

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
 Data File : VU034754.D
 Acq On : 23 Sep 2019 23:40
 Operator : JC/SP
 Sample : K4932-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFDJ9

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_U\METHOD\SOMULM091919WMA.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	1.177	21	27	33	rBV	10955	11133	0.29%	0.036%
2	1.334	69	76	77	rBV6	5357	6156	0.16%	0.020%
3	1.392	85	94	107	rBV	342922	376258	9.80%	1.210%
4	1.540	133	140	151	rVB3	46776	64887	1.69%	0.209%
5	1.672	173	181	195	rVB	276084	353183	9.20%	1.136%
6	2.257	352	363	371	rBV	515340	789423	20.56%	2.539%
7	2.305	371	378	395	rVB	233452	375827	9.79%	1.209%
8	2.498	436	438	443	rVB4	2020	1327	0.03%	0.004%
9	2.601	465	470	482	rVB	5146	8095	0.21%	0.026%
10	2.685	482	496	514	rBV	151075	273989	7.14%	0.881%
11	2.923	566	570	575	rVV3	2143	2956	0.08%	0.010%
12	2.948	576	578	581	rVB4	3077	1778	0.05%	0.006%
13	2.997	590	593	595	rBV3	2339	1736	0.05%	0.006%
14	3.084	618	620	629	rVB7	2011	2154	0.06%	0.007%
15	3.154	634	642	645	rBV6	1329	1480	0.04%	0.005%
16	3.579	769	774	775	rBV5	1558	1304	0.03%	0.004%
17	3.855	856	860	863	rBV3	1514	1394	0.04%	0.004%
18	3.952	886	890	894	rBV4	1605	1738	0.05%	0.006%
19	4.151	938	952	972	rBV	281044	727701	18.95%	2.340%
20	4.624	1077	1099	1121	rBV	410639	984873	25.65%	3.167%
21	4.971	1202	1207	1213	rVB7	2573	3155	0.08%	0.010%
22	5.315	1290	1314	1351	rBV2	852590	2562320	66.73%	8.241%
23	5.530	1368	1381	1407	rVB	100987	222624	5.80%	0.716%
24	5.862	1471	1484	1506	rBV	685239	1401690	36.50%	4.508%
25	6.308	1607	1623	1640	rBV	597672	1202105	31.31%	3.866%
26	6.405	1647	1653	1670	rVV3	12429	30129	0.78%	0.097%
27	6.479	1674	1676	1679	rVB3	3318	1710	0.04%	0.005%
28	6.524	1686	1690	1693	rBV3	2480	2499	0.07%	0.008%
29	6.562	1694	1702	1705	rBV	17408	24086	0.63%	0.077%
30	6.685	1727	1740	1767	rBV	2111549	3839977	100.00%	12.349%
31	7.206	1890	1902	1926	rBV	343048	641555	16.71%	2.063%
32	7.398	1950	1962	1971	rBV	265923	459726	11.97%	1.478%
33	7.543	1993	2007	2023	rBV	1166862	2041318	53.16%	6.565%
34	7.617	2024	2030	2044	rVB2	30568	59675	1.55%	0.192%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
 Title : VOC Analysis

35	7.755	2068	2073	2084	rVB9	5870	9969	0.26%	0.032%
36	7.829	2084	2096	2113	rBV2	297918	550128	14.33%	1.769%
37	8.135	2177	2191	2205	rBV	484619	820063	21.36%	2.637%
38	8.212	2207	2215	2218	rVV3	33373	50037	1.30%	0.161%
39	8.241	2219	2224	2229	rVV2	61891	90340	2.35%	0.291%
40	8.289	2229	2239	2262	rVV	1099494	1983623	51.66%	6.379%
41	8.395	2262	2272	2299	rVV	789538	1299457	33.84%	4.179%
42	8.508	2300	2307	2321	rVB2	33247	60368	1.57%	0.194%
43	8.813	2398	2402	2404	rBV4	2312	1456	0.04%	0.005%
44	8.884	2411	2424	2446	rBV	1019237	1614581	42.05%	5.193%
45	9.067	2462	2481	2507	rVV	1036697	1784159	46.46%	5.738%
46	9.231	2525	2532	2540	rVB5	9672	16653	0.43%	0.054%
47	9.267	2540	2543	2545	rBV4	2295	1553	0.04%	0.005%
48	9.324	2552	2561	2562	rBV7	3985	4773	0.12%	0.015%
49	9.353	2562	2570	2585	rVB2	26448	49978	1.30%	0.161%
50	9.469	2601	2606	2608	rBV5	1922	1597	0.04%	0.005%
51	9.482	2608	2610	2616	rVB5	2057	1774	0.05%	0.006%
52	9.585	2634	2642	2660	rVB	28259	52247	1.36%	0.168%
53	9.749	2682	2693	2707	rBV3	118273	212799	5.54%	0.684%
54	9.919	2733	2746	2748	rBV8	7557	11921	0.31%	0.038%
55	9.967	2757	2761	2768	rVB8	4040	4657	0.12%	0.015%
56	10.154	2812	2819	2827	rVB8	4099	5720	0.15%	0.018%
57	10.202	2827	2834	2835	rBV6	1762	1722	0.04%	0.006%
58	10.302	2860	2865	2868	rBV7	2161	2202	0.06%	0.007%
59	10.408	2882	2898	2924	rBV	775265	1269249	33.05%	4.082%
60	10.543	2936	2940	2941	rBV3	3033	1825	0.05%	0.006%
61	10.569	2943	2948	2950	rBV6	2660	2534	0.07%	0.008%
62	10.662	2964	2977	2981	rBV4	8334	14725	0.38%	0.047%
63	10.755	2999	3006	3012	rVB8	6056	7747	0.20%	0.025%
64	10.797	3012	3019	3022	rBV8	3467	4361	0.11%	0.014%
65	10.903	3043	3052	3057	rBV9	4966	8644	0.23%	0.028%
66	10.955	3065	3068	3080	rVB9	5134	8554	0.22%	0.028%
67	11.025	3087	3090	3094	rVB5	2501	2264	0.06%	0.007%
68	11.057	3094	3100	3101	rBV6	2375	1853	0.05%	0.006%
69	11.128	3112	3122	3137	rVB4	23593	43084	1.12%	0.139%
70	11.234	3150	3155	3156	rBV5	3553	2690	0.07%	0.009%
71	11.250	3156	3160	3162	rBV4	3868	3317	0.09%	0.011%

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72	11.318	3171	3181	3201	rBV7	60062	174348	4.54%	0.561%
73	11.459	3213	3225	3246	rBV	1015261	1622994	42.27%	5.220%
74	11.736	3300	3311	3324	rBV	146662	329870	8.59%	1.061%
75	11.836	3331	3342	3365	rVB	1113676	1993909	51.93%	6.412%
76	12.945	3681	3687	3693	rBV8	6936	11397	0.30%	0.037%
77	13.077	3720	3728	3730	rBV9	10102	13569	0.35%	0.044%
78	13.170	3753	3757	3761	rBV6	2989	3519	0.09%	0.011%
79	13.273	3783	3789	3794	rBV10	6613	10533	0.27%	0.034%
80	13.373	3814	3820	3822	rBV7	5785	6647	0.17%	0.021%
81	13.601	3881	3891	3904	rBV7	23781	58160	1.51%	0.187%
82	13.723	3921	3929	3942	rVB2	79544	130856	3.41%	0.421%
83	13.839	3961	3965	3968	rBV5	8146	8478	0.22%	0.027%
84	14.437	4148	4151	4153	rBV3	2543	1737	0.05%	0.006%
85	14.514	4170	4175	4188	rVB3	9368	15604	0.41%	0.050%
86	14.562	4188	4190	4195	rVB6	3453	2633	0.07%	0.008%
87	14.594	4196	4200	4202	rBV5	3065	2205	0.06%	0.007%
88	14.620	4202	4208	4215	rBV10	8094	11762	0.31%	0.038%
89	14.858	4273	4282	4296	rBV	59467	100107	2.61%	0.322%
90	15.009	4327	4329	4331	rBV3	2609	1533	0.04%	0.005%
91	15.041	4331	4339	4357	rVB3	38491	70921	1.85%	0.228%
92	15.266	4407	4409	4411	rBV3	2388	1367	0.04%	0.004%
93	15.324	4424	4427	4428	rBV3	2983	1752	0.05%	0.006%
94	15.414	4452	4455	4458	rBV6	2889	1861	0.05%	0.006%
95	15.517	4485	4487	4490	rBV4	2095	1487	0.04%	0.005%
96	15.784	4565	4570	4571	rBV5	3762	3151	0.08%	0.010%
97	16.022	4641	4644	4645	rBV3	4392	2538	0.07%	0.008%
98	16.041	4645	4650	4657	rVV9	9780	14960	0.39%	0.048%
99	16.167	4686	4689	4691	rBV4	2647	1962	0.05%	0.006%
100	17.334	5044	5052	5056	rBV4	6688	11749	0.31%	0.038%

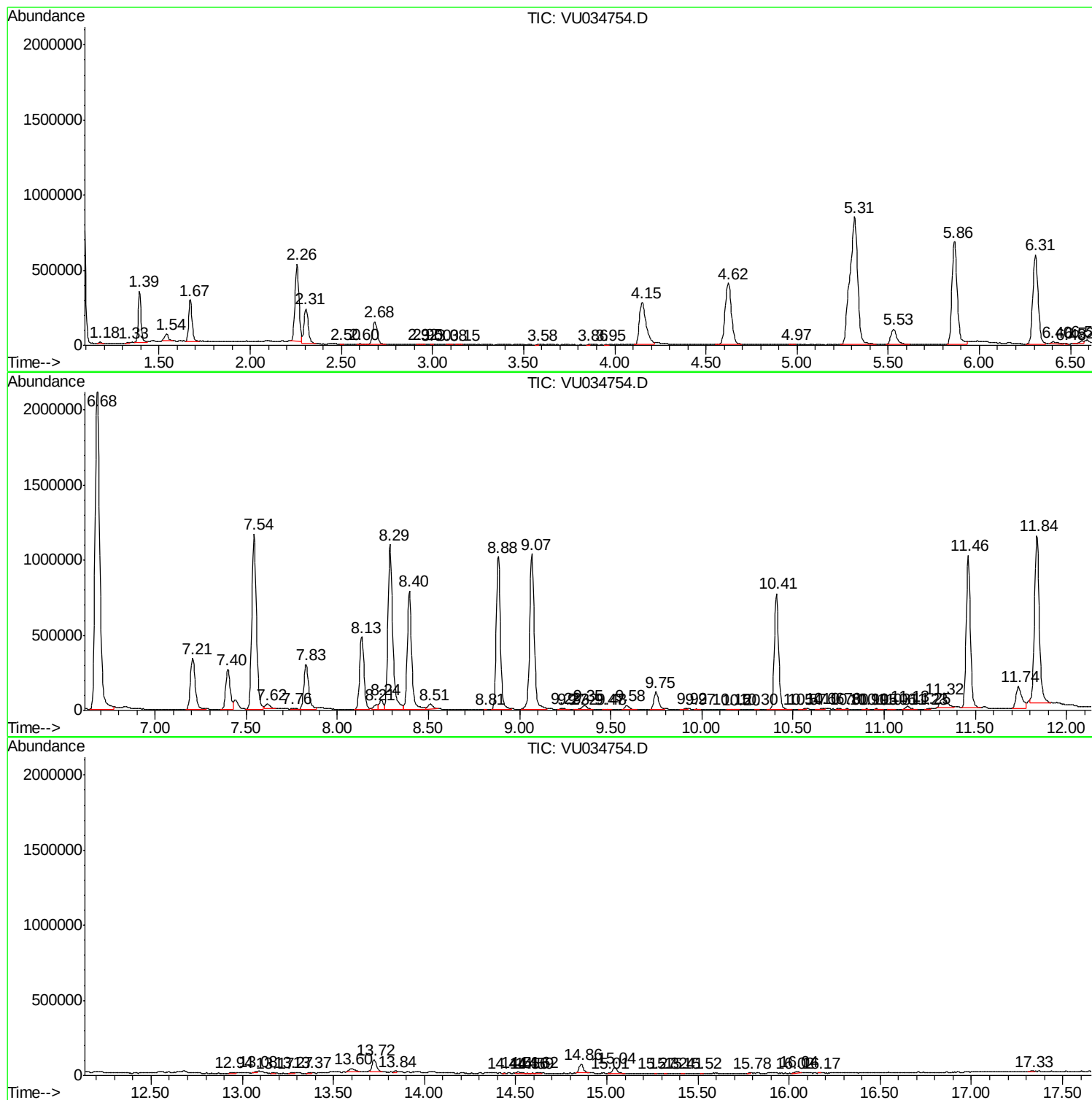
Sum of corrected areas: 31094194

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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

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



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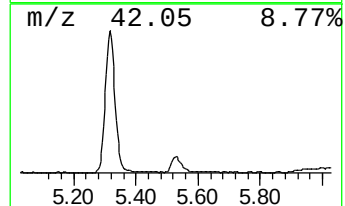
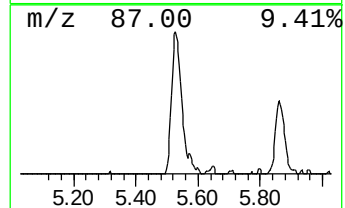
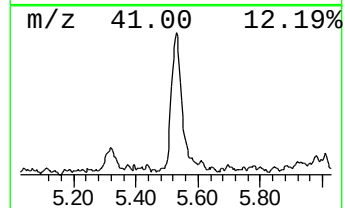
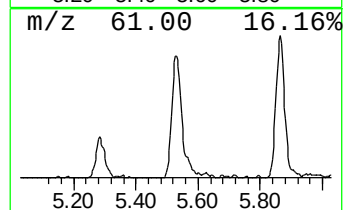
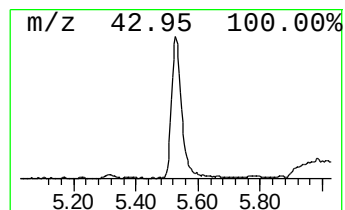
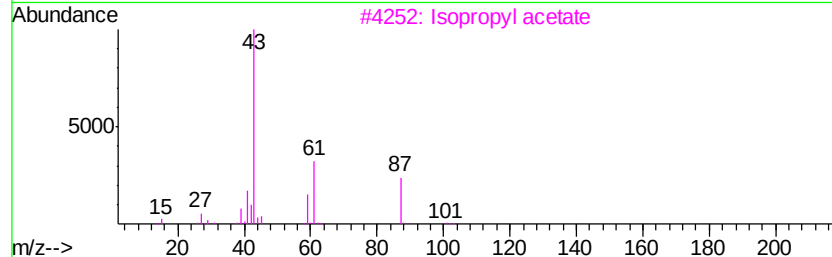
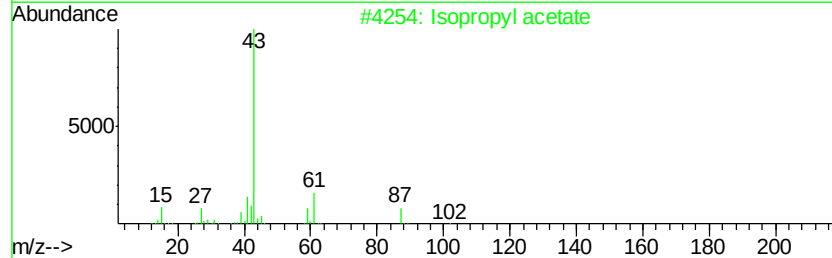
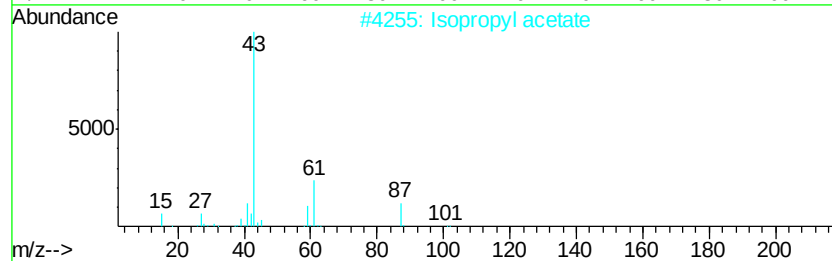
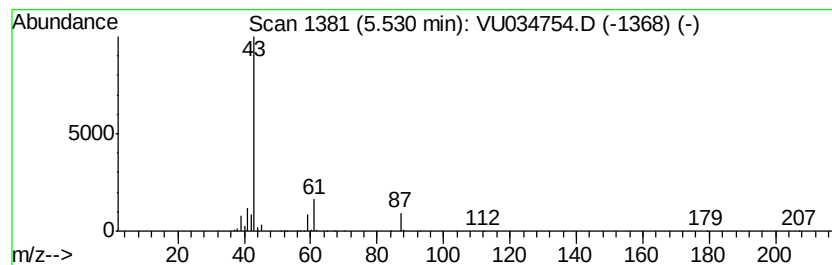
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

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.
5.53	7.94 ug/L	222624	1,4-Difluorobenzene	5.86

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	64
3		Isopropyl acetate	102	C5H10O2	000108-21-4	45
4		Isopropyl acetate	102	C5H10O2	000108-21-4	42
5		Acetaldehyde, O-ethyloxime-	87	C4H9NO	1000288-34-6	5



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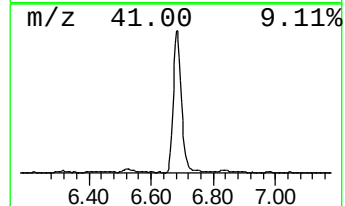
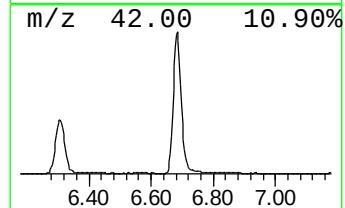
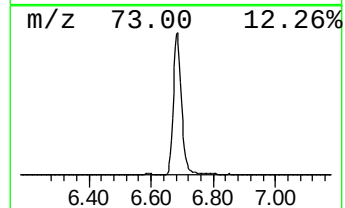
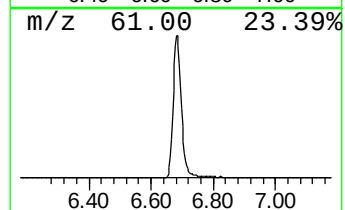
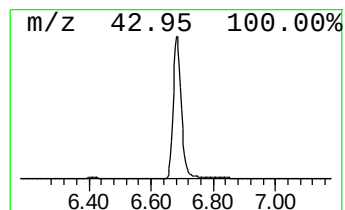
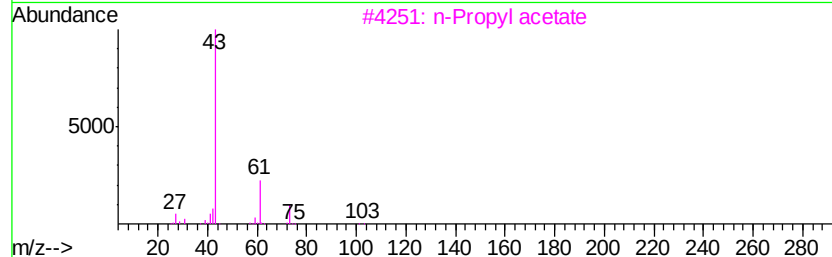
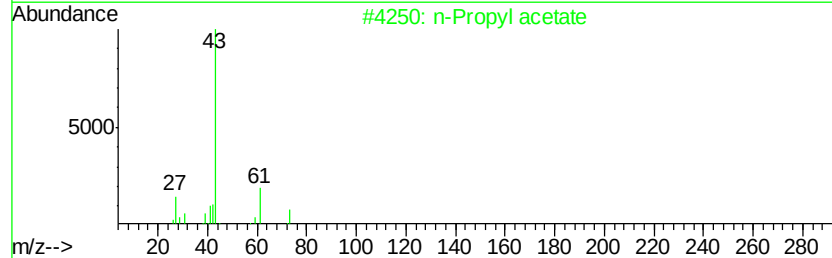
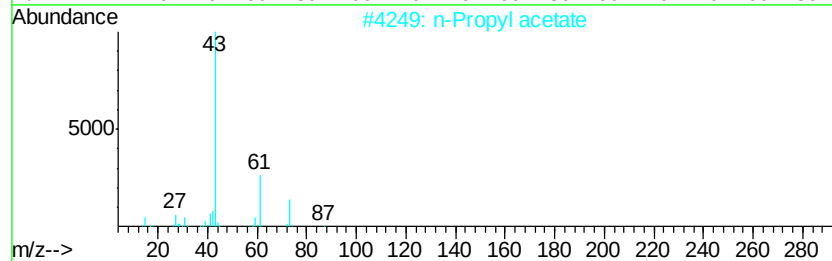
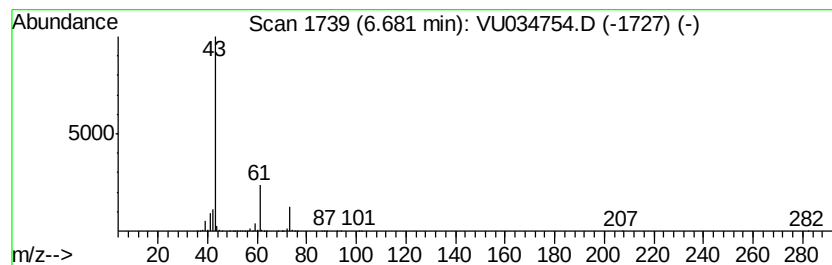
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Quant Title : VOC Analysis

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.
6.68	136.98 ug/L	3839980	1,4-Difluorobenzene	5.86

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	78
3		n-Propyl acetate	102	C5H10O2	000109-60-4	78
4		Isobutylamine	73	C4H11N	000078-81-9	7
5		Isobutylamine	73	C4H11N	000078-81-9	5



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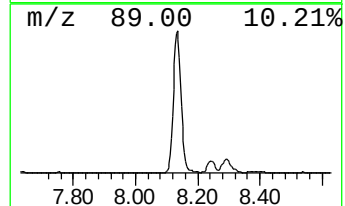
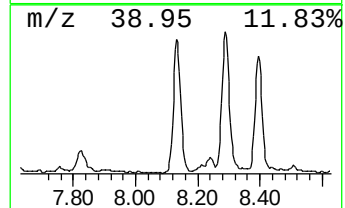
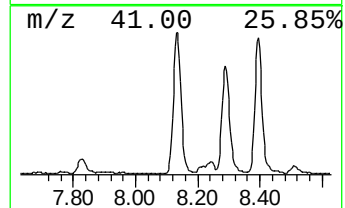
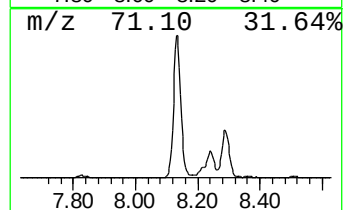
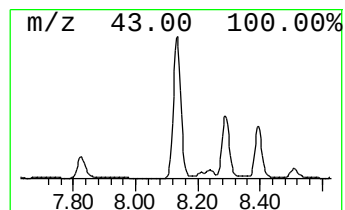
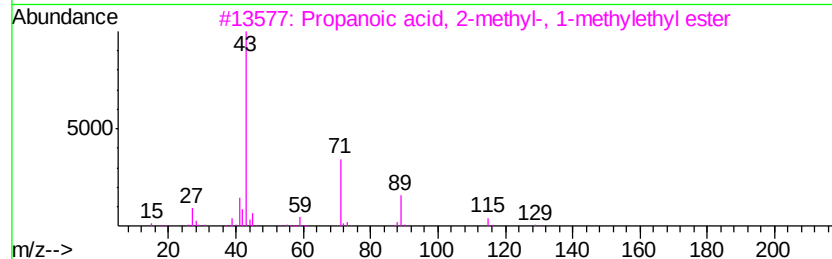
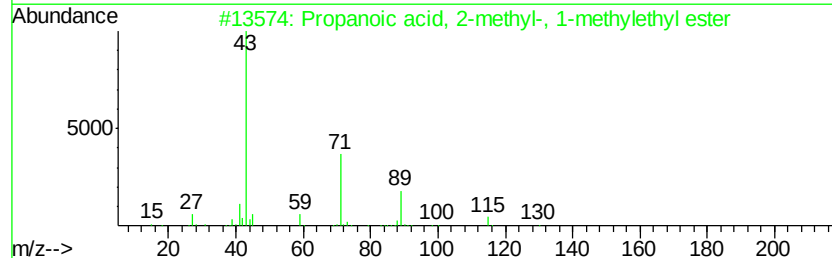
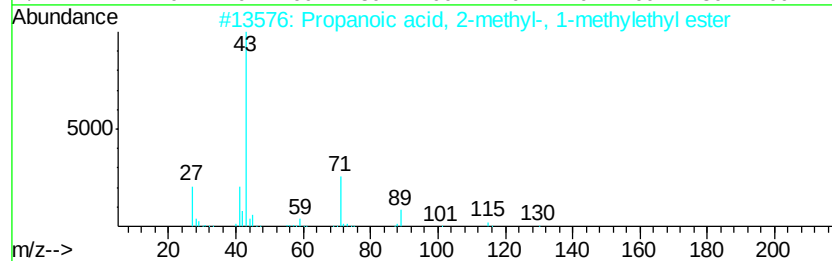
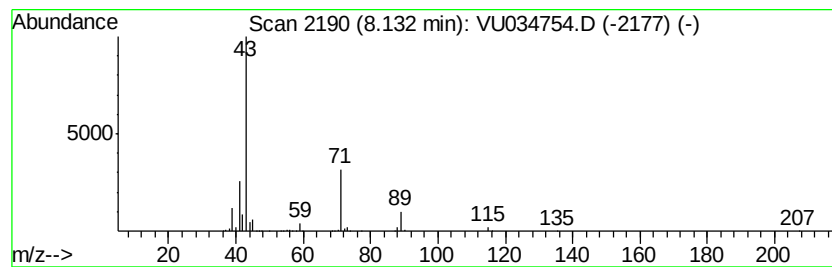
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Propanoic acid, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.13	22.98 ug/L	820063	Chlorobenzene-d5	9.07

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		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
4		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	59
5		3-Pentanone, 2,4-dimethyl-	114	C7H14O	000565-80-0	33



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
Data File : VU034754.D
Acq On : 23 Sep 2019 23:40
Operator : JC/SP
Sample : K4932-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFDJ9

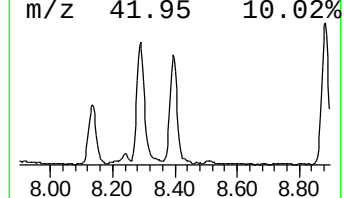
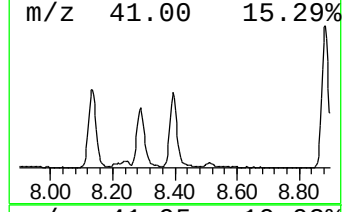
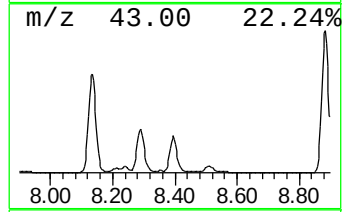
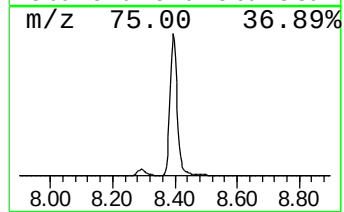
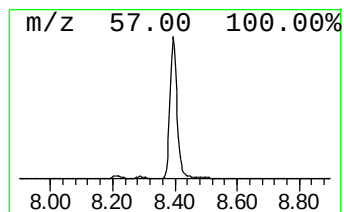
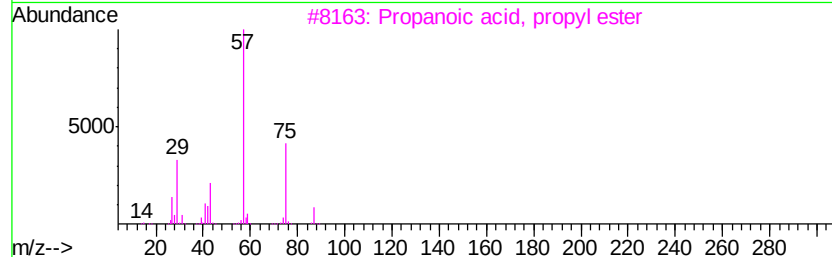
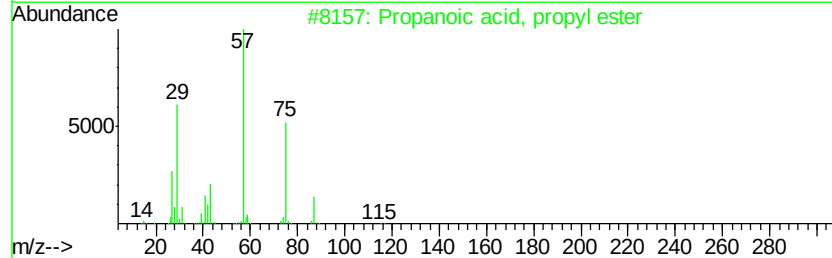
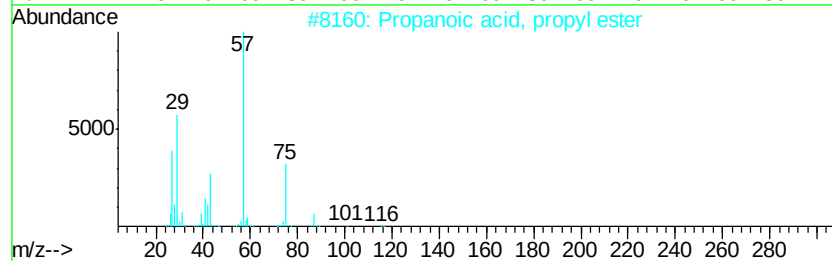
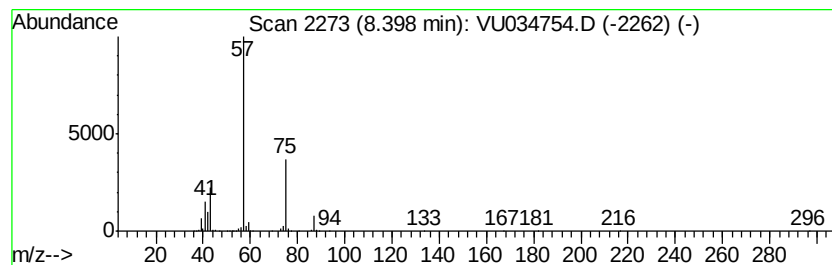
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Quant Title : VOC Analysis

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

Peak Number 5 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	36.42 ug/L	1299460	Chlorobenzene-d5	9.07

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	78
3		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	72
5		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	36



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
Data File : VU034754.D
Acq On : 23 Sep 2019 23:40
Operator : JC/SP
Sample : K4932-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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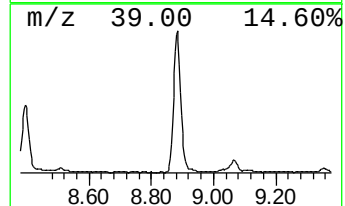
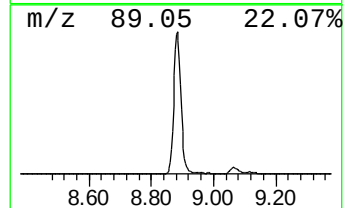
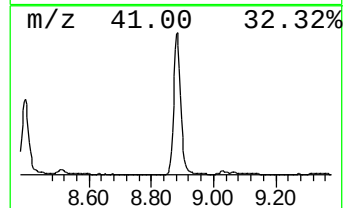
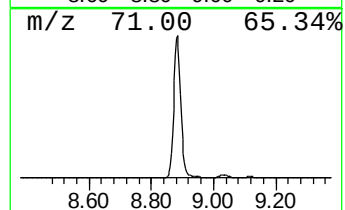
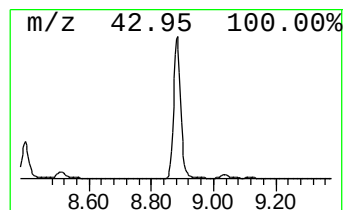
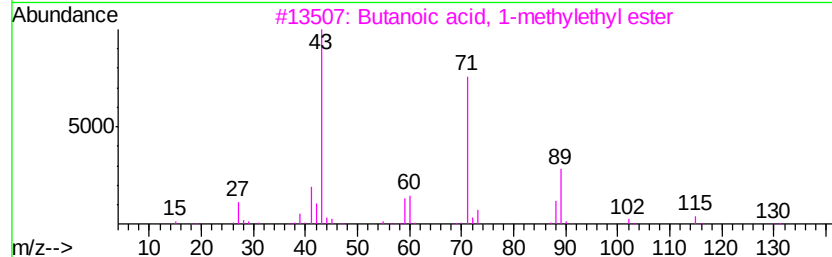
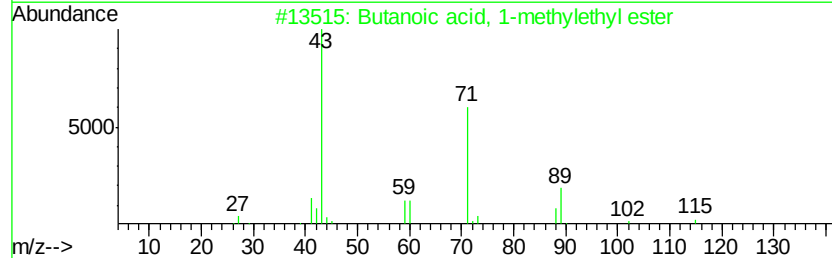
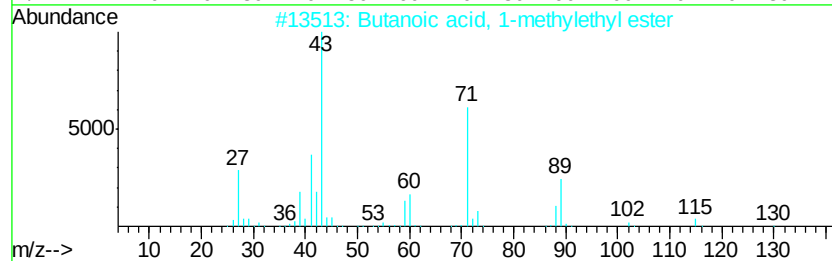
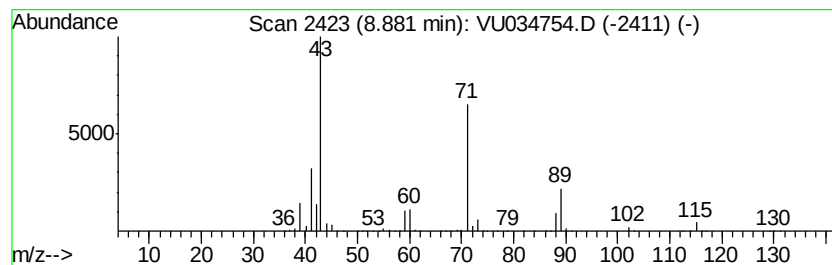
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Quant Title : VOC Analysis

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

Peak Number 6 Butanoic acid, 1-methylethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.88	45.25 ug/L	1614580	Chlorobenzene-d5	9.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
4		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	86
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
Data File : VU034754.D
Acq On : 23 Sep 2019 23:40
Operator : JC/SP
Sample : K4932-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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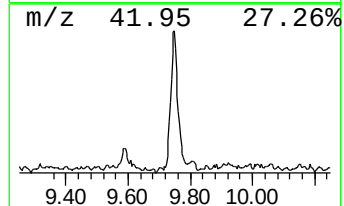
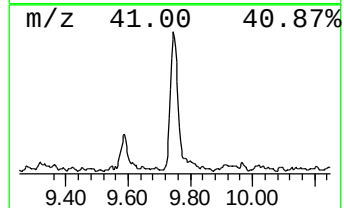
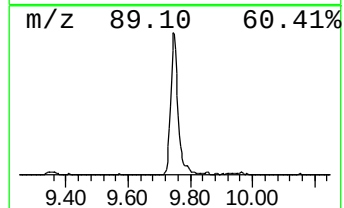
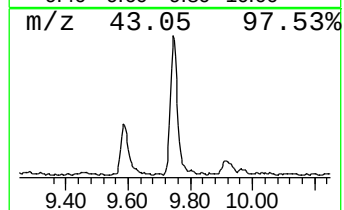
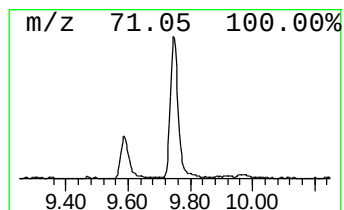
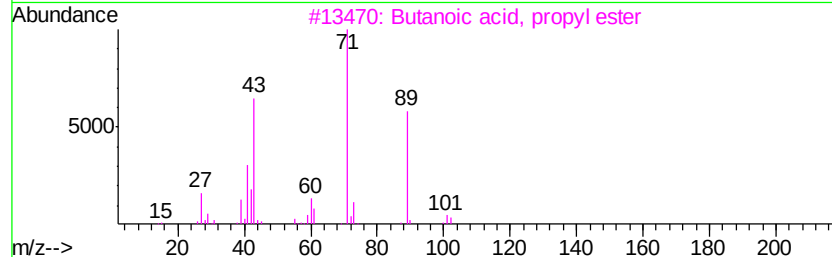
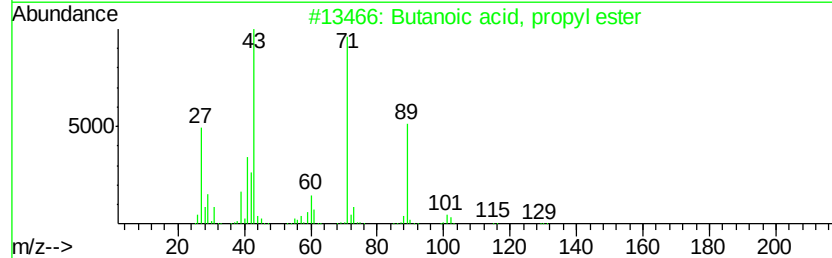
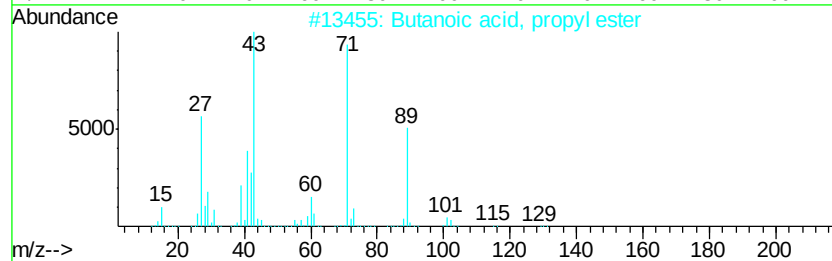
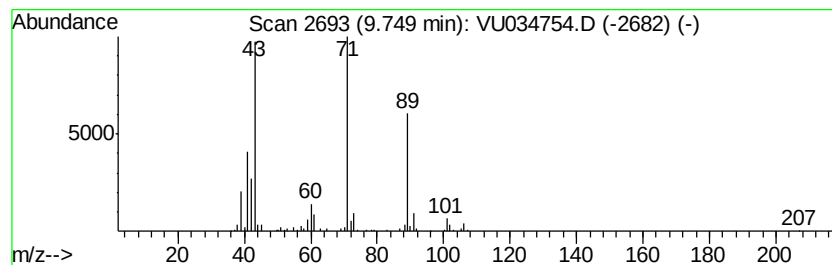
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Quant Title : VOC Analysis

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

Peak Number 7 Butanoic acid, propyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.75	5.96 ug/L	212799	Chlorobenzene-d5	9.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
2		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
3		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	86
4		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	83
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
Data File : VU034754.D
Acq On : 23 Sep 2019 23:40
Operator : JC/SP
Sample : K4932-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

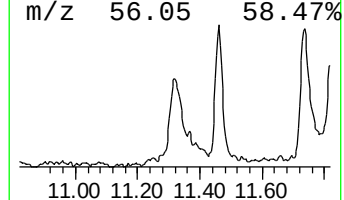
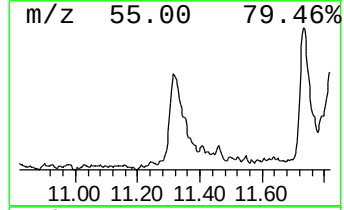
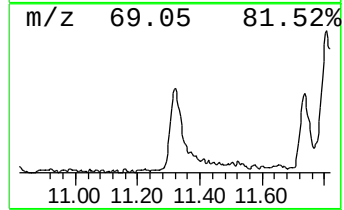
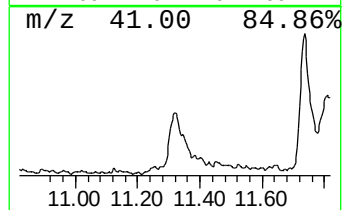
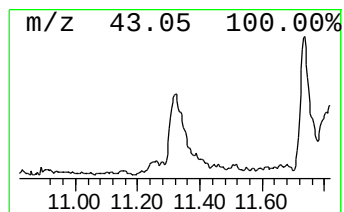
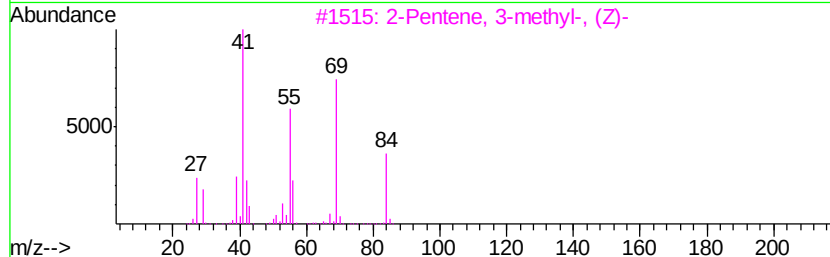
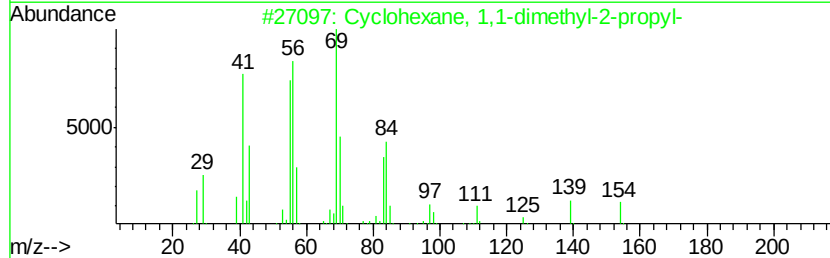
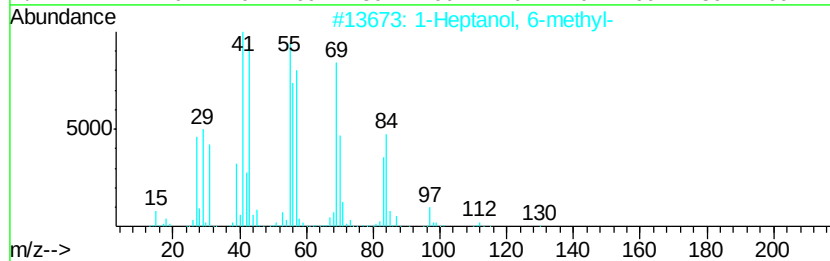
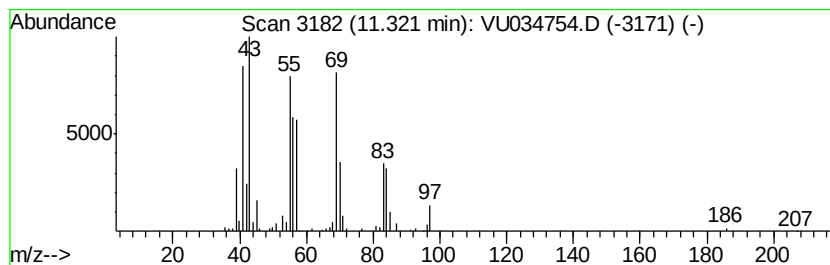
Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

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

Peak Number 8 1-Heptanol, 6-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.32	5.37 ug/L	174348	1,4-Dichlorobenzene-d4	11.46	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptanol, 6-methyl-	130	C8H18O	001653-40-3	91
2	Cyclohexane, 1,1-dimethyl-2-propyl-	154	C11H22	081983-71-3	78
3	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	49
4	Decane, 3-chloro-	176	C10H21Cl	001002-11-5	47
5	1-Heptanol, 6-methyl-	130	C8H18O	001653-40-3	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
Data File : VU034754.D
Acq On : 23 Sep 2019 23:40
Operator : JC/SP
Sample : K4932-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

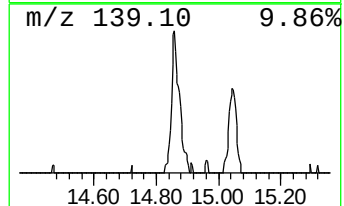
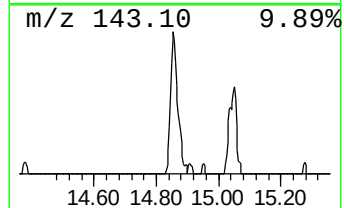
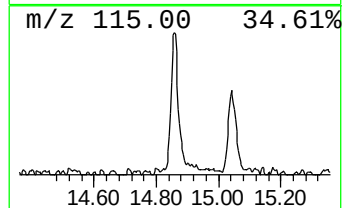
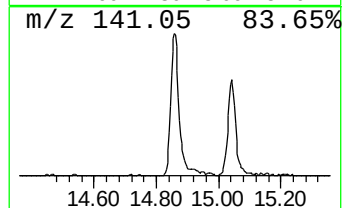
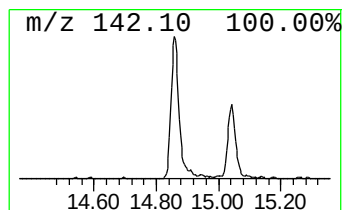
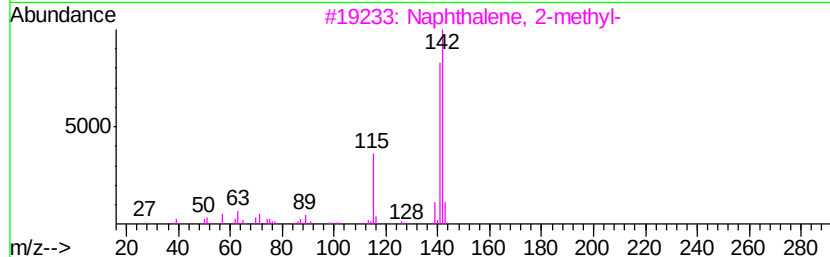
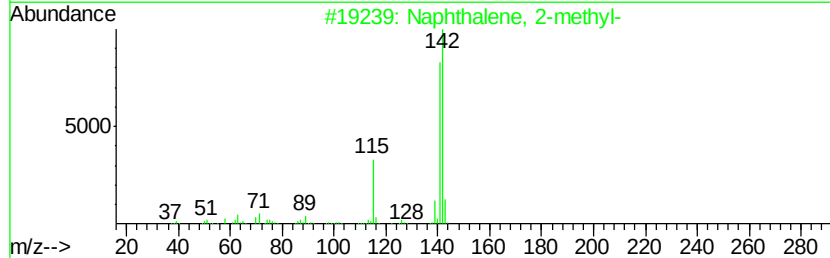
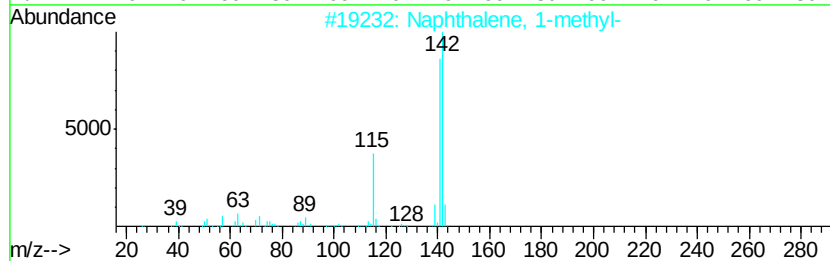
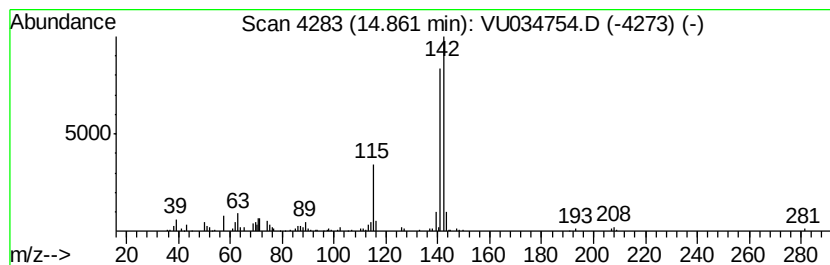
Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

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

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.86	3.08 ug/L	100107	1,4-Dichlorobenzene-d4	11.46	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092319\
Data File : VU034754.D
Acq On : 23 Sep 2019 23:40
Operator : JC/SP
Sample : K4932-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFDJ9

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.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	5.53	7.9	ug/L	222624	1	5.86	1401690	50.0
n-Propyl acetate	6.68	137.0	ug/L	3839980	1	5.86	1401690	50.0
Propanoic acid, 2...	8.13	23.0	ug/L	820063	2	9.07	1784160	50.0
Propanoic acid, p...	8.40	36.4	ug/L	1299460	2	9.07	1784160	50.0
Butanoic acid, 1-...	8.88	45.3	ug/L	1614580	2	9.07	1784160	50.0
Butanoic acid, pr...	9.75	6.0	ug/L	212799	2	9.07	1784160	50.0
1-Heptanol, 6-met...	11.32	5.4	ug/L	174348	3	11.46	1622990	50.0
Naphthalene, 1-me...	14.86	3.1	ug/L	100107	3	11.46	1622990	50.0