

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY051722\
 Data File : VY008786.D
 Acq On : 17 May 2022 21:05
 Operator : KP/MD
 Sample : N2795-01
 Misc : 5.10g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VNJ-239

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_Y\methods\82Y051722S.M
 Title : SW846 8260

Signal : TIC: VY008786.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.028	31	34	37	rVB2	454	575	1.00%	0.160%
2	2.101	43	46	47	rBV2	992	1088	1.90%	0.302%
3	2.119	47	49	50	rBV2	1193	1076	1.87%	0.299%
4	2.168	55	57	60	rBV2	716	954	1.66%	0.265%
5	2.205	60	63	65	rBV4	1059	1063	1.85%	0.296%
6	2.308	78	80	81	rBV	915	584	1.02%	0.162%
7	2.375	88	91	92	rBV3	876	998	1.74%	0.277%
8	2.510	109	113	118	rVB5	1733	3142	5.47%	0.874%
9	2.625	128	132	133	rBV3	1031	1195	2.08%	0.332%
10	2.644	133	135	136	rVV2	1051	814	1.42%	0.226%
11	2.662	136	138	145	rVB5	2157	2823	4.92%	0.785%
12	2.735	148	150	153	rBV4	648	887	1.55%	0.247%
13	2.759	153	154	157	rBV3	962	626	1.09%	0.174%
14	2.802	159	161	162	rBV	705	579	1.01%	0.161%
15	3.009	193	195	200	rVB3	477	624	1.09%	0.173%
16	3.052	200	202	208	rVB2	490	761	1.33%	0.212%
17	3.936	343	347	352	rBV3	342	795	1.38%	0.221%
18	4.692	469	471	472	rBV	835	693	1.21%	0.193%
19	4.716	472	475	480	rVB3	1130	1847	3.22%	0.513%
20	6.984	837	847	853	rBV2	3564	11782	20.52%	3.276%
21	7.301	898	899	904	rVB3	466	634	1.10%	0.176%
22	7.716	959	967	973	rBV3	9898	22975	40.02%	6.387%
23	7.789	973	979	988	rVB3	15479	33165	57.77%	9.220%
24	8.137	1027	1036	1045	rBV2	9722	22026	38.37%	6.124%
25	8.691	1118	1127	1137	rVB	22457	44363	77.28%	12.334%
26	10.179	1364	1371	1378	rBV	34068	57408	100.00%	15.960%
27	10.575	1431	1436	1442	rBV3	2974	6164	10.74%	1.714%
28	10.953	1493	1498	1505	rBV3	3144	6218	10.83%	1.729%
29	11.325	1554	1559	1562	rVB2	638	1089	1.90%	0.303%
30	11.490	1579	1586	1593	rBV	31811	50746	88.40%	14.108%
31	12.483	1743	1749	1756	rBV2	23864	36971	64.40%	10.278%
32	13.422	1898	1903	1910	rVB	26433	41342	72.01%	11.494%
33	13.697	1945	1948	1952	rVB2	720	868	1.51%	0.241%
34	13.965	1988	1992	1995	rVB3	579	888	1.55%	0.247%
35	14.483	2073	2077	2081	rBV4	602	1113	1.94%	0.309%
36	16.062	2333	2336	2339	rBV3	577	818	1.42%	0.227%

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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_Y\methods\82Y051722S.M
Title : SW846 8260

Sum of corrected areas: 359694

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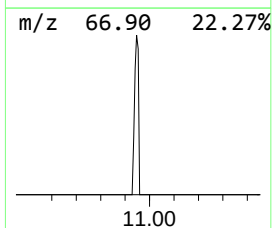
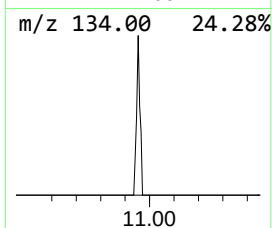
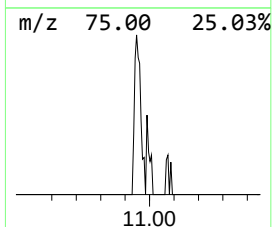
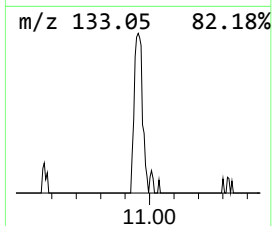
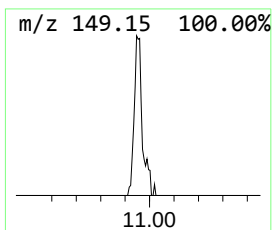
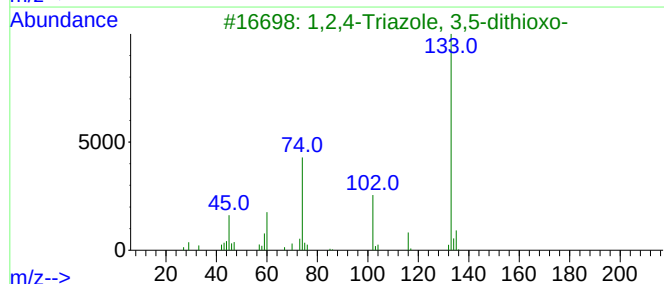
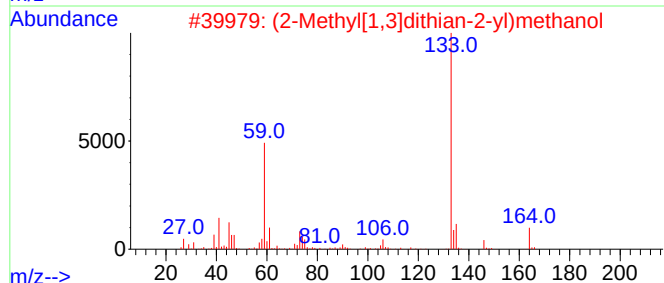
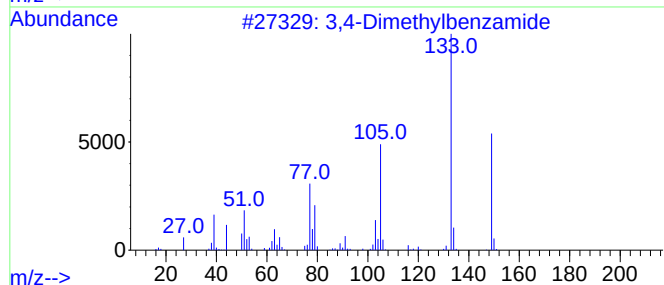
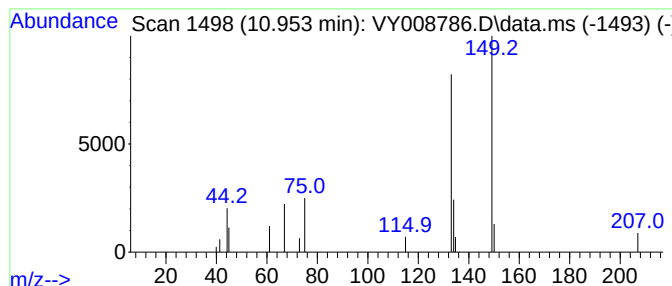
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y051722S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown10.953 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.953	6.13 ug/l	6218	Chlorobenzene-d5	11.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dimethylbenzamide	149	C9H11NO	005580-33-6	17
2			(2-Methyl[1,3]dithian-2-yl)methanol	164	C6H12OS2	1000191-37-4	10
3			1,2,4-Triazole, 3,5-dithio-	133	C2H3N3S2	005650-03-3	10
4			Benzothiazole, 2-methyl-	149	C8H7NS	000120-75-2	9
5			Pyridine, pentamethyl-	149	C10H15N	003748-83-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown10.953	10.953	6.1	ug/l	6218	3	11.490	50746	50.0