

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
 Data File : VW026951.D
 Acq On : 29 Aug 2023 15:23
 Operator : SY/MD
 Sample : 04175-22
 Misc : 5.44g/10mL/MSVOA_W/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EZYW5

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: VW026951.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.247	50	59	62	rBV10	12732	40027	53.29%	7.326%
2	2.503	99	101	106	rVB4	3518	3994	5.32%	0.731%
3	2.588	113	115	117	rVB3	1585	1163	1.55%	0.213%
4	2.808	149	151	156	rVB6	1921	2267	3.02%	0.415%
5	2.899	161	166	172	rVB	5933	12244	16.30%	2.241%
6	3.021	182	186	187	rBV4	2225	2033	2.71%	0.372%
7	3.033	187	188	190	rBV2	1773	1253	1.67%	0.229%
8	3.052	190	191	194	rVB3	2537	2222	2.96%	0.407%
9	3.076	194	195	197	rBV2	2271	1674	2.23%	0.306%
10	3.235	217	221	223	rBV5	1814	2335	3.11%	0.427%
11	3.314	232	234	238	rBV5	1876	2717	3.62%	0.497%
12	3.357	240	241	243	rVV2	1414	1405	1.87%	0.257%
13	3.393	245	247	252	rVB5	1255	1795	2.39%	0.329%
14	3.478	256	261	263	rVB6	1653	2036	2.71%	0.373%
15	3.509	263	266	268	rBV4	1078	1434	1.91%	0.262%
16	3.667	289	292	293	rBV3	1316	1228	1.63%	0.225%
17	3.710	296	299	300	rBV3	1465	1439	1.92%	0.263%
18	3.741	302	304	307	rVB4	2467	2682	3.57%	0.491%
19	3.777	307	310	312	rBV4	1172	1444	1.92%	0.264%
20	3.802	312	314	317	rVV4	1176	1315	1.75%	0.241%
21	4.027	343	351	358	rBV5	8120	21997	29.29%	4.026%
22	4.125	364	367	370	rBV5	3291	5110	6.80%	0.935%
23	4.515	428	431	433	rBV4	1535	1518	2.02%	0.278%
24	4.545	433	436	438	rVB4	1580	1886	2.51%	0.345%
25	4.582	441	442	444	rBV2	1734	1268	1.69%	0.232%
26	4.643	449	452	454	rVV4	1346	1636	2.18%	0.299%
27	4.698	459	461	462	rVV2	1502	1237	1.65%	0.226%
28	4.734	466	467	470	rBV3	1124	1294	1.72%	0.237%
29	4.759	470	471	475	rVB4	1318	1304	1.74%	0.239%
30	4.923	492	498	499	rBV6	2175	4001	5.33%	0.732%
31	4.942	499	501	506	rVV6	2393	3863	5.14%	0.707%
32	5.039	515	517	518	rBV2	1794	1337	1.78%	0.245%
33	5.112	526	529	530	rBV3	1484	1951	2.60%	0.357%
34	5.173	536	539	541	rBV4	2319	2851	3.80%	0.522%
35	5.228	545	548	549	rBV3	1683	1227	1.63%	0.225%
36	5.399	573	576	578	rVB4	1731	1879	2.50%	0.344%

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Integrator: RTE

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

37	5.435	578	582	584	rBV5	1841	3230	4.30%	0.591%
38	5.460	584	586	587	rVB2	2583	1606	2.14%	0.294%
39	5.484	587	590	591	rBV3	1556	1737	2.31%	0.318%
40	5.503	591	593	594	rBV2	1540	1230	1.64%	0.225%
41	5.521	594	596	599	rVB4	1300	1273	1.69%	0.233%
42	5.551	599	601	603	rVB3	1378	1221	1.63%	0.223%
43	5.600	607	609	613	rVB5	1716	1921	2.56%	0.352%
44	5.643	613	616	617	rBV2	1932	1801	2.40%	0.330%
45	5.692	621	624	625	rBV3	1922	1955	2.60%	0.358%
46	5.795	638	641	642	rBV3	2477	2619	3.49%	0.479%
47	5.813	642	644	646	rVV3	1633	1314	1.75%	0.240%
48	5.856	646	651	654	rVB6	1505	2171	2.89%	0.397%
49	5.923	661	662	663	rBV	1976	1168	1.56%	0.214%
50	6.021	676	678	680	rBV3	1942	1558	2.07%	0.285%
51	6.076	683	687	688	rBV4	1425	1485	1.98%	0.272%
52	6.100	688	691	694	rVB5	1136	1468	1.95%	0.269%
53	6.143	694	698	701	rBV6	1854	3193	4.25%	0.584%
54	6.185	703	705	706	rVV2	1949	1462	1.95%	0.268%
55	6.258	715	717	720	rVB4	2245	2293	3.05%	0.420%
56	6.307	723	725	728	rBV4	1143	1445	1.92%	0.264%
57	6.472	748	752	756	rVB7	1069	1870	2.49%	0.342%
58	6.533	760	762	765	rBV4	1726	1660	2.21%	0.304%
59	6.575	767	769	771	rVB3	1738	1358	1.81%	0.249%
60	6.697	786	789	792	rBV4	2105	2991	3.98%	0.547%
61	6.893	819	821	824	rVV4	1969	2269	3.02%	0.415%
62	6.953	828	831	833	rBV3	1718	2178	2.90%	0.399%
63	6.984	833	836	839	rVB5	1666	2132	2.84%	0.390%
64	7.051	846	847	850	rBV3	1232	1204	1.60%	0.220%
65	7.088	850	853	854	rBV3	1187	1256	1.67%	0.230%
66	7.167	864	866	868	rBV3	1975	1307	1.74%	0.239%
67	7.246	877	879	883	rVB5	2204	2467	3.28%	0.452%
68	7.295	883	887	889	rBV5	1797	1783	2.37%	0.326%
69	7.325	889	892	896	rVB6	922	1402	1.87%	0.257%
70	7.362	896	898	901	rVB4	2021	1428	1.90%	0.261%
71	7.399	901	904	905	rBV3	1411	1170	1.56%	0.214%
72	7.484	914	918	920	rVB5	1665	1989	2.65%	0.364%
73	7.508	920	922	924	rBV3	1706	1402	1.87%	0.257%
74	7.600	934	937	939	rBV4	1901	2496	3.32%	0.457%
75	7.648	939	945	953	rVV2	13351	33289	44.32%	6.093%
76	7.770	963	965	968	rVB4	1449	1518	2.02%	0.278%

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 Title : SFAM01.0

77	7.801	968	970	972	rVB3	1409	1353	1.80%	0.248%
78	7.874	980	982	990	rVB8	2830	4942	6.58%	0.905%
79	7.941	990	993	994	rBV3	2636	2812	3.74%	0.515%
80	8.002	1001	1003	1006	rVB4	1664	1637	2.18%	0.300%
81	8.057	1009	1012	1013	rVB3	1748	1513	2.01%	0.277%
82	8.106	1016	1020	1022	rBV5	2100	2462	3.28%	0.451%
83	8.136	1022	1025	1026	rBV3	1128	1317	1.75%	0.241%
84	8.276	1040	1048	1061	rBV3	23278	75108	100.00%	13.747%
85	8.417	1070	1071	1072	rBV	2077	1213	1.62%	0.222%
86	8.709	1117	1119	1122	rVB4	1358	1562	2.08%	0.286%
87	8.843	1133	1141	1149	rBV2	24016	53706	71.51%	9.829%
88	8.953	1158	1159	1160	rBV	1894	1189	1.58%	0.218%
89	9.081	1177	1180	1181	rBV3	2287	2320	3.09%	0.425%
90	9.142	1188	1190	1194	rVB5	2301	2676	3.56%	0.490%
91	9.179	1194	1196	1197	rBV2	2240	1394	1.86%	0.255%
92	9.221	1201	1203	1206	rVV4	1541	2297	3.06%	0.420%
93	9.276	1206	1212	1220	rVB2	18985	40482	53.90%	7.409%
94	9.471	1243	1244	1247	rVB3	2030	1638	2.18%	0.300%
95	9.526	1251	1253	1255	rVB3	2374	1831	2.44%	0.335%
96	9.551	1255	1257	1258	rBV2	1547	1467	1.95%	0.268%
97	10.038	1335	1337	1340	rBV3	3932	3996	5.32%	0.731%
98	10.325	1379	1384	1390	rVB	24501	41160	54.80%	7.533%
99	10.581	1423	1426	1427	rBV3	5089	5281	7.03%	0.967%
100	11.629	1594	1598	1603	rBV2	28762	48636	64.75%	8.902%

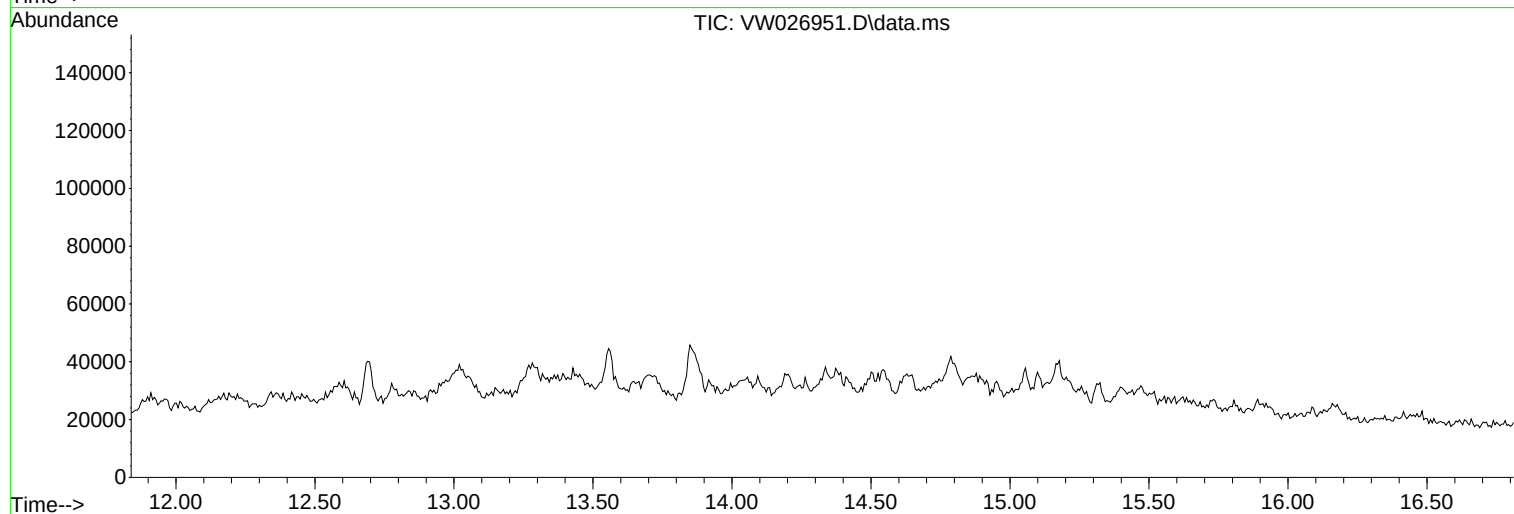
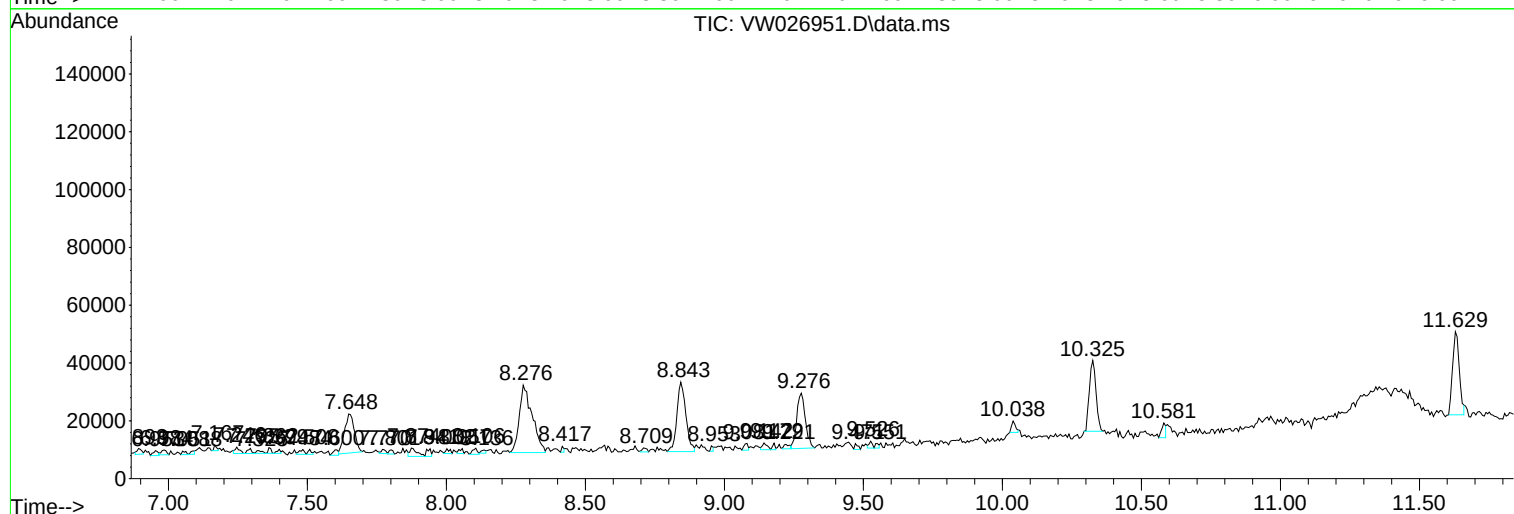
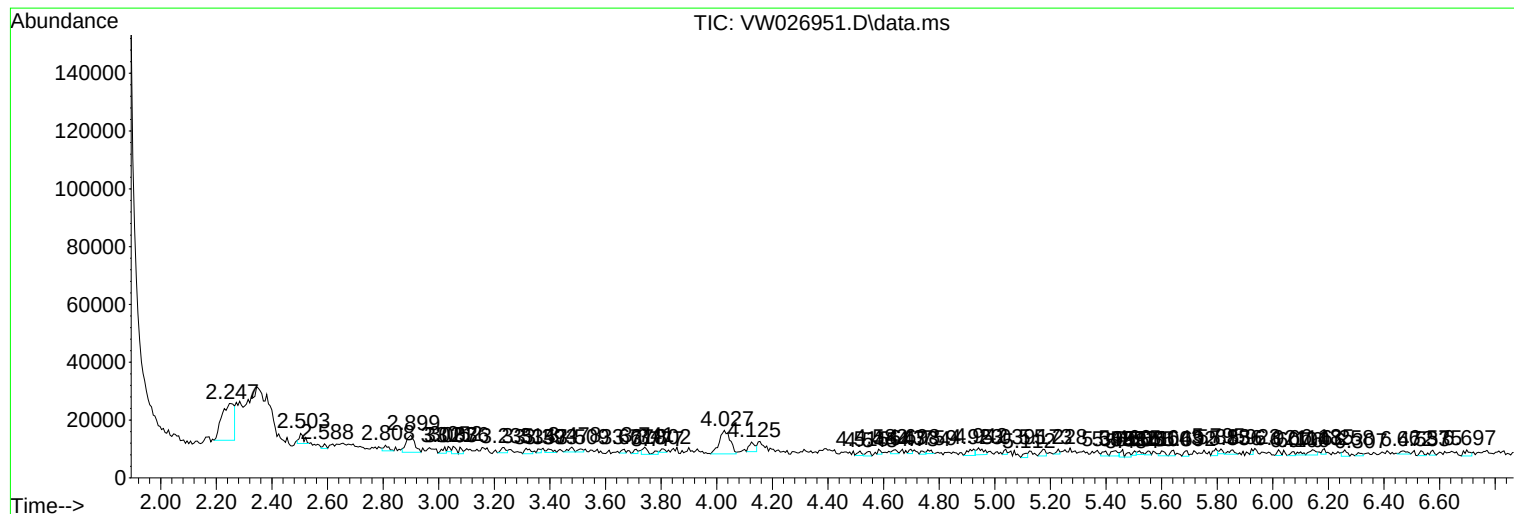
Sum of corrected areas: 546377

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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



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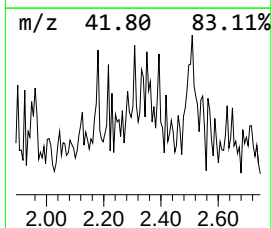
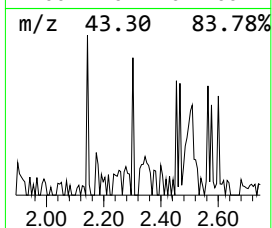
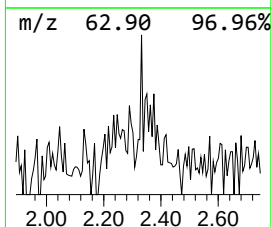
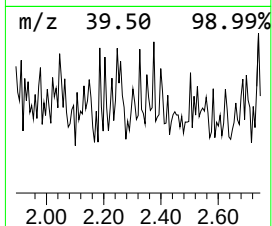
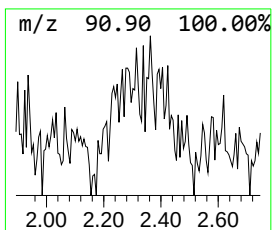
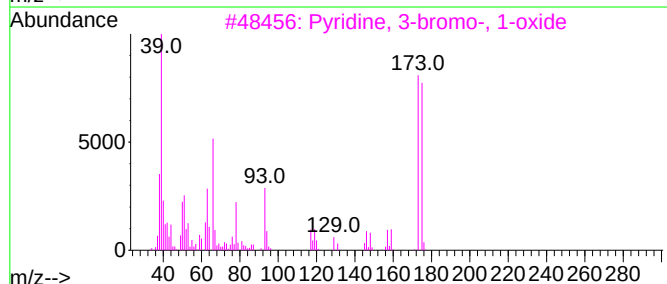
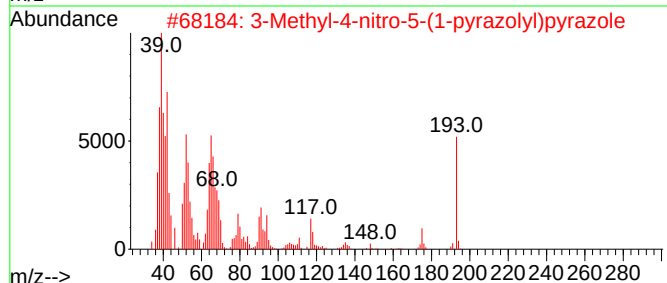
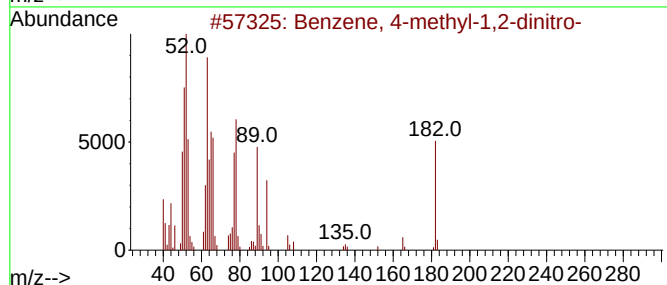
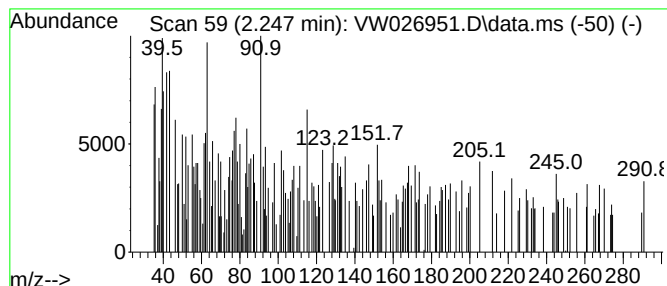
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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.
2.247	18.63 ug/L	40027	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-methyl-1,2-dinitro-	182	C7H6N2O4	000610-39-9	43
2			3-Methyl-4-nitro-5-(1-pyrazolyl)...	193	C7H7N5O2	292855-42-6	38
3			Pyridine, 3-bromo-, 1-oxide	173	C5H4BrNO	002402-97-3	38
4			1-Methoxy-2,5-dinitrobenzene	198	C7H6N2O5	1000362-87-5	35
5			5-Amino-2H-1-benzofuran-3-one	149	C8H7NO2	1174298-02-2	30



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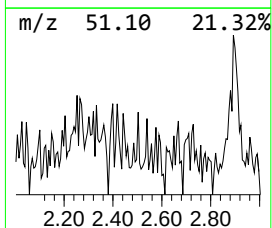
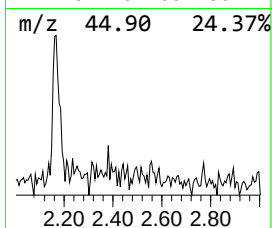
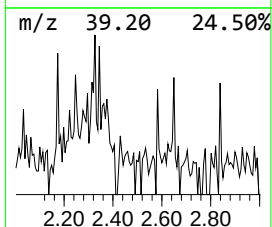
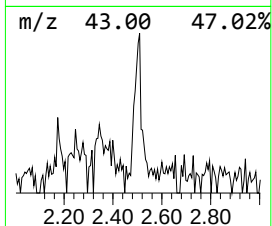
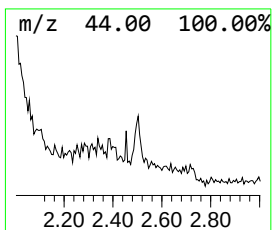
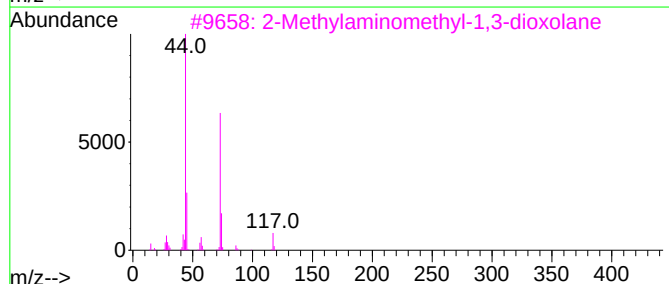
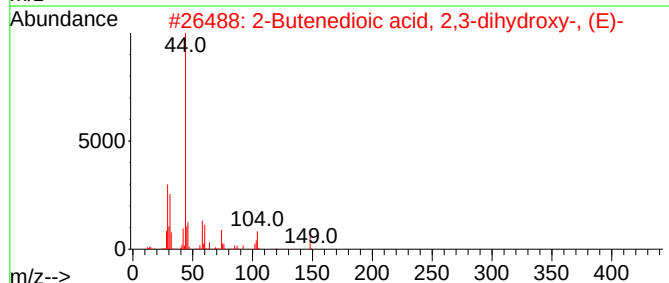
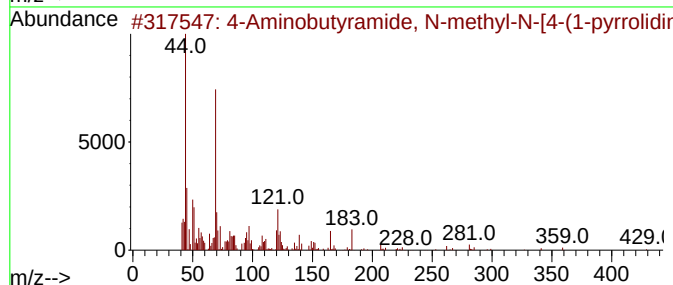
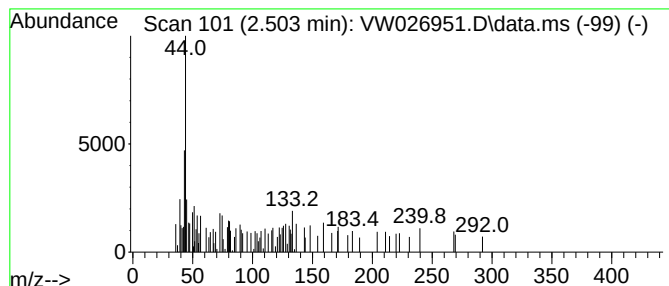
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081023SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.503	1.86 ug/L	3994	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Aminobutyramide, N-methyl-N-[4...	429	C17H21F6N3O3	1000124-10-5	38
2			2-Butenedioic acid, 2,3-dihydrox...	148	C4H4O6	000133-38-0	38
3			2-Methylaminomethyl-1,3-dioxolane	117	C5H11NO2	057366-77-5	38
4			Cyclobutylcarboxamide, N-methyl-...	183	C11H21NO	1000421-30-5	38
5			1,3,5-Trioxepane	104	C4H8O3	005981-06-6	35



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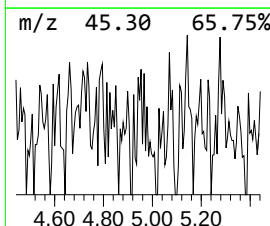
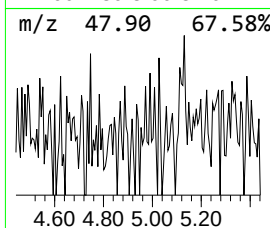
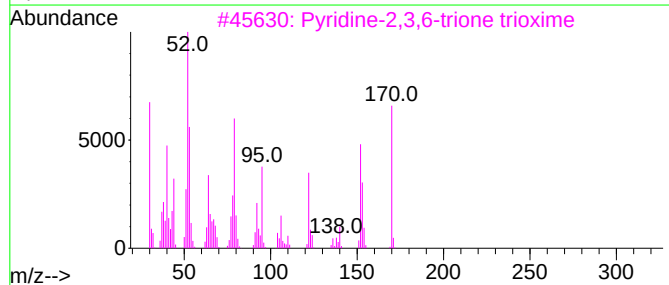
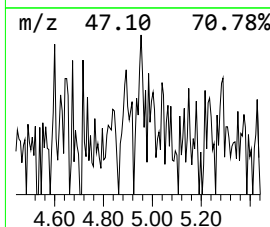
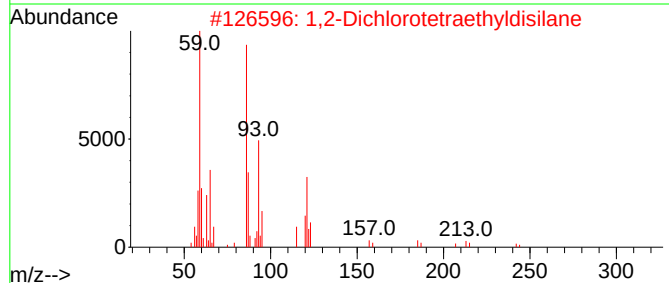
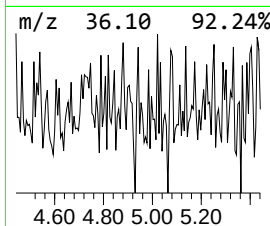
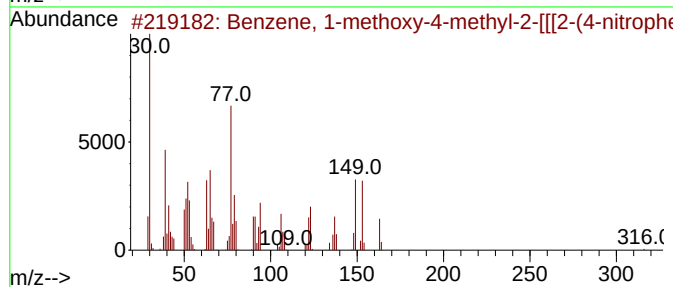
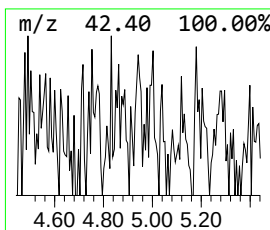
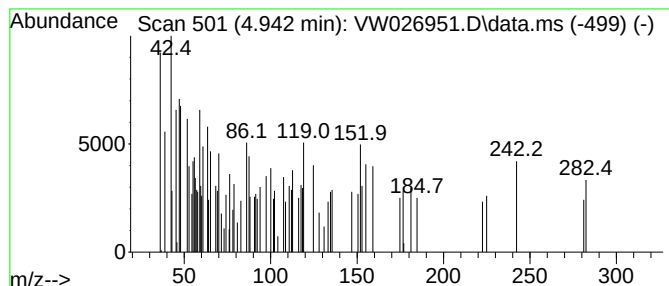
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Quant Title : SFAM01.0

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

Peak Number 4 unknown-03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.942	1.80 ug/L	3863	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methoxy-4-methyl-2-[[...]	316	C15H16N4O4	1000327-64-6	38
2			1,2-Dichlorotetraethylidisilane	242	C8H20Cl2Si2	1000434-03-3	25
3			Pyridine-2,3,6-trione trioxime	170	C5H6N4O3	1000318-31-7	12
4			9,10-Diazatetracyclo[6.2.0(2,7)...	148	C8H8N2O	034098-80-1	12
5			2,6,7-Trimethyl-(1,2,4)-triazolo...	163	C7H9N5	061139-77-3	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYW5

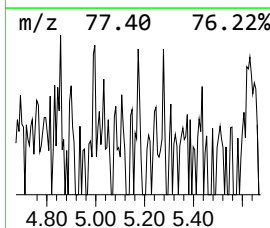
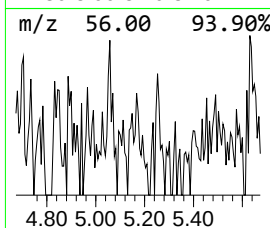
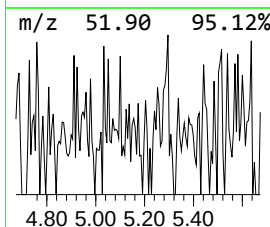
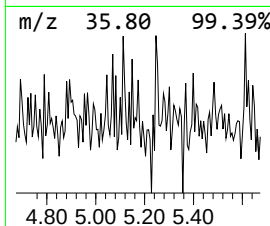
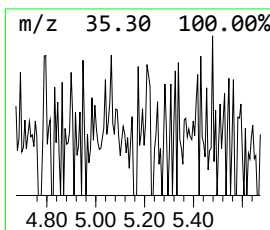
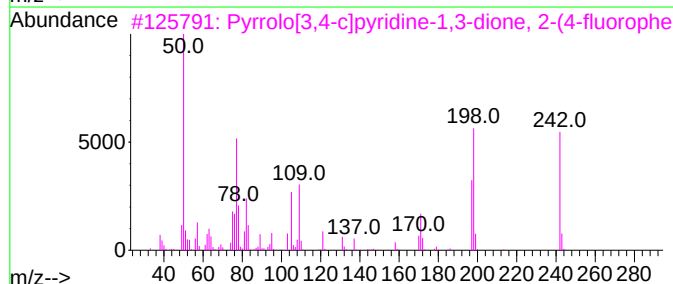
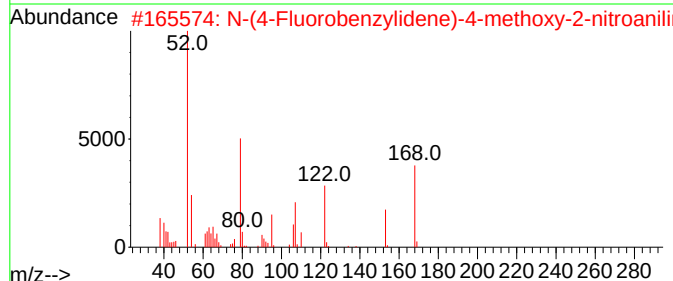
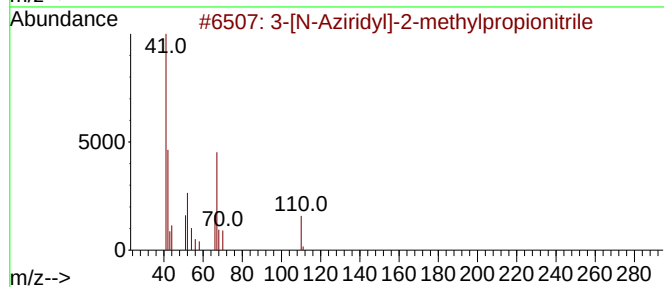
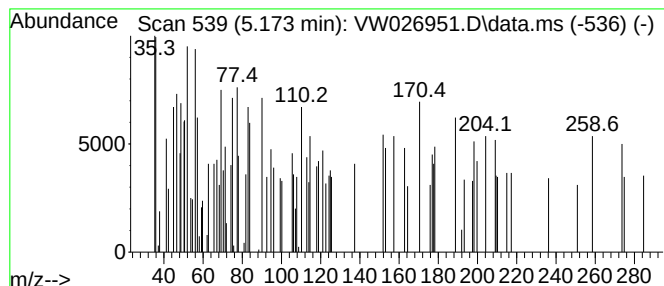
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-04 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.173	1.33 ug/L	2851	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		[N-Aziridyl]-2-methylpropionit...	110	C6H10N2	004078-20-0	14
2	N-(4-Fluorobenzylidene)-4-methox...	274	C14H11FN2O3	1000223-61-1	10		
3	Pyrrolo[3,4-c]pyridine-1,3-dione...	242	C13H7FN2O2	1000310-60-5	10		
4	1,3-Benzenediamine, N,N'-bis(sul...	200	C6H4N2O2S2	017420-01-8	9		
5	Benzene, 1,4-dinitro-	168	C6H4N2O4	000100-25-4	9		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYW5

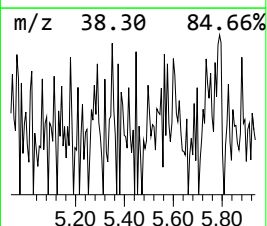
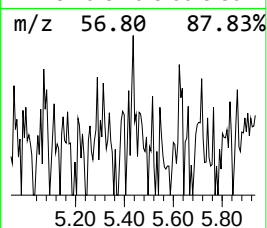
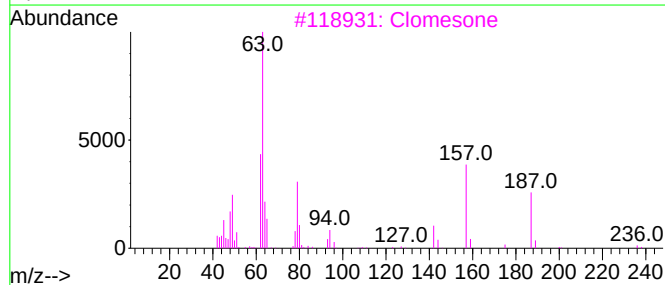
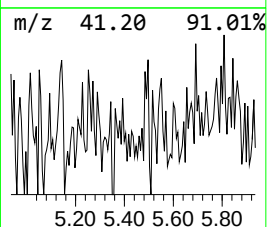
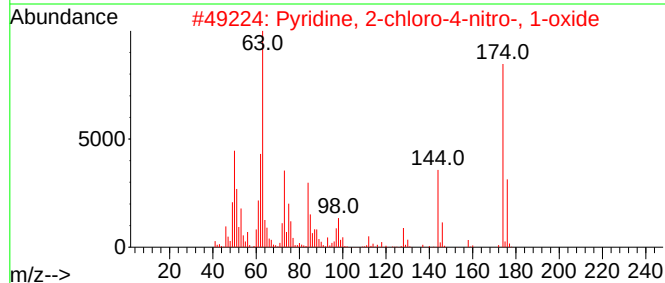
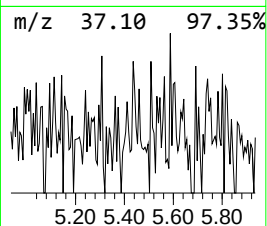
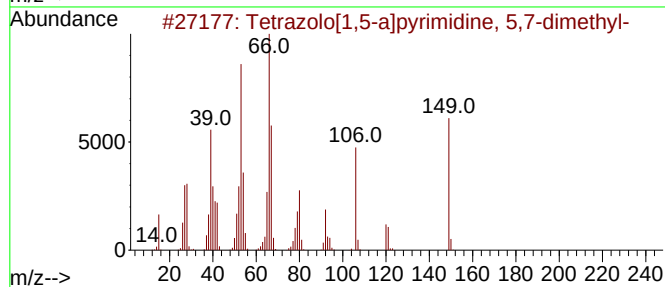
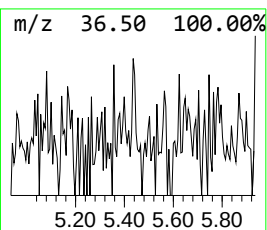
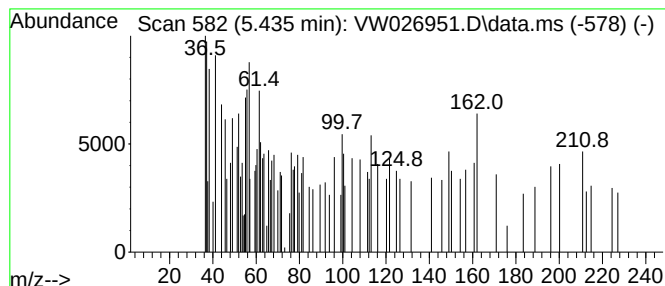
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Quant Title : SFAM01.0

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

Peak Number 6 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.435	1.50 ug/L	3230	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrazolo[1,5-a]pyrimidine, 5,7-...	149	C6H7N5	003210-44-4	35
2			Pyridine, 2-chloro-4-nitro-, 1-o...	174	C5H3ClN2O3	014432-16-7	22
3			Clomesone	236	C4H9ClO5S2	088343-72-0	16
4			Tetrazolo[1,5-a]pyrimidine, 5,7-...	149	C6H7N5	003210-44-4	14
5			Propanenitrile, 3-chloro-2,3-dih...	147	C3H2ClN3O2	139912-33-7	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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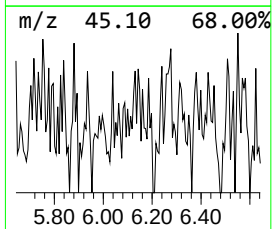
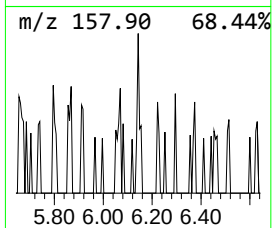
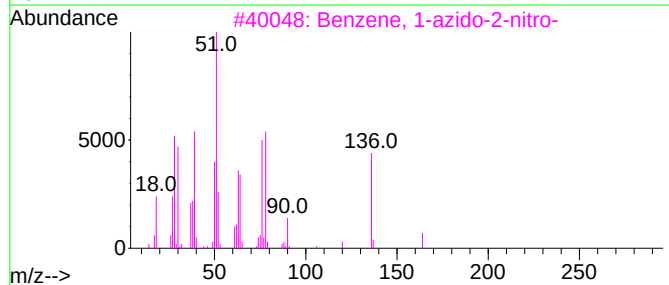
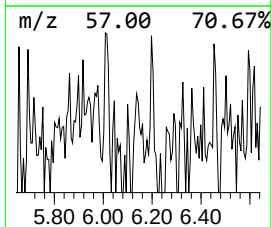
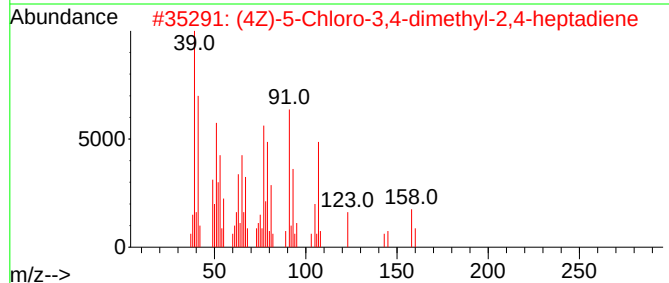
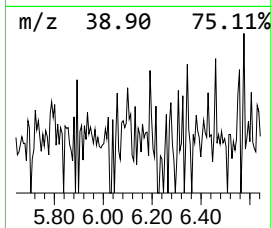
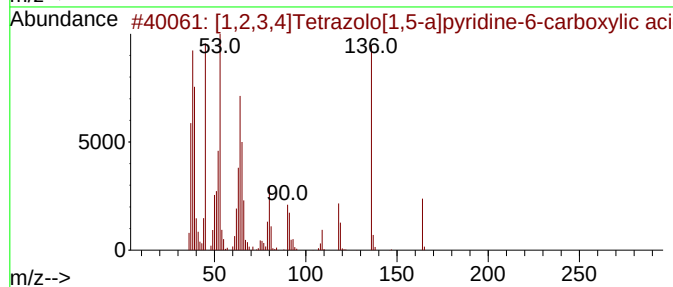
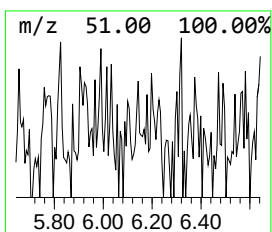
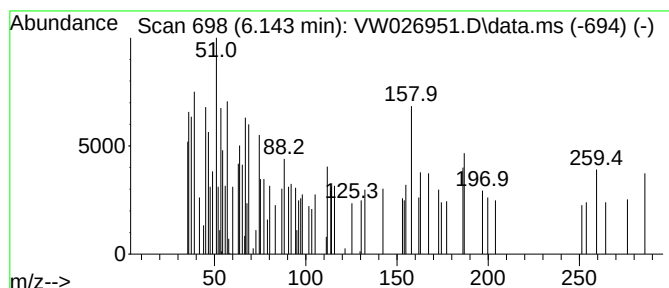
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.143	1.49 ug/L	3193	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,2,3,4]Tetrazolo[1,5-a]pyridin...	164	C6H4N4O2	1000362-60-7	32		
2	(4Z)-5-Chloro-3,4-dimethyl-2,4-h...	158	C9H15Cl	105949-75-5	12		
3	Benzene, 1-azido-2-nitro-	164	C6H4N4O2	001516-58-1	12		
4	N-[(2-Hydroxyphenyl)imino]-2-nit...	259	C12H9N3O4	1000387-20-2	10		
5	3-Methyl-2-propylcyclopent-2-en-...	138	C9H14O	1000423-46-9	10		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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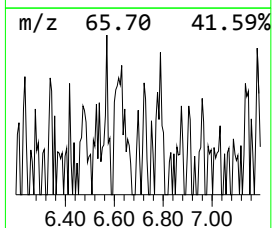
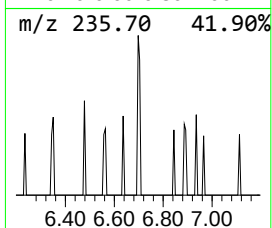
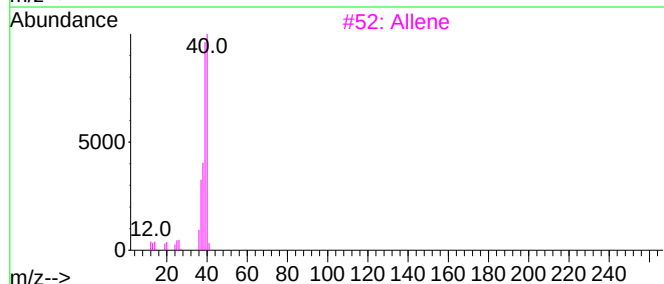
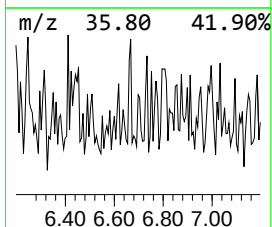
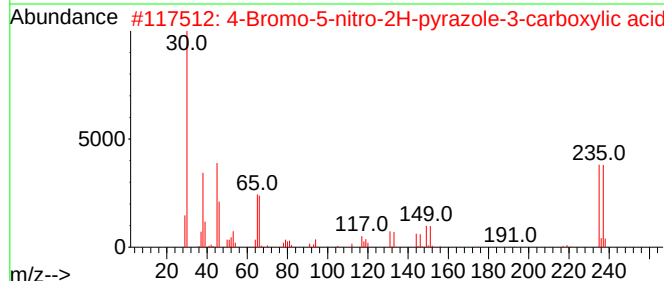
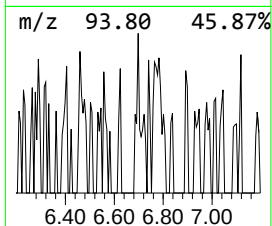
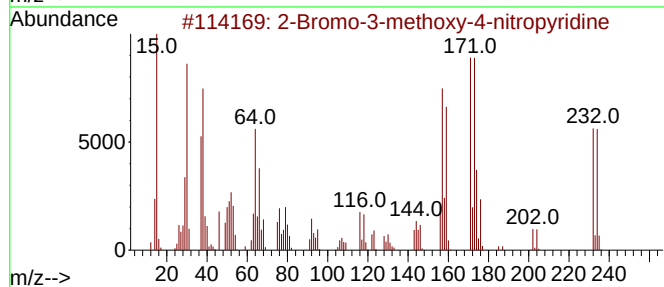
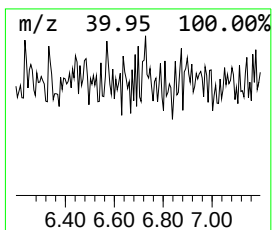
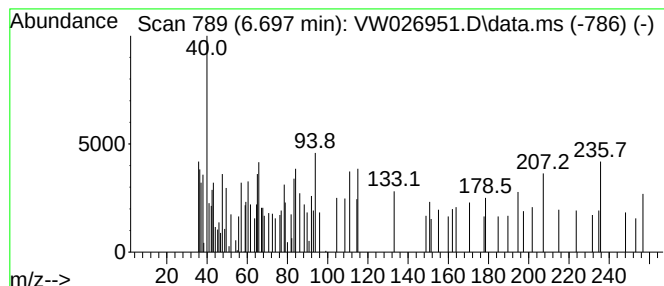
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.697	1.39 ug/L	2991	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo-3-methoxy-4-nitropyridine	232	C6H5BrN2O3	1000444-83-2	27		
2	4-Bromo-5-nitro-2H-pyrazole-3-ca...	235	C4H2BrN3O4	1000459-87-6	10		
3	Allene	40	C3H4	000463-49-0	10		
4	Ethanol, 2,2-dichloro-	114	C2H4Cl2O	000598-38-9	9		
5	1,2-Difluoroethane	66	C2H4F2	000624-72-6	9		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYW5

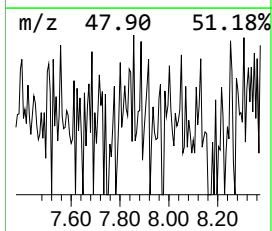
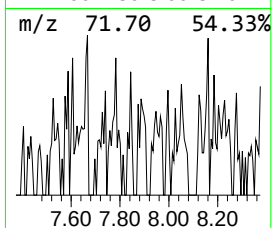
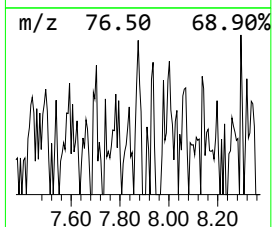
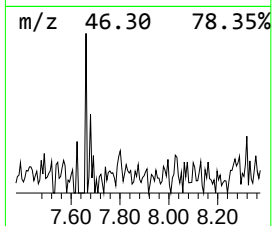
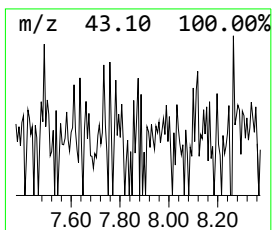
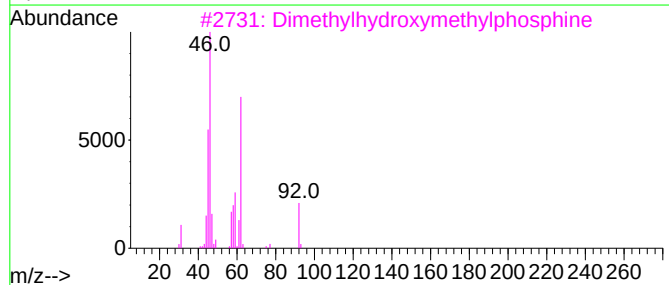
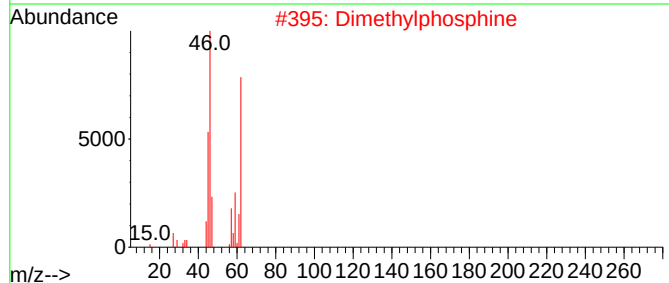
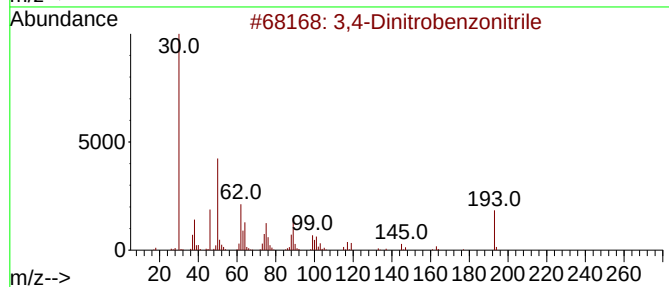
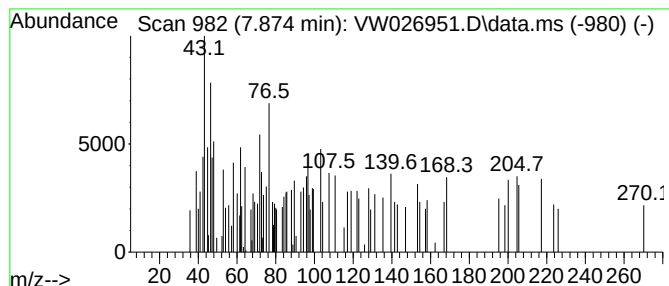
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-08 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.874	2.30 ug/L	4942	1,4-Difluorobenzene	8.843

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,4-Dinitrobenzonitrile	193	C7H3N3O4	004248-33-3	17
2		Dimethylphosphine	62	C2H7P	000676-59-5	14
3		Dimethylhydroxymethylphosphine	92	C3H9OP	033796-25-7	14
4		6-Nitropiperonal	195	C8H5NO5	000712-97-0	12
5		Pyridine-4-aldehyde, N-ethoxycar...	193	C9H11N3O2	083710-34-3	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYW5

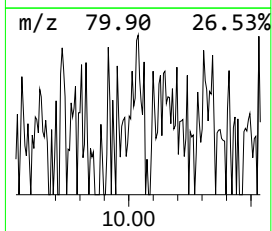
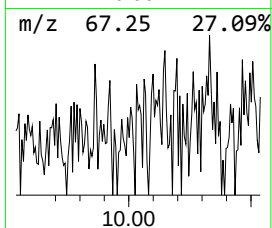
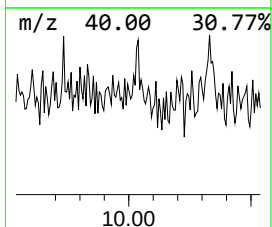
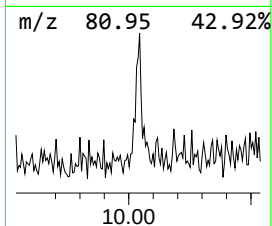
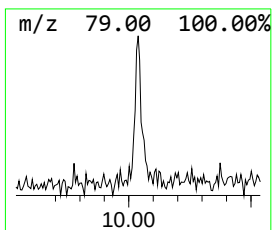
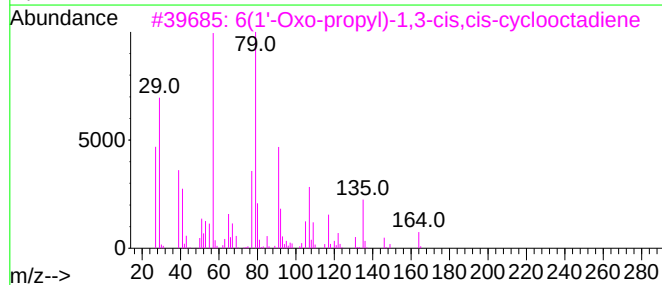
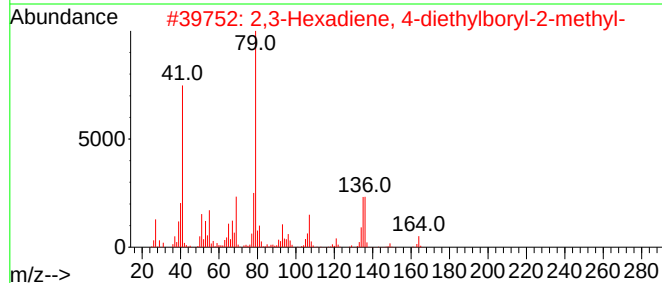
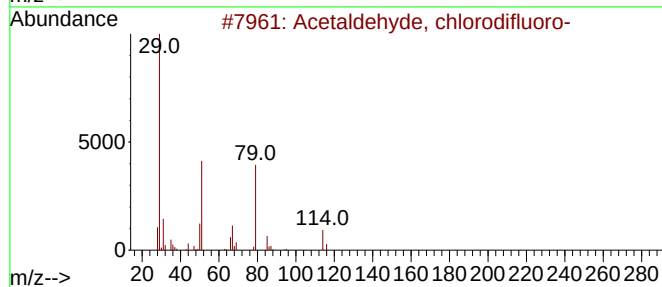
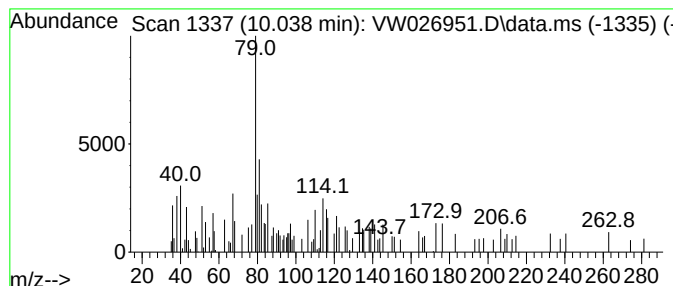
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-09 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.038	1.86 ug/L	3996	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetaldehyde, chlorodifluoro-	114	C2HClF2O	000811-96-1	27
2			2,3-Hexadiene, 4-diethylboryl-2-...	164	C11H21B	1000150-14-8	22
3			6(1'-Oxo-propyl)-1,3-cis,cis-cyc...	164	C11H16O	138146-07-3	22
4			Methyl 4,7,10,13-hexadecatetraen...	262	C17H26O2	1000336-36-7	16
5			Cyclopropane, 1-ethenyl-2-hexeny...	150	C11H18	022822-99-7	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082923\
Data File : VW026951.D
Acq On : 29 Aug 2023 15:23
Operator : SY/MD
Sample : 04175-22
Misc : 5.44g/10mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYW5

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
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unknown-02	2.503	1.9	ug/L	3994	1	8.843	53706	25.0
unknown-03	4.942	1.8	ug/L	3863	1	8.843	53706	25.0
unknown-04	5.173	1.3	ug/L	2851	1	8.843	53706	25.0
unknown-05	5.435	1.5	ug/L	3230	1	8.843	53706	25.0
unknown-06	6.143	1.5	ug/L	3193	1	8.843	53706	25.0
unknown-07	6.697	1.4	ug/L	2991	1	8.843	53706	25.0
unknown-08	7.874	2.3	ug/L	4942	1	8.843	53706	25.0
unknown-09	10.038	1.9	ug/L	3996	1	8.843	53706	25.0