

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
 Data File : VD077806.D
 Acq On : 11 Dec 2023 11:23
 Operator : JC/SY
 Sample : 05750-02
 Misc : 4.24G/10ml/MSVOA_D/SOIL
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 E0LQ5

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_D\Method\SFAMDLM112923SMA.M
 Title : SFAM01.0

Signal : TIC: VD077806.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.687	11	17	20	rBV	11702	20120	1.29%	0.142%
2	1.746	20	27	38	rVB	396767	661001	42.44%	4.666%
3	2.005	69	71	76	rBV2	1664	2931	0.19%	0.021%
4	2.075	79	83	85	rVB2	2745	3429	0.22%	0.024%
5	2.128	90	92	93	rBV2	1289	773	0.05%	0.005%
6	2.275	112	117	124	rBV	257219	394986	25.36%	2.788%
7	2.805	201	207	219	rVB	209916	392825	25.22%	2.773%
8	3.187	270	272	273	rBV	1715	1021	0.07%	0.007%
9	3.264	283	285	287	rBV	1687	1169	0.08%	0.008%
10	3.311	291	293	295	rBV2	1384	1267	0.08%	0.009%
11	3.328	295	296	299	rVB	2455	1897	0.12%	0.013%
12	3.352	299	300	301	rBV	1769	1018	0.07%	0.007%
13	3.417	308	311	314	rBV2	1308	1499	0.10%	0.011%
14	3.570	336	337	341	rBV	1583	2339	0.15%	0.017%
15	3.646	349	350	352	rBV2	954	948	0.06%	0.007%
16	3.805	375	377	378	rBV	1077	786	0.05%	0.006%
17	3.922	387	397	409	rBV	216861	559945	35.96%	3.953%
18	4.234	448	450	453	rVB	1721	1285	0.08%	0.009%
19	4.269	453	456	457	rBV3	1433	1540	0.10%	0.011%
20	4.352	466	470	471	rBV2	949	1016	0.07%	0.007%
21	4.517	496	498	499	rVV2	1197	925	0.06%	0.007%
22	4.528	499	500	502	rVB	2078	1129	0.07%	0.008%
23	4.552	502	504	506	rBV	1379	1287	0.08%	0.009%
24	4.605	511	513	515	rVV2	1257	755	0.05%	0.005%
25	4.769	539	541	542	rBV	2098	1299	0.08%	0.009%
26	4.905	563	564	567	rVB3	1511	1490	0.10%	0.011%
27	4.934	567	569	571	rVV2	1209	867	0.06%	0.006%
28	4.958	571	573	574	rVB2	1585	964	0.06%	0.007%
29	5.040	586	587	590	rBV2	1528	1075	0.07%	0.008%
30	5.099	595	597	598	rVB	1835	1011	0.06%	0.007%
31	5.322	633	635	637	rVB2	1567	1069	0.07%	0.008%
32	5.552	669	674	677	rVV2	1522	2154	0.14%	0.015%
33	5.646	688	690	691	rBV2	1987	1098	0.07%	0.008%
34	5.734	702	705	706	rBV	1463	1446	0.09%	0.010%
35	5.758	706	709	712	rVB	1602	1788	0.11%	0.013%
36	5.781	712	713	715	rVB2	1667	1144	0.07%	0.008%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
 Data File : VD077806.D
 Acq On : 11 Dec 2023 11:23
 Operator : JC/SY
 Sample : 05750-02
 Misc : 4.24G/10ml/MSVOA_D/SOIL
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 E0LQ5

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_D\Method\SFAMDLM112923SMA.M
 Title : SFAM01.0

37	5.816	715	719	720	rBV2	2813	2905	0.19%	0.021%
38	5.887	729	731	734	rVB2	1826	1300	0.08%	0.009%
39	5.946	740	741	743	rBV	2275	1238	0.08%	0.009%
40	5.975	745	746	749	rVB	1907	1118	0.07%	0.008%

41	6.005	749	751	754	rBV2	1256	1616	0.10%	0.011%
42	6.111	768	769	771	rVB	1854	751	0.05%	0.005%
43	6.269	794	796	799	rVB2	1041	933	0.06%	0.007%
44	6.293	799	800	803	rBV	1429	1615	0.10%	0.011%
45	6.387	812	816	818	rBV2	1228	1786	0.11%	0.013%

46	6.452	825	827	828	rBV	1589	1031	0.07%	0.007%
47	6.505	834	836	839	rBV2	1535	1853	0.12%	0.013%
48	6.599	849	852	854	rBV	1444	1631	0.10%	0.012%
49	6.987	910	918	926	rBV2	30772	87430	5.61%	0.617%
50	7.205	953	955	957	rBV	1696	1902	0.12%	0.013%

51	7.246	961	962	966	rVB	2171	1355	0.09%	0.010%
52	7.310	971	973	974	rBV2	1356	1114	0.07%	0.008%
53	7.440	994	995	998	rBV2	1660	1383	0.09%	0.010%
54	7.569	1007	1017	1029	rBV	262938	664824	42.69%	4.693%
55	7.781	1051	1053	1056	rVB	1837	1420	0.09%	0.010%

56	7.810	1056	1058	1060	rVB2	1340	1108	0.07%	0.008%
57	7.834	1060	1062	1063	rBV	1641	797	0.05%	0.006%
58	8.205	1114	1125	1139	rBV2	541157	1557315	100.00%	10.994%
59	8.599	1191	1192	1193	rBV	2457	1137	0.07%	0.008%
60	8.781	1214	1223	1234	rBV	513059	1042591	66.95%	7.360%

61	8.969	1254	1255	1256	rBV	2092	1080	0.07%	0.008%
62	9.034	1259	1266	1272	rVV2	55829	114883	7.38%	0.811%
63	9.210	1288	1296	1308	rBV	330074	699781	44.94%	4.940%
64	9.610	1360	1364	1368	rVV3	2062	3238	0.21%	0.023%
65	9.810	1397	1398	1400	rBV	1360	824	0.05%	0.006%

66	9.852	1403	1405	1408	rBV	1801	1036	0.07%	0.007%
67	9.981	1420	1427	1437	rVB	184609	345199	22.17%	2.437%
68	10.210	1464	1466	1467	rBV2	1991	862	0.06%	0.006%
69	10.269	1469	1476	1485	rVV	750081	1336402	85.81%	9.434%
70	10.528	1512	1520	1530	rBV	116906	205957	13.23%	1.454%

71	10.810	1561	1568	1575	rBV	537166	962966	61.84%	6.798%
72	10.881	1575	1580	1589	rVB	128798	224113	14.39%	1.582%
73	11.087	1613	1615	1619	rBV3	2481	3008	0.19%	0.021%
74	11.263	1644	1645	1651	rBV2	1350	2038	0.13%	0.014%
75	11.581	1692	1699	1708	rBV	680748	1179874	75.76%	8.329%

76	11.687	1713	1717	1720	rVB5	6928	9741	0.63%	0.069%
----	--------	------	------	------	------	------	------	-------	--------

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
 Data File : VD077806.D
 Acq On : 11 Dec 2023 11:23
 Operator : JC/SY
 Sample : 05750-02
 Misc : 4.24G/10ml/MSVOA_D/SOIL
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 E0LQ5

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_D\Method\SFAMDLM112923SMA.M
 Title : SFAM01.0

77	11.793	1729	1735	1741	rBV4	22400	43305	2.78%	0.306%
78	12.051	1776	1779	1780	rBV2	1580	1764	0.11%	0.012%
79	12.122	1786	1791	1797	rVB3	17598	32640	2.10%	0.230%
80	12.169	1797	1799	1803	rBV4	1515	1379	0.09%	0.010%
81	12.257	1812	1814	1816	rVB2	1664	1139	0.07%	0.008%
82	12.422	1838	1842	1845	rVB3	3028	4704	0.30%	0.033%
83	12.581	1865	1869	1873	rBV4	5913	8892	0.57%	0.063%
84	12.651	1874	1881	1888	rVB	199201	339686	21.81%	2.398%
85	12.828	1906	1911	1915	rVB3	39083	68446	4.40%	0.483%
86	12.904	1920	1924	1930	rVB3	35724	58673	3.77%	0.414%
87	13.063	1947	1951	1957	rVB	21319	34336	2.20%	0.242%
88	13.122	1959	1961	1962	rBV2	5487	4130	0.27%	0.029%
89	13.216	1971	1977	1983	rVB2	84567	136512	8.77%	0.964%
90	13.457	2015	2018	2021	rBV4	9283	12817	0.82%	0.090%
91	13.522	2022	2029	2039	rVB	602463	1051607	67.53%	7.424%
92	13.716	2059	2062	2066	rBV2	44198	62926	4.04%	0.444%
93	13.816	2073	2079	2089	rVB	540600	961981	61.77%	6.791%
94	14.063	2117	2121	2126	rVB	47402	73948	4.75%	0.522%
95	14.175	2135	2140	2145	rVB3	50678	83445	5.36%	0.589%
96	14.469	2187	2190	2194	rBV2	28461	36936	2.37%	0.261%
97	14.657	2218	2222	2226	rBV3	56682	103895	6.67%	0.733%
98	14.775	2235	2242	2247	rBV4	120849	271277	17.42%	1.915%
99	14.928	2263	2268	2275	rVB	89624	154935	9.95%	1.094%
100	15.392	2343	2347	2351	rBV2	48965	73628	4.73%	0.520%

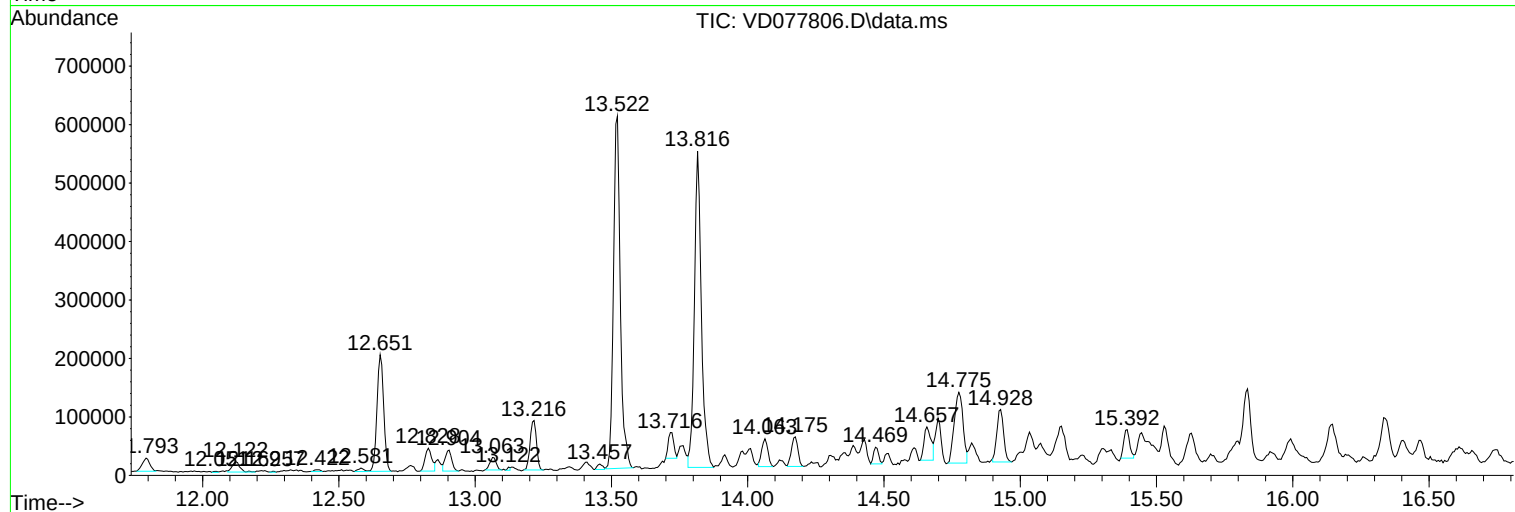
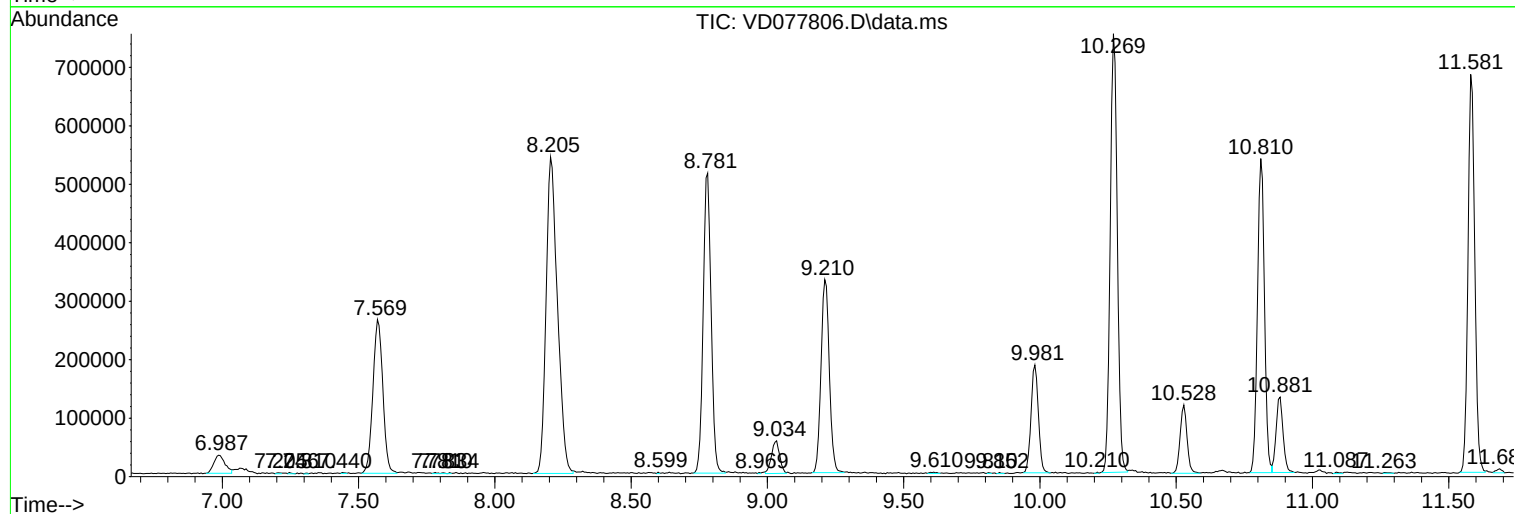
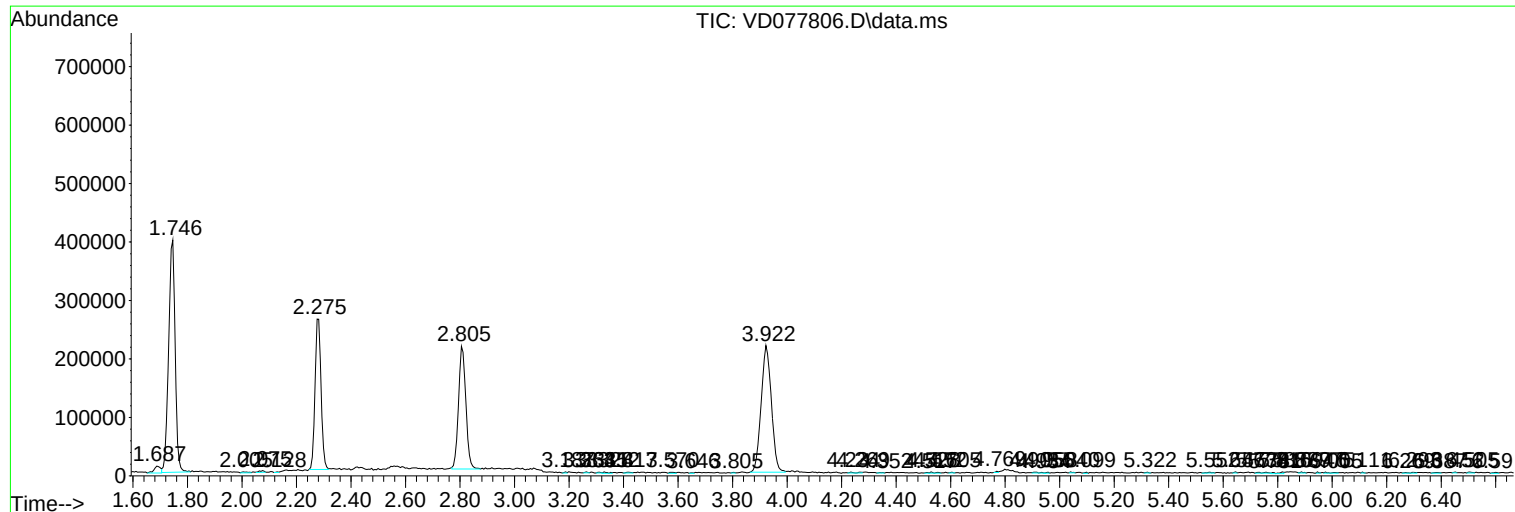
Sum of corrected areas: 14165452

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

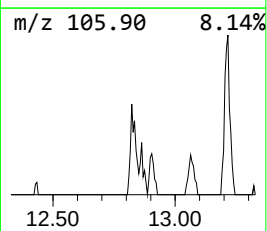
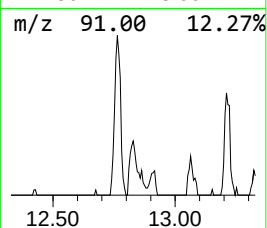
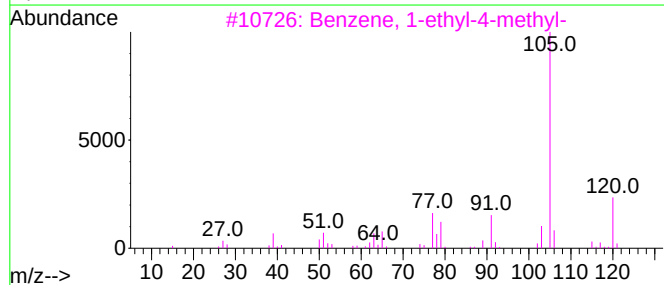
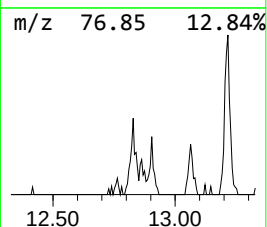
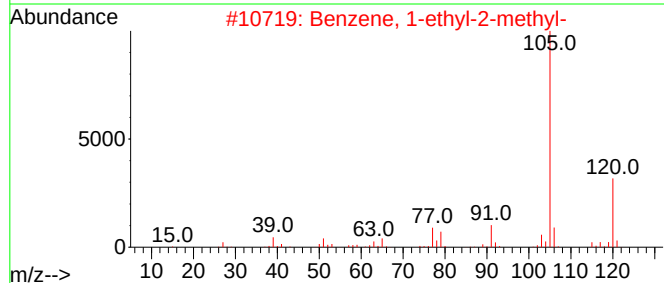
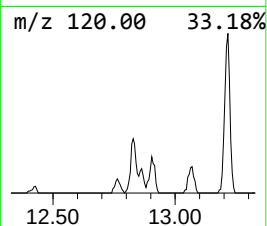
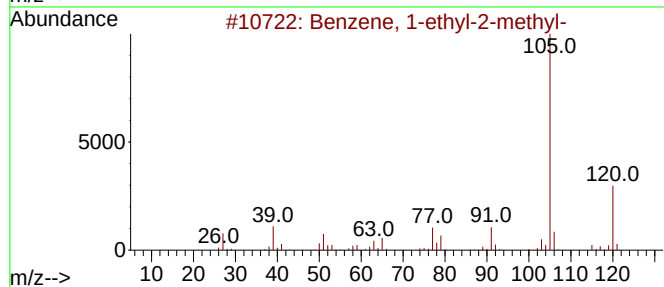
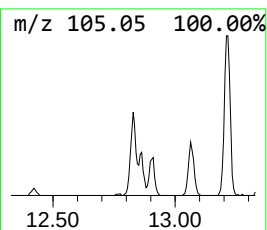
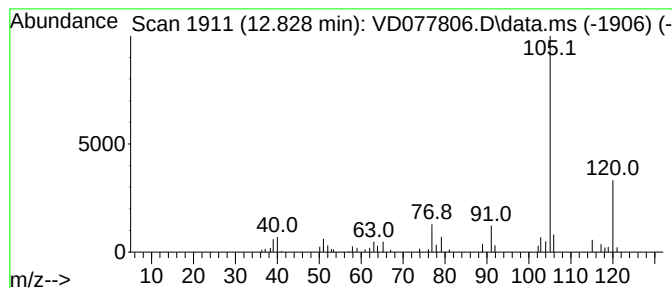
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.828	1.63 ug/L	68446	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

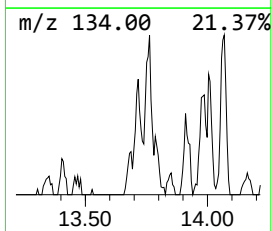
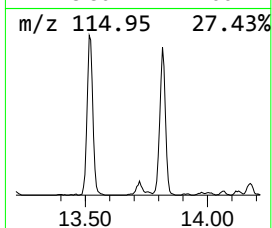
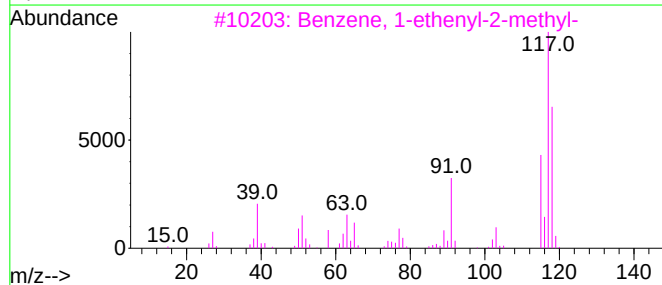
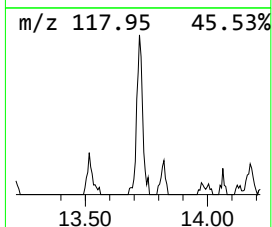
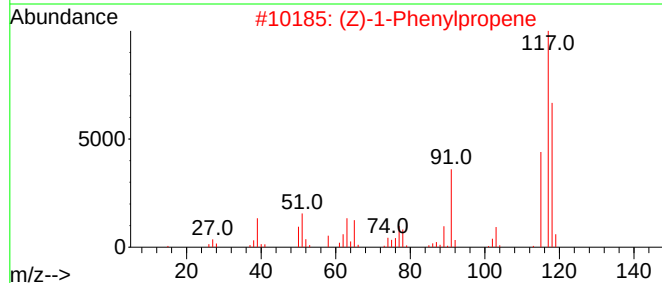
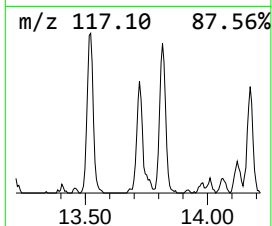
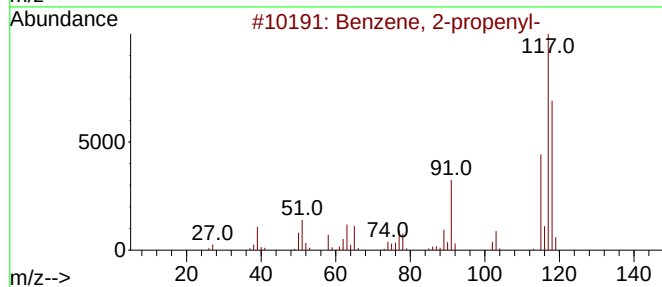
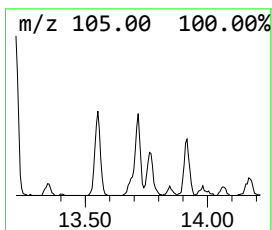
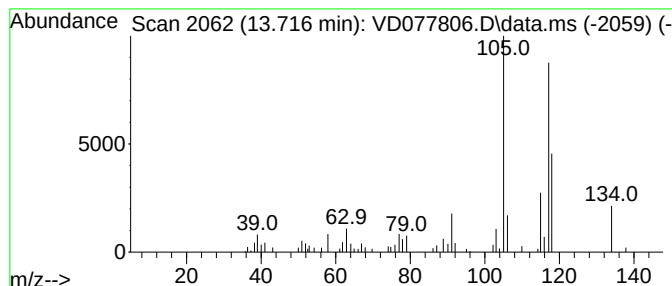
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 Benzene, 2-propenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	1.50 ug/L	62926	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
2			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	60
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	58
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

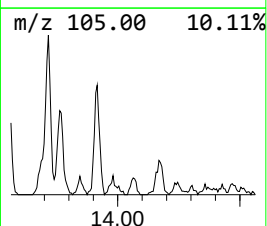
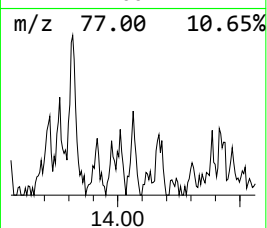
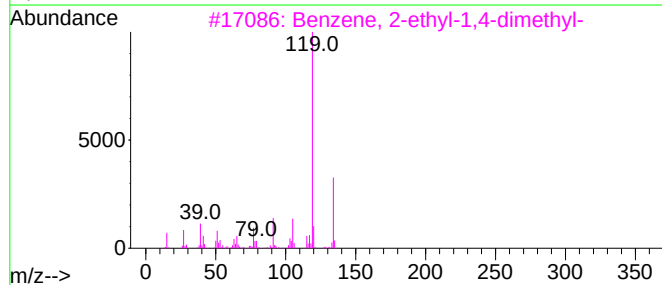
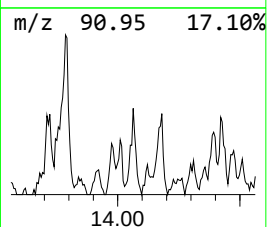
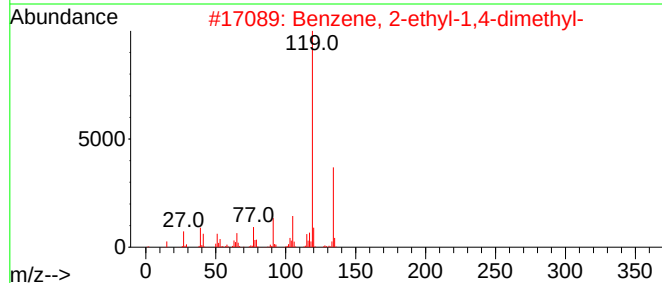
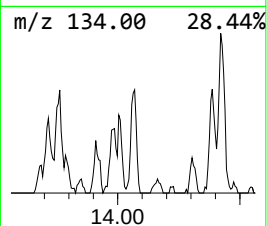
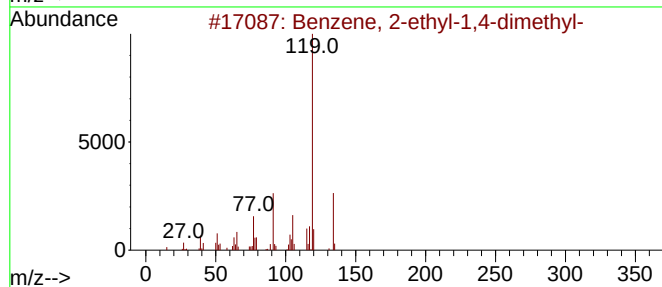
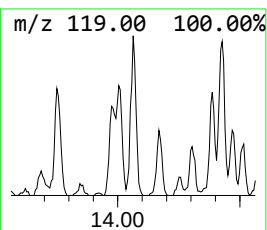
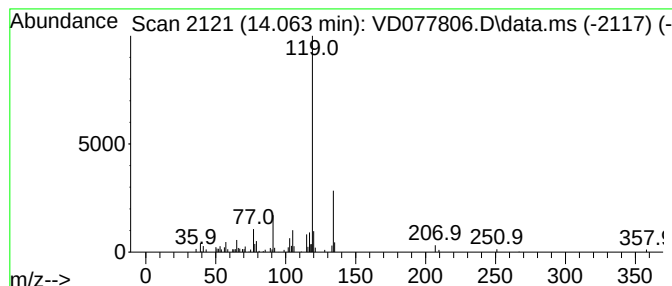
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.063	1.76 ug/L	73948	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

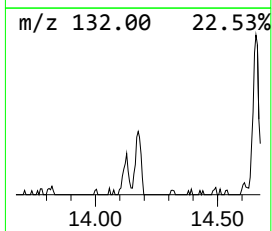
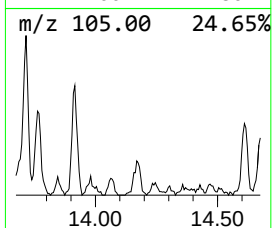
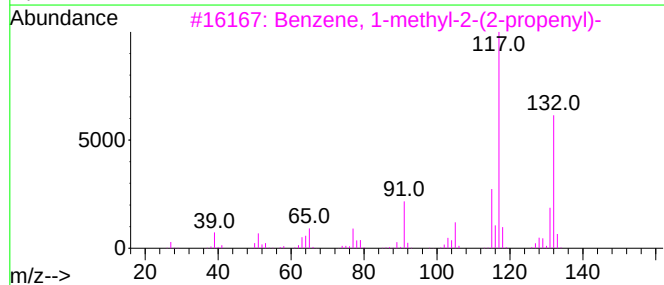
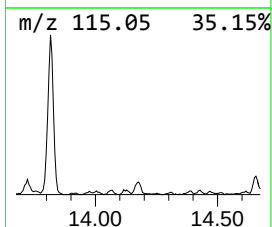
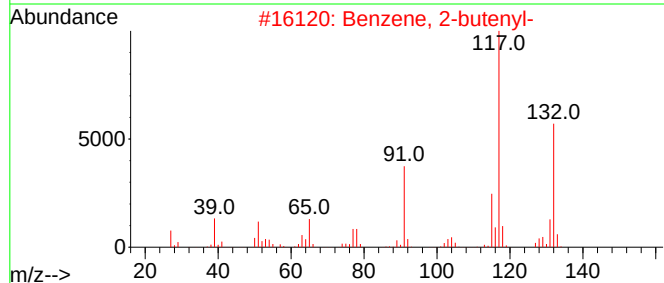
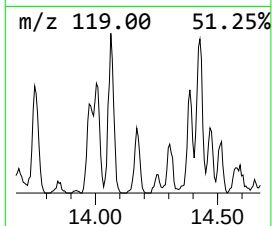
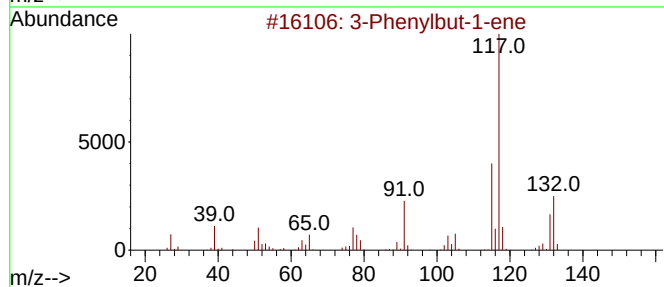
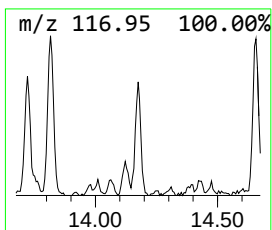
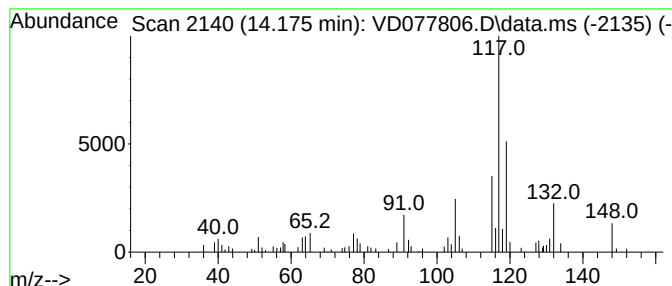
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 6 3-Phenylbut-1-ene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.175	1.98 ug/L	83445	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	58
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	58
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	58
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

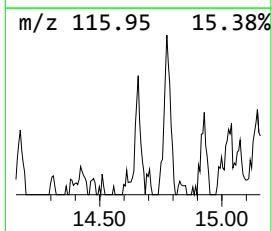
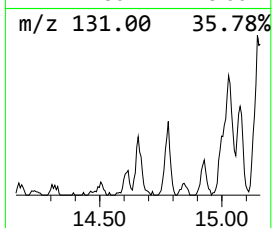
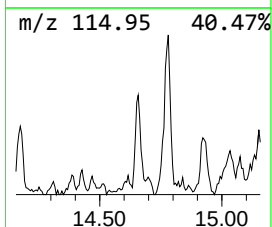
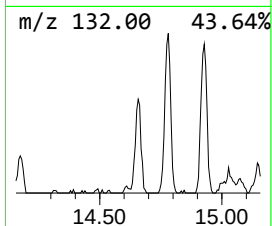
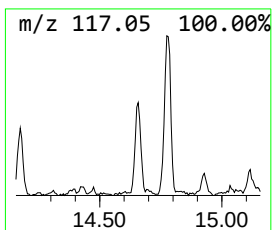
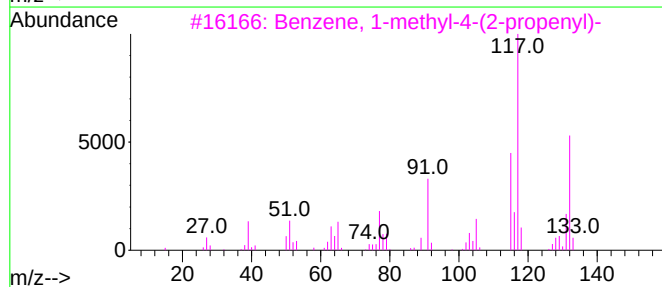
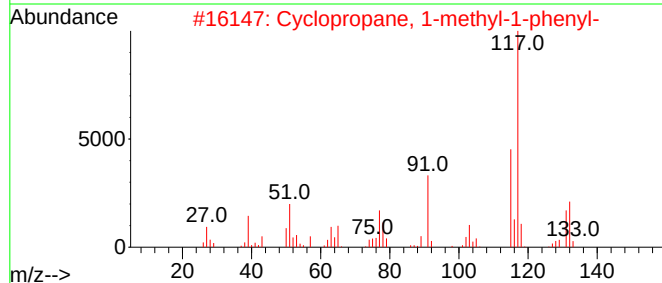
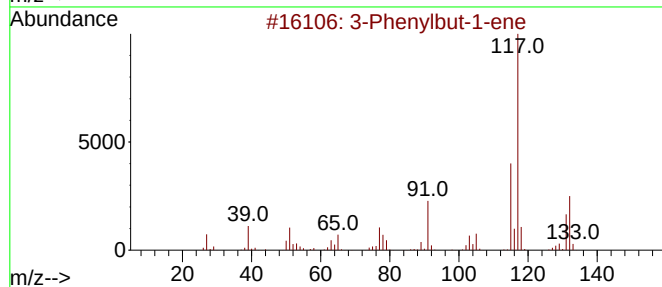
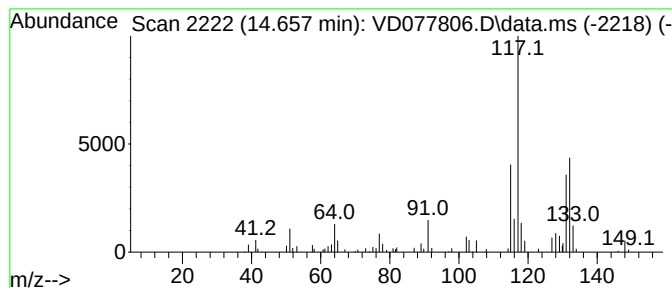
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 7 Cyclopropane, 1-methyl-1-ph... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.657	2.47 ug/L	103895	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
2			Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	87
3			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	86
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

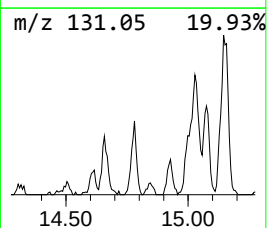
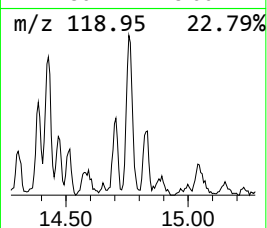
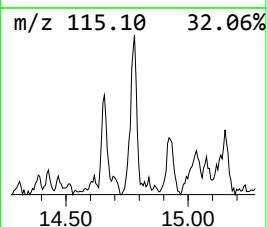
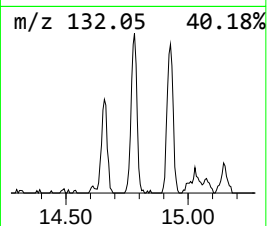
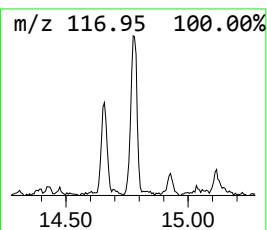
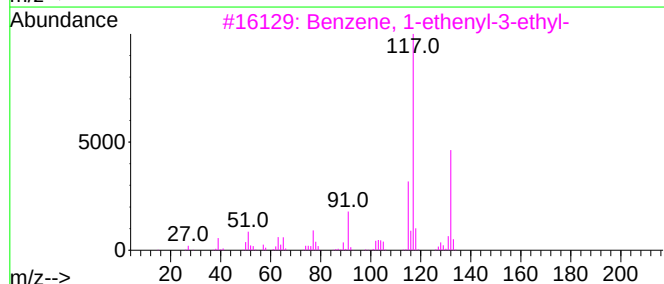
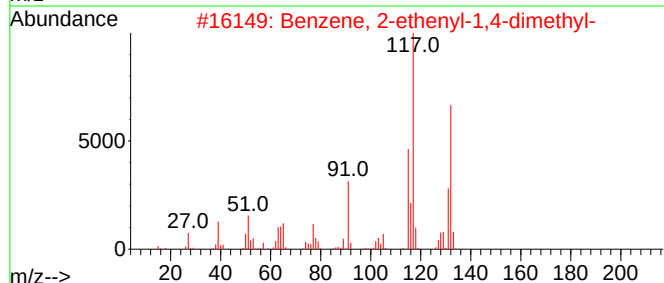
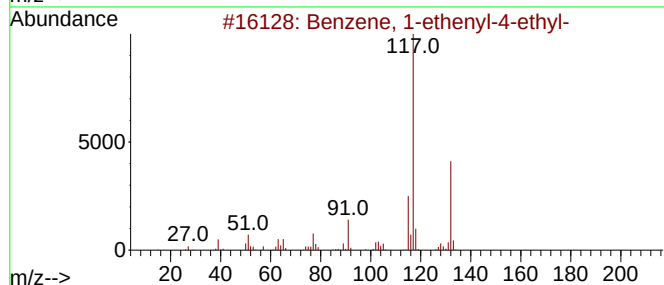
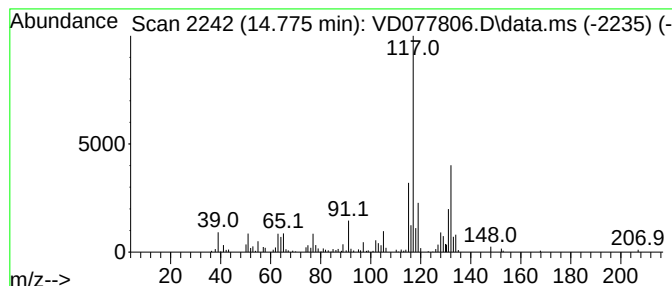
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.775	6.45 ug/L	271277	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

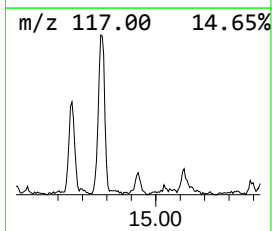
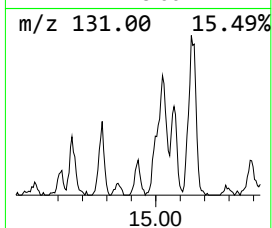
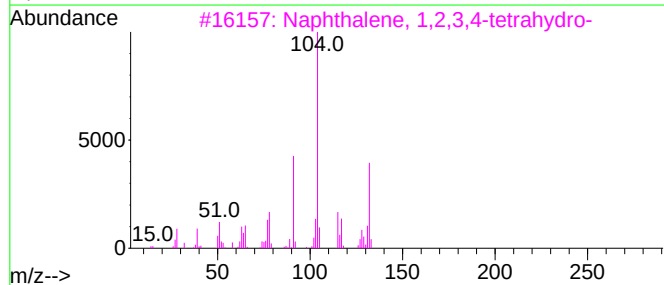
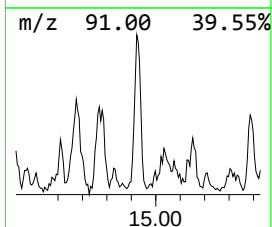
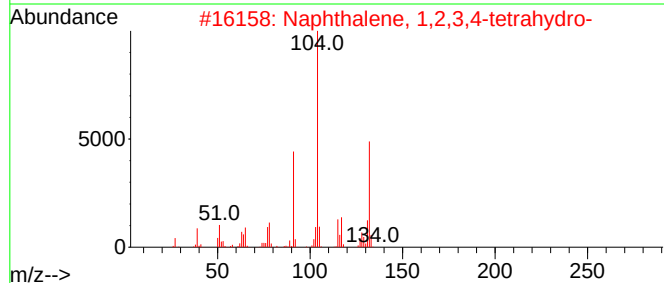
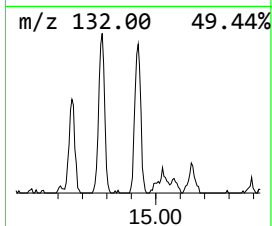
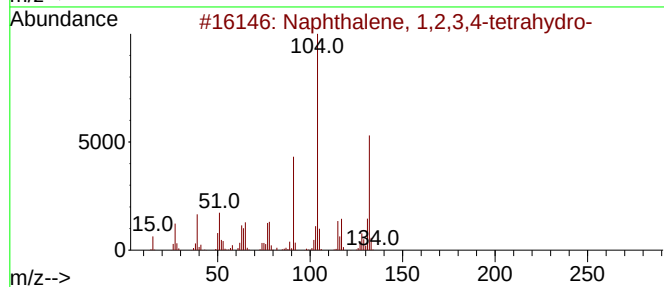
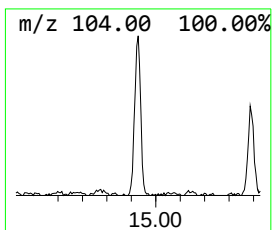
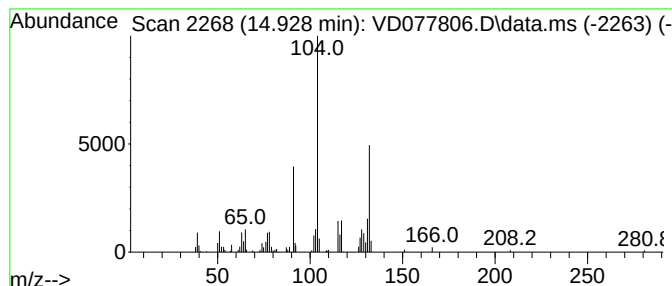
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.928	3.68 ug/L	154935	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	87
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	72
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	64
5			.delta.-6,7-Octalin, 1.beta.,4.b...	168	C10H16O2	051921-13-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

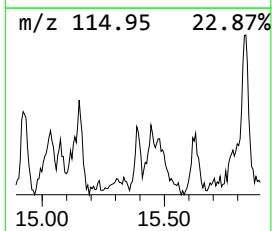
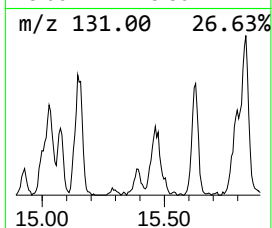
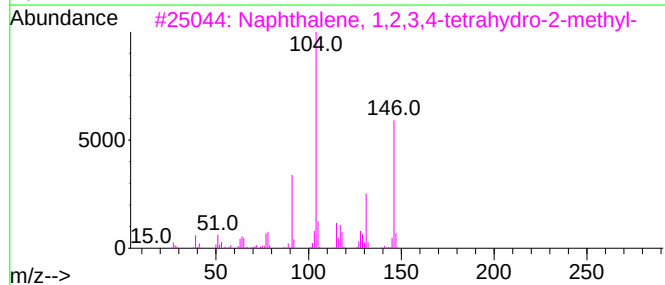
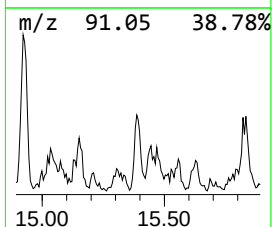
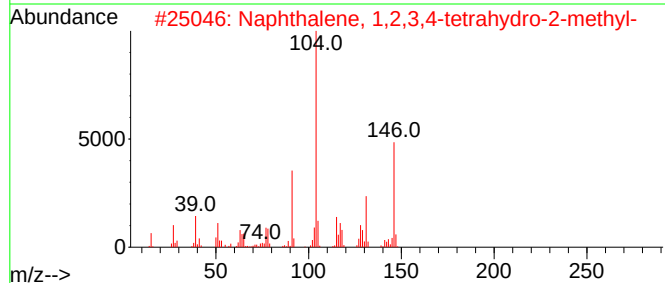
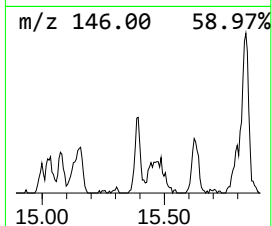
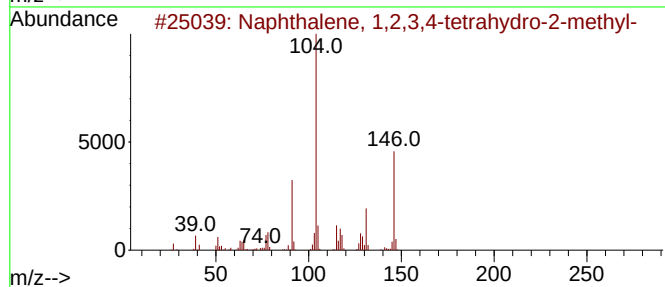
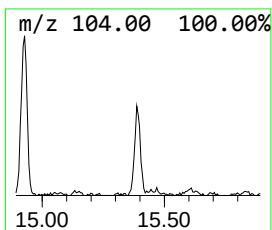
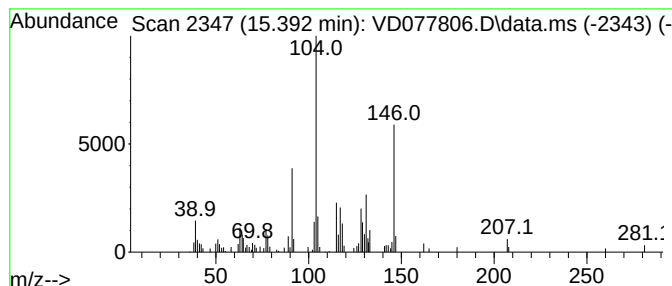
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.M
Quant Title : SFAM01.0

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.392	1.75 ug/L	73628	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	93
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	70
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	64
5			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121123\
Data File : VD077806.D
Acq On : 11 Dec 2023 11:23
Operator : JC/SY
Sample : 05750-02
Misc : 4.24G/10ml/MSVOA_D/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
E0LQ5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM112923SMA.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
Benzene, 1-ethy...	12.828	1.6	ug/L	68446	3	13.522	1051610	25.0
Benzene, 2-prop...	13.716	1.5	ug/L	62926	3	13.522	1051610	25.0
Benzene, 2-ethy...	14.063	1.8	ug/L	73948	3	13.522	1051610	25.0
3-Phenylbut-1-ene	14.175	2.0	ug/L	83445	3	13.522	1051610	25.0
Cyclopropane, 1...	14.657	2.5	ug/L	103895	3	13.522	1051610	25.0
Benzene, 1-ethe...	14.775	6.5	ug/L	271277	3	13.522	1051610	25.0
Naphthalene, 1,...	14.928	3.7	ug/L	154935	3	13.522	1051610	25.0
Naphthalene, 1,...	15.392	1.8	ug/L	73628	3	13.522	1051610	25.0