

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV020222\
 Data File : VV024530.D
 Acq On : 02 Feb 2022 17:22
 Operator : SY/MD
 Sample : N1353-04MEDL 10X
 Misc : 2.91g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 COVE5MEDL

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: VV024530.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	76	83	93	rVB	188895	192508	14.80%	1.711%
2	1.568	157	164	178	rVB	167631	187495	14.41%	1.666%
3	2.105	323	331	343	rVB	348375	483311	37.15%	4.295%
4	3.371	723	725	728	rBV2	325	278	0.02%	0.002%
5	3.439	743	746	749	rBV5	461	301	0.02%	0.003%
6	3.503	763	766	768	rBV3	625	439	0.03%	0.004%
7	3.529	772	774	776	rVB2	450	163	0.01%	0.001%
8	3.548	776	780	781	rBV4	538	405	0.03%	0.004%
9	3.577	786	789	791	rBV3	517	304	0.02%	0.003%
10	3.613	797	800	802	rBV3	637	400	0.03%	0.004%
11	3.838	866	870	872	rBV2	343	285	0.02%	0.003%
12	3.912	879	893	934	rBV	54411	259150	19.92%	2.303%
13	4.349	1014	1029	1052	rBV	202980	504576	38.79%	4.484%
14	4.847	1179	1184	1188	rBV4	773	750	0.06%	0.007%
15	4.973	1219	1223	1224	rBV2	489	356	0.03%	0.003%
16	5.043	1229	1245	1277	rBV2	510128	1300950	100.00%	11.562%
17	5.616	1410	1423	1448	rBV	424519	853980	65.64%	7.590%
18	5.905	1502	1513	1529	rBV2	29900	60341	4.64%	0.536%
19	6.069	1544	1564	1593	rBV	302669	678308	52.14%	6.028%
20	6.889	1817	1819	1821	rBV3	367	206	0.02%	0.002%
21	6.985	1838	1849	1872	rBV	180568	329543	25.33%	2.929%
22	7.227	1913	1924	1941	rBV	519486	975094	74.95%	8.666%
23	7.313	1942	1951	1969	rVB	619386	1064746	81.84%	9.463%
24	7.619	2037	2046	2072	rBV	123592	217583	16.72%	1.934%
25	7.899	2129	2133	2134	rBV2	379	323	0.02%	0.003%
26	7.924	2139	2141	2145	rBV4	456	403	0.03%	0.004%
27	8.040	2174	2177	2179	rBV2	543	396	0.03%	0.004%
28	8.088	2182	2192	2227	rBV	316740	677551	52.08%	6.022%
29	8.500	2318	2320	2323	rVB3	610	292	0.02%	0.003%
30	8.538	2329	2332	2336	rBV3	569	490	0.04%	0.004%
31	8.567	2336	2341	2342	rBV4	676	563	0.04%	0.005%
32	8.638	2359	2363	2364	rBV2	398	251	0.02%	0.002%
33	8.673	2369	2374	2377	rBV5	722	673	0.05%	0.006%
34	8.689	2377	2379	2382	rBV2	405	195	0.01%	0.002%
35	8.751	2395	2398	2400	rBV3	401	235	0.02%	0.002%
36	8.850	2417	2429	2449	rBV	630287	1016794	78.16%	9.037%

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

37	9.442	2610	2613	2618	rBV4	443	440	0.03%	0.004%
38	9.519	2629	2637	2639	rBV5	2205	2964	0.23%	0.026%
39	9.709	2690	2696	2712	rVB	13074	21531	1.66%	0.191%
40	9.850	2738	2740	2742	rBV3	583	335	0.03%	0.003%
41	9.937	2762	2767	2768	rBV3	786	657	0.05%	0.006%
42	9.966	2774	2776	2779	rBV3	550	350	0.03%	0.003%
43	10.053	2801	2803	2806	rVB3	410	201	0.02%	0.002%
44	10.124	2818	2825	2835	rVB3	4534	6963	0.54%	0.062%
45	10.214	2841	2853	2872	rBV	345430	548433	42.16%	4.874%
46	10.538	2949	2954	2955	rBV3	1434	1120	0.09%	0.010%
47	10.654	2986	2990	2991	rBV2	580	405	0.03%	0.004%
48	10.696	2999	3003	3006	rVB2	310	208	0.02%	0.002%
49	10.709	3006	3007	3009	rBV	440	210	0.02%	0.002%
50	10.770	3025	3026	3030	rBV3	333	206	0.02%	0.002%
51	10.985	3090	3093	3094	rBV2	231	129	0.01%	0.001%
52	10.995	3094	3096	3101	rVB5	704	562	0.04%	0.005%
53	11.033	3101	3108	3111	rBV4	1145	1347	0.10%	0.012%
54	11.091	3124	3126	3127	rBV	543	235	0.02%	0.002%
55	11.246	3163	3174	3195	rBV	618070	942126	72.42%	8.373%
56	11.522	3257	3260	3262	rBV2	592	425	0.03%	0.004%
57	11.622	3279	3291	3315	rVB	583831	900819	69.24%	8.006%
58	11.783	3335	3341	3342	rBV4	1519	1587	0.12%	0.014%
59	11.911	3378	3381	3385	rVB3	784	554	0.04%	0.005%
60	11.959	3385	3396	3399	rBV5	692	1370	0.11%	0.012%
61	12.390	3526	3530	3532	rBV3	723	607	0.05%	0.005%
62	12.455	3547	3550	3552	rBV2	429	250	0.02%	0.002%
63	12.644	3607	3609	3610	rBV2	360	169	0.01%	0.002%
64	12.680	3618	3620	3621	rBV	611	248	0.02%	0.002%
65	12.718	3628	3632	3633	rBV2	594	340	0.03%	0.003%
66	12.744	3638	3640	3641	rBV	330	150	0.01%	0.001%
67	12.924	3694	3696	3697	rBV	264	111	0.01%	0.001%
68	12.972	3706	3711	3715	rBV5	857	807	0.06%	0.007%
69	13.014	3722	3724	3728	rBV4	484	344	0.03%	0.003%
70	13.091	3746	3748	3752	rVB2	394	301	0.02%	0.003%
71	13.111	3752	3754	3758	rVB3	570	390	0.03%	0.003%
72	13.136	3759	3762	3764	rBV2	491	328	0.03%	0.003%
73	13.165	3769	3771	3775	rBV3	595	511	0.04%	0.005%
74	13.258	3798	3800	3802	rBV2	349	164	0.01%	0.001%
75	13.368	3832	3834	3837	rVB	323	140	0.01%	0.001%
76	13.393	3837	3842	3846	rBV6	621	630	0.05%	0.006%

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Title : VOC Analysis

77	13.432	3852	3854	3857	rBV2	500	345	0.03%	0.003%
78	13.590	3900	3903	3905	rBV3	508	317	0.02%	0.003%
79	13.676	3927	3930	3931	rBV2	591	323	0.02%	0.003%
80	13.718	3937	3943	3944	rBV5	716	748	0.06%	0.007%
81	13.744	3948	3951	3958	rVB4	734	655	0.05%	0.006%
82	13.776	3958	3961	3964	rBV5	652	508	0.04%	0.005%
83	13.837	3977	3980	3986	rVB5	639	617	0.05%	0.005%
84	14.448	4168	4170	4171	rBV2	495	188	0.01%	0.002%

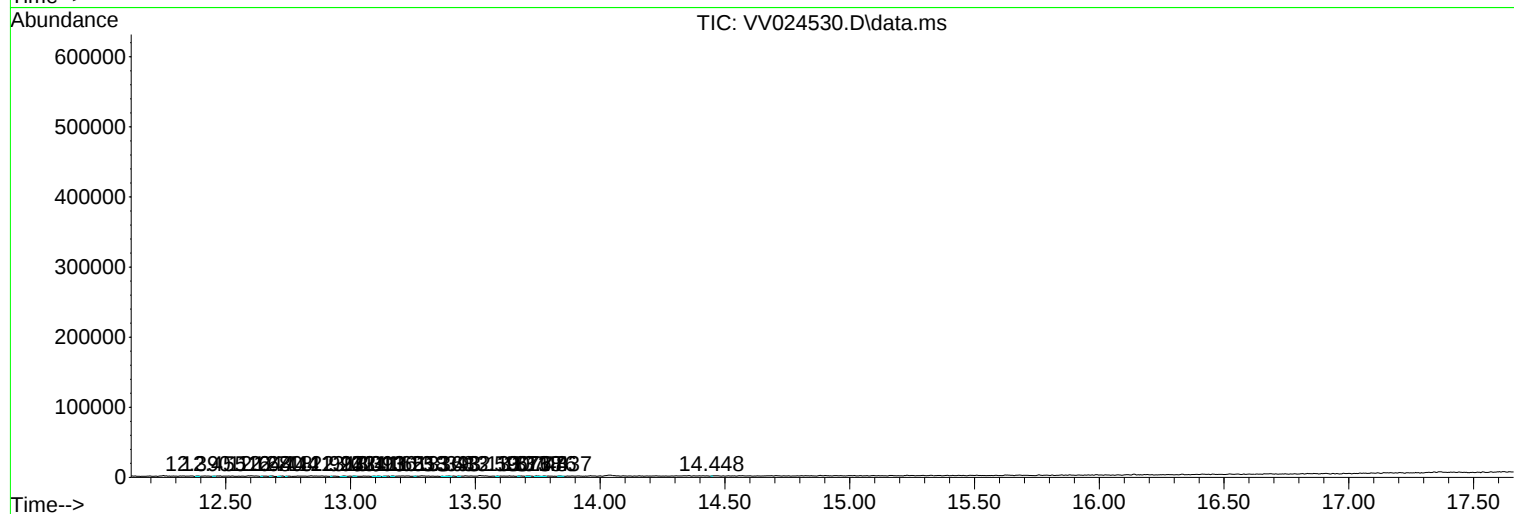
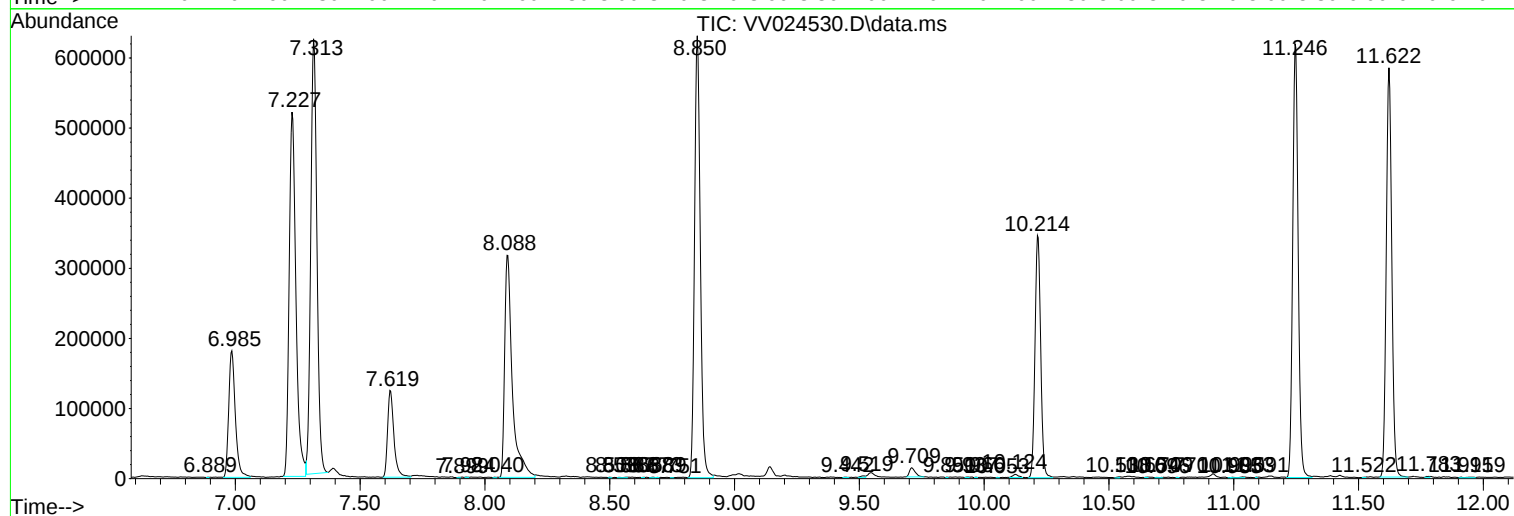
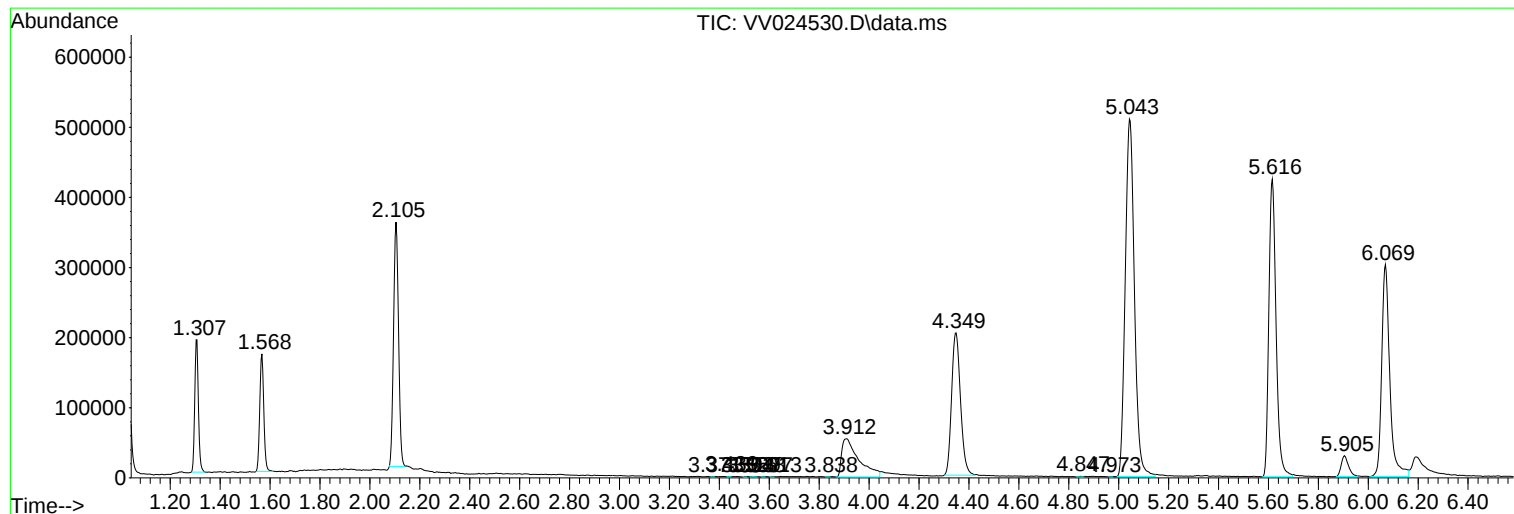
Sum of corrected areas: 11252006

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Instrument :
MSVOA_V
ClientSampleId :
C0VE5MEDL

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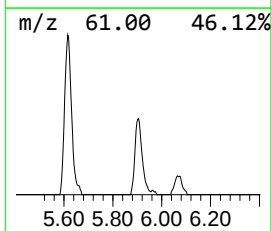
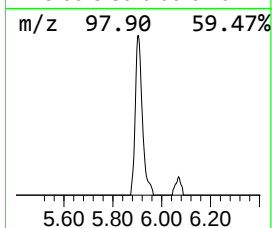
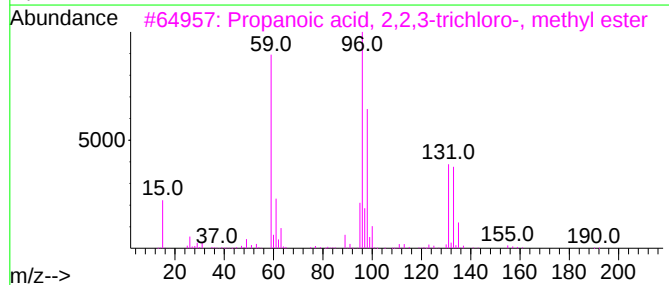
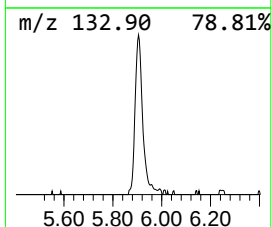
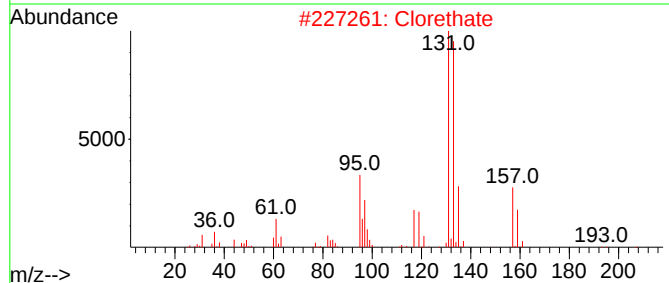
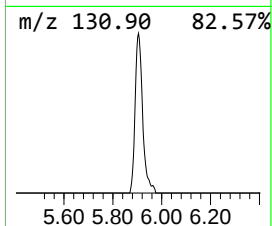
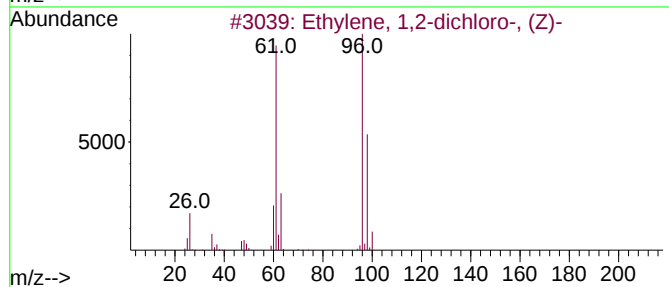
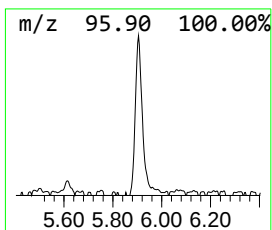
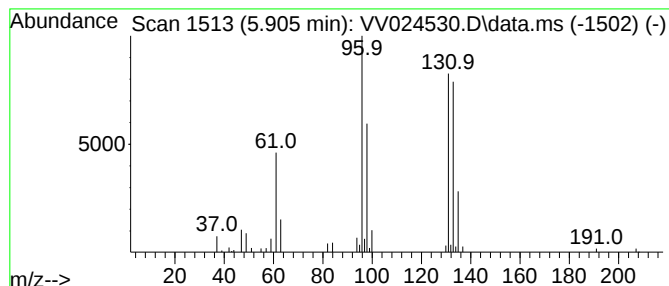
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM020222WMA.M
Quant Title : VOC Analysis

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	3.53 ug/L	60341	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylene, 1,2-dichloro-, (Z)-	96	C2H2Cl2	000156-59-2	38
2			Clorethate	322	C5H4Cl6O3	005634-37-7	28
3			Propanoic acid, 2,2,3-trichloro-...	190	C4H5Cl3O2	004749-35-3	25
4			Dihydro-4,5-dichloro-2(3H)furanone	154	C4H4Cl2O2	114434-99-0	25
5			2-Chloro-5-methylthiazole	133	C4H4ClNS	033342-65-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	5.905	3.5	ug/L	60341	1	5.616	853980	50.0