

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW100219\
 Data File : VW013400.D
 Acq On : 03 Oct 2019 02:27
 Operator : SY/VA
 Sample : VW1002SBL02
 Misc : 5.00G/5ML/MSVOA W/SOIL
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 VW1002SBL02

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.367	231	240	252	rVB2	12023	30667	2.41%	0.346%
2	4.903	479	492	493	rBV3	12246	26582	2.09%	0.300%
3	5.117	511	527	544	rBV	42069	155914	12.27%	1.758%
4	5.745	616	630	641	rBV5	7577	25492	2.01%	0.287%
5	5.921	649	659	670	rBV3	7469	22919	1.80%	0.258%
6	7.147	848	860	872	rBV	57524	176027	13.85%	1.985%
7	7.878	971	980	985	rBV	164615	390732	30.75%	4.406%
8	7.945	985	991	1000	rVB	348095	774948	60.99%	8.738%
9	8.250	1031	1041	1046	rBV	339675	865940	68.15%	9.764%
10	8.299	1047	1049	1061	rVB	162946	290916	22.89%	3.280%
11	8.842	1127	1138	1148	rBV	467656	943913	74.28%	10.643%
12	9.756	1279	1288	1297	rBV2	18794	36619	2.88%	0.413%
13	9.890	1300	1310	1318	rBV	53016	110582	8.70%	1.247%
14	9.988	1318	1326	1332	rVB3	10940	24744	1.95%	0.279%
15	10.244	1358	1368	1373	rBV	27610	53237	4.19%	0.600%
16	10.323	1373	1381	1389	rVB	713038	1270680	100.00%	14.327%
17	10.701	1436	1443	1450	rBV	127174	235886	18.56%	2.660%
18	10.847	1461	1467	1472	rVB	10293	18148	1.43%	0.205%
19	11.061	1494	1502	1509	rBV	63909	110541	8.70%	1.246%
20	11.177	1515	1521	1527	rVB2	23042	38327	3.02%	0.432%
21	11.439	1552	1564	1574	rVB	92527	168850	13.29%	1.904%
22	11.628	1586	1595	1604	rBV	622566	1062386	83.61%	11.979%
23	12.469	1726	1733	1740	rBV2	18414	32335	2.54%	0.365%
24	12.615	1747	1757	1767	rBV2	562018	939089	73.90%	10.588%
25	12.932	1804	1809	1814	rBV3	10075	15688	1.23%	0.177%
26	13.292	1859	1868	1875	rVB7	7864	17590	1.38%	0.198%
27	13.554	1905	1911	1918	rBV	584883	948373	74.64%	10.693%
28	14.072	1991	1996	2003	rVB	23219	36865	2.90%	0.416%
29	15.438	2212	2220	2225	rVB2	26319	45053	3.55%	0.508%

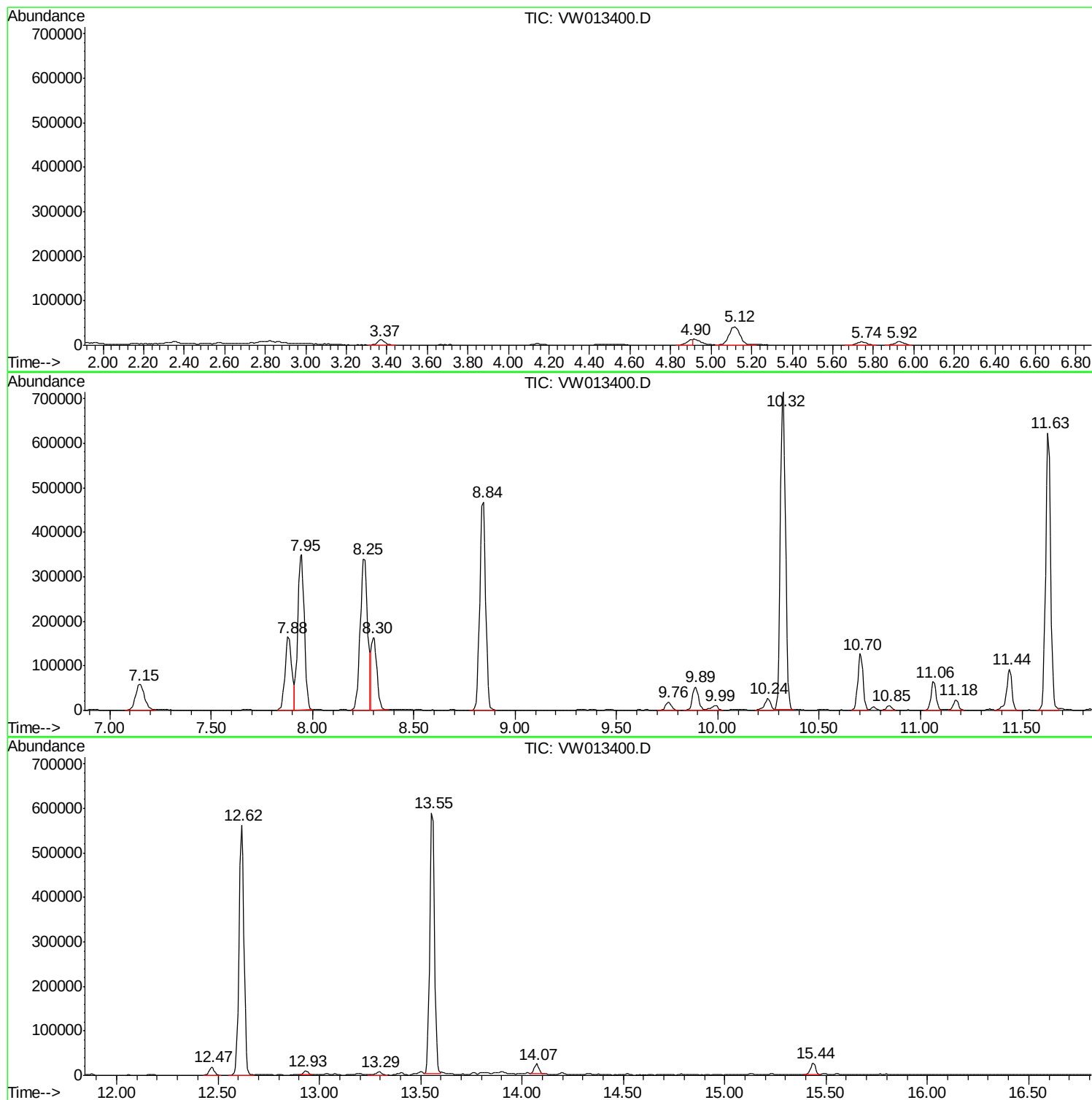
Sum of corrected areas: 8869043

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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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



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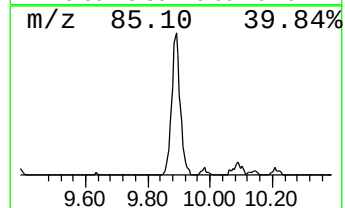
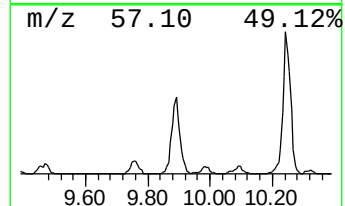
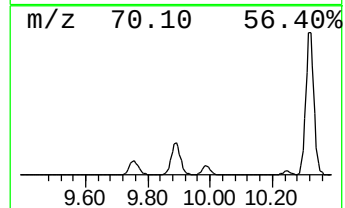
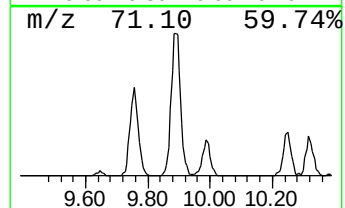
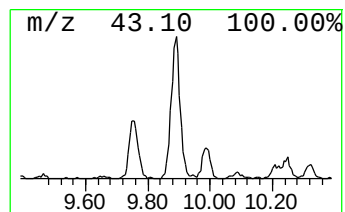
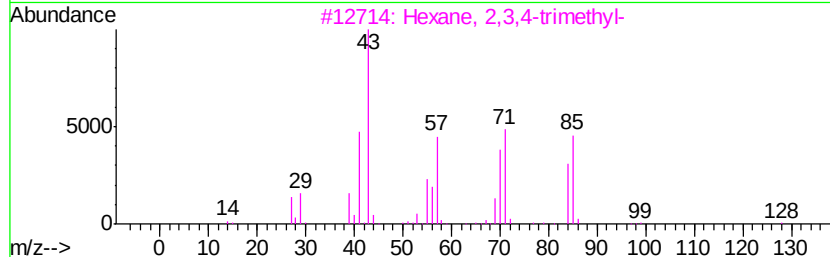
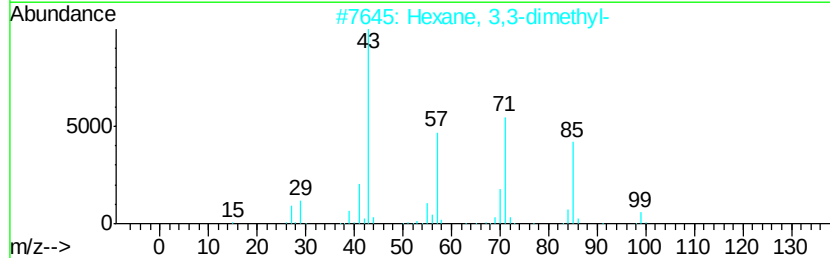
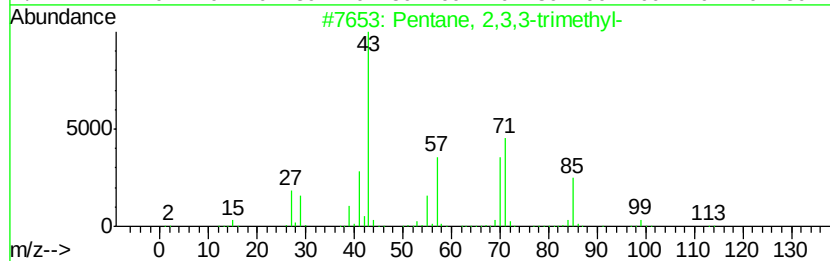
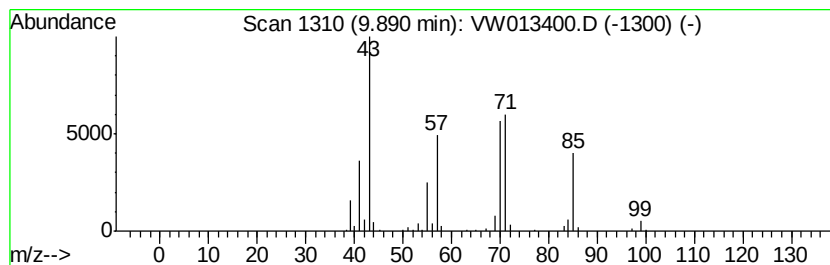
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	5.86 ug/l	110582	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	53
5		Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	53



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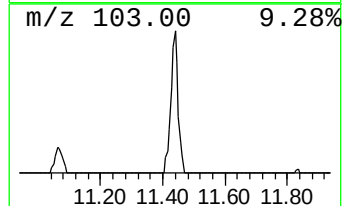
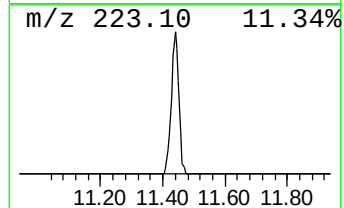
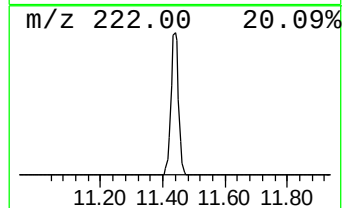
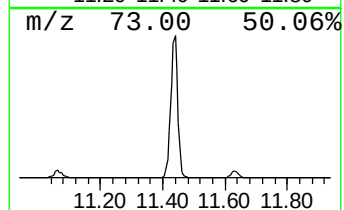
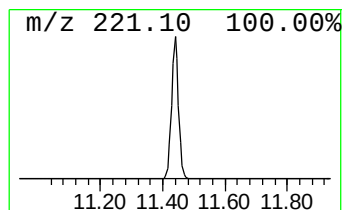
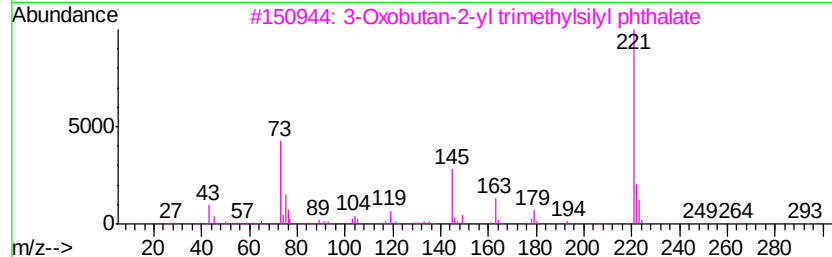
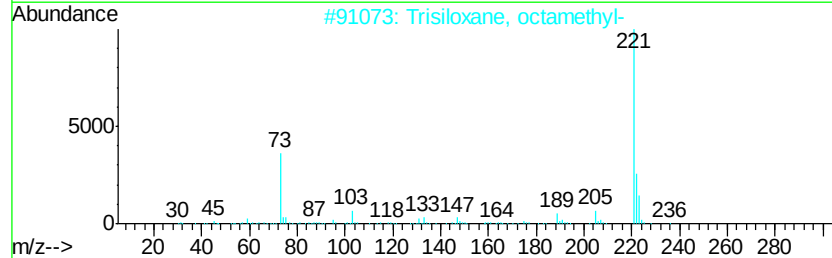
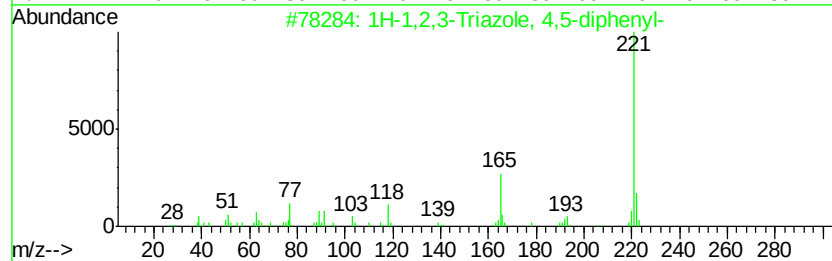
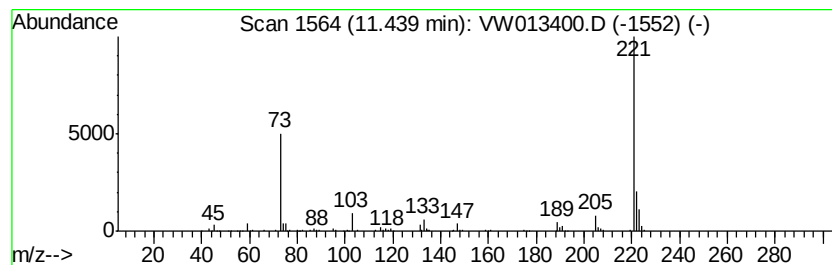
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Peak Number 6 1H-1,2,3-Triazole, 4,5-diph... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.44	7.95 ug/l	168850	Chlorobenzene-d5	11.63		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-1,2,3-Triazole, 4,5-diphenyl-	221	C14H11N3	005533-73-3	59
2		Trisiloxane, octamethyl-	236	C8H24O2Si3	000107-51-7	53
3		3-Oxobutan-2-yl trimethylsilyl p...	308	C15H20O5Si	1000373-49-5	50
4		2-Methyl-thieno[2',3':3,4]benzo[...	221	C10H7NOS2	294668-48-7	45
5		1-Aza-2-sila-5-boracyclopent-3-e...	221	C12H24BNSi	081620-70-4	40



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 2,3,3-tr...	9.89	5.9	ug/l	110582	2	8.84	943913	50.0
1H-1,2,3-Triazole...	11.44	8.0	ug/l	168850	3	11.63	1062390	50.0