

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV120919\
 Data File : VV013962.D
 Acq On : 09 Dec 2019 14:56
 Operator : SY/MD
 Sample : K6146-06
 Misc : 5.0mL/MSVOA V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBBT6

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM120219WMA.M
 Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.209	31	37	40	rBV	40790	49711	2.32%	0.226%
2	1.318	63	71	82	rBV	290703	298119	13.94%	1.357%
3	1.556	140	145	160	rVB	287925	346126	16.19%	1.575%
4	2.122	312	321	333	rBV	596938	825412	38.60%	3.756%
5	2.460	418	426	436	rVV3	13326	22186	1.04%	0.101%
6	2.531	439	448	458	rVB	25605	41836	1.96%	0.190%
7	2.679	491	494	497	rBV4	1288	855	0.04%	0.004%
8	2.926	565	571	575	rVB7	1276	1498	0.07%	0.007%
9	3.177	647	649	655	rVB6	975	830	0.04%	0.004%
10	3.225	659	664	665	rBV3	1651	1137	0.05%	0.005%
11	3.482	742	744	750	rVB6	1684	1417	0.07%	0.006%
12	3.540	758	762	768	rVB6	1375	1329	0.06%	0.006%
13	3.576	768	773	775	rBV4	1453	1272	0.06%	0.006%
14	3.617	781	786	788	rBV5	1016	842	0.04%	0.004%
15	3.659	795	799	802	rVV6	1245	1092	0.05%	0.005%
16	3.730	818	821	827	rVB6	991	924	0.04%	0.004%
17	3.762	827	831	835	rBV6	1280	1193	0.06%	0.005%
18	3.807	844	845	851	rVB4	1704	1169	0.05%	0.005%
19	3.833	851	853	855	rBV3	1717	906	0.04%	0.004%
20	3.923	869	881	906	rBV	165226	423042	19.78%	1.925%
21	4.264	983	987	989	rVV4	1389	1175	0.05%	0.005%
22	4.392	1008	1027	1052	rBV	354996	840427	39.30%	3.824%
23	4.627	1096	1100	1103	rBV5	1471	856	0.04%	0.004%
24	4.836	1162	1165	1173	rVB9	1109	1253	0.06%	0.006%
25	4.945	1196	1199	1201	rBV4	1369	856	0.04%	0.004%
26	5.090	1225	1244	1275	rBV2	824445	2138361	100.00%	9.730%
27	5.322	1305	1316	1337	rBV2	56026	125382	5.86%	0.571%
28	5.460	1356	1359	1363	rBV4	1689	955	0.04%	0.004%
29	5.579	1386	1396	1397	rBV2	18236	18975	0.89%	0.086%
30	5.659	1409	1421	1437	rBV	730197	1426610	66.72%	6.491%
31	6.113	1549	1562	1587	rVB	443221	894487	41.83%	4.070%
32	6.405	1637	1653	1664	rVB4	20084	57716	2.70%	0.263%
33	6.498	1671	1682	1707	rBV	1230236	2130126	99.61%	9.693%
34	7.022	1833	1845	1861	rBV	247330	437464	20.46%	1.991%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM120219WMA.M
 Title : VOC Analysis

35	7.222	1897	1907	1920	rBV	177048	297329	13.90%	1.353%
36	7.354	1935	1948	1969	rBV	917155	1607522	75.18%	7.315%
37	7.656	2030	2042	2063	rBV2	201491	362041	16.93%	1.647%
38	7.961	2125	2137	2154	rBV	313298	522423	24.43%	2.377%
39	8.125	2178	2188	2210	rBV2	634134	1123704	52.55%	5.113%
40	8.228	2210	2220	2234	rVB	571545	874158	40.88%	3.978%
41	8.338	2250	2254	2266	rVB3	20318	32635	1.53%	0.148%
42	8.530	2309	2314	2316	rBV5	1285	1048	0.05%	0.005%
43	8.717	2361	2372	2401	rBV	679457	1065177	49.81%	4.847%
44	8.891	2412	2426	2440	rBV	1101586	1786036	83.52%	8.127%
45	9.180	2510	2516	2524	rVB4	9651	14312	0.67%	0.065%
46	9.325	2557	2561	2566	rBV7	1419	1725	0.08%	0.008%
47	9.424	2584	2592	2600	rBV2	15569	26112	1.22%	0.119%
48	9.582	2632	2641	2656	rBV	78571	126799	5.93%	0.577%
49	9.916	2742	2745	2748	rBV3	1246	927	0.04%	0.004%
50	9.952	2751	2756	2758	rBV6	1753	1542	0.07%	0.007%
51	10.051	2784	2787	2791	rVB5	2003	1312	0.06%	0.006%
52	10.257	2833	2851	2873	rBV	688772	1079191	50.47%	4.911%
53	10.337	2873	2876	2878	rBV3	1565	1209	0.06%	0.006%
54	10.392	2890	2893	2894	rBV3	1577	903	0.04%	0.004%
55	10.460	2910	2914	2917	rVB4	1699	1510	0.07%	0.007%
56	10.482	2917	2921	2925	rBV5	2001	2173	0.10%	0.010%
57	10.505	2925	2928	2933	rVB7	1534	1575	0.07%	0.007%
58	10.527	2933	2935	2939	rVB4	1368	906	0.04%	0.004%
59	10.569	2946	2948	2950	rBV2	1939	1198	0.06%	0.005%
60	10.691	2982	2986	2990	rBV6	2112	1861	0.09%	0.008%
61	10.743	2999	3002	3005	rVV5	1589	1104	0.05%	0.005%
62	10.887	3044	3047	3050	rBV4	1634	1288	0.06%	0.006%
63	10.958	3063	3069	3074	rBV8	3132	3863	0.18%	0.018%
64	11.289	3159	3172	3190	rBV	981749	1489452	69.65%	6.777%
65	11.434	3214	3217	3221	rVB4	2123	1276	0.06%	0.006%
66	11.540	3248	3250	3256	rVB6	1633	864	0.04%	0.004%
67	11.665	3277	3289	3306	rBV	905637	1406587	65.78%	6.400%
68	11.842	3342	3344	3350	rVB6	1453	934	0.04%	0.004%
69	11.878	3353	3355	3359	rBV5	1331	1072	0.05%	0.005%
70	11.968	3380	3383	3387	rBV3	1084	913	0.04%	0.004%
71	12.122	3427	3431	3433	rBV2	1157	903	0.04%	0.004%

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72	12.177	3446	3448	3451	rBV3	1886	1266	0.06%	0.006%
73	12.257	3470	3473	3476	rBV3	1439	1046	0.05%	0.005%
74	12.402	3509	3518	3522	rBV3	2405	3372	0.16%	0.015%
75	12.427	3522	3526	3532	rBV8	1646	2039	0.10%	0.009%
76	12.501	3546	3549	3553	rVV3	1493	986	0.05%	0.004%
77	12.591	3575	3577	3579	rBV2	1396	901	0.04%	0.004%
78	12.688	3602	3607	3608	rBV4	1274	868	0.04%	0.004%
79	12.707	3610	3613	3617	rVV6	2120	1795	0.08%	0.008%
80	12.839	3649	3654	3658	rBV6	1516	1479	0.07%	0.007%
81	12.861	3658	3661	3667	rVB6	1279	1297	0.06%	0.006%
82	12.890	3667	3670	3674	rBV5	1885	1527	0.07%	0.007%
83	13.058	3718	3722	3726	rVB7	2128	2156	0.10%	0.010%
84	13.202	3763	3767	3770	rBV5	1626	1430	0.07%	0.007%
85	13.247	3779	3781	3784	rBV4	1222	836	0.04%	0.004%
86	13.421	3831	3835	3838	rBV5	1252	1188	0.06%	0.005%
87	13.472	3848	3851	3854	rBV3	1408	872	0.04%	0.004%
88	13.492	3854	3857	3860	rVV4	1582	1092	0.05%	0.005%
89	13.559	3876	3878	3881	rVV3	1220	899	0.04%	0.004%
90	13.617	3892	3896	3899	rBV5	1540	1196	0.06%	0.005%
91	13.778	3943	3946	3948	rBV3	1472	1212	0.06%	0.006%
92	14.061	4029	4034	4035	rBV4	1157	999	0.05%	0.005%
93	14.164	4064	4066	4069	rVV4	1297	893	0.04%	0.004%
94	14.257	4092	4095	4097	rBV4	1873	1177	0.06%	0.005%
95	14.363	4124	4128	4131	rVB5	1630	1240	0.06%	0.006%
96	14.533	4177	4181	4183	rBV4	1199	868	0.04%	0.004%
97	14.594	4196	4200	4202	rBV4	1319	913	0.04%	0.004%
98	14.849	4276	4279	4281	rBV3	1508	915	0.04%	0.004%
99	15.025	4331	4334	4338	rVB5	2474	1600	0.07%	0.007%
100	15.141	4366	4370	4372	rBV3	2279	1784	0.08%	0.008%

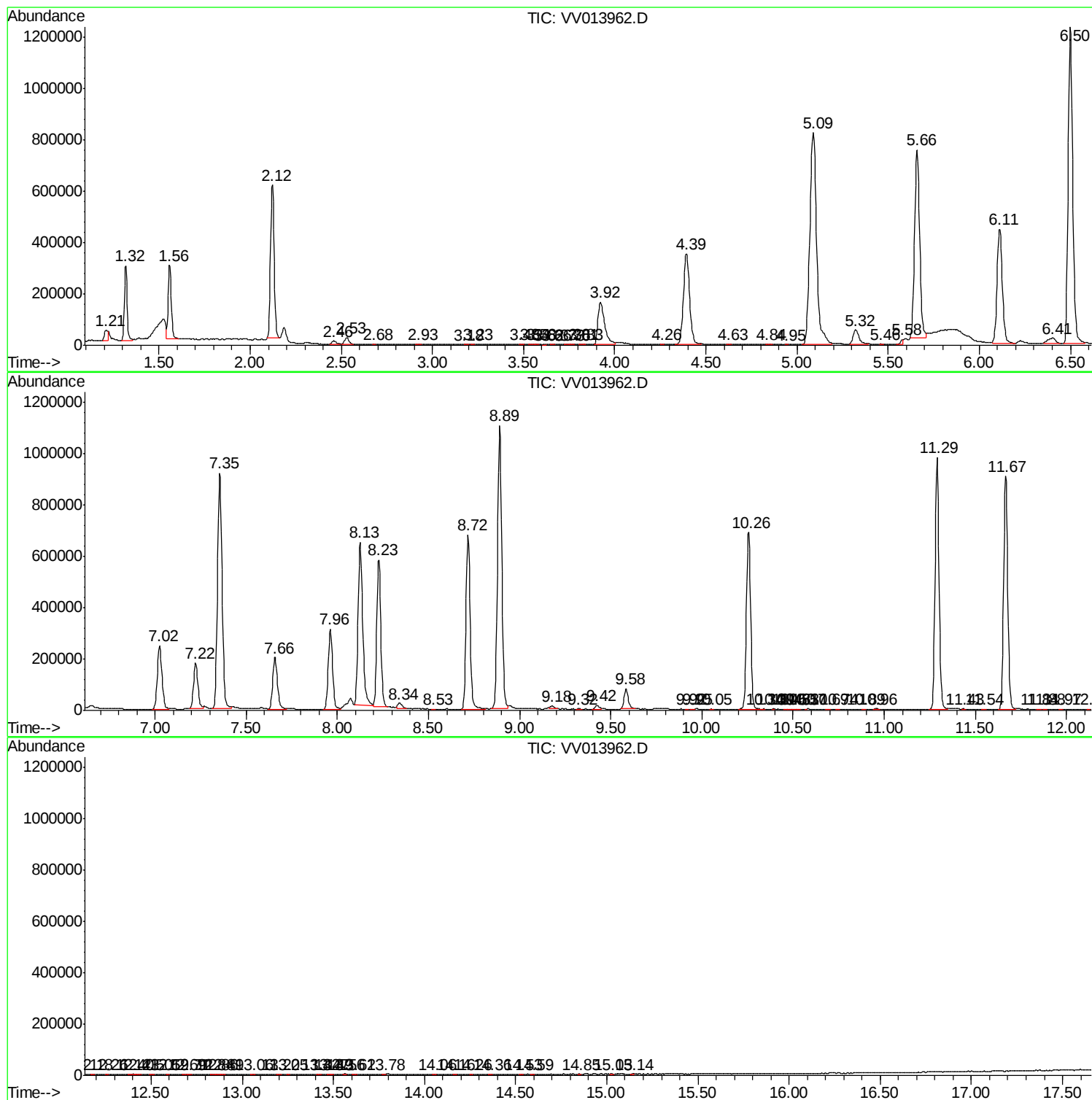
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM120219WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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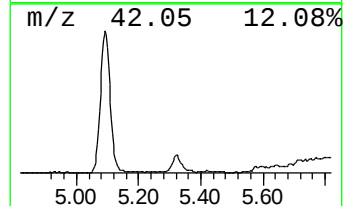
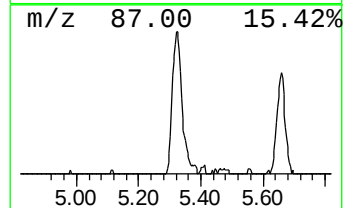
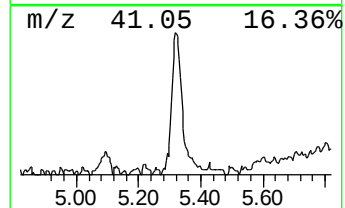
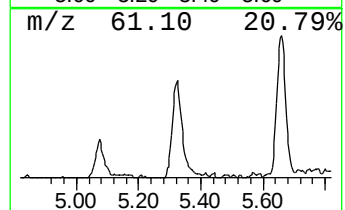
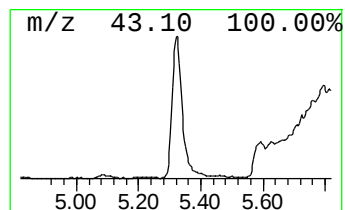
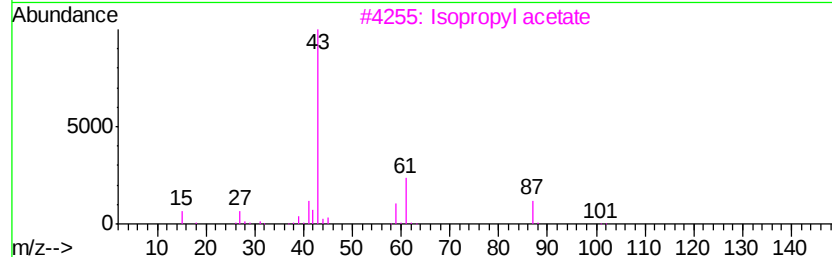
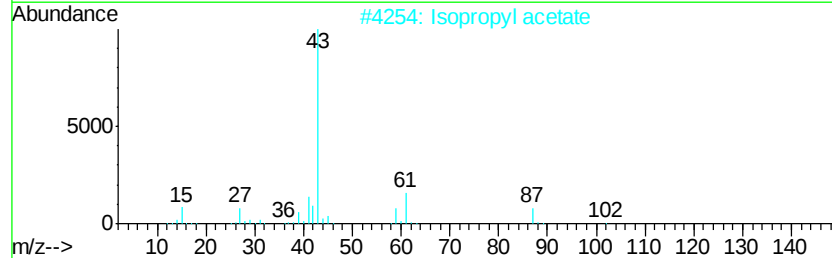
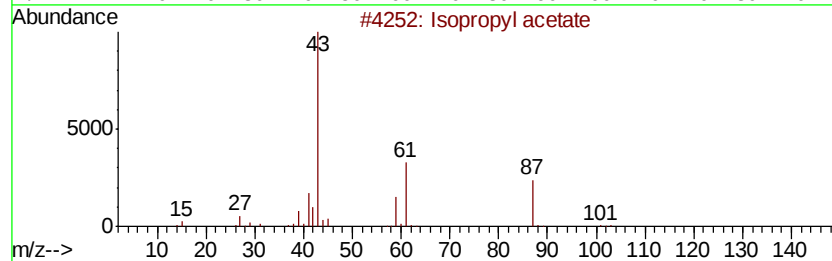
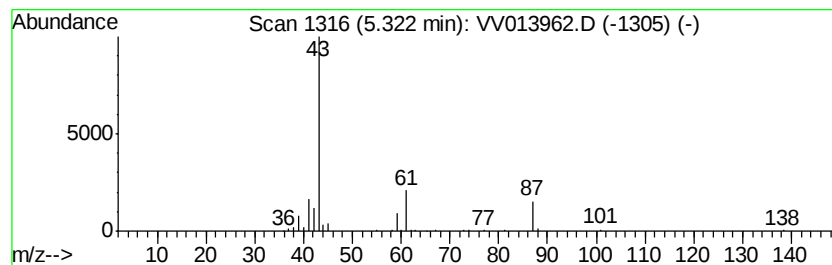
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM120219WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Isopropyl acetate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	4.39 ug/L	125382	1,4-Difluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl acetate	102	C5H10O2	000108-21-4	72
2		Isopropyl acetate	102	C5H10O2	000108-21-4	56
3		Isopropyl acetate	102	C5H10O2	000108-21-4	56
4		Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	28
5		Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	28



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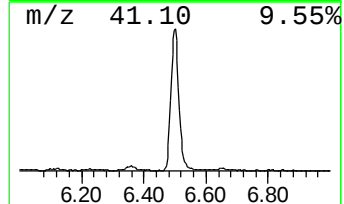
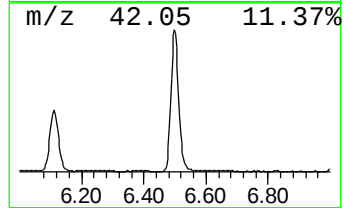
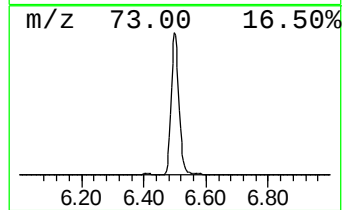
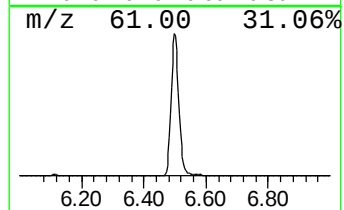
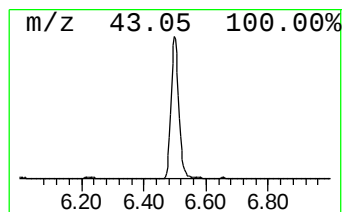
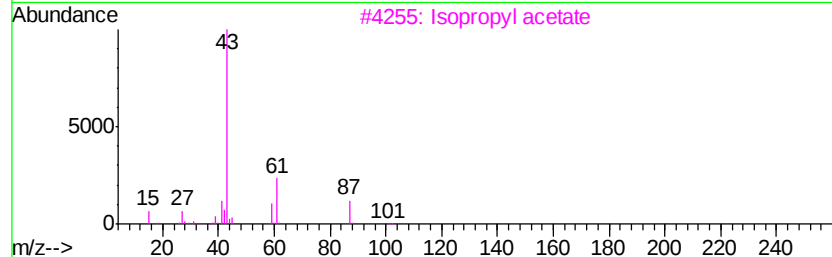
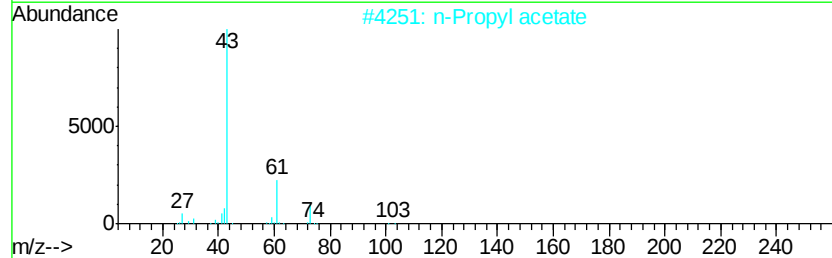
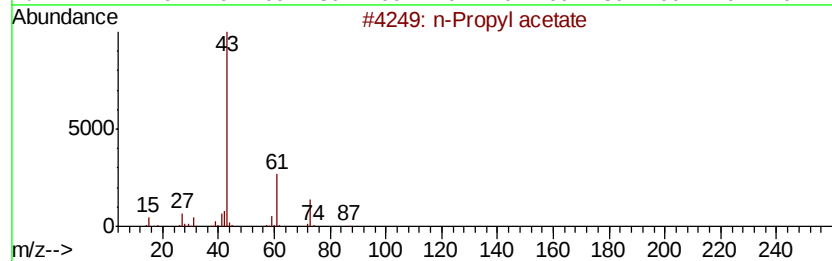
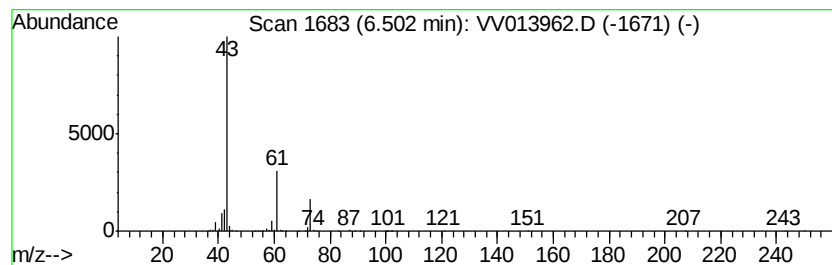
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	74.66 ug/L	2130130	1,4-Difluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2		n-Propyl acetate	102	C5H10O2	000109-60-4	83
3		Isopropyl acetate	102	C5H10O2	000108-21-4	36
4		n-Propyl acetate	102	C5H10O2	000109-60-4	9
5		Isobutyl acetate	116	C6H12O2	000110-19-0	9



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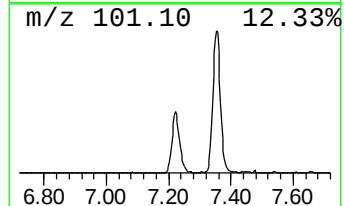
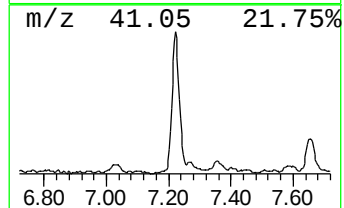
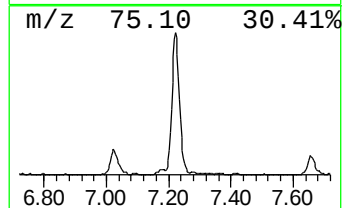
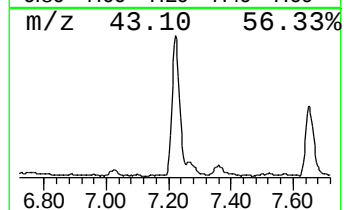
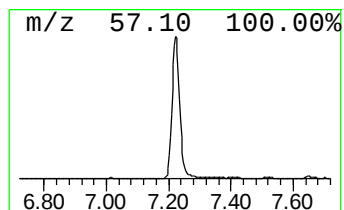
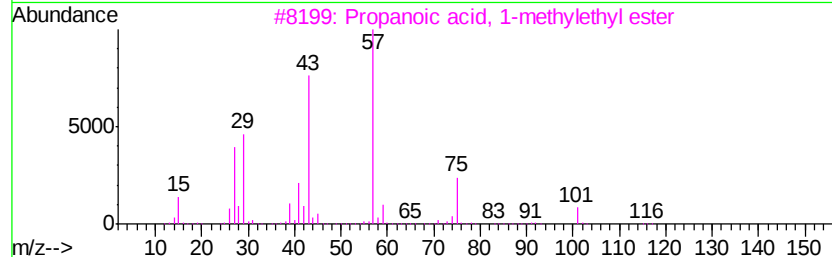
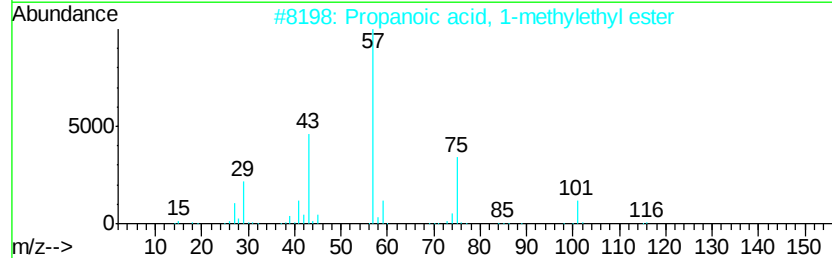
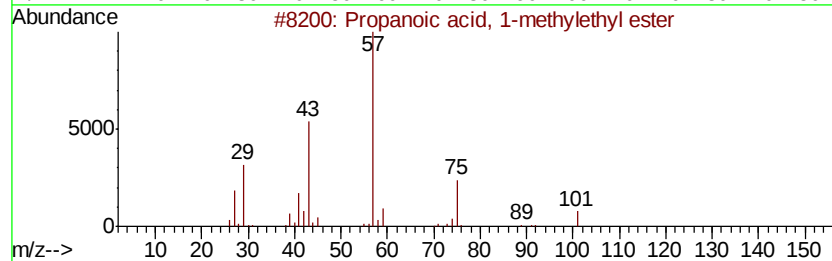
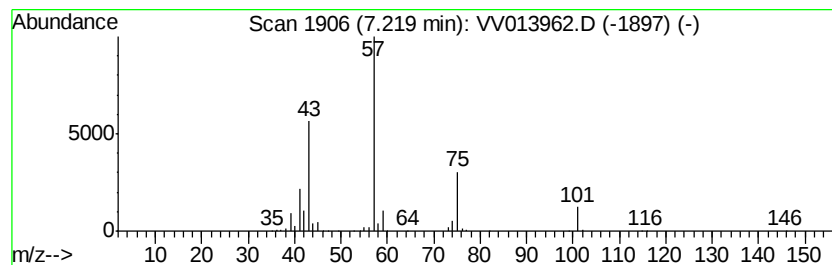
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Quant Title : VOC Analysis

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TIC Integration Parameters: LSCINT.P

Peak Number 4 Propanoic acid, 1-methyleth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	10.42 ug/L	297329	1,4-Difluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
2		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	72
3		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	64
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	38
5		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV120919\
Data File : VV013962.D
Acq On : 09 Dec 2019 14:56
Operator : SY/MD
Sample : K6146-06
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBBT6

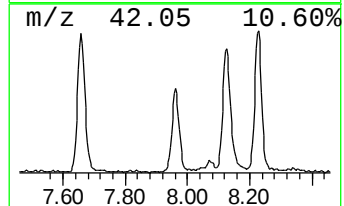
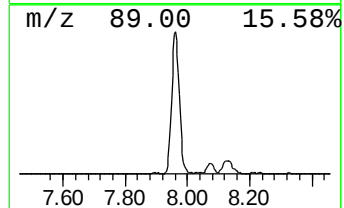
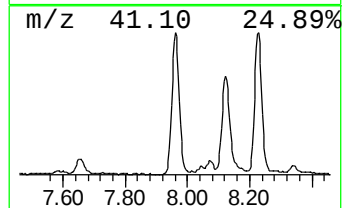
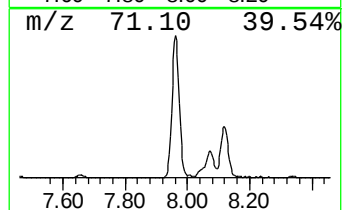
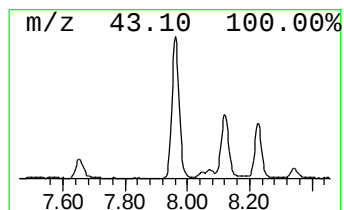
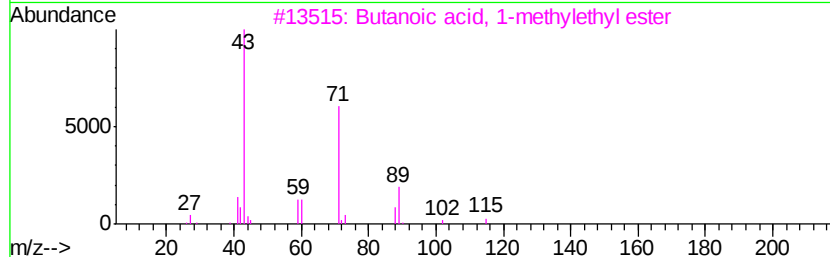
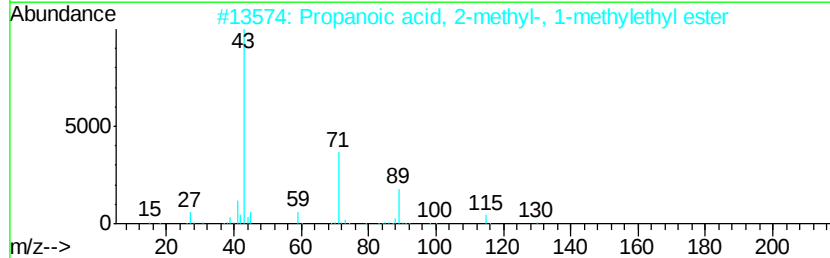
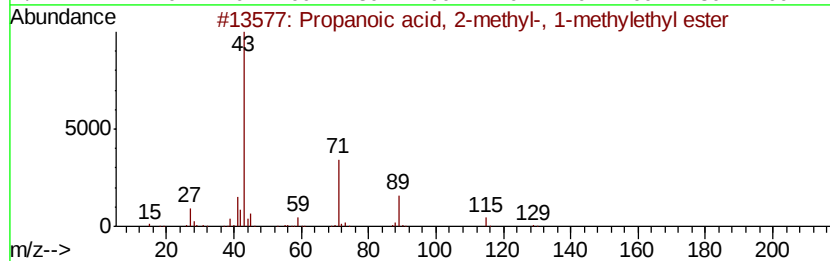
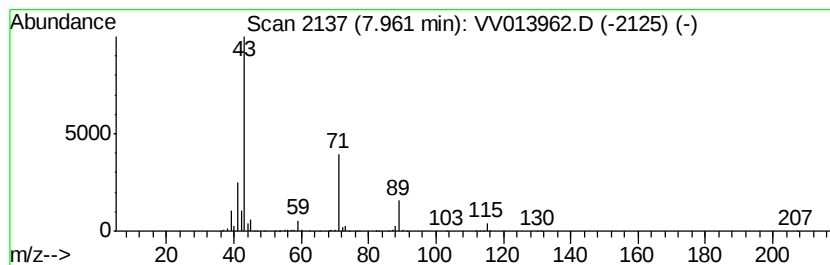
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Propanoic acid, 2-methyl-, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.96	14.63 ug/L	522423	Chlorobenzene-d5	8.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
2		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
3		Butanoic acid, 1-methyl-ethyl ester	130	C7H14O2	000638-11-9	72
4		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV120919\
Data File : VV013962.D
Acq On : 09 Dec 2019 14:56
Operator : SY/MD
Sample : K6146-06
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBBT6

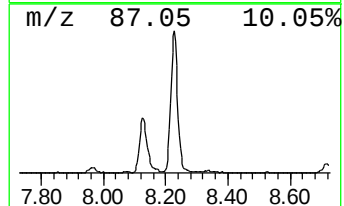
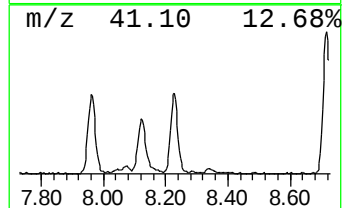
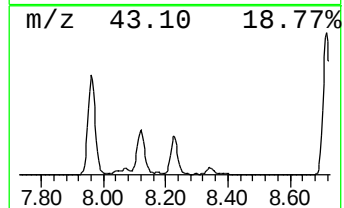
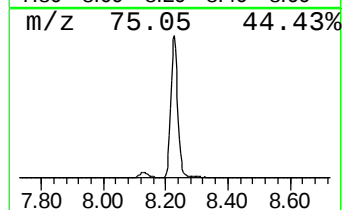
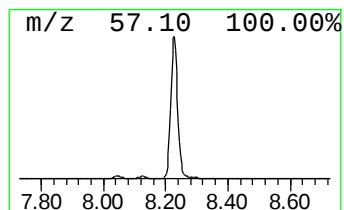
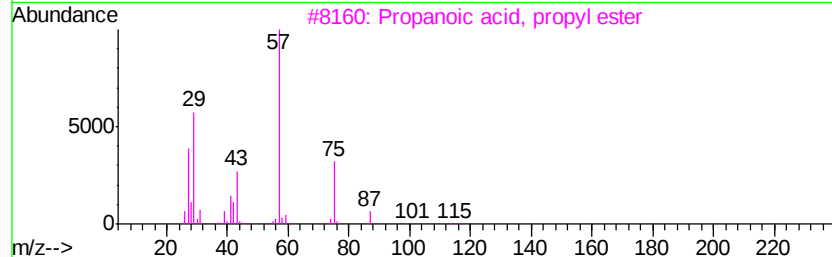
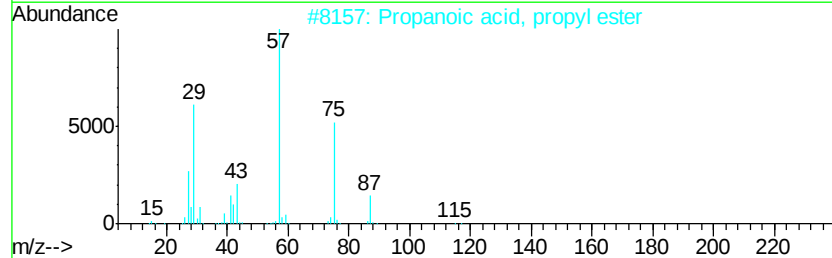
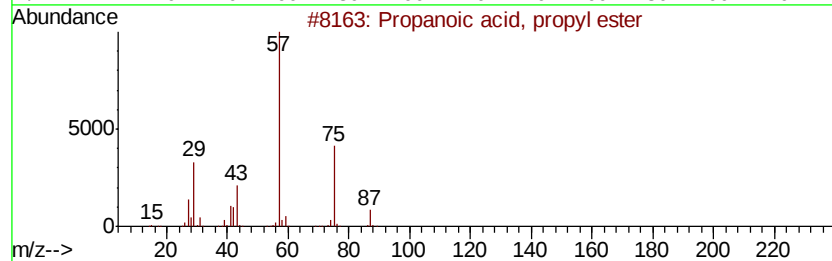
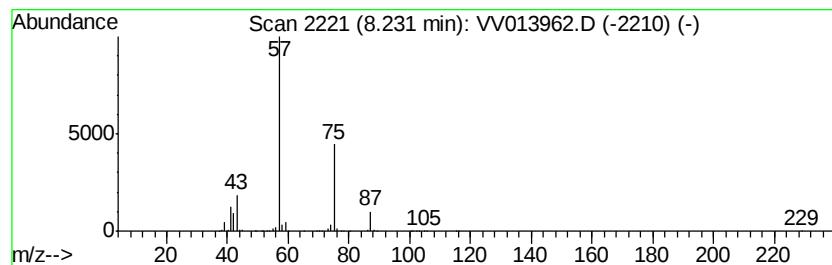
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.23	24.47 ug/L	874158	Chlorobenzene-d5	8.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
3		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
5		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	33



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV120919\
Data File : VV013962.D
Acq On : 09 Dec 2019 14:56
Operator : SY/MD
Sample : K6146-06
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBBT6

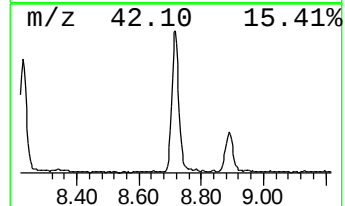
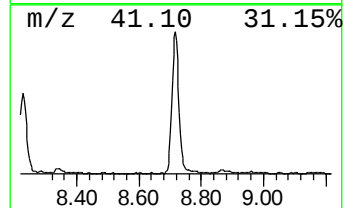
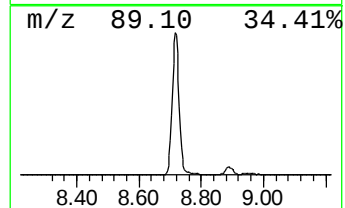
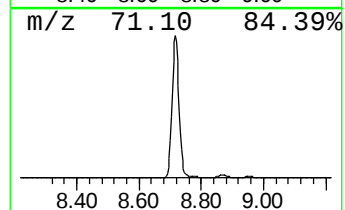
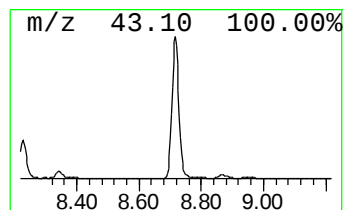
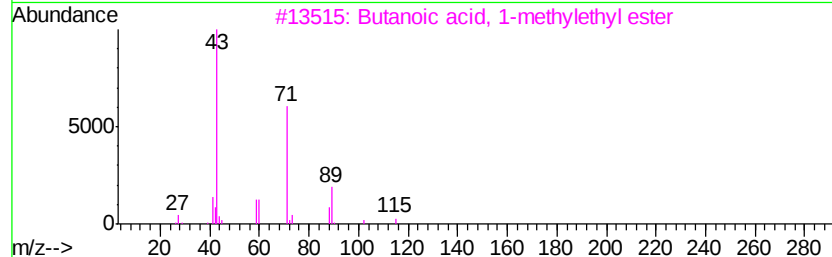
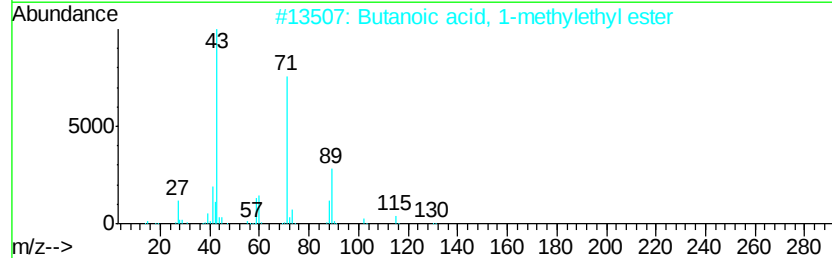
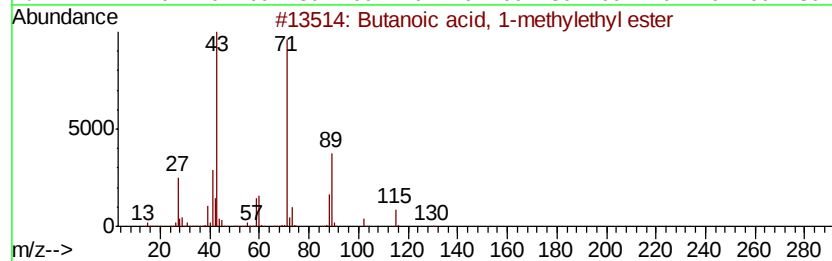
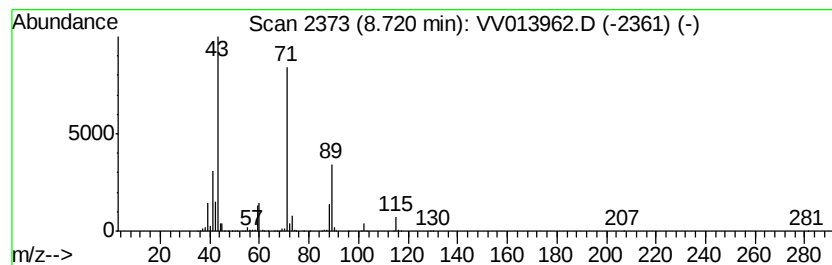
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Butanoic acid, 1-methylethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	29.82 ug/L	1065180	Chlorobenzene-d5	8.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	91
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	83
4		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	72
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV120919\
Data File : VV013962.D
Acq On : 09 Dec 2019 14:56
Operator : SY/MD
Sample : K6146-06
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBBT6

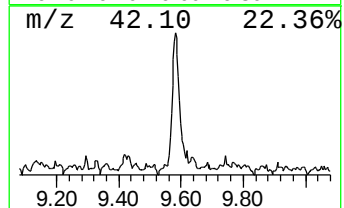
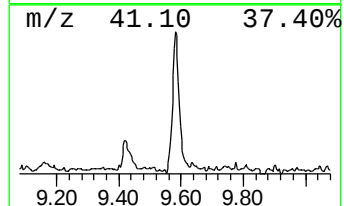
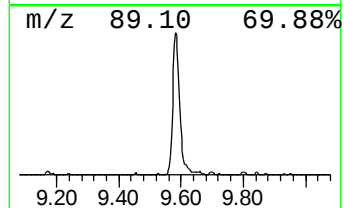
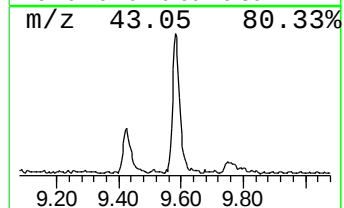
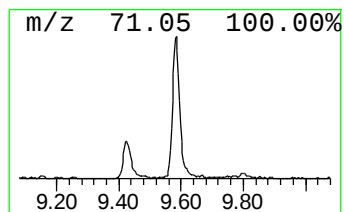
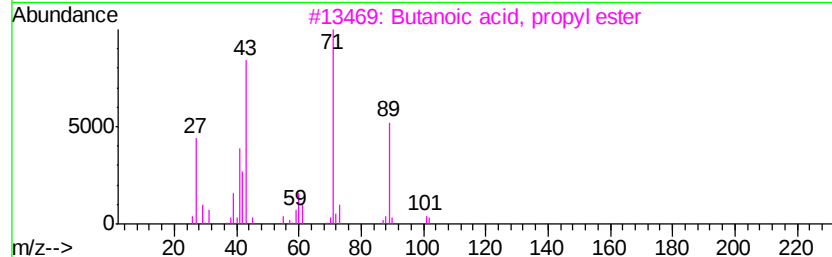
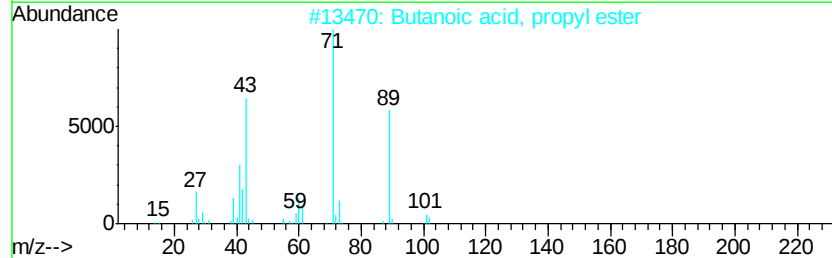
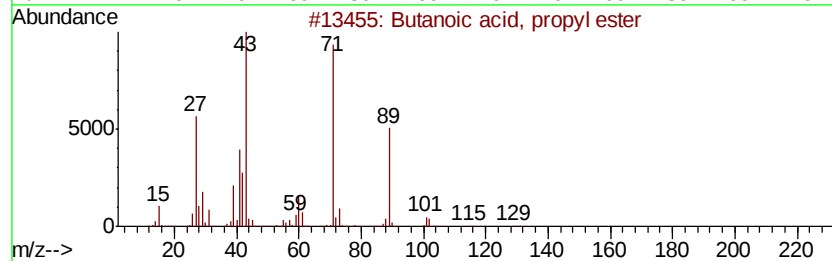
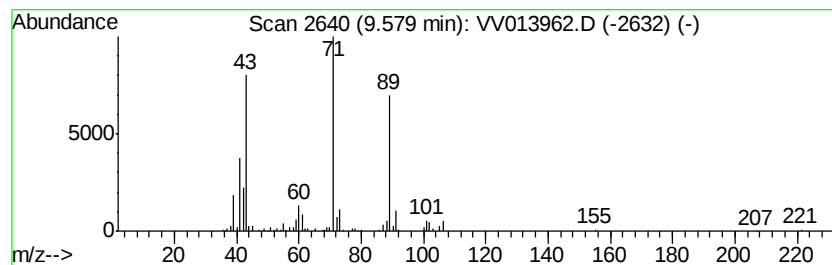
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM120219WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Butanoic acid, propyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.58	3.55 ug/L	126799	Chlorobenzene-d5	8.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	86
2		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	86
3		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	86
4		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	86
5		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	72



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV120919\
Data File : VV013962.D
Acq On : 09 Dec 2019 14:56
Operator : SY/MD
Sample : K6146-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBBT6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM120219WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl acetate	5.32	4.4	ug/L	125382	1	5.66	1426610	50.0
n-Propyl acetate	6.50	74.7	ug/L	2130130	1	5.66	1426610	50.0
Propanoic acid, 1...	7.22	10.4	ug/L	297329	1	5.66	1426610	50.0
Propanoic acid, 2...	7.96	14.6	ug/L	522423	2	8.89	1786040	50.0
Propanoic acid, p...	8.23	24.5	ug/L	874158	2	8.89	1786040	50.0
Butanoic acid, 1-...	8.72	29.8	ug/L	1065180	2	8.89	1786040	50.0
Butanoic acid, pr...	9.58	3.5	ug/L	126799	2	8.89	1786040	50.0