

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\
 Data File : VY006932.D
 Acq On : 04 Dec 2021 11:46
 Operator : SY/MD
 Sample : M4884-02
 Misc : 3.17g/10.0mL/MSVOA_Y/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 EW9H8

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_Y\methods\SFAMYL120321SMA.M
 Title : VOC Analysis

Signal : TIC: VY006932.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.253	66	71	80	rVB	24922	42839	11.90%	1.671%
2	2.772	148	156	166	rBV	17340	40749	11.32%	1.589%
3	3.089	206	208	211	rVB3	456	483	0.13%	0.019%
4	3.229	228	231	232	rBV2	243	316	0.09%	0.012%
5	3.259	234	236	241	rVB4	681	934	0.26%	0.036%
6	3.302	241	243	248	rVB2	313	372	0.10%	0.015%
7	3.351	248	251	252	rBV2	336	296	0.08%	0.012%
8	3.540	281	282	286	rVB3	455	429	0.12%	0.017%
9	3.595	289	291	299	rVB2	325	571	0.16%	0.022%
10	3.741	309	315	316	rBV2	358	358	0.10%	0.014%
11	3.857	324	334	344	rBV2	39635	103781	28.83%	4.047%
12	4.198	381	390	399	rBV2	2995	7949	2.21%	0.310%
13	4.442	426	430	432	rVB	276	330	0.09%	0.013%
14	4.472	432	435	436	rBV	351	315	0.09%	0.012%
15	4.716	466	475	488	rBV2	13038	38925	10.81%	1.518%
16	4.918	501	508	512	rBV4	865	2137	0.59%	0.083%
17	4.985	517	519	521	rVV	470	359	0.10%	0.014%
18	5.259	560	564	567	rVV2	249	417	0.12%	0.016%
19	5.344	575	578	582	rVB2	254	325	0.09%	0.013%
20	5.741	638	643	644	rVV3	1679	2161	0.60%	0.084%
21	5.759	644	646	653	rVV3	1803	3021	0.84%	0.118%
22	5.875	661	665	668	rBV	413	455	0.13%	0.018%
23	6.277	728	731	734	rVB	274	332	0.09%	0.013%
24	6.466	760	762	765	rBV2	247	305	0.08%	0.012%
25	6.679	795	797	799	rVB	471	326	0.09%	0.013%
26	6.722	799	804	807	rBV3	336	647	0.18%	0.025%
27	6.789	811	815	816	rBV2	399	482	0.13%	0.019%
28	6.887	822	831	841	rBV	21149	61833	17.18%	2.411%
29	6.984	841	847	856	rVB2	6877	19943	5.54%	0.778%
30	7.484	918	929	942	rVB	58407	144847	40.24%	5.648%
31	7.624	949	952	956	rVB3	227	368	0.10%	0.014%
32	7.746	968	972	974	rBV	360	429	0.12%	0.017%
33	7.844	984	988	992	rBV	232	404	0.11%	0.016%
34	8.051	1016	1022	1024	rBV2	362	558	0.16%	0.022%
35	8.118	1024	1033	1047	rVV2	115151	359953	100.00%	14.036%
36	8.240	1047	1053	1060	rVB3	5447	11035	3.07%	0.430%

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYL120321SMA.M
 Title : VOC Analysis

37	8.466	1088	1090	1094	rVB2	204	289	0.08%	0.011%
38	8.691	1119	1127	1136	rBV	88658	173690	48.25%	6.773%
39	8.801	1143	1145	1149	rBV2	214	369	0.10%	0.014%
40	8.941	1163	1168	1174	rVB4	1171	2028	0.56%	0.079%
41	9.075	1186	1190	1191	rVV	573	583	0.16%	0.023%
42	9.124	1191	1198	1206	rVV	78299	157666	43.80%	6.148%
43	9.191	1208	1209	1211	rVV	1209	725	0.20%	0.028%
44	9.209	1211	1212	1216	rVV	492	360	0.10%	0.014%
45	9.362	1234	1237	1240	rVV2	354	483	0.13%	0.019%
46	9.514	1258	1262	1265	rVV4	719	1157	0.32%	0.045%
47	9.563	1265	1270	1272	rVB2	318	478	0.13%	0.019%
48	9.746	1296	1300	1306	rVV4	852	1267	0.35%	0.049%
49	9.892	1315	1324	1333	rVB	41497	69714	19.37%	2.718%
50	9.971	1335	1337	1344	rVV3	619	1115	0.31%	0.043%
51	10.039	1344	1348	1349	rVV2	223	344	0.10%	0.013%
52	10.069	1350	1353	1358	rVV2	1199	1764	0.49%	0.069%
53	10.124	1361	1362	1364	rVV	488	413	0.11%	0.016%
54	10.179	1364	1371	1378	rVV	119472	202332	56.21%	7.890%
55	10.246	1378	1382	1388	rVV2	11391	18703	5.20%	0.729%
56	10.441	1407	1414	1420	rBV	25038	42492	11.80%	1.657%
57	10.514	1423	1426	1428	rVB2	241	291	0.08%	0.011%
58	10.581	1428	1437	1446	rBV	38659	71816	19.95%	2.800%
59	10.709	1454	1458	1462	rBV2	703	1222	0.34%	0.048%
60	10.788	1464	1471	1480	rVV	63248	99919	27.76%	3.896%
61	10.862	1482	1483	1487	rVB	350	471	0.13%	0.018%
62	10.947	1490	1497	1507	rBV2	20879	37375	10.38%	1.457%
63	11.038	1510	1512	1518	rVV7	1437	2179	0.61%	0.085%
64	11.319	1547	1558	1565	rVB2	28930	48380	13.44%	1.887%
65	11.489	1579	1586	1593	rVV	95795	153700	42.70%	5.994%
66	11.593	1598	1603	1608	rVB2	1822	2851	0.79%	0.111%
67	11.654	1608	1613	1614	rBV	247	332	0.09%	0.013%
68	11.697	1614	1620	1627	rVV3	5353	9086	2.52%	0.354%
69	12.032	1668	1675	1681	rVB3	3953	7185	2.00%	0.280%
70	12.099	1683	1686	1689	rBV3	259	293	0.08%	0.011%
71	12.203	1700	1703	1707	rBV3	197	308	0.09%	0.012%
72	12.245	1707	1710	1711	rBV	354	296	0.08%	0.012%
73	12.282	1714	1716	1719	rVB2	300	363	0.10%	0.014%
74	12.337	1719	1725	1729	rBV4	1605	2802	0.78%	0.109%
75	12.398	1731	1735	1738	rBV2	453	603	0.17%	0.024%
76	12.489	1741	1750	1756	rBV	132470	223591	62.12%	8.719%

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYLM120321SMA.M
 Title : VOC Analysis

77	12.562	1756	1762	1769	rVB	62155	98341	27.32%	3.835%
78	12.623	1769	1772	1774	rBV2	326	471	0.13%	0.018%
79	12.660	1776	1778	1780	rBV2	461	380	0.11%	0.015%
80	12.739	1788	1791	1793	rBV3	782	824	0.23%	0.032%
81	12.764	1793	1795	1799	rVB3	639	573	0.16%	0.022%
82	12.819	1799	1804	1810	rBV	8601	14850	4.13%	0.579%
83	12.904	1816	1818	1821	rBV2	433	329	0.09%	0.013%
84	12.965	1826	1828	1832	rVB3	651	777	0.22%	0.030%
85	13.068	1840	1845	1850	rVB4	1576	2509	0.70%	0.098%
86	13.117	1850	1853	1857	rBV3	1685	2539	0.71%	0.099%
87	13.184	1859	1864	1873	rVB2	11065	17091	4.75%	0.666%
88	13.276	1873	1879	1885	rBV5	2060	4182	1.16%	0.163%
89	13.373	1892	1895	1896	rVV2	720	666	0.19%	0.026%
90	13.428	1898	1904	1909	rVV	47827	76772	21.33%	2.994%
91	13.483	1909	1913	1917	rVV7	2382	4689	1.30%	0.183%
92	13.532	1917	1921	1926	rVB4	2505	3864	1.07%	0.151%
93	13.611	1932	1934	1935	rBV2	465	288	0.08%	0.011%
94	13.642	1935	1939	1943	rVV6	1944	3282	0.91%	0.128%
95	13.721	1946	1952	1959	rVV	50616	83635	23.23%	3.261%
96	13.965	1983	1992	1999	rBV2	21483	40111	11.14%	1.564%
97	14.702	2111	2113	2115	rBV3	855	523	0.15%	0.020%
98	15.318	2210	2214	2219	rVB3	3228	4839	1.34%	0.189%
99	15.391	2221	2226	2235	rVB	8002	14525	4.04%	0.566%
100	15.727	2280	2281	2282	rBV	847	350	0.10%	0.014%

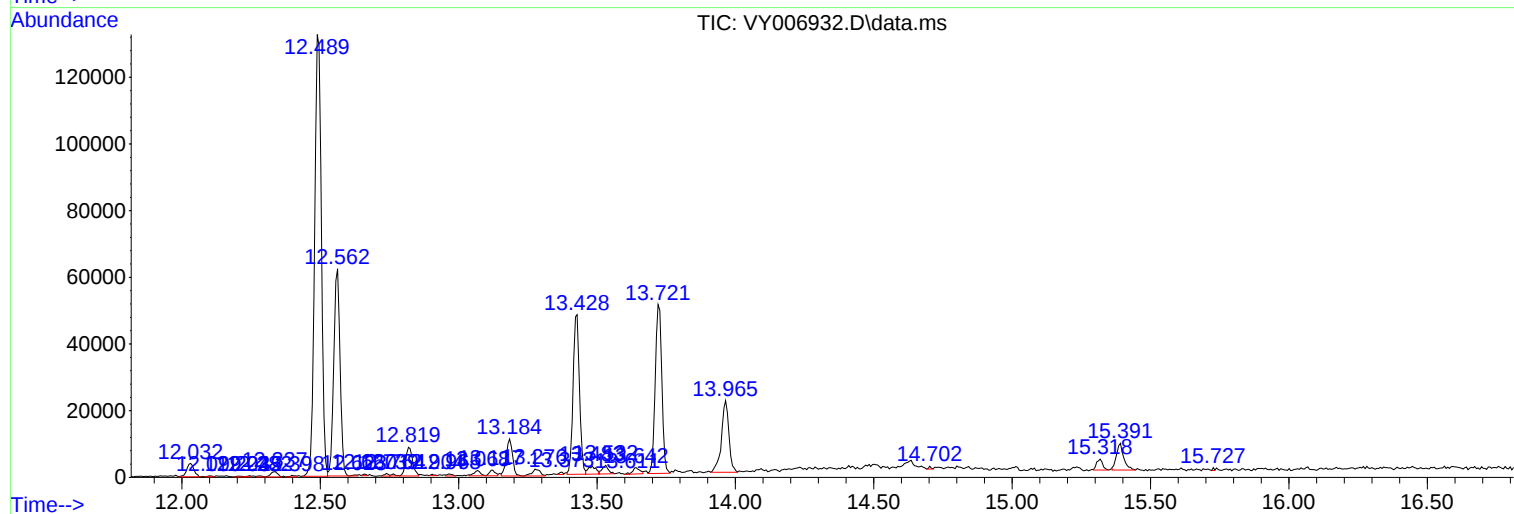
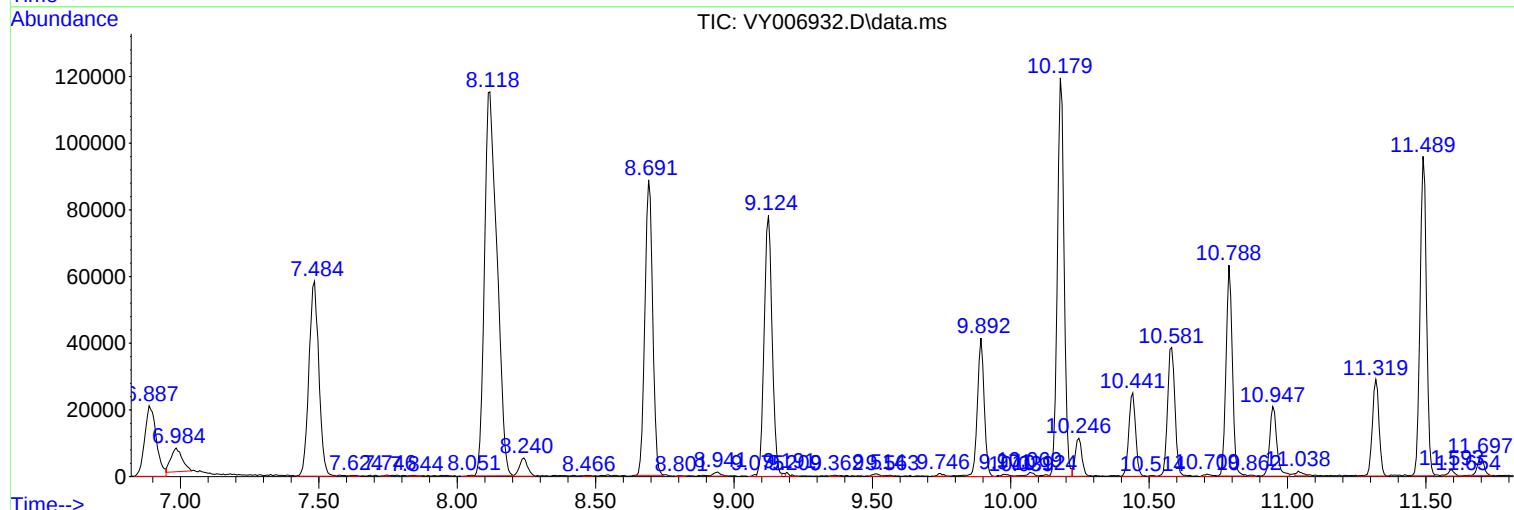
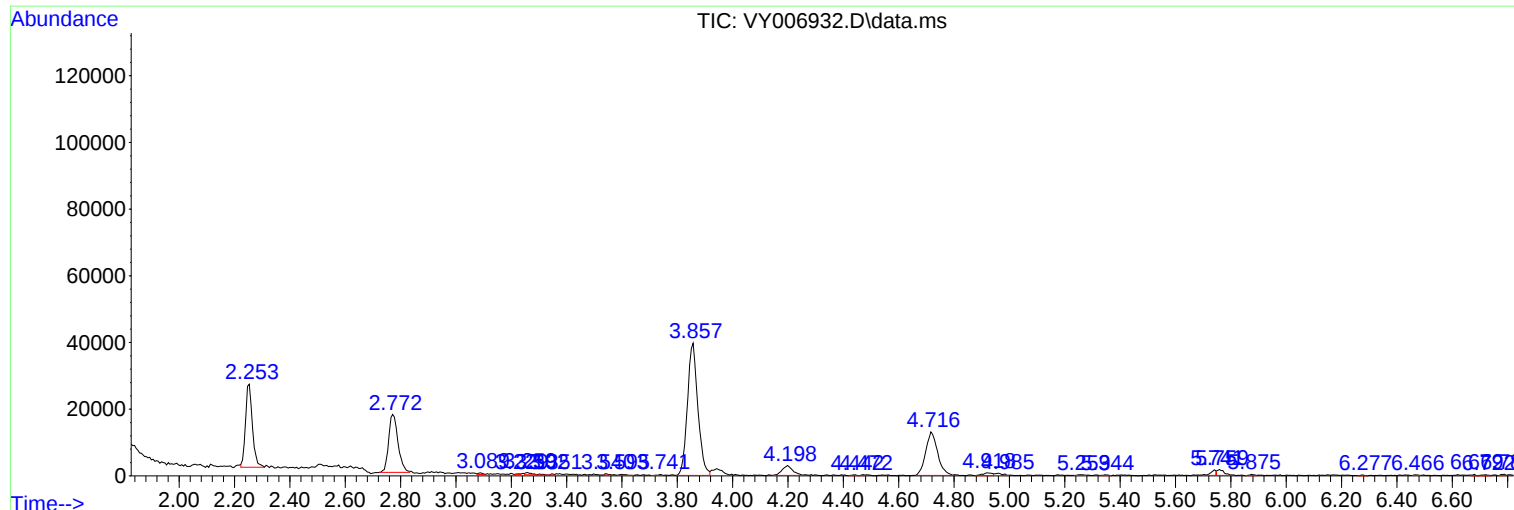
Sum of corrected areas: 2564429

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Instrument :
MSVOA_Y
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYL120321SMA.M
Quant Title : VOC Analysis

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



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Instrument :
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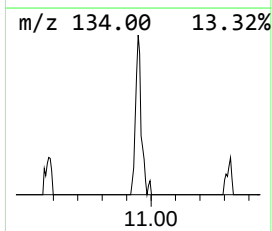
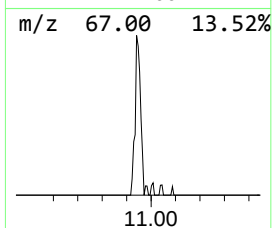
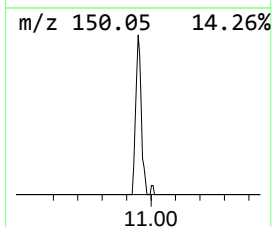
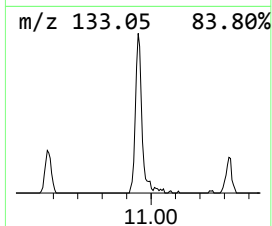
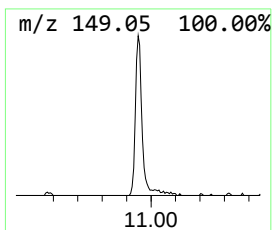
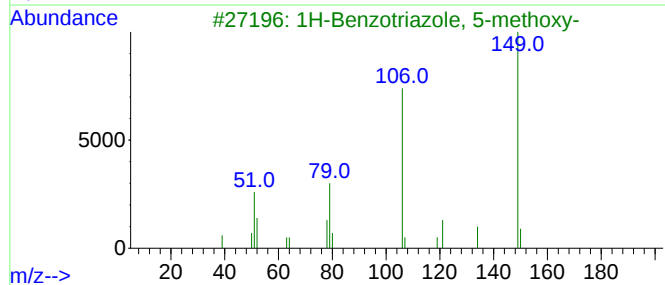
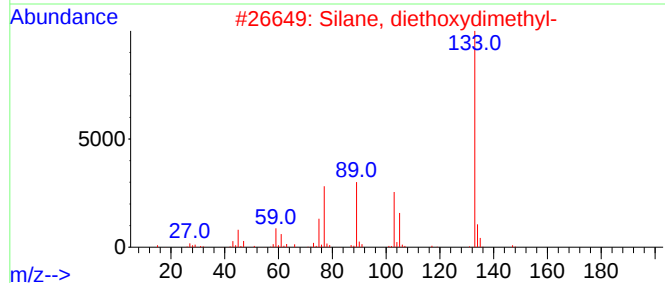
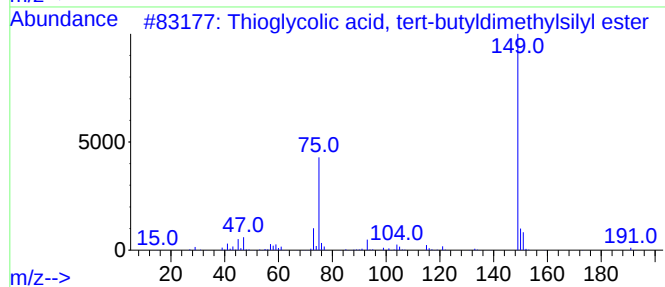
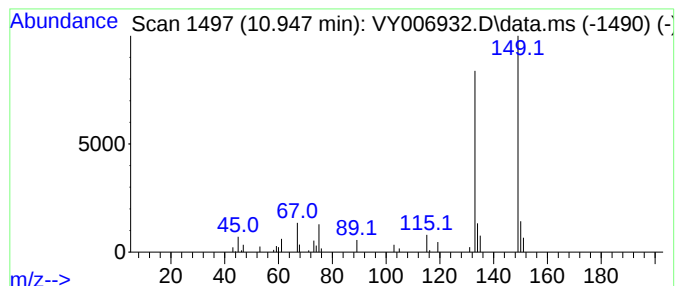
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYL120321SMA.M
Quant Title : VOC Analysis

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

Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.947	6.08 ug/L	37375	Chlorobenzene-d5	11.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thioglycolic acid, tert-butyldim...	206	C8H18O2SSi	1000476-01-2	32
2			Silane, diethoxydimethyl-	148	C6H16O2Si	000078-62-6	25
3			1H-Benzotriazole, 5-methoxy-	149	C7H7N3O	027799-91-3	25
4			Silane, diethoxydimethyl-	148	C6H16O2Si	000078-62-6	25
5			4-Methylthieno[2,3-b]pyridine	149	C8H7NS	013362-81-7	25



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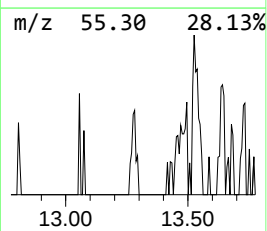
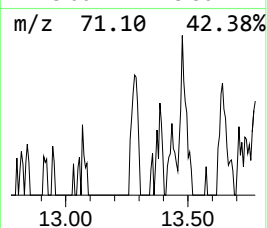
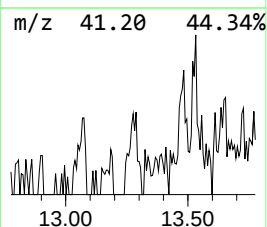
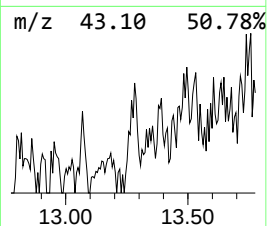
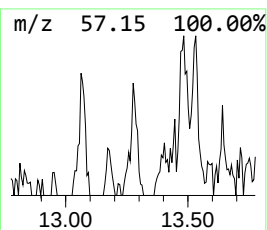
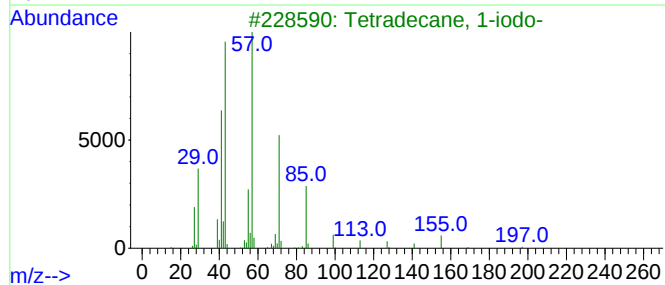
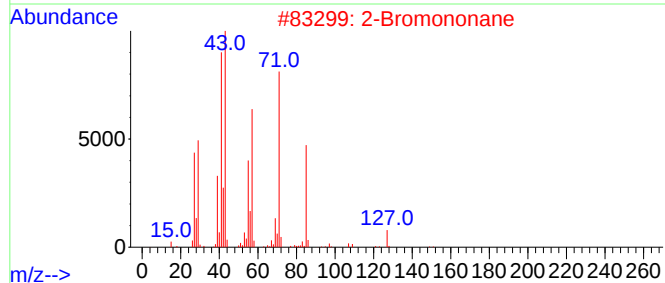
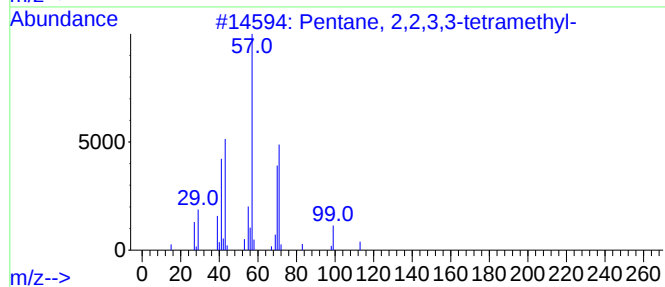
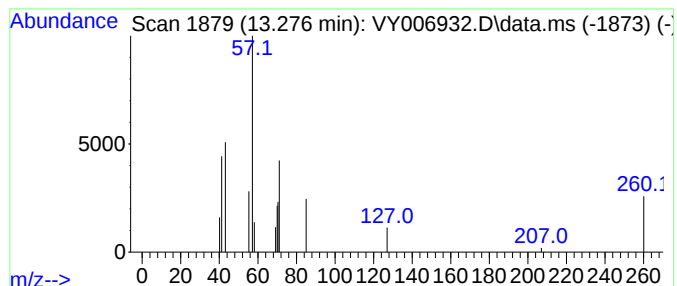
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYLEM120321SMA.M
Quant Title : VOC Analysis

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.276	1.36 ug/L	4182	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	40
2			2-Bromononane	206	C9H19Br	002216-35-5	25
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	25
4			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	25
5			Trichloroacetic acid, 2-methylbu...	232	C7H11Cl3O2	1000330-88-8	25



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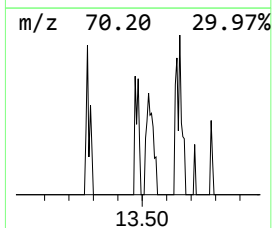
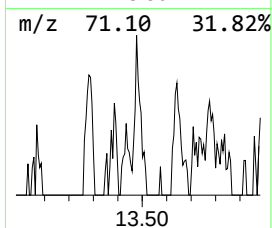
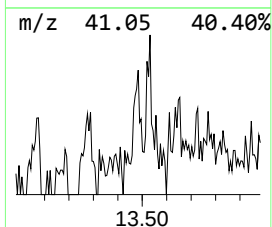
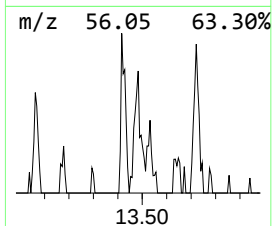
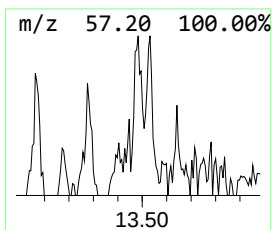
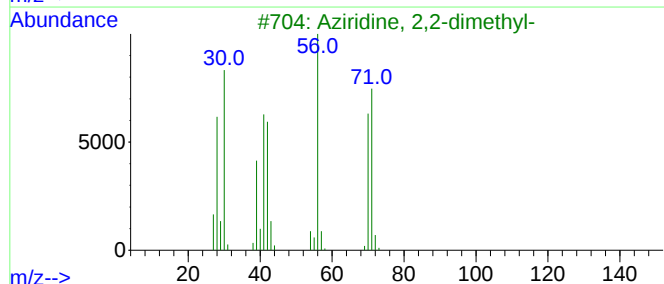
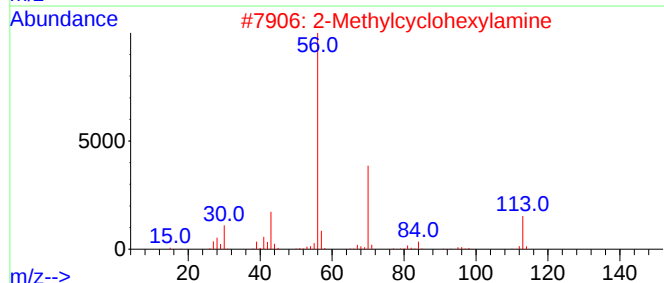
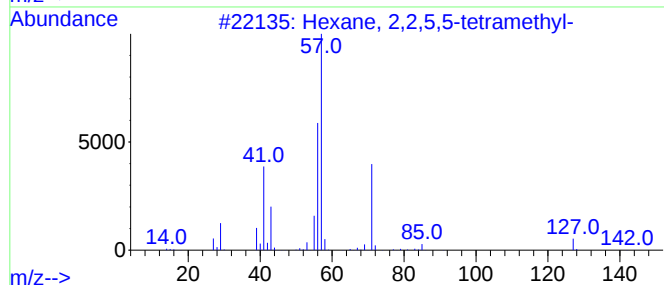
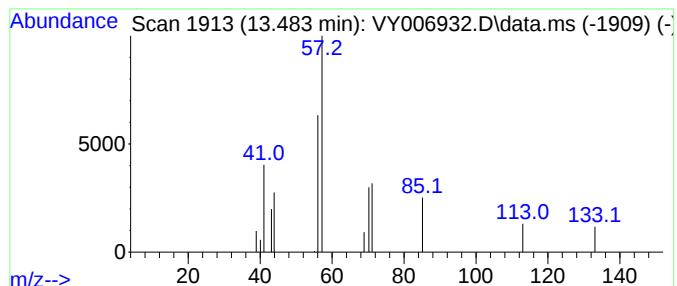
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYL120321SMA.M
Quant Title : VOC Analysis

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.483	1.53 ug/L	4689	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	23
2			2-Methylcyclohexylamine	113	C7H15N	007003-32-9	9
3			Aziridine, 2,2-dimethyl-	71	C4H9N	002658-24-4	9
4			2-Methylcyclohexylamine	113	C7H15N	007003-32-9	9
5			Aziridine, 2,2-dimethyl-	71	C4H9N	002658-24-4	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\
Data File : VY006932.D
Acq On : 04 Dec 2021 11:46
Operator : SY/MD
Sample : M4884-02
Misc : 3.17g/10.0mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EW9H8

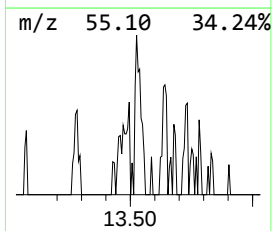
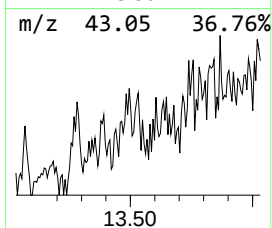
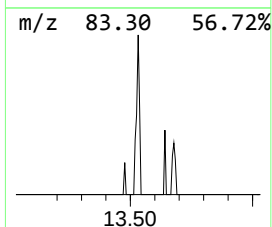
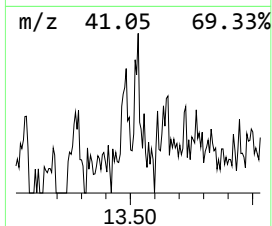
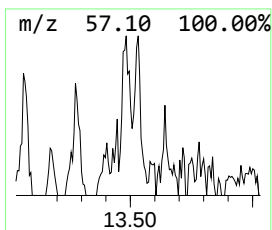
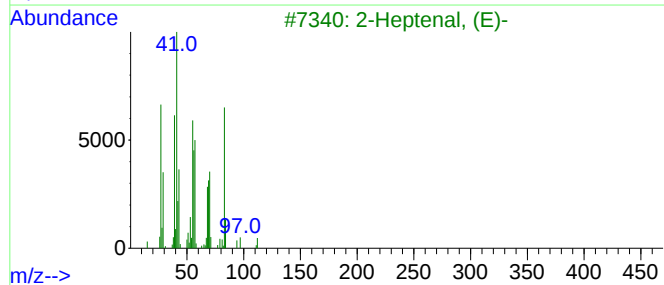
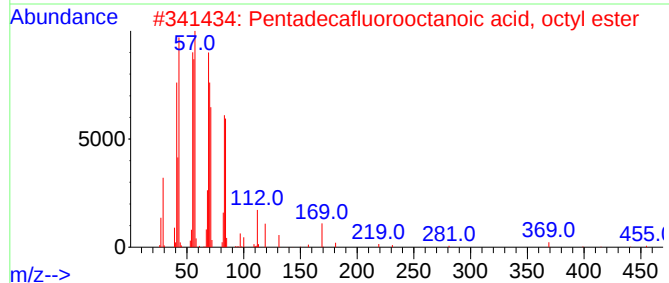
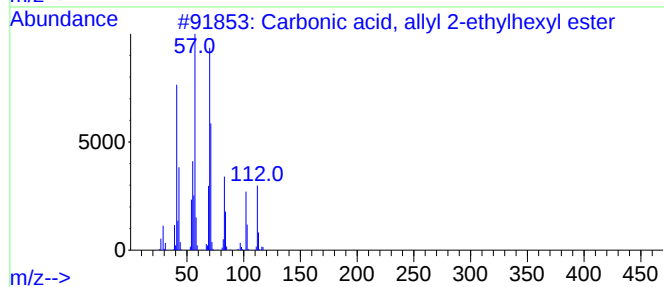
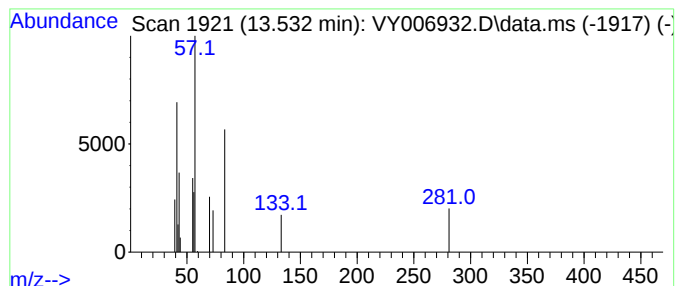
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Quant Title : VOC Analysis

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

Peak Number 11 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.532	1.26 ug/L	3864	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, allyl 2-ethylhexy...	214	C12H22O3	1000357-80-6	17
2			Pentadecafluorooctanoic acid, oc...	526	C16H17F15O2	1000406-03-8	12
3			2-Heptenal, (E)-	112	C7H12O	018829-55-5	9
4			3-Tridecene, (Z)-	182	C13H26	041446-53-1	9
5			1-Propanol, 2,2-dimethyl-	88	C5H12O	000075-84-3	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\
Data File : VY006932.D
Acq On : 04 Dec 2021 11:46
Operator : SY/MD
Sample : M4884-02
Misc : 3.17g/10.0mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EW9H8

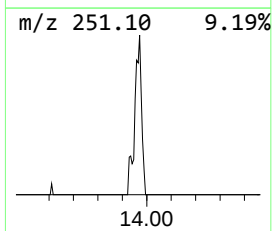
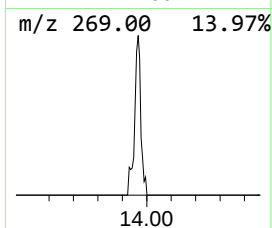
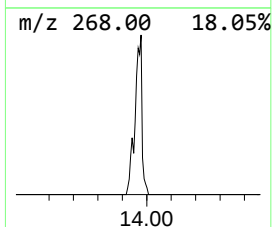
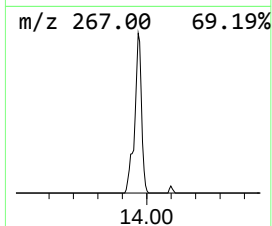
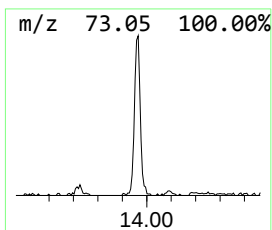
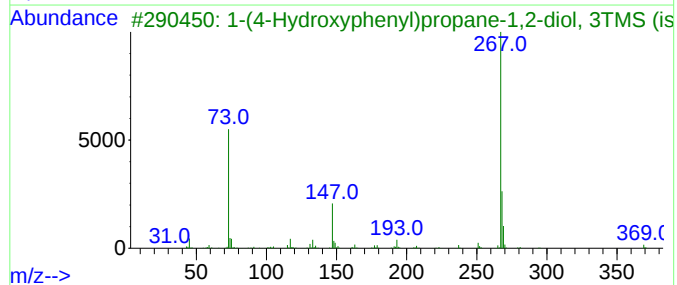
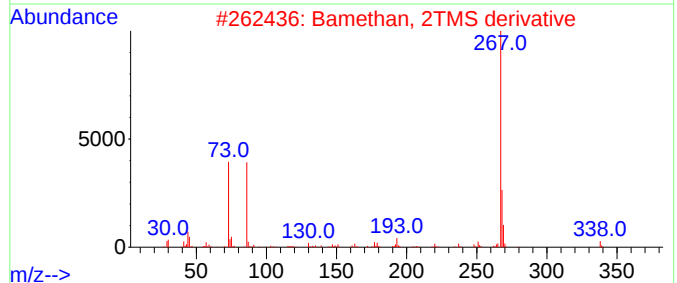
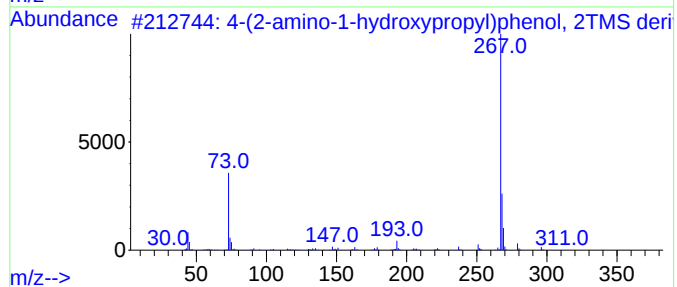
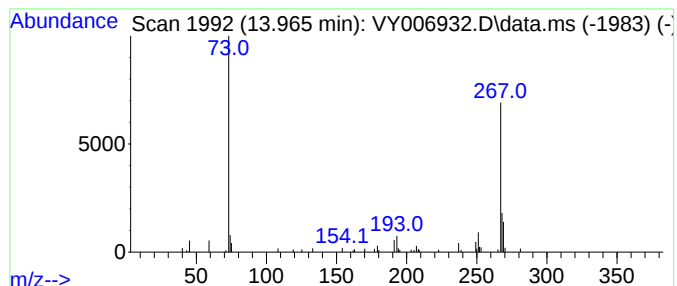
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Quant Title : VOC Analysis

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

Peak Number 12 4-(2-amino-1-hydroxypropyl)... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.965	13.06 ug/L	40111	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(2-amino-1-hydroxypropyl)pheno...	311	C15H29NO2Si2	1000478-98-6	64
2			Bamethan, 2TMS derivative	353	C18H35NO2Si2	040629-66-1	64
3			1-(4-Hydroxyphenyl)propane-1,2-d...	384	C18H36O3Si3	1000495-85-7	59
4			2-Hydroxymandelic acid, ethyl es...	340	C16H28O4Si2	1000071-88-8	59
5			1-(4-Hydroxyphenyl)propane-1,2-d...	384	C18H36O3Si3	1000495-85-6	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\
Data File : VY006932.D
Acq On : 04 Dec 2021 11:46
Operator : SY/MD
Sample : M4884-02
Misc : 3.17g/10.0mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EW9H8

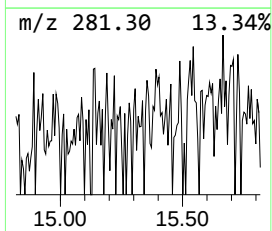
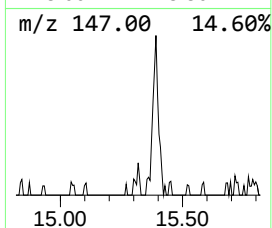
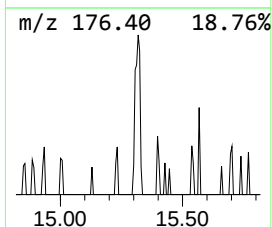
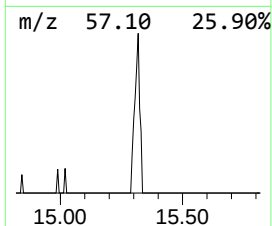
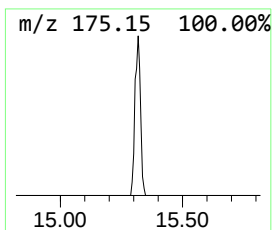
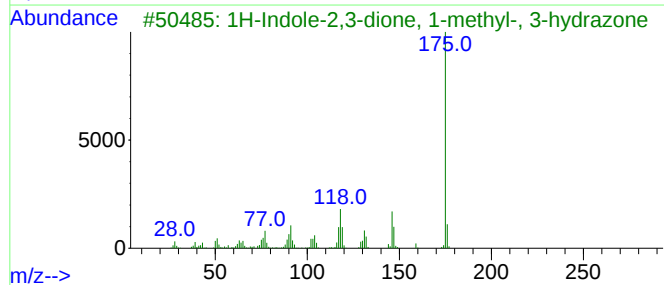
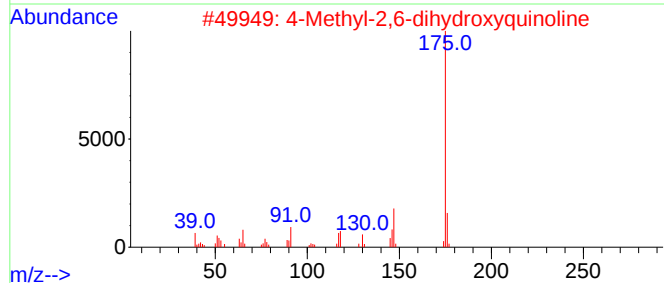
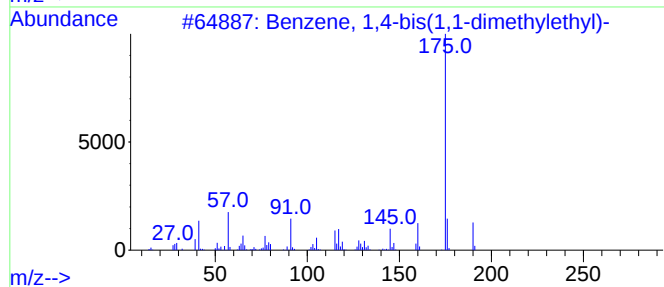
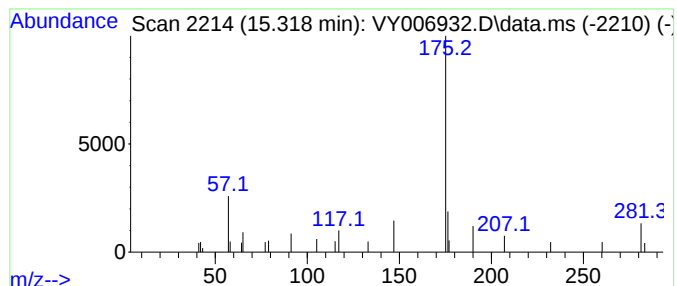
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Quant Title : VOC Analysis

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

Peak Number 13 Benzene, 1,4-bis(1,1-dimeth... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.318	1.58 ug/L	4839	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-bis(1,1-dimethyleth...	190	C14H22	001012-72-2	58
2			4-Methyl-2,6-dihydroxyquinoline	175	C10H9NO2	034982-01-9	53
3			1H-Indole-2,3-dione, 1-methyl-, ...	175	C9H9N3O	003265-23-4	53
4			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	53
5			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\
Data File : VY006932.D
Acq On : 04 Dec 2021 11:46
Operator : SY/MD
Sample : M4884-02
Misc : 3.17g/10.0mL/MSVOA_Y/soil
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EW9H8

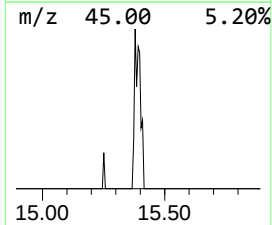
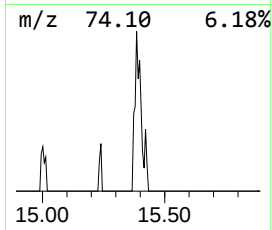
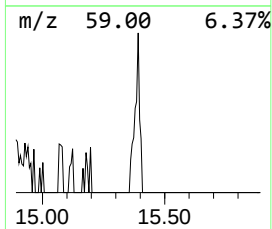
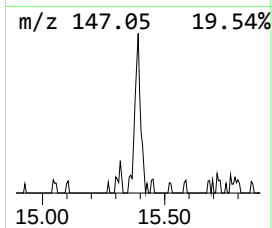
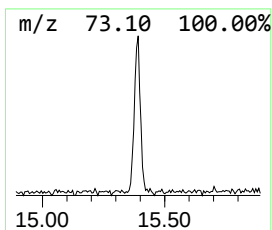
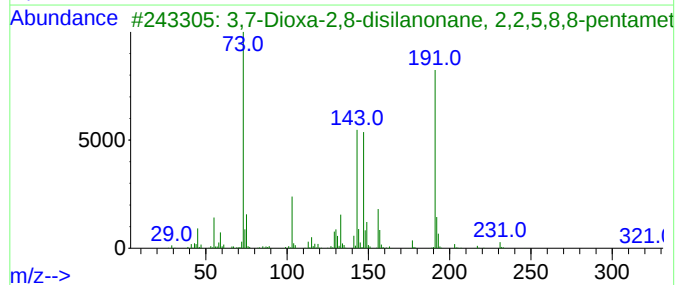
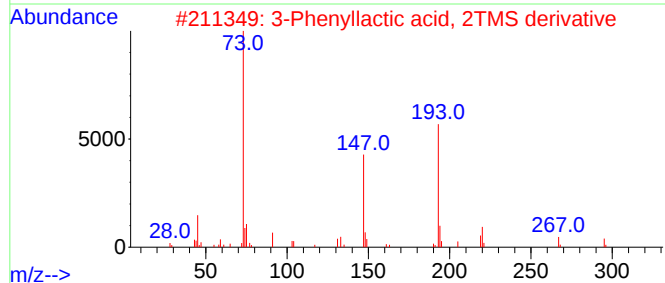
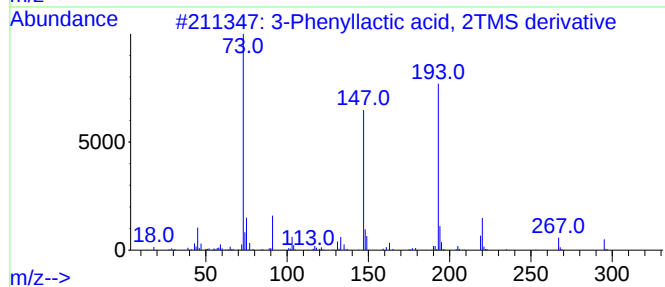
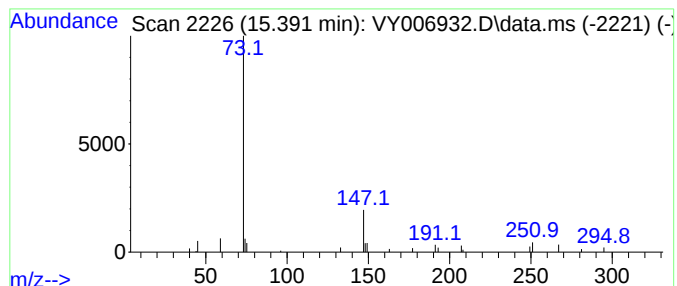
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYL120321SMA.M
Quant Title : VOC Analysis

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

Peak Number 14 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.391	4.73 ug/L	14525	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenyllactic acid, 2TMS deriva...	310	C15H26O3Si2	027750-45-4	42
2			3-Phenyllactic acid, 2TMS deriva...	310	C15H26O3Si2	027750-45-4	36
3			3,7-Dioxa-2,8-disilanonane, 2,2,...	336	C14H36O3Si3	017866-29-4	33
4			2-Hydroxybutyric acid, 2TMS	248	C10H24O3Si2	055133-93-2	9
5			Acetic acid, bis[(trimethylsilyl)...	308	C11H28O4Si3	055517-38-9	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\
Data File : VY006932.D
Acq On : 04 Dec 2021 11:46
Operator : SY/MD
Sample : M4884-02
Misc : 3.17g/10.0mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EW9H8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYL120321SMA.M
Quant Title : VOC Analysis

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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(DEL) Alkane: S...	13.276	1.4	ug/L	4182	3	13.428	76772	25.0
(DEL) Alkane: S...	13.483	1.5	ug/L	4689	3	13.428	76772	25.0
unknown-02	13.532	1.3	ug/L	3864	3	13.428	76772	25.0
4-(2-amino-1-hy...	13.965	13.1	ug/L	40111	3	13.428	76772	25.0
Benzene, 1,4-bi...	15.318	1.6	ug/L	4839	3	13.428	76772	25.0
unknown-03	15.391	4.7	ug/L	14525	3	13.428	76772	25.0