

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD071421\
 Data File : VD069694.D
 Acq On : 14 Jul 2021 18:08
 Operator : VA/SY
 Sample : M2997-01RE
 Misc : 5.08G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 RBR-200029RE

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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_D\Method\82D071321S.M
 Title : SW846 8260

Signal : TIC: VD069694.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.221	45	51	52	rBV3	7300	12335	3.70%	0.593%
2	2.268	57	59	61	rBV3	3852	3846	1.15%	0.185%
3	2.626	117	120	121	rBV3	5577	6056	1.81%	0.291%
4	2.826	152	154	156	rBV2	2814	3470	1.04%	0.167%
5	2.850	156	158	161	rVV4	4918	5915	1.77%	0.284%
6	2.879	161	163	172	rVB6	9495	17753	5.32%	0.853%
7	3.450	257	260	265	rBV2	2056	3561	1.07%	0.171%
8	4.956	508	516	529	rBV4	11303	33756	10.12%	1.622%
9	7.679	974	979	981	rBV2	3385	6280	1.88%	0.302%
10	7.697	981	982	987	rVB3	2791	3699	1.11%	0.178%
11	7.914	1010	1019	1024	rBV3	46234	108809	32.61%	5.229%
12	7.979	1024	1030	1040	rVV2	74357	167271	50.12%	8.039%
13	8.326	1081	1089	1097	rVV2	60730	137799	41.29%	6.623%
14	8.861	1172	1180	1189	rVB2	114463	237391	71.14%	11.409%
15	9.773	1332	1335	1339	rBV2	2340	3443	1.03%	0.165%
16	10.332	1423	1430	1439	rBV	177089	331627	99.37%	15.938%
17	10.973	1535	1539	1544	rBV3	2007	4088	1.23%	0.196%
18	11.638	1646	1652	1658	rBV	190764	333713	100.00%	16.038%
19	12.626	1814	1820	1827	rVV2	133296	226659	67.92%	10.893%
20	12.679	1827	1829	1836	rVB4	3050	5409	1.62%	0.260%
21	13.161	1908	1911	1916	rVB3	2787	3781	1.13%	0.182%
22	13.567	1971	1980	1986	rBV	166213	277362	83.11%	13.330%
23	13.661	1989	1996	2001	rBV4	14807	24038	7.20%	1.155%
24	13.926	2036	2041	2043	rBV2	3152	4183	1.25%	0.201%
25	13.985	2047	2051	2055	rBV	2533	3869	1.16%	0.186%
26	14.255	2092	2097	2106	rVB9	11608	30250	9.06%	1.454%
27	14.390	2116	2120	2123	rVV4	3284	4639	1.39%	0.223%
28	14.426	2123	2126	2129	rVV3	2991	4346	1.30%	0.209%
29	14.467	2129	2133	2137	rVB4	3603	5419	1.62%	0.260%
30	14.573	2149	2151	2157	rVB5	3143	5038	1.51%	0.242%
31	14.708	2168	2174	2178	rBV5	3297	6283	1.88%	0.302%
32	14.761	2181	2183	2187	rVB4	3178	4087	1.22%	0.196%
33	14.808	2187	2191	2194	rBV4	3889	5799	1.74%	0.279%
34	14.873	2198	2202	2205	rBV6	5234	6219	1.86%	0.299%
35	14.932	2209	2212	2216	rBV4	3277	4329	1.30%	0.208%
36	15.143	2245	2248	2253	rVB5	2894	4267	1.28%	0.205%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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37	15.432	2294	2297	2299	rBV3	3221	3573	1.07%	0.172%
38	15.885	2372	2374	2376	rBV2	2978	3463	1.04%	0.166%
39	16.014	2395	2396	2400	rVB3	3921	3709	1.11%	0.178%
40	16.120	2410	2414	2416	rBV3	3124	3499	1.05%	0.168%
41	16.320	2446	2448	2451	rVV3	3130	3365	1.01%	0.162%
42	16.361	2454	2455	2458	rVB3	4585	3727	1.12%	0.179%
43	16.449	2469	2470	2473	rBV4	3210	3647	1.09%	0.175%
44	16.655	2503	2505	2509	rVB3	3850	5426	1.63%	0.261%
45	16.702	2511	2513	2515	rVV3	3882	3508	1.05%	0.169%

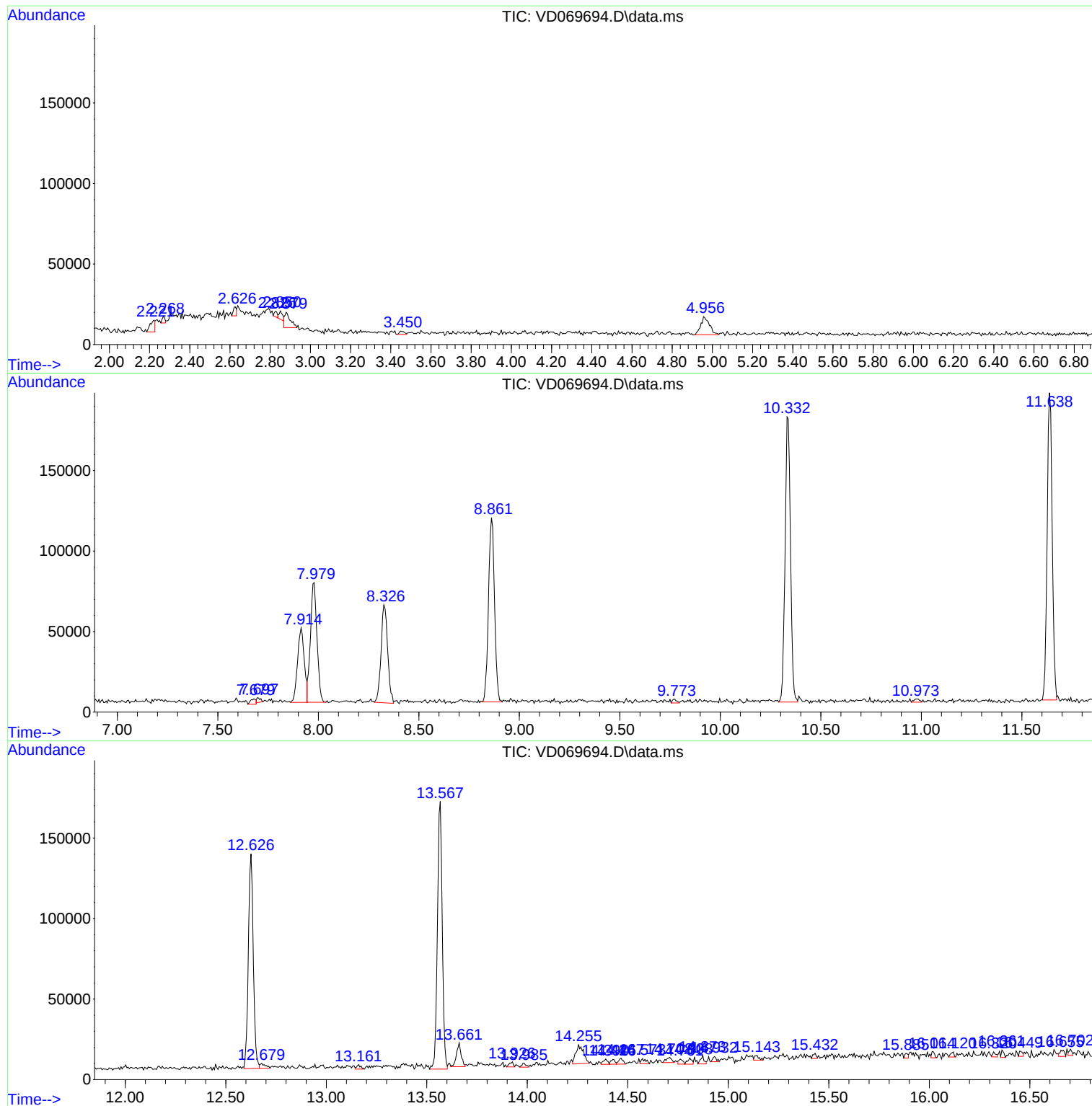
Sum of corrected areas: 2080706

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D071321S.M
Quant Title : SW846 8260

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



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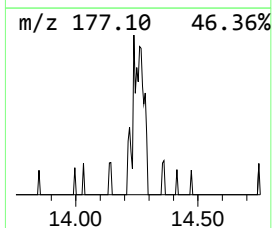
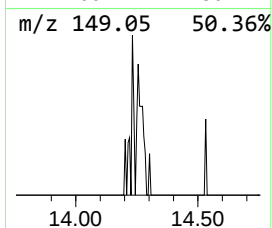
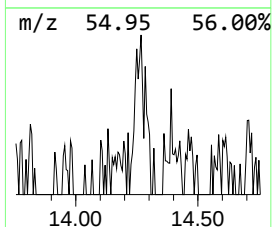
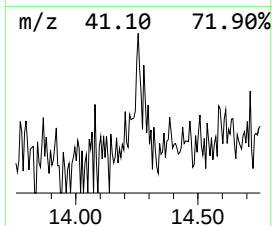
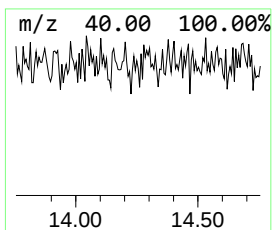
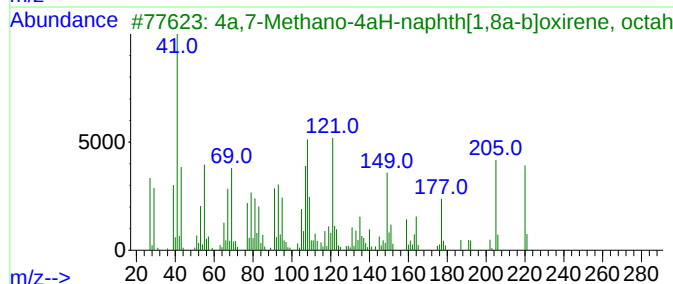
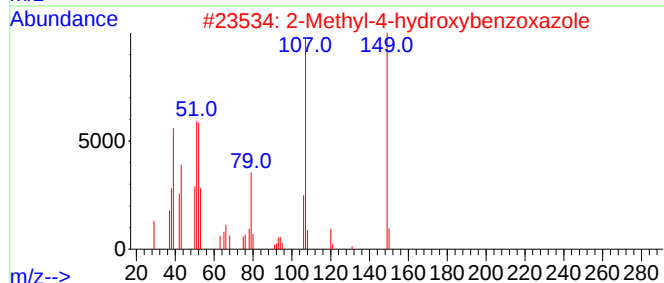
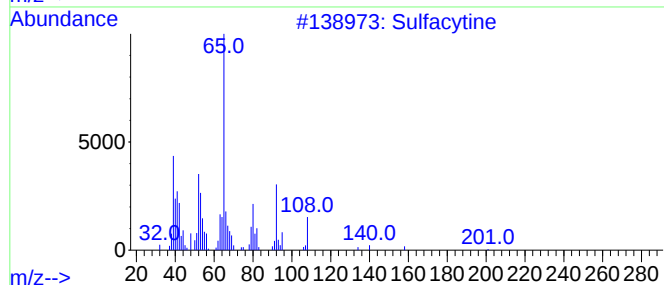
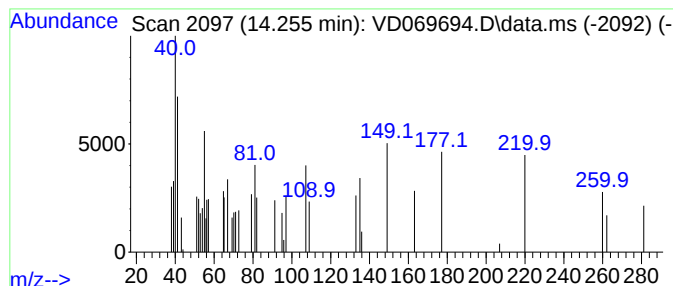
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Peak Number 3 unknown14.255 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.255	5.45 ug/l	30250	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfacytine	294	C12H14N4O3S	017784-12-2	17
2			2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	14
3			4a,7-Methano-4aH-naphth[1,8a-b]o...	220	C15H24O	067999-56-8	10
4			Phenol, 2-amino-4-(1H-1,2,3,4-te...	177	C7H7NO	1000362-63-1	9
5			Phenoxyacetoneitrile	133	C8H7NO	003598-14-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown14.255	14.255	5.5	ug/l	30250	4	13.561	277362	50.0