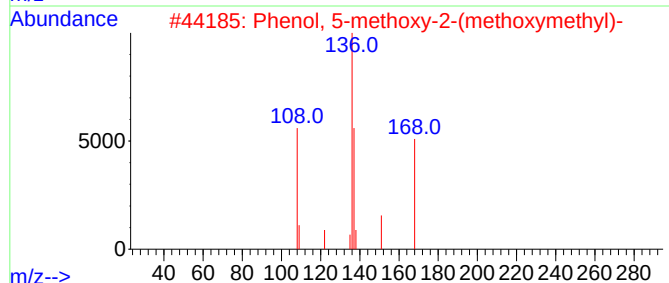
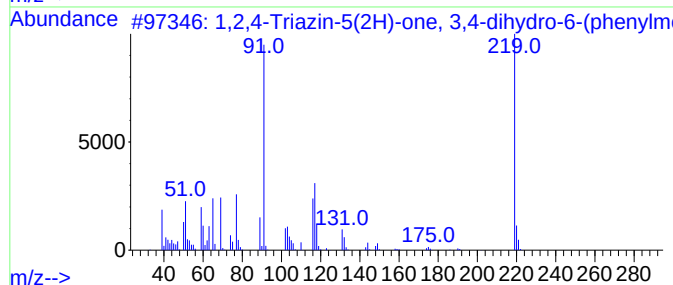
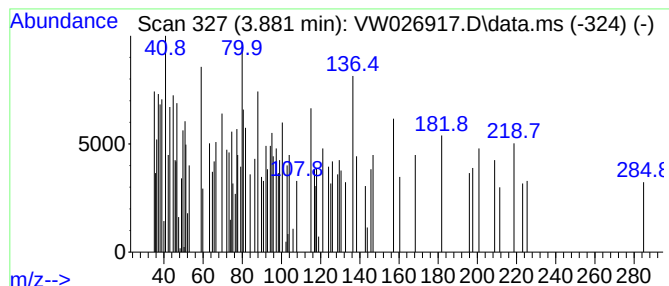
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.881	3.82 ug/L	3820	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,4-Triazin-5(2H)-one, 3,4-dih...	219	C10H9N3OS	1000319-59-8	35
2			Phenol, 5-methoxy-2-(methoxymeth...	168	C9H12O3	062849-09-6	32
3			2-Penten-4-yne, 2-methyl-	80	C6H8	001595-53-5	32
4			3-Butenenitrile, 2-methyl-	81	C5H7N	016529-56-9	27
5			2-(2-Hydroxyphenoxy)-1-phenyleth...	230	C14H14O3	328104-89-8	16



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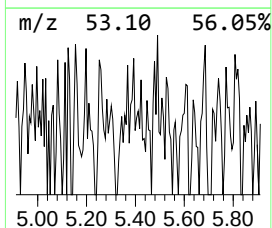
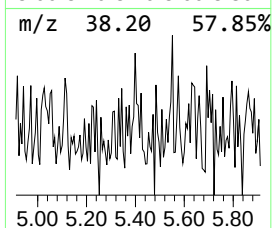
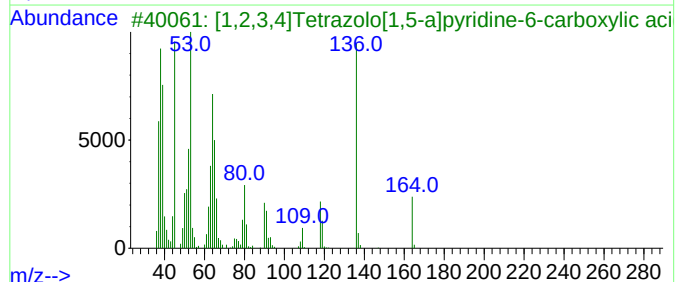
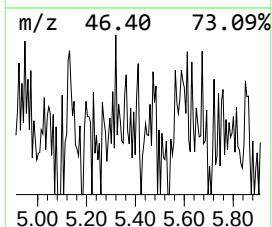
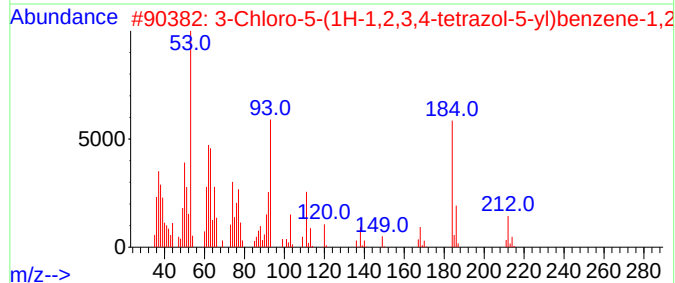
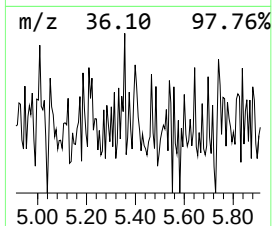
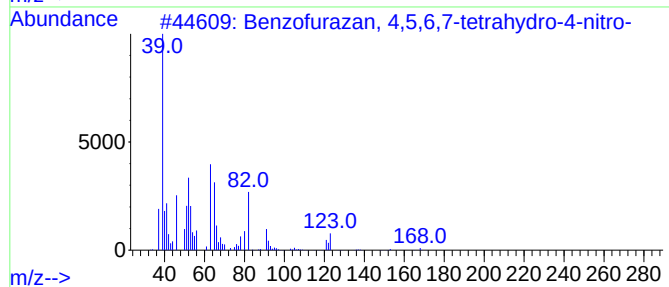
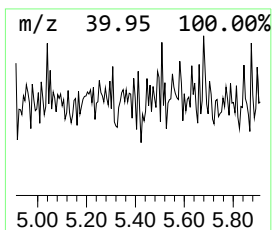
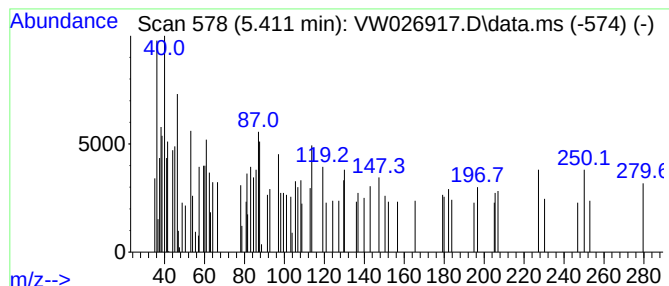
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.411	3.33 ug/L	3334	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofurazan, 4,5,6,7-tetrahydro...	169	C6H7N3O3	1000262-66-9	23
2			3-Chloro-5-(1H-1,2,3,4-tetrazol-...	212	C7H5ClN4O2	1000409-86-2	22
3			[1,2,3,4]Tetrazolo[1,5-a]pyridin...	164	C6H4N4O2	1000362-60-7	14
4			Benzoic acid, 2,5-dichloro-3,6-d...	264	C10H10Cl2O4	062059-39-6	10
5			1-(4-Amino-furazan-3-yl)-5-methy...	250	C9H10N6O3	1000300-81-0	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYX0

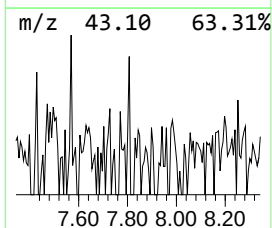
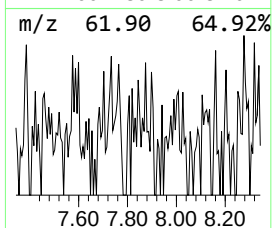
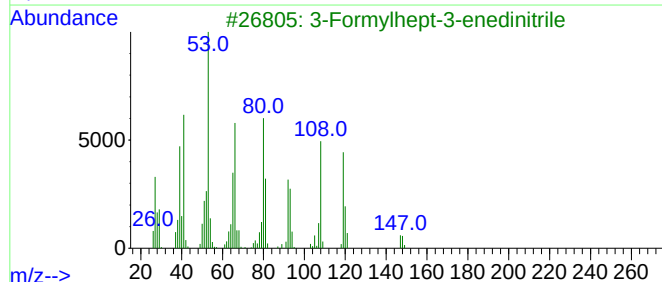
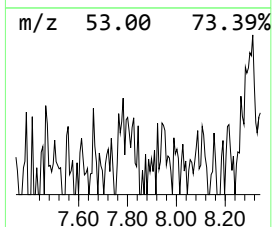
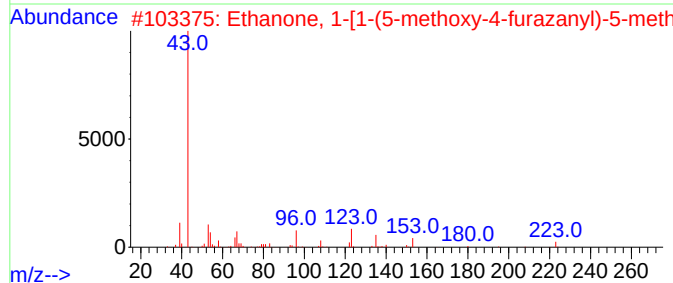
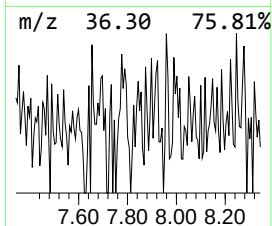
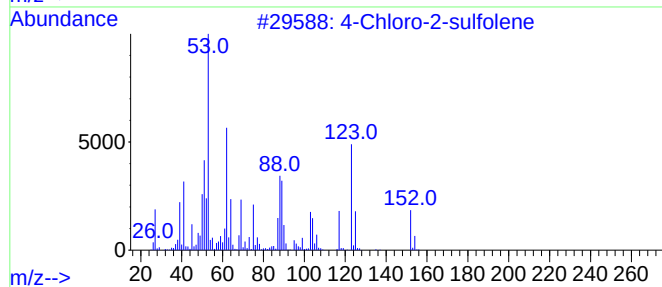
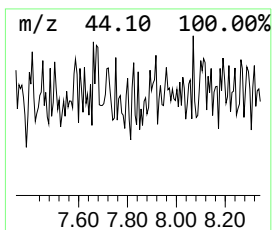
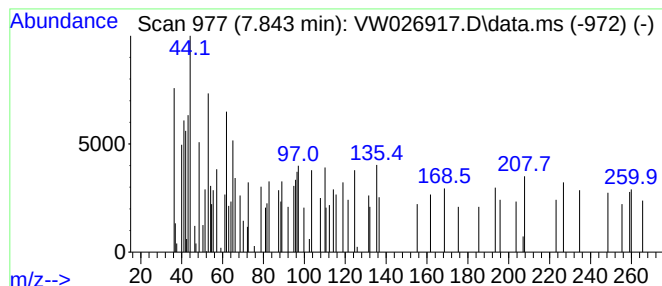
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-04 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.843	3.51 ug/L	3509	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-2-sulfolene	152	C4H5ClO2S	042925-42-8	43
2			Ethanone, 1-[1-(5-methoxy-4-fura...	223	C8H9N5O3	298685-74-2	35
3			3-Formylhept-3-enedinitrile	148	C8H8N2O	1000192-13-5	25
4			2-Butenal, 2-ethenyl-	96	C6H8O	020521-42-0	25
5			Phenol, 2,6-dinitro-4-(trifluoro...	252	C7H3F3N2O5	000393-77-1	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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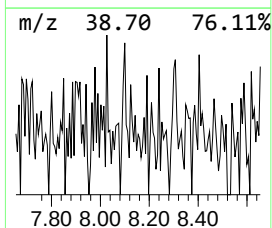
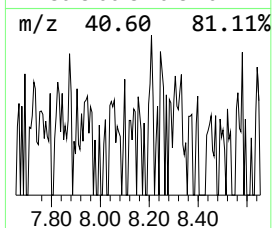
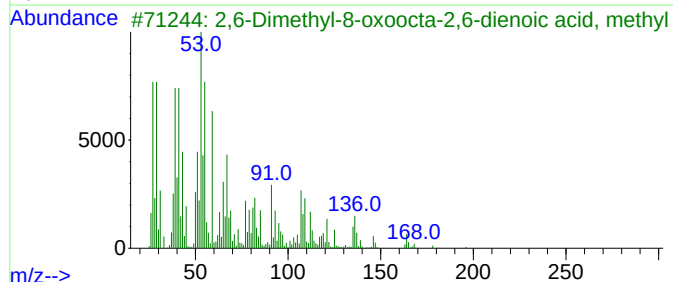
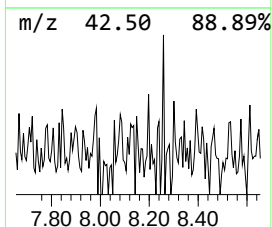
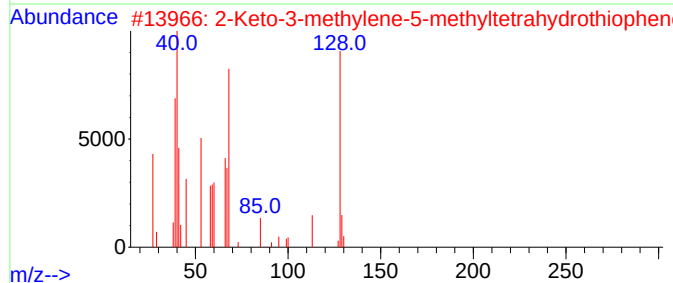
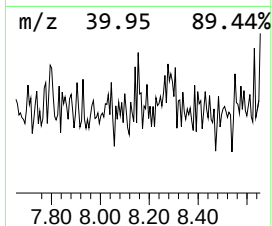
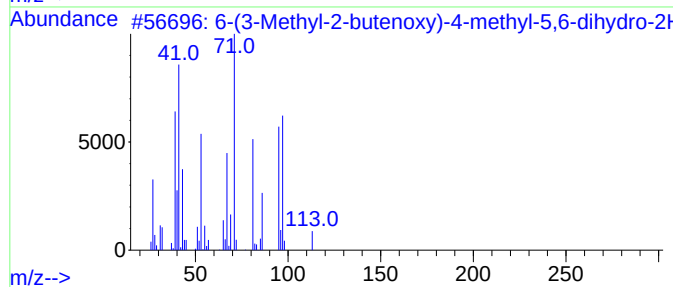
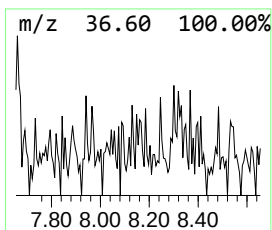
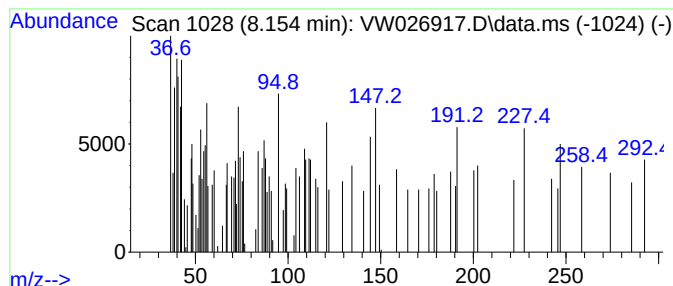
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.154	3.78 ug/L	3781	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-(3-Methyl-2-butenoxy)-4-methyl...	182	C11H18O2	1000432-15-5	33		
2	2-Keto-3-methylene-5-methyltetra...	128	C6H8OS	072359-73-0	32		
3	2,6-Dimethyl-8-oxoocta-2,6-dieno...	196	C11H16O3	1000195-67-9	28		
4	Ethyl 2-heptynoate	154	C9H14O2	016930-95-3	25		
5	4-Chloro-N-(1,2-dihydro-2-pyridy...	268	C11H9ClN2O2S	001213-38-3	22		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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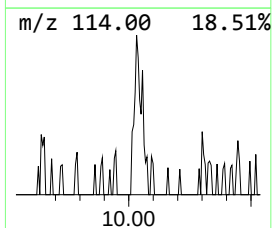
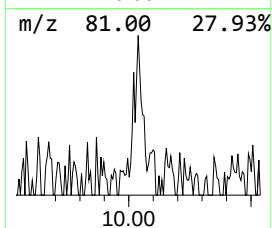
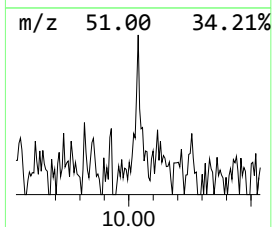
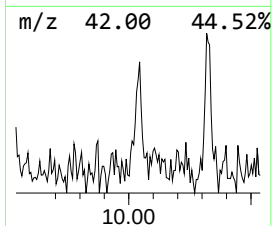
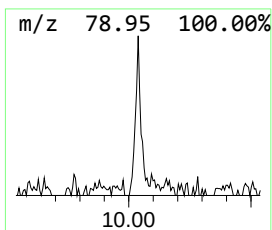
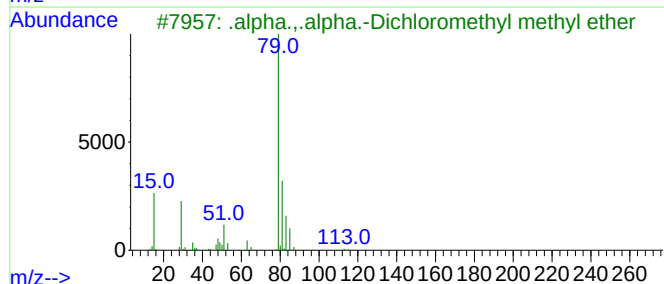
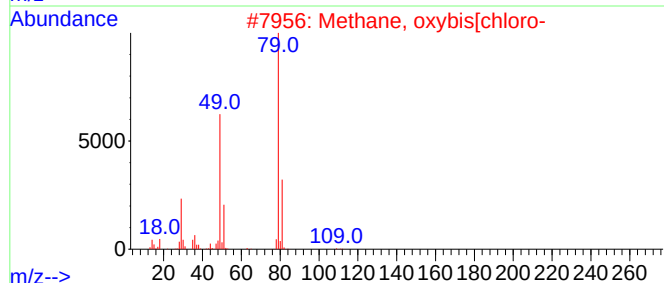
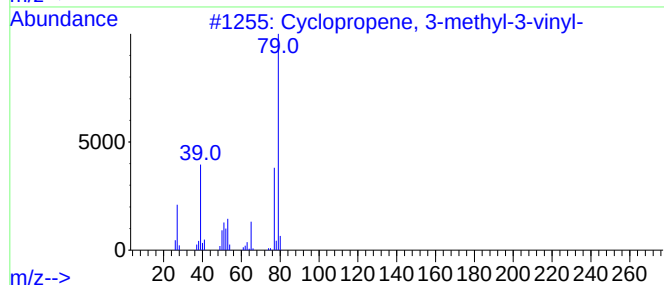
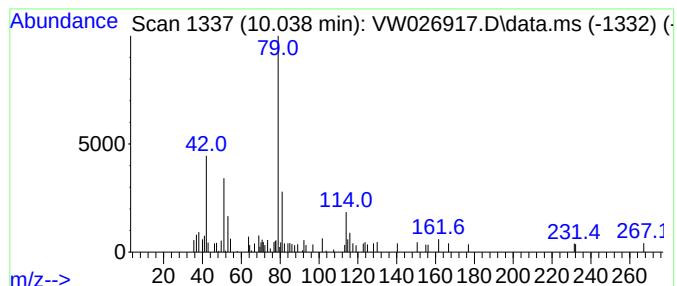
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Quant Title : SFAM01.0

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

Peak Number 6 unknown-06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.038	5.62 ug/L	5624	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropene, 3-methyl-3-vinyl-	80	C6H8	071153-30-5	32
2			Methane, oxybis[chloro-	114	C2H4Cl2O	000542-88-1	25
3			.alpha.,.alpha.-Dichloromethyl m...	114	C2H4Cl2O	004885-02-3	25
4			2-Hexen-4-yne	80	C6H8	014092-20-7	23
5			Dicyclopropylmethanol, chlorodif...	224	C9H11ClF2O2	1000376-25-1	16



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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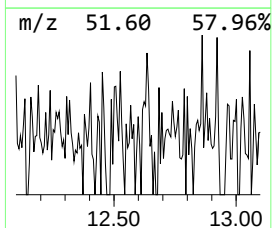
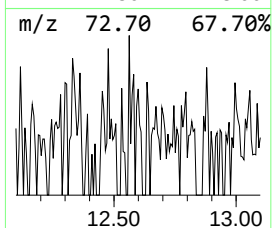
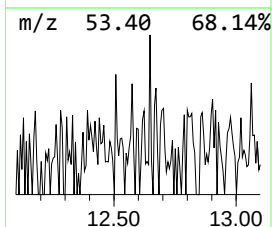
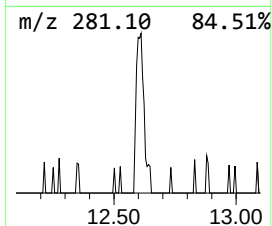
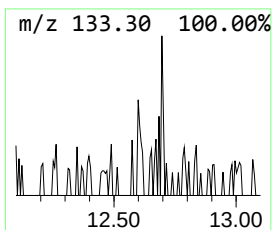
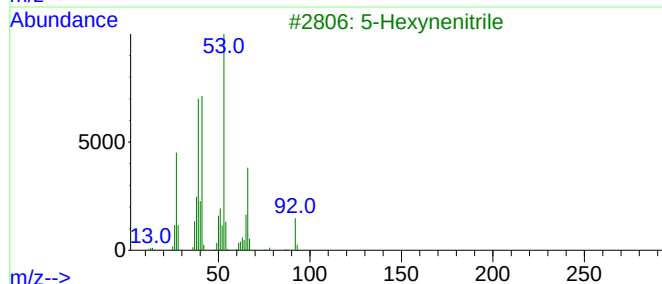
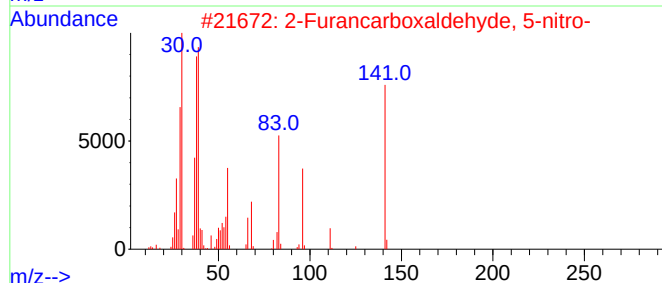
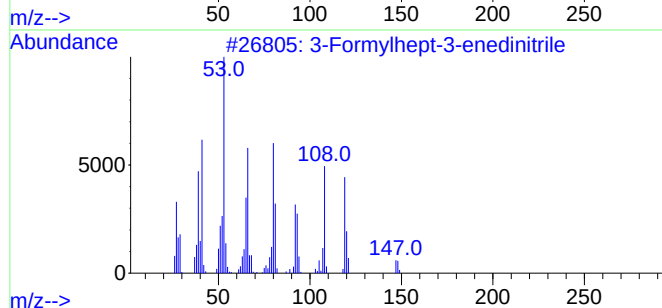
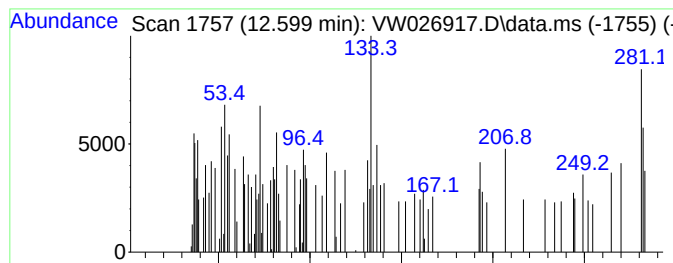
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-07 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.599	4.51 ug/L	3052	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Formylhept-3-enedinitrile	148	C8H8N2O	1000192-13-5	10
2			2-Furancarboxaldehyde, 5-nitro-	141	C5H3NO4	000698-63-5	10
3			5-Hexynenitrile	93	C6H7N	014918-21-9	10
4			2-Pentenedinitrile	92	C5H4N2	007717-24-0	10
5			1,3-Butadiene-1-carboxylic acid	98	C5H6O2	000626-99-3	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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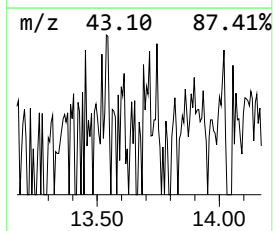
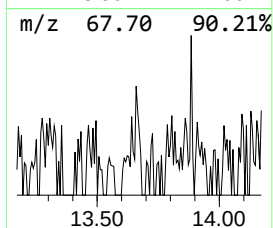
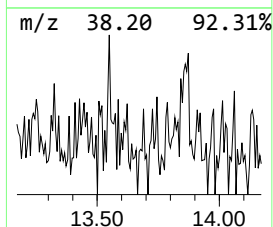
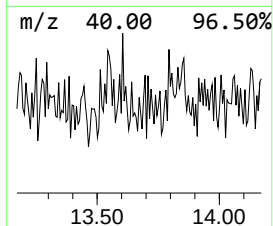
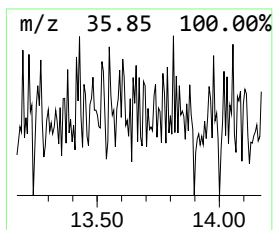
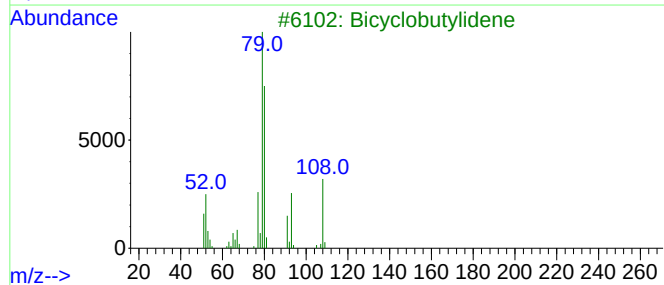
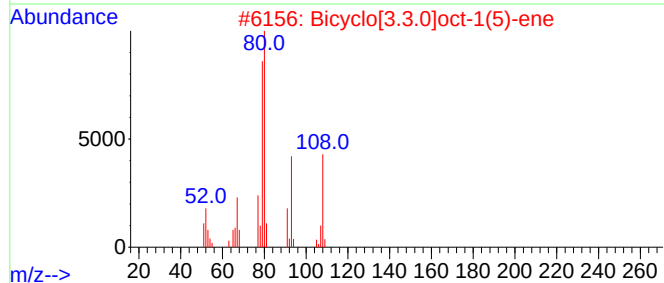
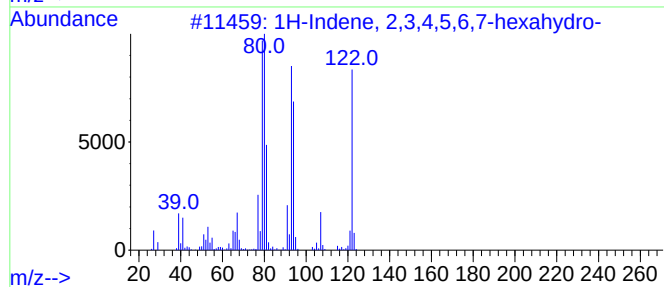
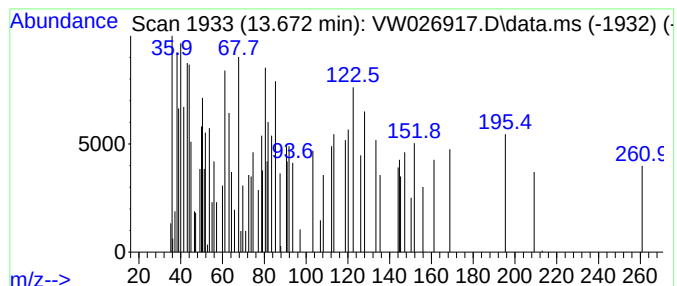
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.672	5.40 ug/L	3653	1,4-Dichlorobenzene-d4	13.556

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3,4,5,6,7-hexahydro-	122	C9H14	000695-90-9	35
2		Bicyclo[3.3.0]oct-1(5)-ene	108	C8H12	006491-93-6	25
3		Bicyclobutylidene	108	C8H12	006708-14-1	25
4		(1H)-3a,4,5,6,7,7a-Hexahydroindene	122	C9H14	1000196-59-5	18
5		Bicyclobutylidene	108	C8H12	006708-14-1	15



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

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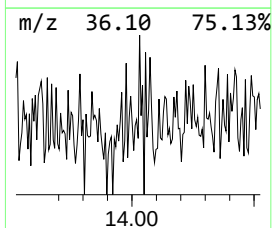
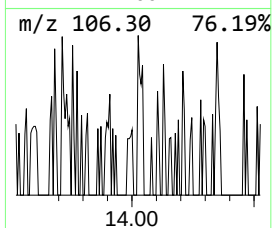
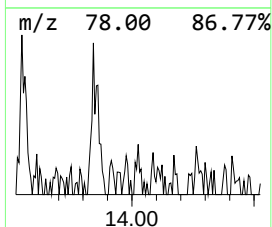
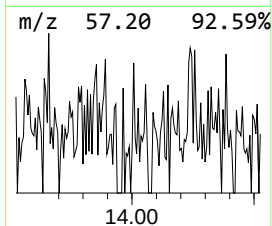
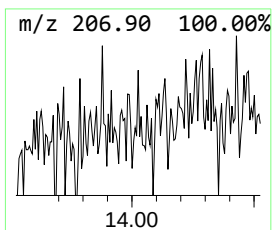
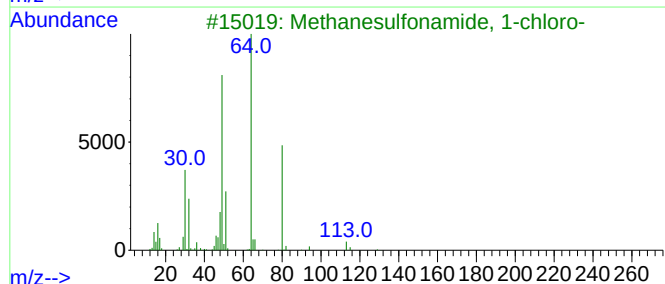
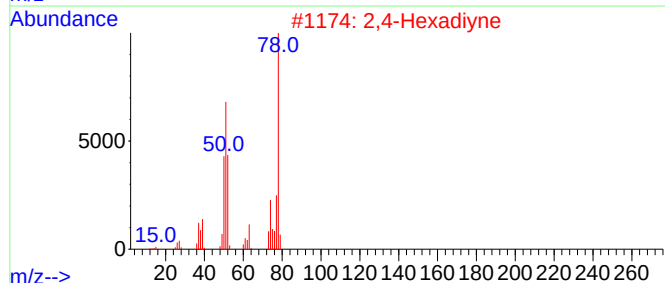
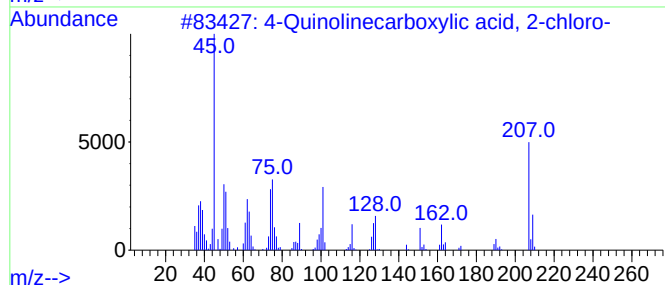
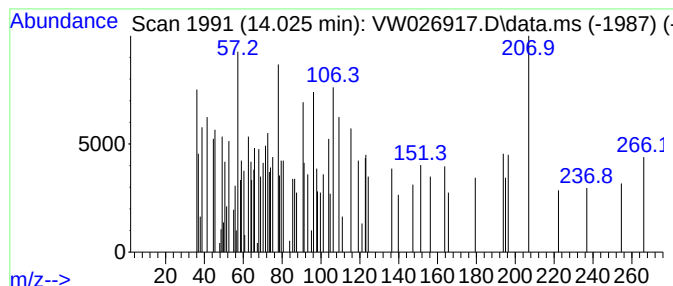
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-09 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.025	5.32 ug/L	3594	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Quinolinecarboxylic acid, 2-ch...	207	C10H6ClNO2	005467-57-2	14
2			2,4-Hexadiyne	78	C6H6	002809-69-0	10
3			Methanesulfonamide, 1-chloro-	129	CH4ClNO2S	021335-43-3	10
4			1-(Dicyanomethyl)eth-1-ene-1,2,2...	221	C8H7N5O3	1000388-01-7	10
5			1-Ethynylcyclohex-2-en-1-ol	122	C8H10O	038693-58-2	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

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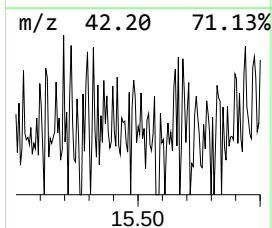
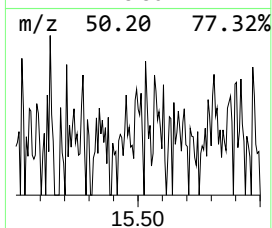
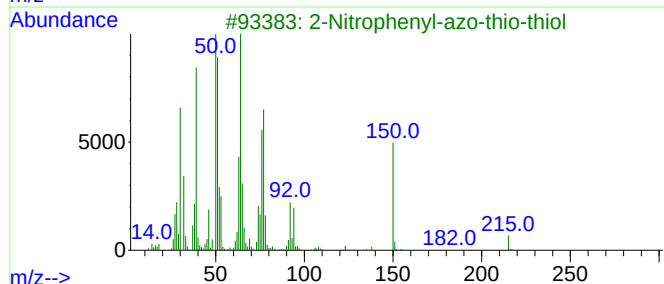
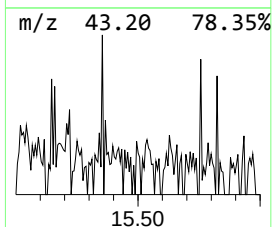
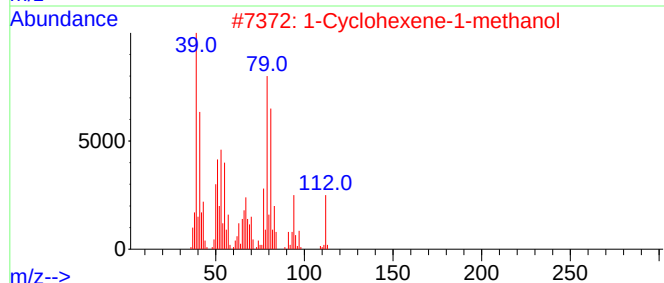
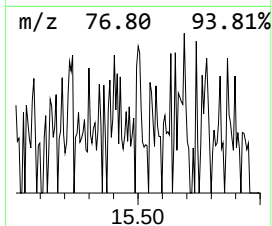
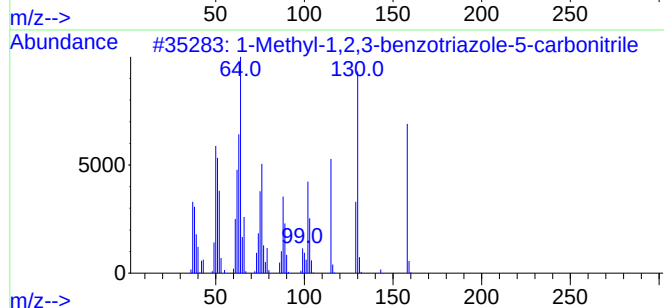
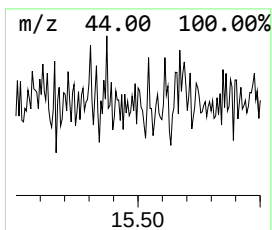
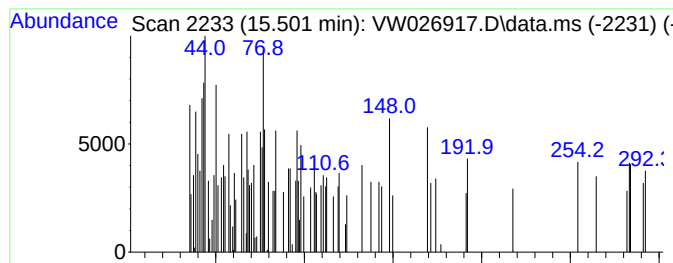
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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

Peak Number 10 unknown-10 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.501	4.86 ug/L	3287	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-1,2,3-benzotriazole-5-c...	158	C8H6N4	1000435-82-2	47
2			1-Cyclohexene-1-methanol	112	C7H12O	004845-04-9	35
3			2-Nitrophenyl-azo-thio-thiol	215	C6H5N3O2S2	1000286-39-5	25
4			1,3-Butadiyne	50	C4H2	000460-12-8	25
5			2-(3-Methyl-4-nitro-5-pyrazolyla...	262	C11H10N4O4	1000223-64-7	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082523\
Data File : VW026917.D
Acq On : 26 Aug 2023 01:03
Operator : SY/MD
Sample : 04175-05
Misc : 5.76g/10mL/MSVOA_W/SOIL
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYX0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.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
unknown-01	2.241	14.9	ug/L	14900	1	8.843	25009	25.0
unknown-02	3.881	3.8	ug/L	3820	1	8.843	25009	25.0
unknown-03	5.411	3.3	ug/L	3334	1	8.843	25009	25.0
unknown-04	7.843	3.5	ug/L	3509	1	8.843	25009	25.0
unknown-05	8.154	3.8	ug/L	3781	1	8.843	25009	25.0
unknown-06	10.038	5.6	ug/L	5624	1	8.843	25009	25.0
unknown-07	12.599	4.5	ug/L	3052	3	13.556	16905	25.0
unknown-08	13.672	5.4	ug/L	3653	3	13.556	16905	25.0
unknown-09	14.025	5.3	ug/L	3594	3	13.556	16905	25.0
unknown-10	15.501	4.9	ug/L	3287	3	13.556	16905	25.0