

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
 Data File : VW026111.D
 Acq On : 06 Jun 2023 16:49
 Operator : SY/MD
 Sample : 03062-07
 Misc : 6.19g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EZEL3

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\SFAMWLM053023SMA.M
 Title : SFAM01.0

Signal : TIC: VW026111.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.844	155	157	158	rBV2	1998	1269	10.12%	0.812%
2	3.606	278	282	285	rBV6	1766	2604	20.76%	1.666%
3	3.692	293	296	297	rBV3	1476	1769	14.10%	1.132%
4	3.948	336	338	339	rBV2	1372	613	4.89%	0.392%
5	4.362	404	406	408	rBV3	1512	820	6.54%	0.525%
6	4.570	438	440	441	rVB2	2115	1283	10.23%	0.821%
7	4.588	441	443	444	rBV2	1292	904	7.21%	0.578%
8	5.057	518	520	522	rBV3	1373	1352	10.78%	0.865%
9	5.899	656	658	661	rVB4	2003	1699	13.54%	1.087%
10	6.222	709	711	712	rBV2	1594	1461	11.65%	0.935%
11	6.685	784	787	791	rBV6	2059	3108	24.77%	1.988%
12	6.746	795	797	798	rBV2	1576	809	6.45%	0.518%
13	6.966	831	833	835	rBV3	1551	1702	13.57%	1.089%
14	7.142	860	862	865	rVB4	1491	1309	10.43%	0.837%
15	7.533	924	926	929	rBV4	1464	1411	11.25%	0.903%
16	7.630	940	942	943	rBV2	1714	1350	10.76%	0.864%
17	7.813	970	972	977	rVB6	1607	1722	13.73%	1.102%
18	7.874	980	982	985	rBV4	1813	2189	17.45%	1.400%
19	7.923	988	990	992	rBV3	1607	1299	10.35%	0.831%
20	8.002	1001	1003	1006	rBV4	1384	1445	11.52%	0.924%
21	8.228	1038	1040	1043	rBV4	1920	2229	17.77%	1.426%
22	8.264	1043	1046	1047	rBV2	5226	4940	39.38%	3.160%
23	8.673	1110	1113	1116	rBV5	2699	4830	38.50%	3.090%
24	8.843	1137	1141	1148	rVB4	5900	12545	100.00%	8.025%
25	9.038	1172	1173	1174	rBV	521	194	1.55%	0.124%
26	9.075	1174	1179	1180	rBV5	1040	1511	12.04%	0.967%
27	9.124	1186	1187	1190	rBV2	1333	1165	9.29%	0.745%
28	9.197	1193	1199	1201	rBV7	1455	2530	20.17%	1.619%
29	9.276	1206	1212	1217	rVV5	4274	8395	66.92%	5.371%
30	9.362	1225	1226	1228	rBV2	1157	794	6.33%	0.508%
31	9.404	1230	1233	1235	rVB4	2718	2968	23.66%	1.899%
32	9.435	1235	1238	1239	rBV3	463	320	2.55%	0.205%
33	9.453	1239	1241	1244	rBV4	480	478	3.81%	0.306%
34	9.502	1247	1249	1252	rVB4	630	755	6.02%	0.483%
35	9.538	1252	1255	1258	rBV2	512	842	6.71%	0.539%
36	9.581	1258	1262	1265	rBV4	829	1355	10.80%	0.867%

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Integrator: RTE

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 Title : SFAM01.0

37	9.605	1265	1266	1270	rVB3	694	536	4.27%	0.343%
38	9.666	1273	1276	1283	rBV6	1000	2499	19.92%	1.599%
39	9.715	1283	1284	1285	rBV	283	148	1.18%	0.095%
40	9.733	1285	1287	1288	rVB3	570	289	2.30%	0.185%
41	9.746	1288	1289	1291	rVB2	666	407	3.24%	0.260%
42	9.776	1291	1294	1296	rBV2	390	346	2.76%	0.221%
43	9.819	1296	1301	1302	rBV4	891	815	6.50%	0.521%
44	9.837	1302	1304	1308	rVB4	1148	1403	11.18%	0.898%
45	9.868	1308	1309	1310	rBV	621	340	2.71%	0.218%
46	9.892	1310	1313	1315	rVB4	518	494	3.94%	0.316%
47	9.910	1315	1316	1317	rBV	503	186	1.48%	0.119%
48	9.935	1317	1320	1323	rBV5	837	1283	10.23%	0.821%
49	9.977	1323	1327	1329	rVB3	866	836	6.66%	0.535%
50	10.032	1329	1336	1337	rBV7	1351	2068	16.48%	1.323%
51	10.105	1344	1348	1350	rVB4	1041	1049	8.36%	0.671%
52	10.130	1350	1352	1354	rVB3	1164	867	6.91%	0.555%
53	10.154	1354	1356	1358	rBV2	1295	1355	10.80%	0.867%
54	10.178	1358	1360	1361	rVB2	576	384	3.06%	0.246%
55	10.197	1361	1363	1366	rBV4	606	779	6.21%	0.498%
56	10.227	1366	1368	1369	rBV2	815	585	4.66%	0.374%
57	10.325	1376	1384	1390	rBV4	5543	12357	98.50%	7.905%
58	10.404	1395	1397	1399	rBV2	973	788	6.28%	0.504%
59	10.434	1399	1402	1403	rBV3	939	789	6.29%	0.505%
60	10.453	1403	1405	1406	rVB	577	341	2.72%	0.218%
61	10.471	1406	1408	1410	rBV3	633	435	3.47%	0.278%
62	10.495	1410	1412	1416	rVB5	996	1152	9.18%	0.737%
63	10.538	1416	1419	1420	rBV3	772	826	6.58%	0.528%
64	10.605	1427	1430	1431	rBV3	1858	2188	17.44%	1.400%
65	10.971	1488	1490	1491	rBV2	1063	629	5.01%	0.402%
66	11.233	1531	1533	1536	rVB4	1966	1728	13.77%	1.105%
67	11.270	1536	1539	1540	rBV3	1305	1427	11.38%	0.913%
68	11.373	1554	1556	1557	rBV2	1715	945	7.53%	0.605%
69	11.514	1577	1579	1580	rBV2	1200	695	5.54%	0.445%
70	11.629	1594	1598	1603	rBV5	6760	12244	97.60%	7.833%
71	11.977	1654	1655	1657	rVB2	1448	920	7.33%	0.589%
72	12.166	1684	1686	1689	rBV4	2026	1828	14.57%	1.169%
73	12.391	1721	1723	1724	rBV2	1210	891	7.10%	0.570%
74	12.928	1807	1811	1814	rBV6	1523	2644	21.08%	1.691%
75	13.562	1910	1915	1921	rBV9	3786	7987	63.67%	5.110%
76	13.958	1978	1980	1982	rBV3	1427	1120	8.93%	0.717%

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Peak separation: 5

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Title : SFAM01.0

77	14.050	1993	1995	1998	rBV4	1577	1821	14.52%	1.165%
78	14.287	2033	2034	2037	rBV3	1586	1549	12.35%	0.991%
79	14.684	2096	2099	2101	rBV4	1705	2660	21.20%	1.702%
80	14.751	2108	2110	2112	rBV3	1736	1165	9.29%	0.745%
81	14.928	2137	2139	2141	rBV3	1598	1409	11.23%	0.901%
82	16.092	2328	2330	2333	rBV4	1645	1247	9.94%	0.798%
83	16.757	2437	2439	2441	rBV3	2025	1892	15.08%	1.210%

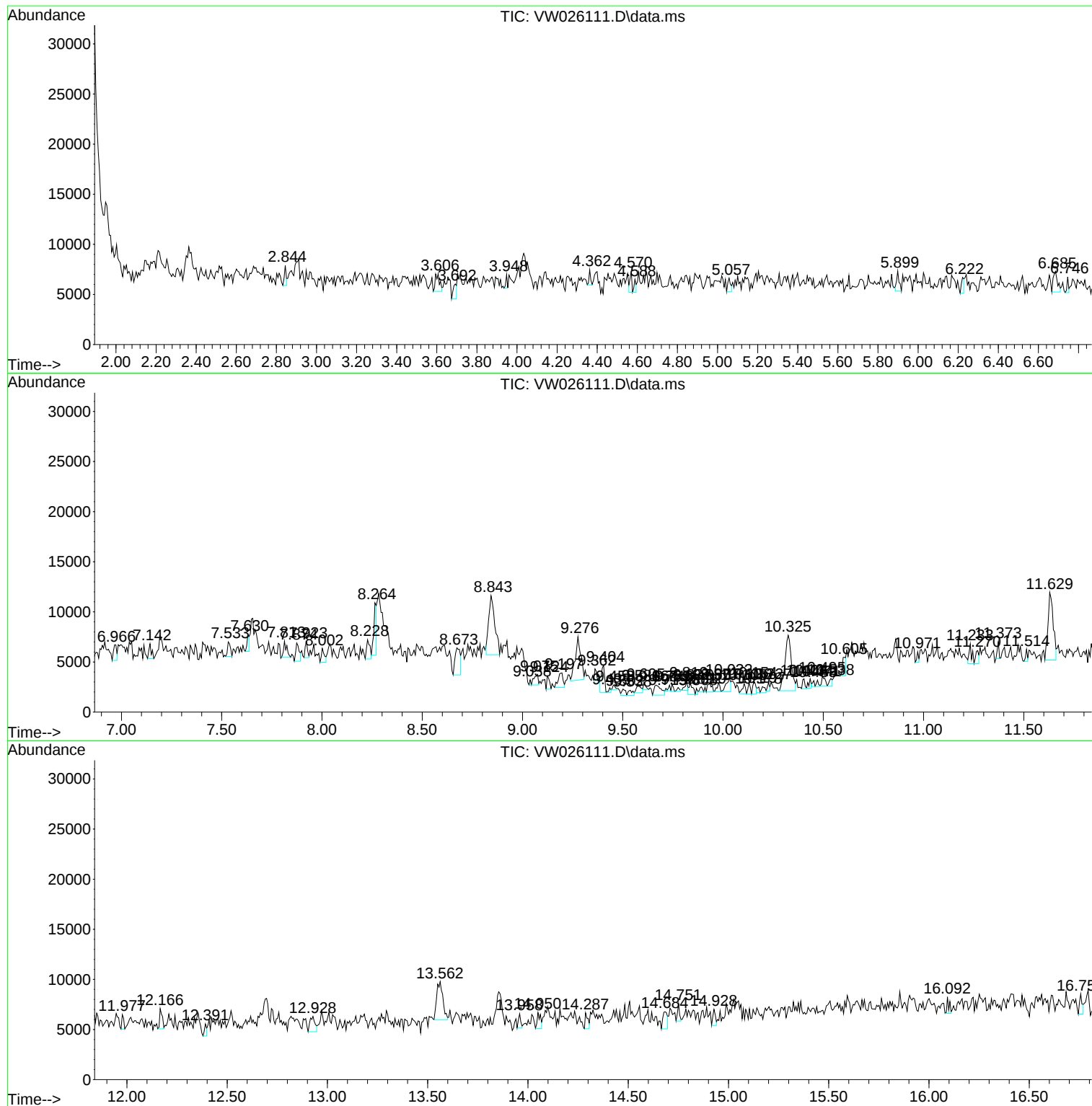
Sum of corrected areas: 156315

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MSVOA_W
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EZEL3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM053023SMA.M
Quant Title : SFAM01.0

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



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Instrument :
MSVOA_W
ClientSampleId :
EZEL3

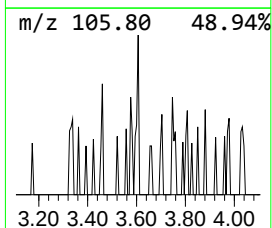
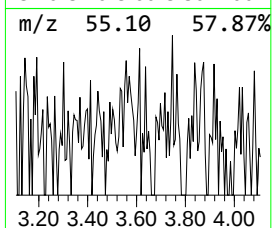
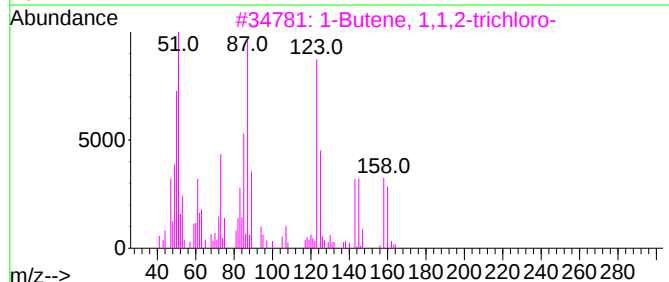
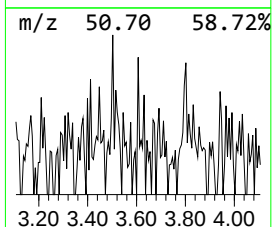
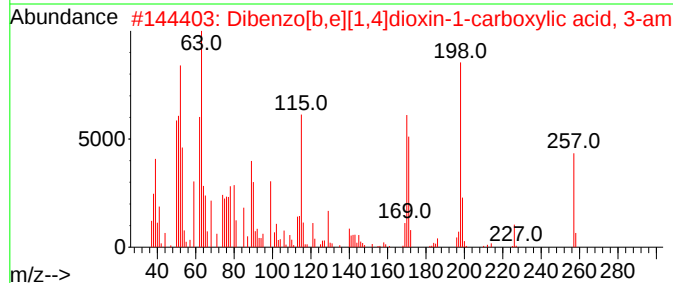
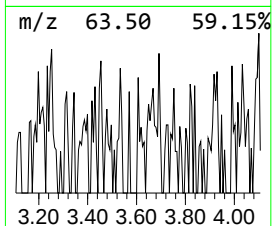
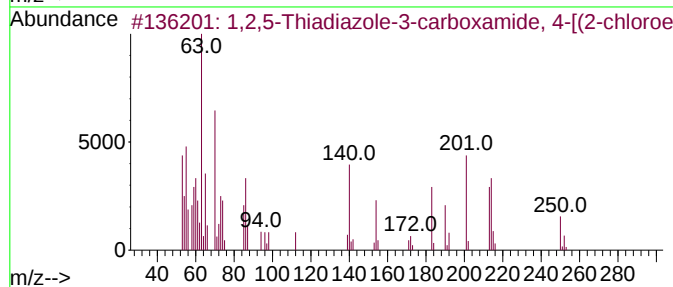
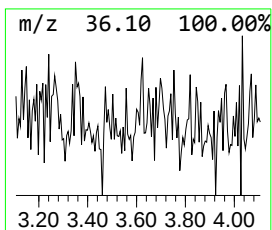
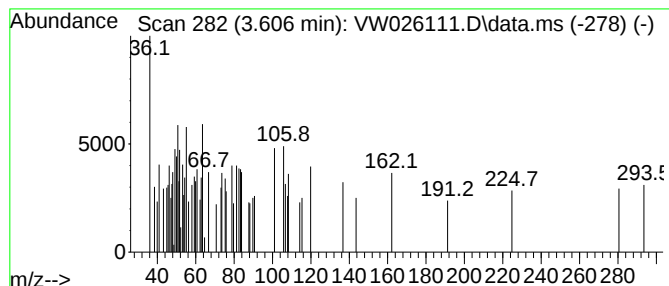
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Quant Title : SFAM01.0

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

Peak Number 1 1,2,5-Thiadiazole-3-carboxa... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.606	5.19 ug/L	2604	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,5-Thiadiazole-3-carboxamide,...	250	C7H11ClN4O2S	147194-45-4	64		
2	Dibenzo[b,e][1,4]dioxin-1-carbox...	257	C14H11NO4	1000362-82-4	25		
3	1-Butene, 1,1,2-trichloro-	158	C4H5Cl3	042860-89-9	17		
4	Carmustine	213	C5H9Cl2N3O2	000154-93-8	16		
5	2-Chloro-6-nitrophenol	173	C6H4ClNO3	000603-86-1	14		



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Instrument :
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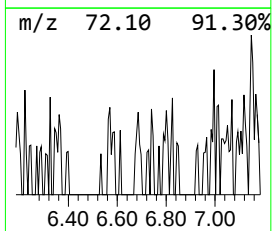
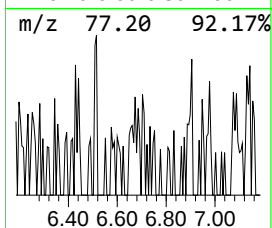
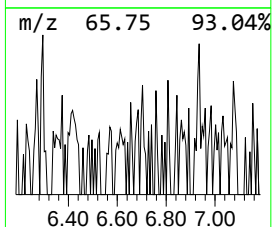
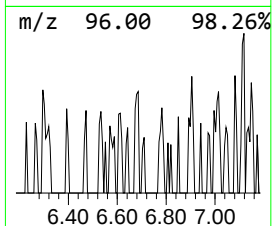
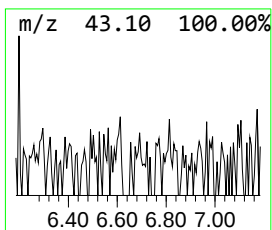
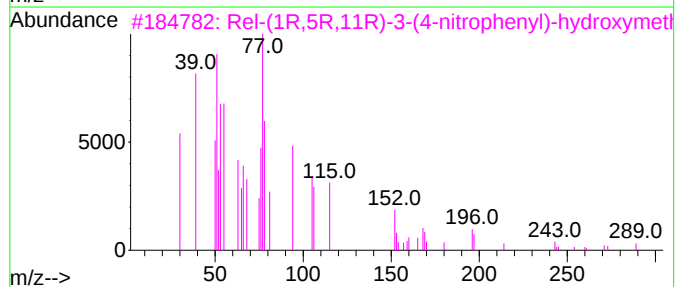
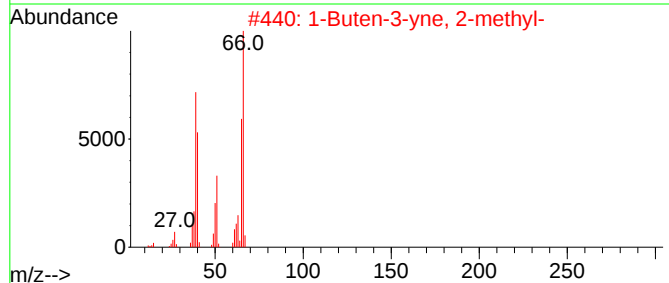
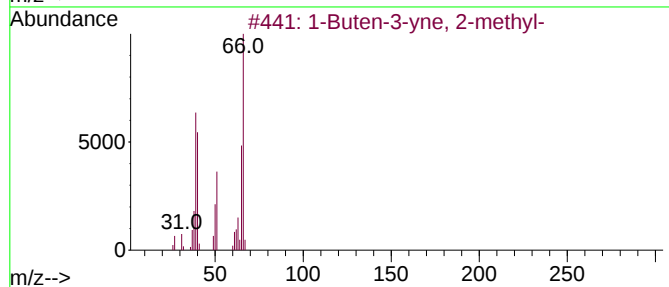
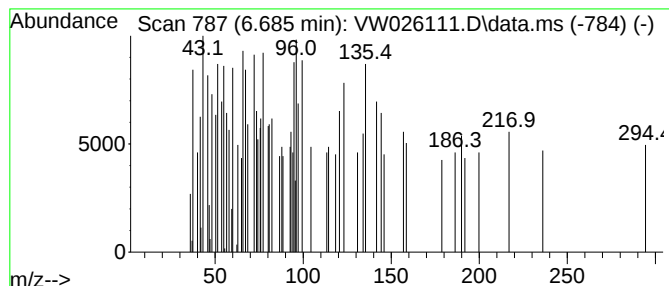
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM053023SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.685	6.19 ug/L	3108	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Buten-3-yne, 2-methyl-	66	C5H6	000078-80-8	35
2			1-Buten-3-yne, 2-methyl-	66	C5H6	000078-80-8	25
3			Rel-(1R,5R,11R)-3-(4-nitrophenyl...	289	C14H11NO6	1000433-34-4	14
4			2-Propen-1-one, 3-(4-fluoropheny...	226	C15H11FO	001608-51-1	12
5			Bicyclo[2.2.2]oct-5-en-2-one, 7-...	138	C8H10O2	1000153-14-8	10



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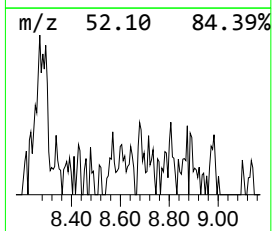
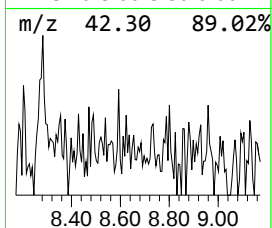
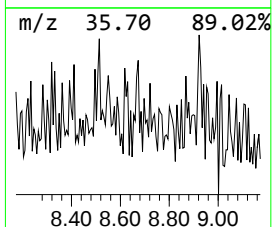
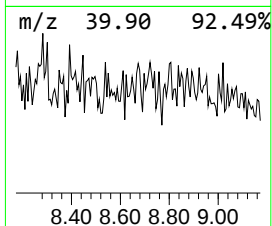
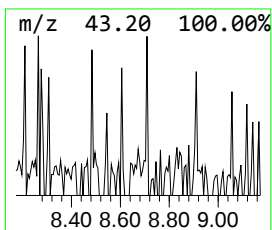
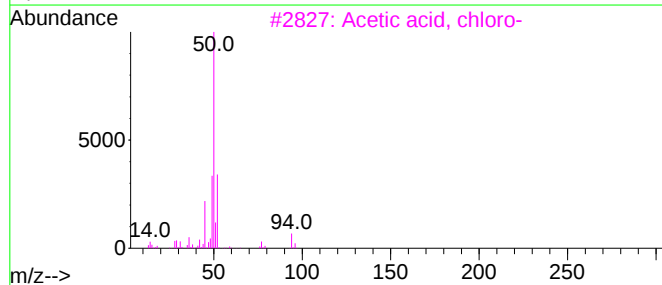
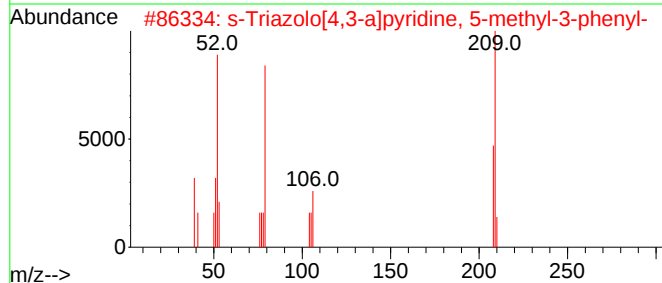
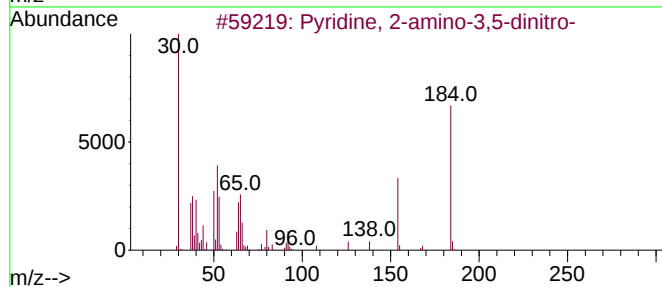
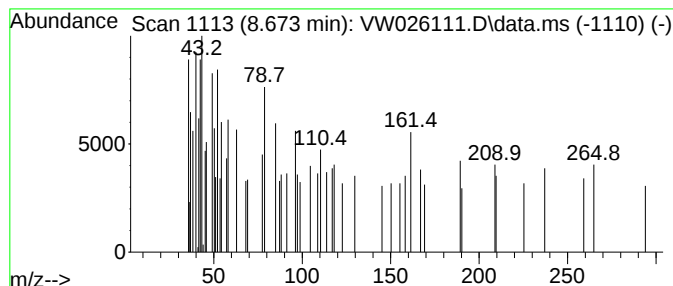
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM053023SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.673	9.63 ug/L	4830	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2-amino-3,5-dinitro-	184	C5H4N4O4	003073-30-1	25		
2	s-Triazolo[4,3-a]pyridine, 5-met...	209	C13H11N3	004926-14-1	22		
3	Acetic acid, chloro-	94	C2H3ClO2	000079-11-8	12		
4	Cyclopropanecarboxamide, 1,2,2-t...	160	C7H4N4O	1000351-24-4	10		
5	1,2-Propadiene, 1-bromo-	118	C3H3Br	010024-18-7	10		



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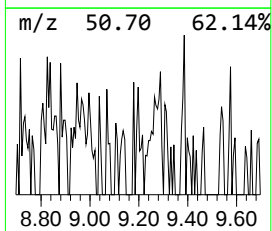
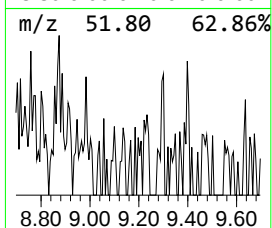
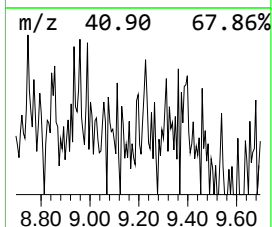
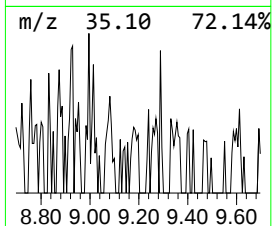
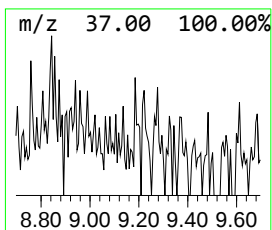
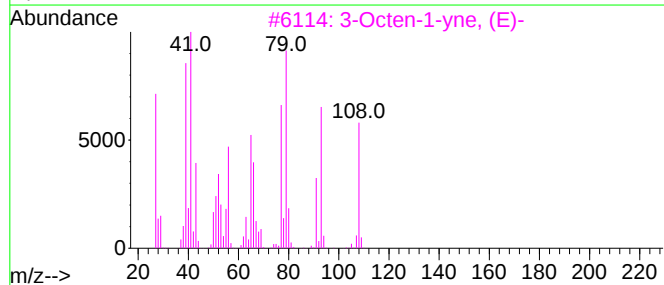
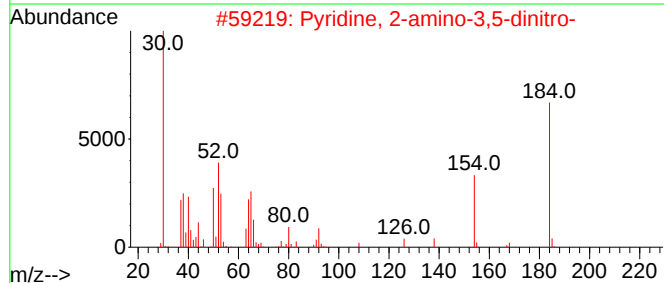
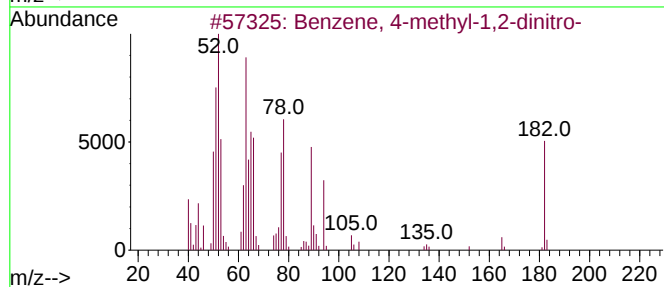
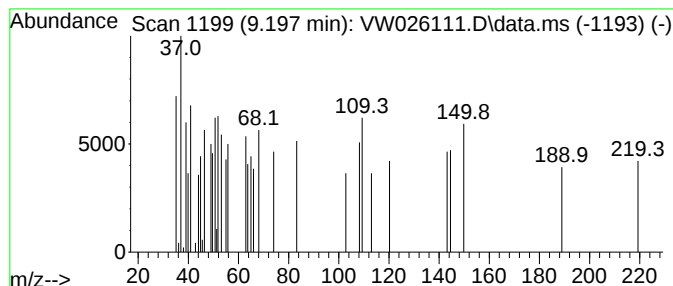
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Quant Title : SFAM01.0

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

Peak Number 4 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.197	5.04 ug/L	2530	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	38
2			Pyridine, 2-amino-3,5-dinitro-	184	C5H4N4O4	003073-30-1	25
3			3-Octen-1-yne, (E)-	108	C8H12	042104-42-7	11
4			Carbonochloridic acid, chloromet...	128	C2H2Cl2O2	022128-62-7	10
5			2,4-Dinitrobenzoic acid	212	C7H4N2O6	000610-30-0	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

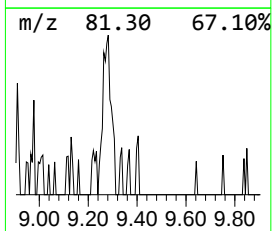
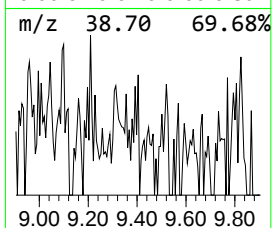
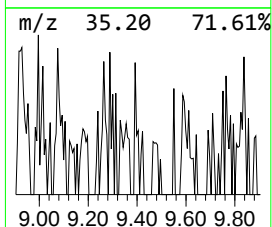
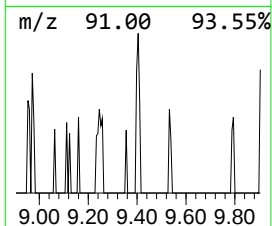
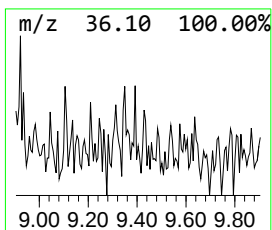
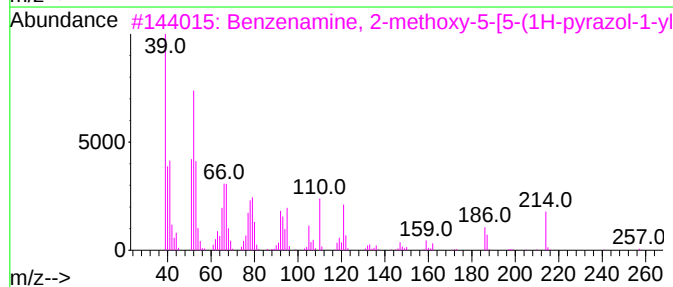
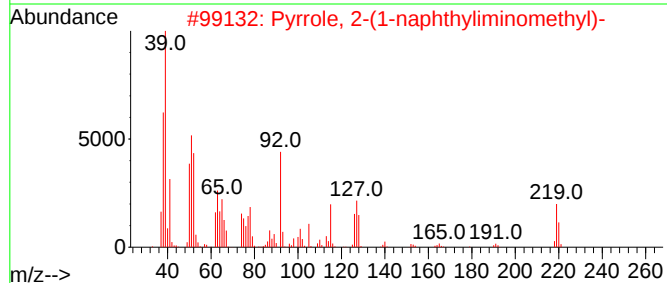
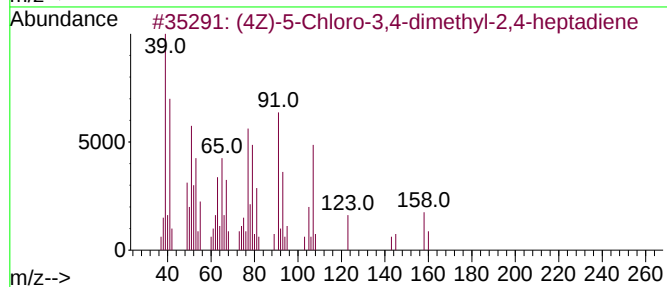
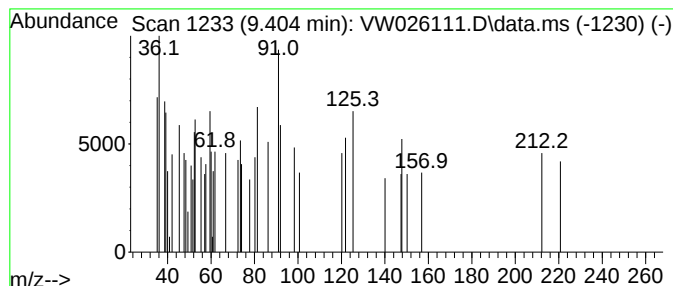
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-04 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.404	5.91 ug/L	2968	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(4Z)-5-Chloro-3,4-dimethyl-2,4-h...	158	C9H15Cl	105949-75-5	39		
2	Pyrrole, 2-(1-naphthyliminomethyl)-	220	C15H12N2	212759-16-5	38		
3	Benzenamine, 2-methoxy-5-[5-(1H-...	257	C11H11N7O	1000362-82-7	38		
4	3-Methyl-2-propylcyclopent-2-en-...	138	C9H14O	1000423-46-9	37		
5	(2E,4E)-5-Chloro-3,4-dimethyl-2,...	158	C9H15Cl	1000112-49-7	33		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

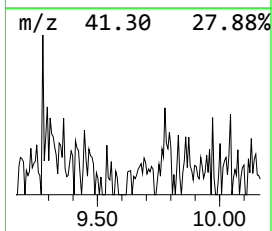
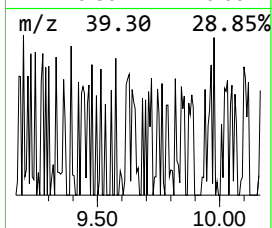
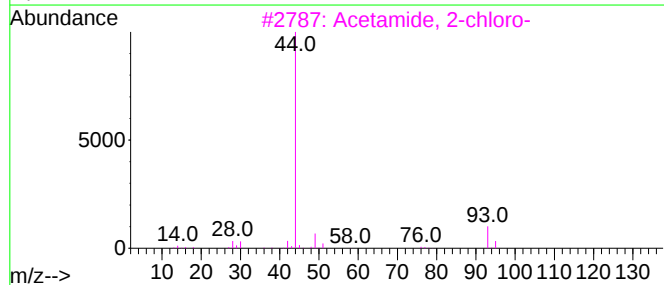
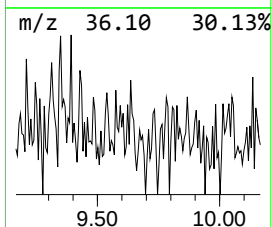
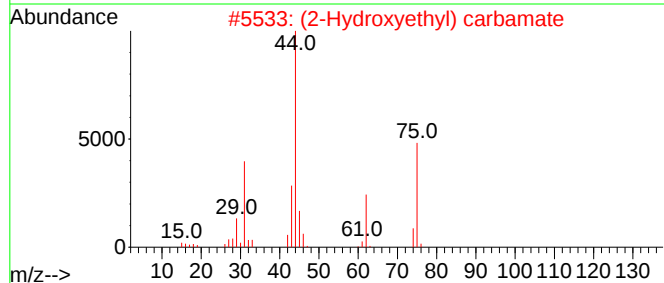
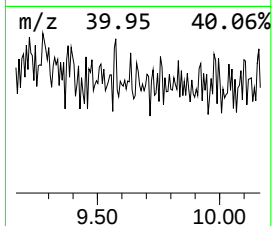
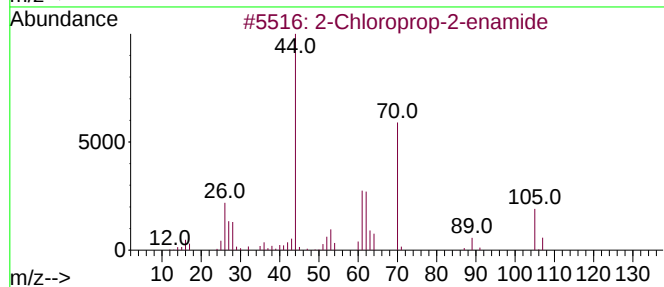
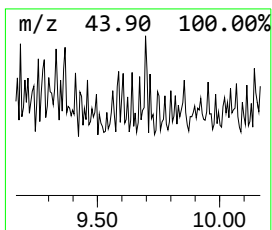
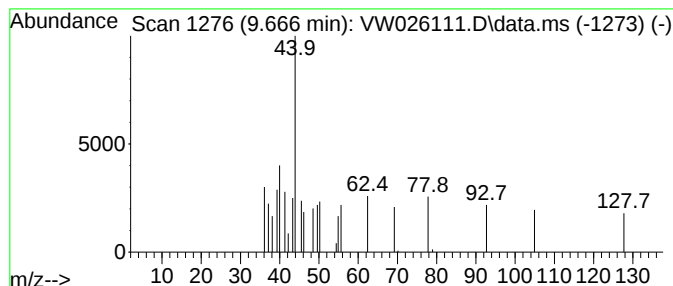
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Quant Title : SFAM01.0

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

Peak Number 6 unknown-05 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.666	4.98 ug/L	2499	1,4-Difluorobenzene	8.843

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Chloroprop-2-enamide	105	C3H4ClNO	1000411-34-7	9
2		(2-Hydroxyethyl) carbamate	105	C3H7NO3	005395-01-7	9
3		Acetamide, 2-chloro-	93	C2H4ClNO	000079-07-2	4
4		n-Propyl fluoride	62	C3H7F	000460-13-9	4
5		Cyclopropanecarboxamide	85	C4H7NO	006228-73-5	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

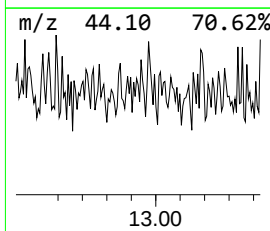
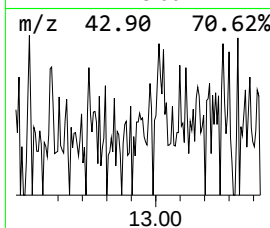
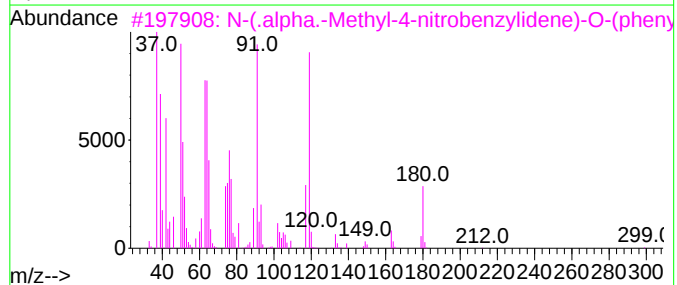
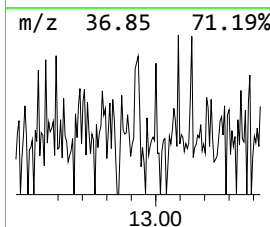
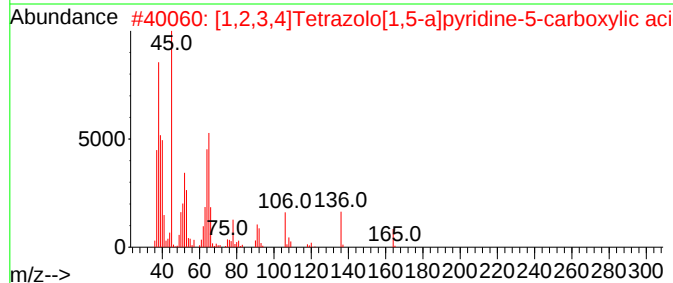
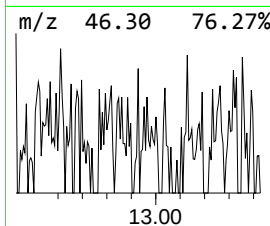
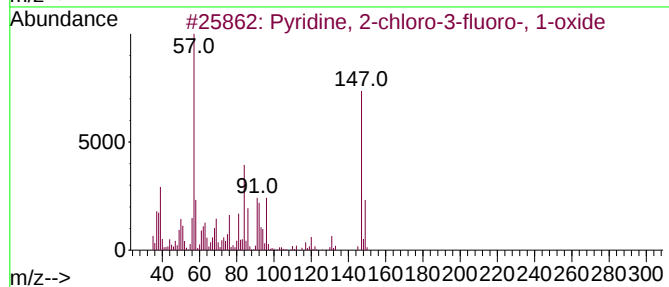
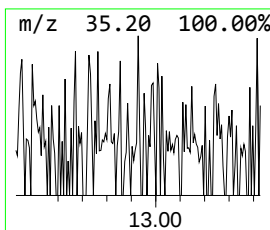
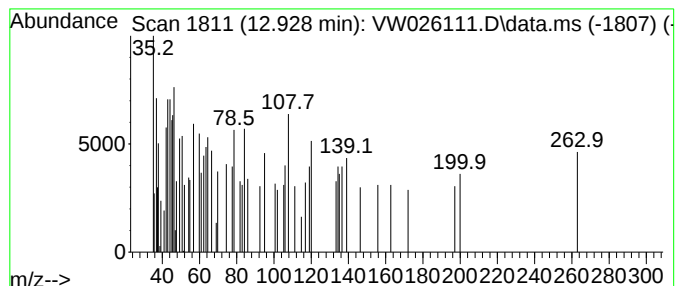
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.928	8.28 ug/L	2644	1,4-Dichlorobenzene-d4	13.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine, 2-chloro-3-fluoro-, 1-...	147	C5H3ClFNO	1000327-47-4	35
2			[1,2,3,4]Tetrazolo[1,5-a]pyridin...	164	C6H4N4O2	1010386-90-2	17
3			N-(.alpha.-Methyl-4-nitrobenzyl)...	299	C15H13N3O4	1000223-36-2	16
4			[.alpha.-Acetylmethylene-4-metho...	240	C11H11BF2O3	1000223-62-2	12
5			Ethanedial,1,2-dichloro-, bis(0-...	240	C6H6Cl2N2O4	089715-76-4	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

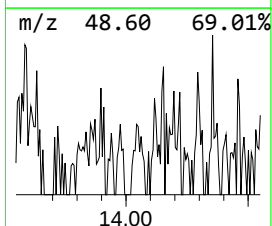
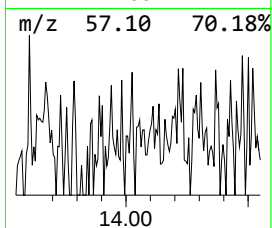
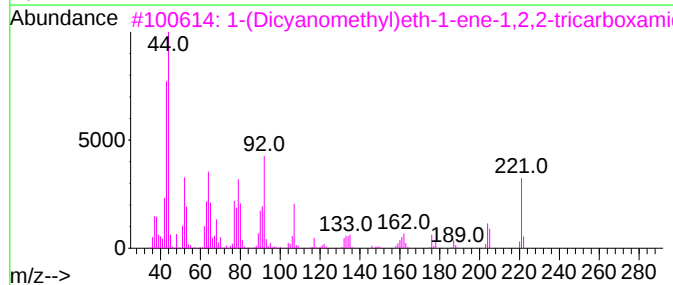
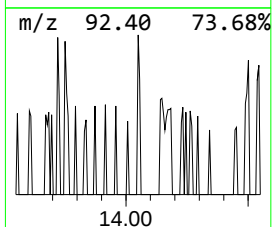
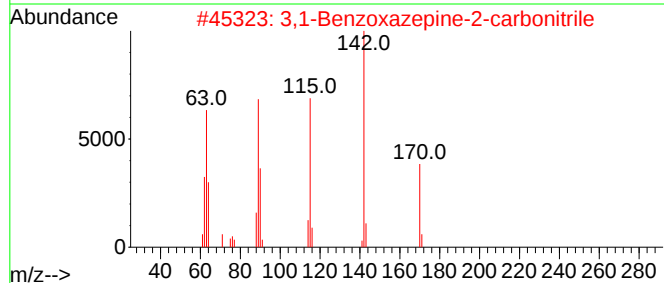
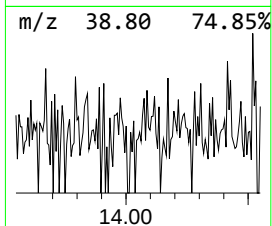
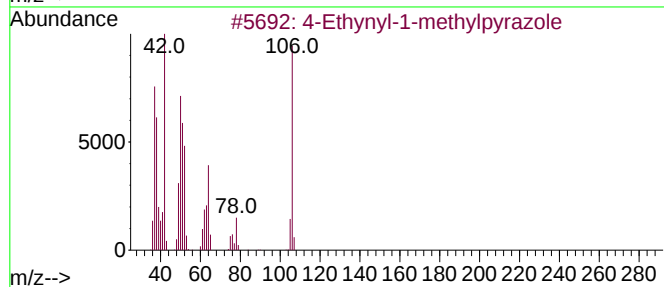
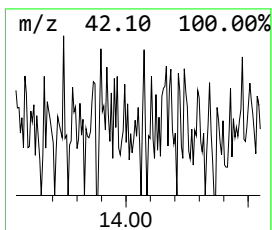
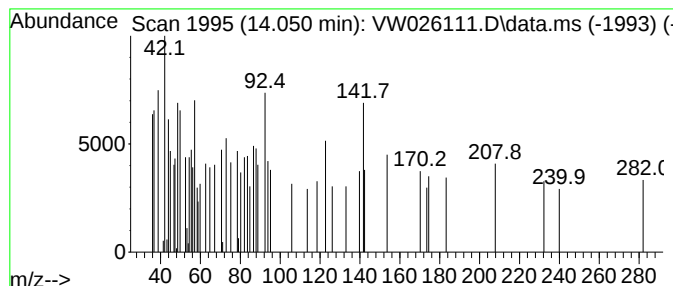
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-07 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.050	5.70 ug/L	1821	1,4-Dichlorobenzene-d4	13.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Ethynyl-1-methylpyrazole	106	C6H6N2	039806-89-8	35
2			3,1-Benzoxazepine-2-carbonitrile	170	C10H6N2O	018457-80-2	17
3			1-(Dicyanomethyl)eth-1-ene-1,2,2...	221	C8H7N5O3	1000388-01-7	17
4			3,1-Benzoxazepine-2-carbonitrile	170	C10H6N2O	018457-80-2	17
5			N-(.alpha.-Methyl-4-nitrobenzyl)li...	299	C15H13N3O4	1000223-36-2	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

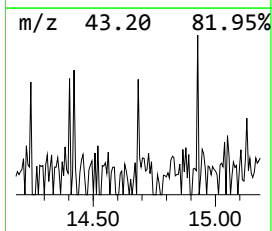
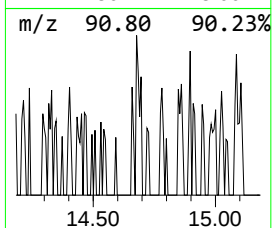
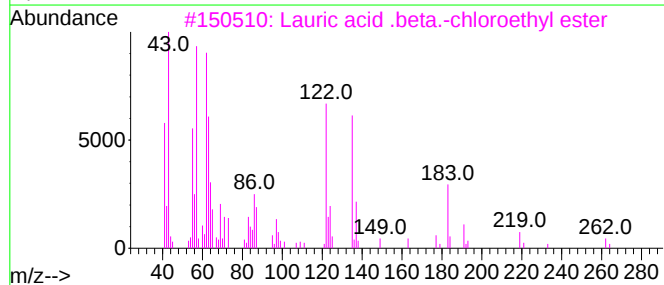
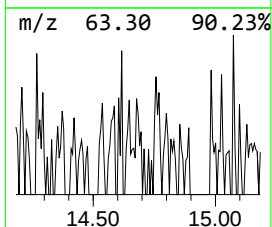
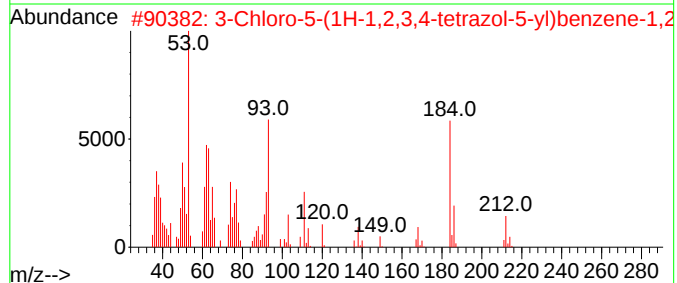
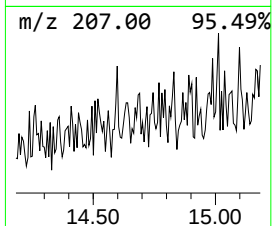
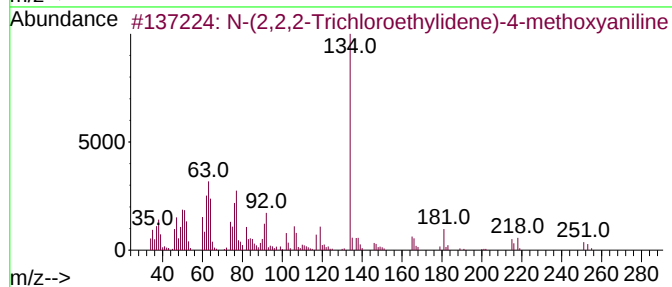
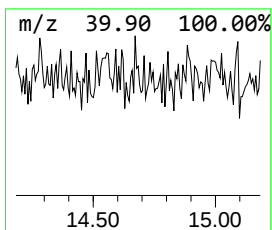
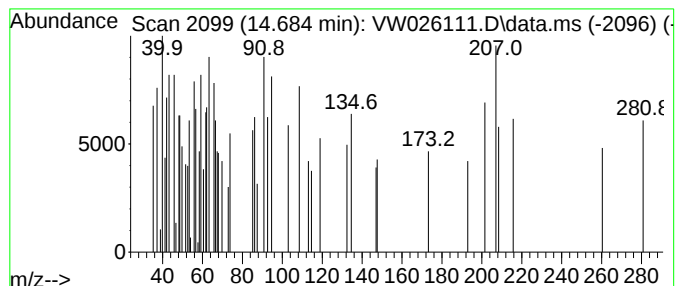
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Quant Title : SFAM01.0

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.
14.684	8.33 ug/L	2660	1,4-Dichlorobenzene-d4	13.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-(2,2,2-Trichloroethylidene)-4-...	251	C9H8Cl3NO	083320-59-6	32
2			3-Chloro-5-(1H-1,2,3,4-tetrazol-...	212	C7H5ClN4O2	1000409-86-2	25
3			Lauric acid .beta.-chloroethyl e...	262	C14H27ClO2	064919-15-9	23
4			2-Chloroethyl caprate	234	C12H23ClO2	015175-04-9	23
5			Phenol, 4-bromo-2,5-dichloro-	240	C6H3BrCl2O	001940-42-7	23



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

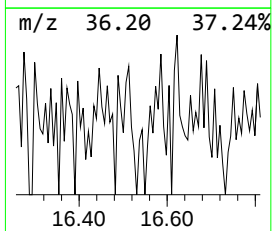
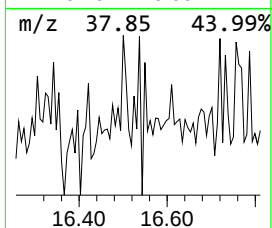
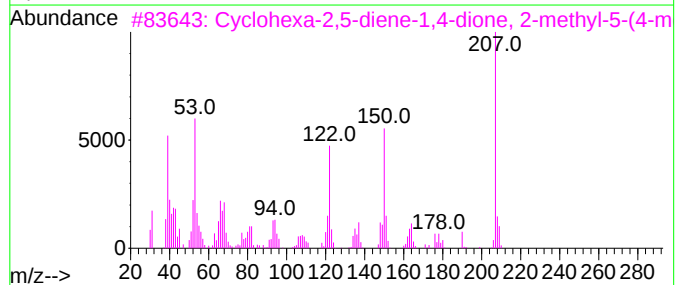
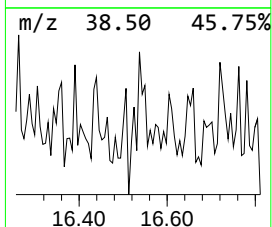
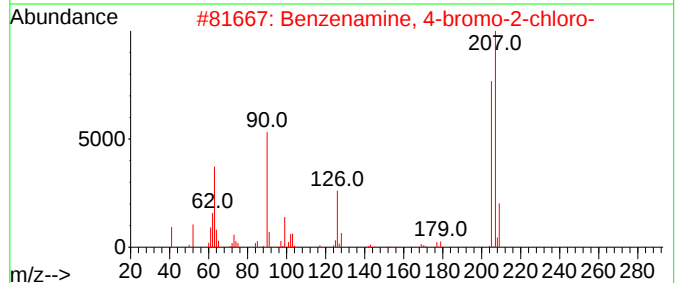
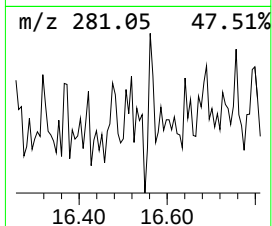
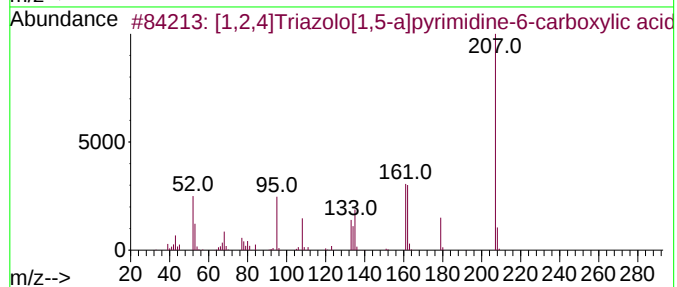
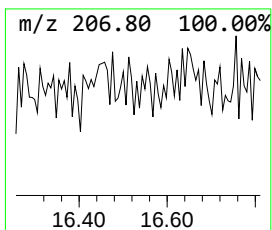
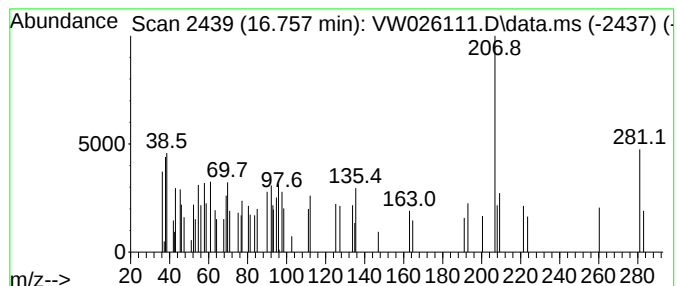
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM053023SMA.M
Quant Title : SFAM01.0

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

Peak Number 10 unknown-09 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.756	5.92 ug/L	1892	1,4-Dichlorobenzene-d4	13.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			[1,2,4]Triazolo[1,5-a]pyrimidine...	207	C8H9N5O2	092673-40-0	38
2			Benzenamine, 4-bromo-2-chloro-	205	C6H5BrClN	038762-41-3	25
3			Cyclohexa-2,5-diene-1,4-dione, 2...	207	C11H13NO3	002158-89-6	10
4			5-Nitro-2-(trifluoromethyl)pyrid...	207	C6H4F3N3O2	438564-36-4	10
5			8-Chloro-5-quinolinecarboxylic acid	207	C10H6ClNO2	121490-68-4	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW060623\
Data File : VW026111.D
Acq On : 06 Jun 2023 16:49
Operator : SY/MD
Sample : 03062-07
Misc : 6.19g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZEL3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM053023SMA.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
1,2,5-Thiadiaz...	3.606	5.2	ug/L	2604	1	8.843	12545	25.0
unknown-01	6.685	6.2	ug/L	3108	1	8.843	12545	25.0
unknown-02	8.673	9.6	ug/L	4830	1	8.843	12545	25.0
unknown-03	9.197	5.0	ug/L	2530	1	8.843	12545	25.0
unknown-04	9.404	5.9	ug/L	2968	1	8.843	12545	25.0
unknown-05	9.666	5.0	ug/L	2499	1	8.843	12545	25.0
unknown-06	12.928	8.3	ug/L	2644	3	13.568	7987	25.0
unknown-07	14.050	5.7	ug/L	1821	3	13.568	7987	25.0
unknown-08	14.684	8.3	ug/L	2660	3	13.568	7987	25.0
unknown-09	16.756	5.9	ug/L	1892	3	13.568	7987	25.0