

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020222\
 Data File : VV024524.D
 Acq On : 02 Feb 2022 15:00
 Operator : SY/MD
 Sample : N1353-06
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 COVE7

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_V\Method\SFAMVLM020222WMA.M

Title : VOC Analysis

Signal : TIC: VV024524.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	77	83	96	rVB	200728	200531	14.54%	1.889%
2	1.568	157	164	176	rBV	193581	215443	15.62%	2.030%
3	2.108	323	332	361	rVB	358809	537018	38.94%	5.060%
4	3.031	615	619	620	rBV3	463	281	0.02%	0.003%
5	3.262	689	691	693	rBV	503	270	0.02%	0.003%
6	3.317	704	708	711	rBV3	860	816	0.06%	0.008%
7	3.513	765	769	773	rBV3	875	803	0.06%	0.008%
8	3.593	792	794	798	rBV2	376	221	0.02%	0.002%
9	3.626	801	804	805	rBV2	533	267	0.02%	0.003%
10	3.786	852	854	858	rBV2	460	419	0.03%	0.004%
11	3.825	864	866	870	rBV3	360	190	0.01%	0.002%
12	3.847	870	873	875	rBV2	505	294	0.02%	0.003%
13	3.886	875	885	923	rBV	88268	284472	20.63%	2.680%
14	4.349	1013	1029	1053	rBV	218685	539696	39.13%	5.085%
15	4.664	1122	1127	1129	rBV4	522	494	0.04%	0.005%
16	4.802	1167	1170	1173	rBV3	498	364	0.03%	0.003%
17	4.834	1178	1180	1185	rBV3	416	282	0.02%	0.003%
18	4.908	1200	1203	1206	rBV2	420	292	0.02%	0.003%
19	5.047	1229	1246	1279	rBV2	532563	1379065	100.00%	12.993%
20	5.616	1411	1423	1449	rBV	434053	871966	63.23%	8.215%
21	5.809	1480	1483	1484	rBV2	743	454	0.03%	0.004%
22	5.905	1501	1513	1533	rBV	33299	72203	5.24%	0.680%
23	6.011	1542	1546	1547	rBV3	789	580	0.04%	0.005%
24	6.069	1551	1564	1595	rVV	315035	633689	45.95%	5.970%
25	6.436	1677	1678	1679	rBV	431	131	0.01%	0.001%
26	6.484	1691	1693	1697	rVB4	570	277	0.02%	0.003%
27	6.728	1766	1769	1774	rVB5	666	637	0.05%	0.006%
28	6.754	1774	1777	1781	rBV3	647	589	0.04%	0.006%
29	6.789	1786	1788	1790	rBV2	318	160	0.01%	0.002%
30	6.841	1801	1804	1806	rVB3	798	403	0.03%	0.004%
31	6.854	1806	1808	1813	rVB4	572	362	0.03%	0.003%
32	6.902	1821	1823	1825	rBV2	311	112	0.01%	0.001%
33	6.918	1825	1828	1830	rBV2	418	267	0.02%	0.003%
34	6.931	1830	1832	1833	rBV	398	208	0.02%	0.002%
35	6.985	1838	1849	1870	rBV	190476	345955	25.09%	3.259%
36	7.313	1940	1951	1970	rBV	652173	1121600	81.33%	10.567%

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Peak Location: TOP

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM020222WMA.M
 Title : VOC Analysis

37	7.619	2036	2046	2070	rBV	132214	229201	16.62%	2.159%
38	7.834	2111	2113	2114	rBV	483	206	0.01%	0.002%
39	7.883	2123	2128	2131	rBV5	934	887	0.06%	0.008%
40	7.940	2137	2146	2154	rBV3	4248	7119	0.52%	0.067%
41	8.088	2182	2192	2224	rBV	344207	667504	48.40%	6.289%
42	8.477	2310	2313	2317	rBV4	631	546	0.04%	0.005%
43	8.558	2335	2338	2341	rBV3	447	465	0.03%	0.004%
44	8.616	2350	2356	2358	rBV4	711	793	0.06%	0.007%
45	8.693	2376	2380	2384	rBV6	822	694	0.05%	0.007%
46	8.792	2409	2411	2413	rBV2	508	321	0.02%	0.003%
47	8.850	2417	2429	2455	rBV	643665	1041195	75.50%	9.810%
48	9.291	2563	2566	2571	rBV4	661	504	0.04%	0.005%
49	9.333	2575	2579	2583	rBV4	669	647	0.05%	0.006%
50	9.403	2598	2601	2602	rBV2	537	302	0.02%	0.003%
51	9.490	2624	2628	2632	rBV3	629	722	0.05%	0.007%
52	9.548	2641	2646	2649	rBV6	568	693	0.05%	0.007%
53	9.590	2657	2659	2663	rBV3	309	140	0.01%	0.001%
54	9.632	2669	2672	2675	rBV4	572	381	0.03%	0.004%
55	9.731	2698	2703	2704	rBV4	473	429	0.03%	0.004%
56	9.915	2758	2760	2762	rBV3	399	194	0.01%	0.002%
57	10.018	2790	2792	2793	rBV	418	162	0.01%	0.002%
58	10.127	2820	2826	2837	rVB3	4864	7873	0.57%	0.074%
59	10.169	2837	2839	2842	rBV2	583	391	0.03%	0.004%
60	10.214	2842	2853	2873	rBV	349773	543976	39.45%	5.125%
61	10.503	2939	2943	2944	rBV3	659	330	0.02%	0.003%
62	10.866	3054	3056	3057	rBV2	505	176	0.01%	0.002%
63	11.095	3123	3127	3130	rVB4	576	471	0.03%	0.004%
64	11.149	3137	3144	3148	rBV4	1834	2694	0.20%	0.025%
65	11.246	3163	3174	3192	rBV	619873	946319	68.62%	8.916%
66	11.532	3255	3263	3279	rBV2	14206	25026	1.81%	0.236%
67	11.622	3280	3291	3316	rVV	589020	918443	66.60%	8.653%
68	11.966	3394	3398	3401	rBV4	851	889	0.06%	0.008%
69	12.204	3469	3472	3476	rBV2	473	389	0.03%	0.004%
70	12.480	3554	3558	3563	rBV4	1108	1258	0.09%	0.012%
71	12.960	3702	3707	3709	rBV4	760	613	0.04%	0.006%
72	13.342	3823	3826	3829	rBV2	556	386	0.03%	0.004%
73	13.677	3928	3930	3938	rVB5	959	878	0.06%	0.008%

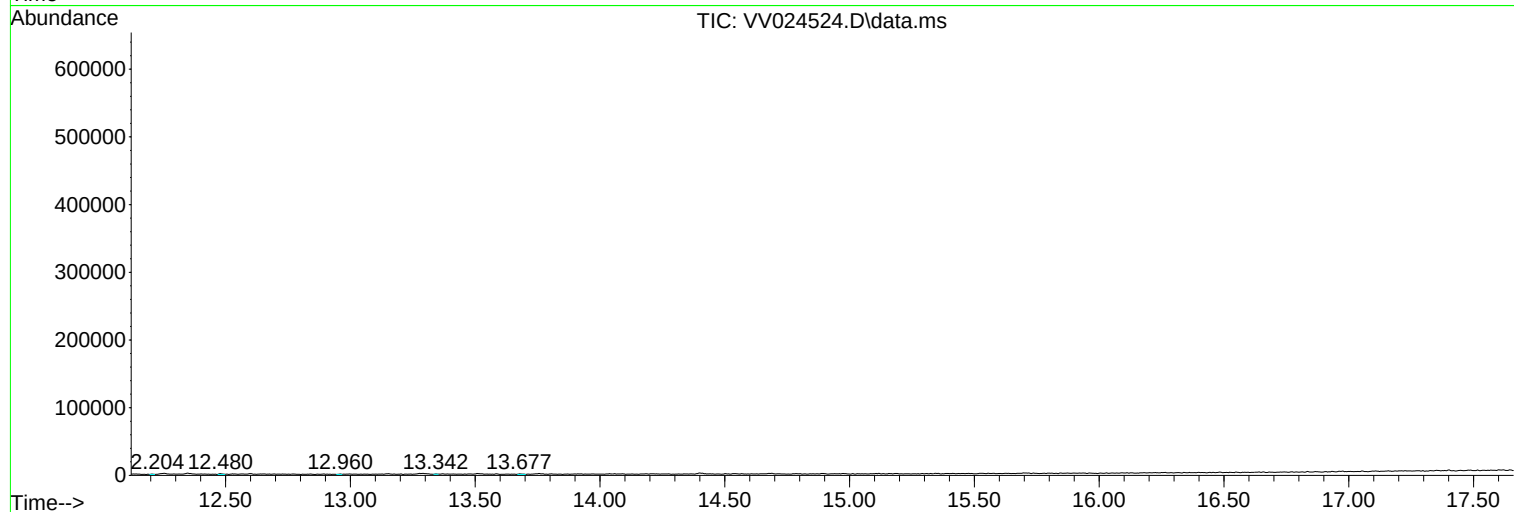
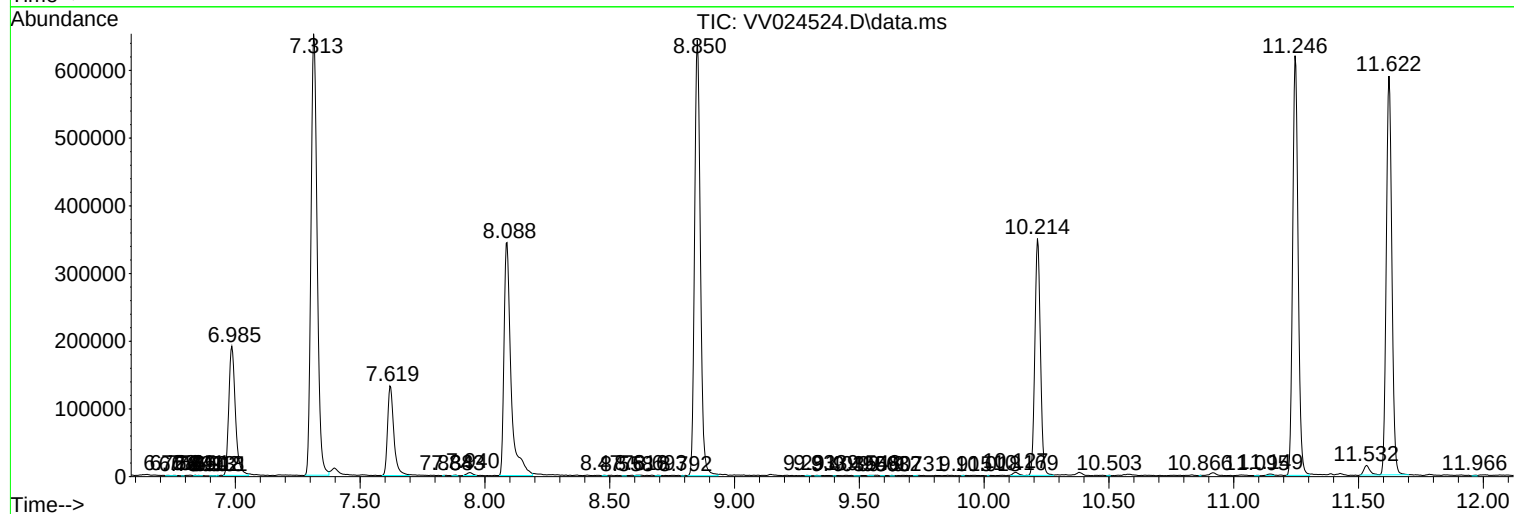
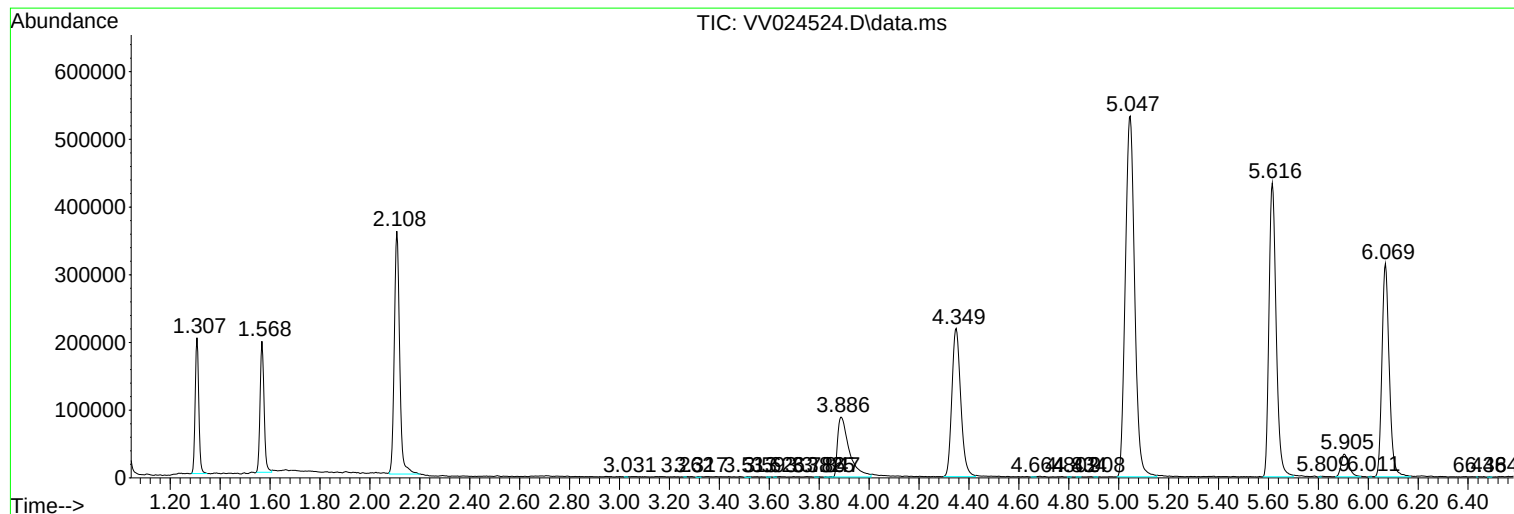
Sum of corrected areas: 10614028

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM020222WMA.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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ClientSampleId :
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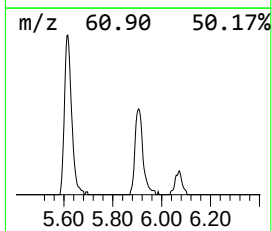
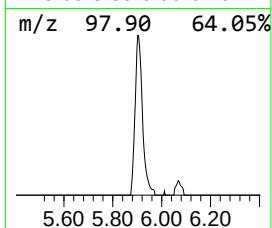
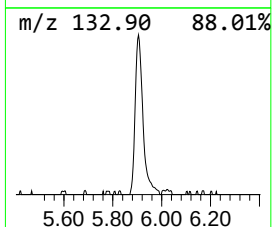
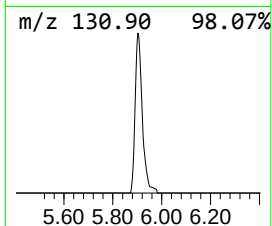
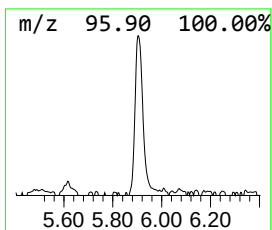
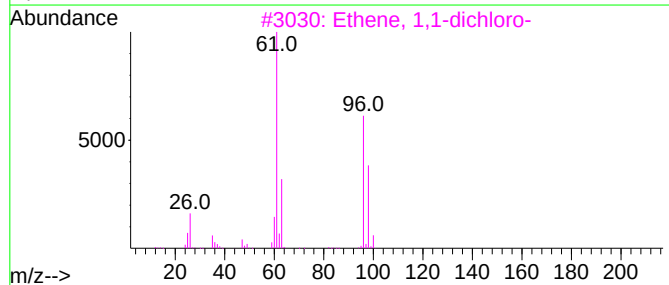
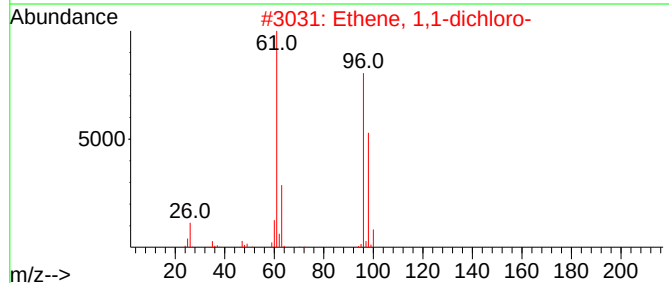
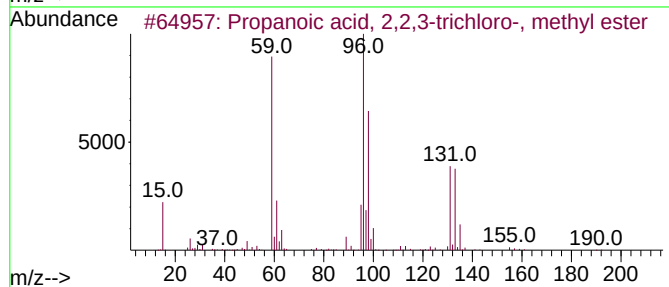
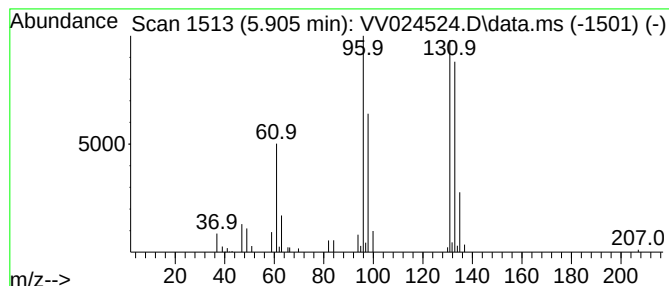
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM020222WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.905	4.14 ug/L	72203	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2,2,3-trichloro-...	190	C4H5Cl3O2	004749-35-3	39
2			Ethene, 1,1-dichloro-	96	C2H2Cl2	000075-35-4	38
3			Ethene, 1,1-dichloro-	96	C2H2Cl2	000075-35-4	38
4			Ethylene, 1,2-dichloro-, (Z)-	96	C2H2Cl2	000156-59-2	35
5			Ethene, 1,1-dichloro-	96	C2H2Cl2	000075-35-4	35



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	5.905	4.1	ug/L	72203	1	5.616	871966	50.0