

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071724\
 Data File : VU059996.D
 Acq On : 17 Jul 2024 15:42
 Operator : MD/SY
 Sample : P3260-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 E08F9

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_U\Method\SFAMULM070224WMA.M
 Title : VOC Analysis

Signal : TIC: VU059996.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.596	87	95	116	rBV	70850	86329	1.92%	0.858%
2	1.808	145	161	177	rBV4	8202	24204	0.54%	0.241%
3	1.904	183	191	205	rVB	59352	82725	1.84%	0.822%
4	2.396	342	344	348	rVB3	574	287	0.01%	0.003%
5	2.560	383	395	408	rVV	119529	194702	4.34%	1.935%
6	2.763	454	458	459	rBV2	451	291	0.01%	0.003%
7	2.779	460	463	464	rBV2	337	218	0.00%	0.002%
8	3.039	537	544	550	rBV4	896	1393	0.03%	0.014%
9	3.084	555	558	563	rBV2	198	208	0.00%	0.002%
10	3.181	578	588	599	rBV	2321	4570	0.10%	0.045%
11	3.354	633	642	645	rBV4	863	1170	0.03%	0.012%
12	3.425	661	664	671	rVB2	207	194	0.00%	0.002%
13	3.731	754	759	763	rBV2	240	263	0.01%	0.003%
14	3.817	779	786	792	rVB	188	189	0.00%	0.002%
15	3.901	806	812	815	rVB2	215	192	0.00%	0.002%
16	4.033	849	853	858	rVB2	217	171	0.00%	0.002%
17	4.065	858	863	870	rBV2	196	349	0.01%	0.003%
18	4.319	939	942	945	rBV2	220	170	0.00%	0.002%
19	4.557	1014	1016	1023	rVB2	197	228	0.01%	0.002%
20	4.618	1023	1035	1061	rBV	72625	190380	4.24%	1.892%
21	5.055	1155	1171	1195	rBV	126355	281979	6.28%	2.802%
22	5.354	1257	1264	1268	rBV2	556	688	0.02%	0.007%
23	5.718	1356	1377	1402	rBV3	244119	668189	14.88%	6.641%
24	5.956	1448	1451	1454	rBV2	243	191	0.00%	0.002%
25	6.113	1497	1500	1503	rVB3	455	342	0.01%	0.003%
26	6.155	1503	1513	1516	rBV2	257	547	0.01%	0.005%
27	6.174	1516	1519	1523	rBV2	378	220	0.00%	0.002%
28	6.242	1526	1540	1562	rBV	248442	476436	10.61%	4.735%
29	6.361	1574	1577	1582	rVB4	934	717	0.02%	0.007%
30	6.470	1610	1611	1619	rVB2	280	282	0.01%	0.003%
31	6.538	1619	1632	1648	rBV4	39785	78323	1.74%	0.778%
32	6.682	1663	1677	1702	rBV	156592	307402	6.85%	3.055%
33	6.827	1719	1722	1724	rVB	441	199	0.00%	0.002%
34	7.103	1799	1808	1817	rVB4	1491	2724	0.06%	0.027%
35	7.155	1820	1824	1828	rVB2	211	167	0.00%	0.002%
36	7.187	1828	1834	1835	rBV2	525	368	0.01%	0.004%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM070224WMA.M
 Title : VOC Analysis

37	7.226	1843	1846	1851	rVB2	328	300	0.01%	0.003%
38	7.348	1879	1884	1887	rVV2	234	214	0.00%	0.002%
39	7.370	1887	1891	1896	rVB2	274	250	0.01%	0.002%
40	7.412	1900	1904	1907	rBV2	274	217	0.00%	0.002%
41	7.560	1938	1950	1976	rBV2	77322	137709	3.07%	1.369%
42	7.708	1990	1996	1999	rBV3	563	489	0.01%	0.005%
43	7.814	2024	2029	2032	rBV2	287	258	0.01%	0.003%
44	7.891	2041	2053	2075	rBV	275695	484173	10.79%	4.812%
45	8.116	2120	2123	2125	rBV2	241	200	0.00%	0.002%
46	8.174	2128	2141	2164	rBV	52049	90553	2.02%	0.900%
47	8.290	2173	2177	2180	rBV2	398	339	0.01%	0.003%
48	8.309	2180	2183	2186	rVV2	251	220	0.00%	0.002%
49	8.325	2186	2188	2194	rVB	380	264	0.01%	0.003%
50	8.467	2222	2232	2242	rVV2	1782	2643	0.06%	0.026%
51	8.547	2242	2257	2273	rVV	2677344	4489078	100.00%	44.613%
52	8.627	2274	2282	2323	rVV	205316	408823	9.11%	4.063%
53	8.769	2324	2326	2327	rVV3	1164	593	0.01%	0.006%
54	8.804	2330	2337	2348	rVV4	5194	8447	0.19%	0.084%
55	8.846	2348	2350	2353	rVB2	416	244	0.01%	0.002%
56	8.872	2353	2358	2361	rBV4	409	451	0.01%	0.004%
57	8.971	2384	2389	2392	rBV2	242	233	0.01%	0.002%
58	8.988	2392	2394	2399	rBV	241	187	0.00%	0.002%
59	9.094	2423	2427	2432	rVB2	250	213	0.00%	0.002%
60	9.152	2440	2445	2448	rBV2	166	179	0.00%	0.002%
61	9.187	2453	2456	2459	rVB2	244	170	0.00%	0.002%
62	9.222	2464	2467	2471	rVB	282	212	0.00%	0.002%
63	9.412	2514	2526	2560	rVV	366630	619338	13.80%	6.155%
64	9.624	2589	2592	2594	rBV	272	182	0.00%	0.002%
65	9.692	2607	2613	2615	rBV	548	463	0.01%	0.005%
66	10.074	2726	2732	2736	rBV2	165	204	0.00%	0.002%
67	10.103	2736	2741	2747	rVV3	734	1051	0.02%	0.010%
68	10.283	2791	2797	2806	rBV4	3149	4727	0.11%	0.047%
69	10.470	2852	2855	2859	rVV	213	194	0.00%	0.002%
70	10.750	2929	2942	2964	rBV	257972	418572	9.32%	4.160%
71	10.891	2976	2986	2992	rBV2	2139	3405	0.08%	0.034%
72	10.987	3012	3016	3018	rBV2	265	208	0.00%	0.002%
73	11.161	3067	3070	3075	rVB	244	174	0.00%	0.002%
74	11.463	3160	3164	3168	rBV2	269	270	0.01%	0.003%
75	11.650	3219	3222	3225	rBV	205	172	0.00%	0.002%
76	11.724	3241	3245	3249	rBV2	270	260	0.01%	0.003%

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77	11.743	3249	3251	3254	rVB3	331	177	0.00%	0.002%
78	11.807	3259	3271	3292	rBV	313283	515235	11.48%	5.120%
79	12.187	3377	3389	3411	rBV	270587	454592	10.13%	4.518%
80	12.711	3549	3552	3557	rVB2	323	247	0.01%	0.002%
81	12.769	3563	3570	3580	rVB2	1556	2380	0.05%	0.024%
82	12.827	3585	3588	3592	rBV2	189	198	0.00%	0.002%
83	13.103	3670	3674	3681	rBV	311	415	0.01%	0.004%
84	13.470	3784	3788	3793	rBV2	264	282	0.01%	0.003%
85	13.495	3793	3796	3800	rVV	257	192	0.00%	0.002%
86	14.450	4091	4093	4096	rBV2	285	210	0.00%	0.002%
87	14.508	4108	4111	4116	rBV3	277	297	0.01%	0.003%
88	14.547	4121	4123	4127	rBV	323	232	0.01%	0.002%
89	14.595	4135	4138	4141	rVB3	333	228	0.01%	0.002%
90	14.614	4141	4144	4147	rBV2	297	221	0.00%	0.002%
91	14.929	4239	4242	4244	rBV2	319	210	0.00%	0.002%
92	15.193	4320	4324	4329	rVV3	487	563	0.01%	0.006%
93	15.254	4341	4343	4346	rBV	372	171	0.00%	0.002%
94	15.563	4436	4439	4440	rBV2	335	185	0.00%	0.002%
95	15.608	4451	4453	4456	rBV4	542	324	0.01%	0.003%
96	15.859	4529	4531	4534	rBV4	565	382	0.01%	0.004%
97	15.917	4547	4549	4551	rBV2	483	228	0.01%	0.002%
98	16.074	4592	4598	4602	rBV4	1079	1156	0.03%	0.011%
99	16.171	4626	4628	4631	rBV2	705	520	0.01%	0.005%
100	16.331	4676	4678	4683	rBV3	657	688	0.02%	0.007%

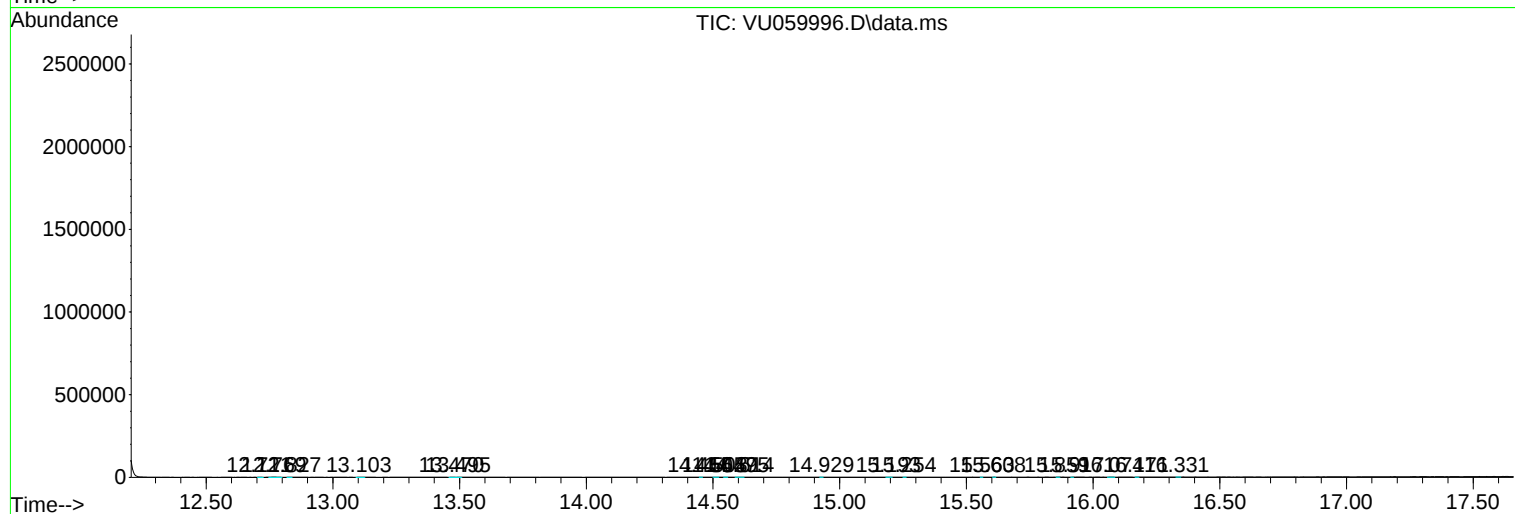
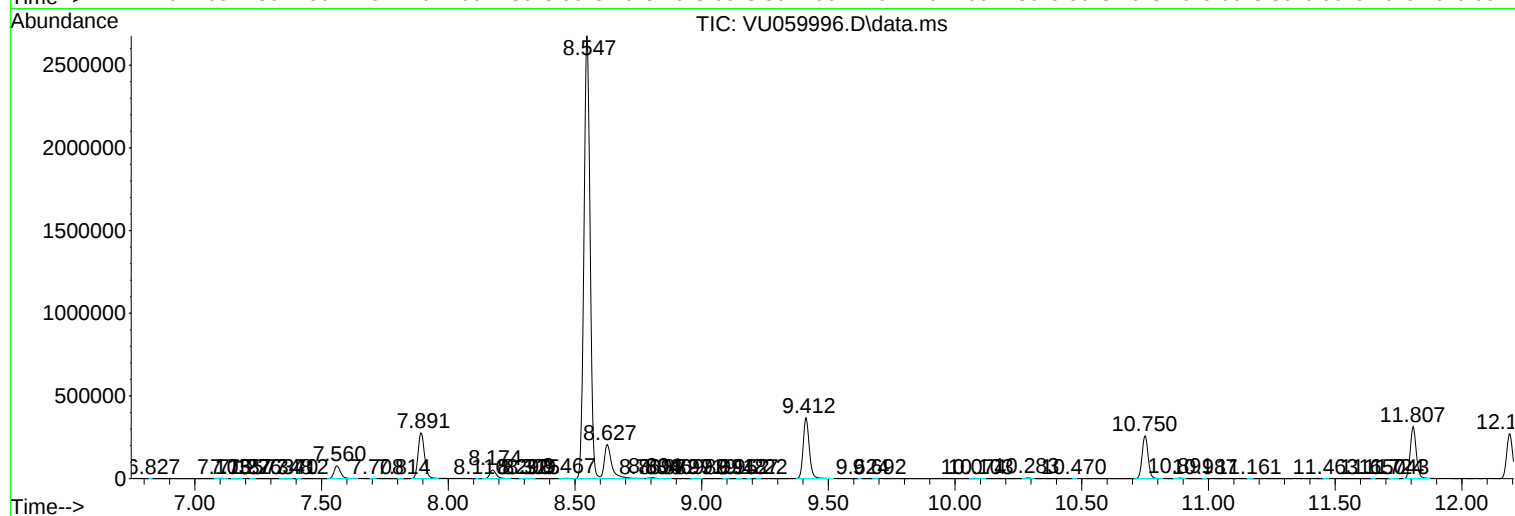
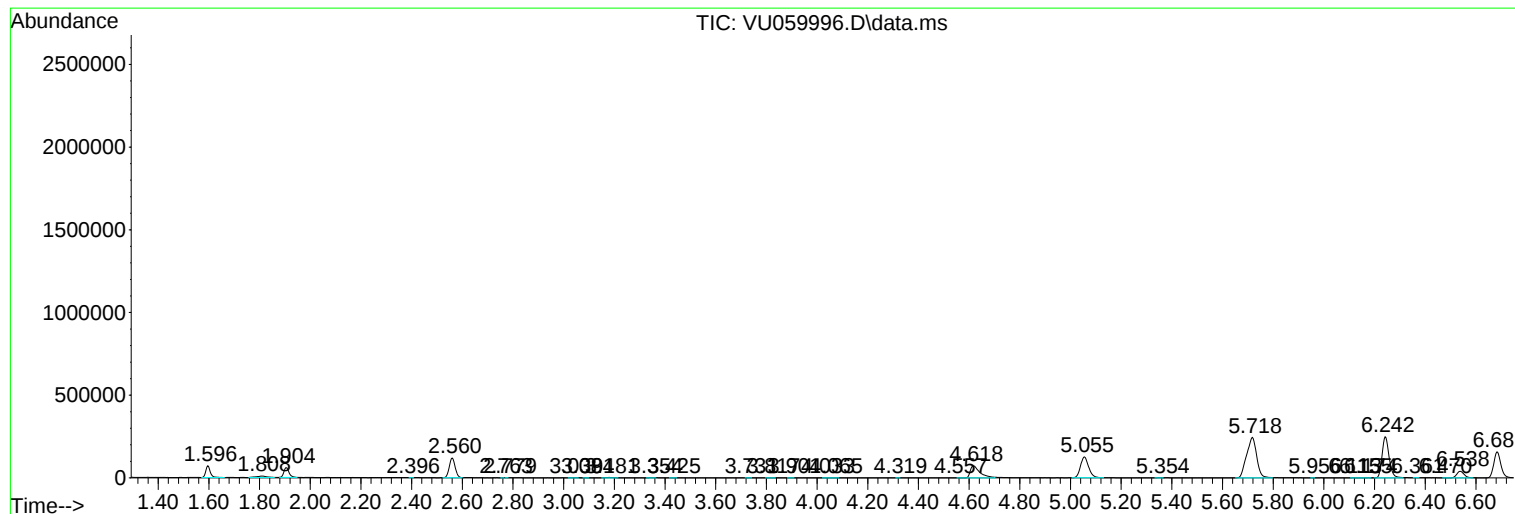
Sum of corrected areas: 10062219

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM070224WMA.M
Quant Title : VOC Analysis

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



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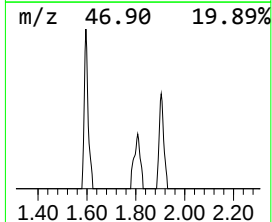
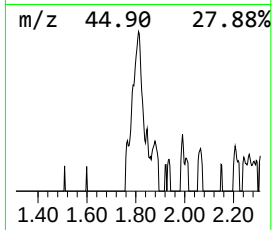
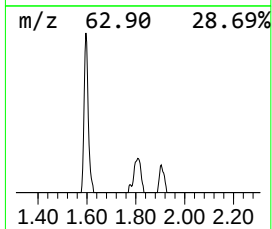
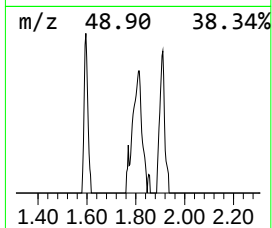
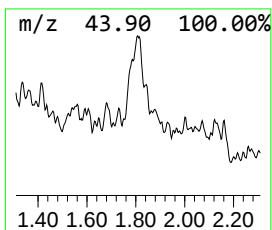
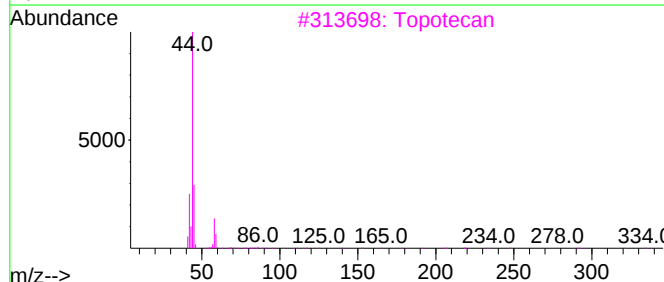
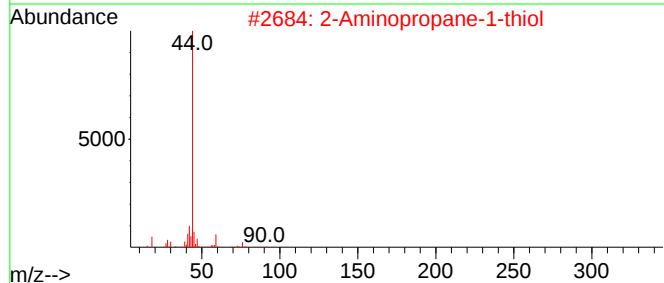
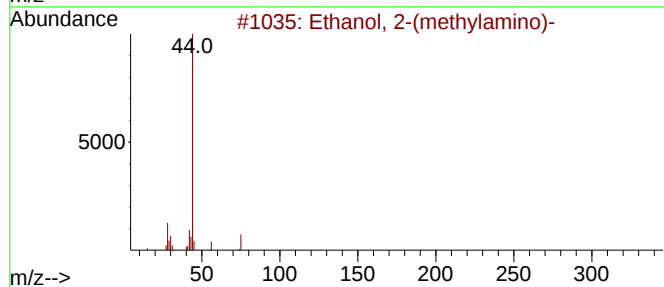
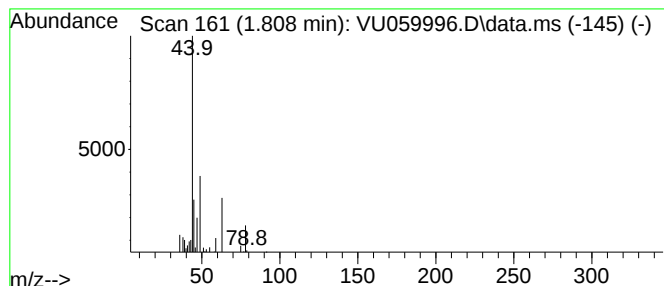
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM070224WMA.M
Quant Title : VOC Analysis

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.
1.808	2.54 ug/L	24204	1,4-Difluorobenzene	6.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(methylamino)-	75	C3H9NO	000109-83-1	9
2			2-Aminopropane-1-thiol	91	C3H9NS	010229-29-5	9
3			Topotecan	421	C23H23N3O5	1000131-70-7	9
4			Ethanol, 2-(methylamino)-	75	C3H9NO	000109-83-1	5
5			Hydrazinecarboxamide	75	CH5N3O	000057-56-7	4



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TIC Library : C:\Database\NIST20.L
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	1.808	2.5	ug/L	24204	1	6.242	476436	50.0