

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU063018\
 Data File : VU025102.D
 Acq On : 30 Jun 2018 20:50
 Operator : MD/SY
 Sample : J3759-15
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WP021MW120-180626

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_U\METHOD\82U061318W.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	1.088	137	155	179	rBV	1290589	1449312	2.65%	1.321%
2	1.397	244	251	290	rBV2	37459121	54607219	100.00%	49.784%
3	4.230	1110	1132	1183	rBV	9441150	22811339	41.77%	20.796%
4	4.992	1353	1369	1390	rVB3	399106	979874	1.79%	0.893%
5	5.394	1480	1494	1515	rVB	981715	2078575	3.81%	1.895%
6	5.889	1632	1648	1680	rBV	382961	762463	1.40%	0.695%
7	7.567	2149	2170	2182	rBV	701475	1234660	2.26%	1.126%
8	7.641	2182	2193	2211	rVB	732543	1250842	2.29%	1.140%
9	9.124	2632	2654	2687	rBV	11921917	20437573	37.43%	18.632%
10	10.310	3011	3023	3040	rBV	562764	903136	1.65%	0.823%
11	11.487	3378	3389	3407	rBV2	686325	1667777	3.05%	1.520%
12	11.882	3498	3512	3534	rVB	541262	906186	1.66%	0.826%
13	12.358	3649	3660	3682	rBV3	259220	599634	1.10%	0.547%

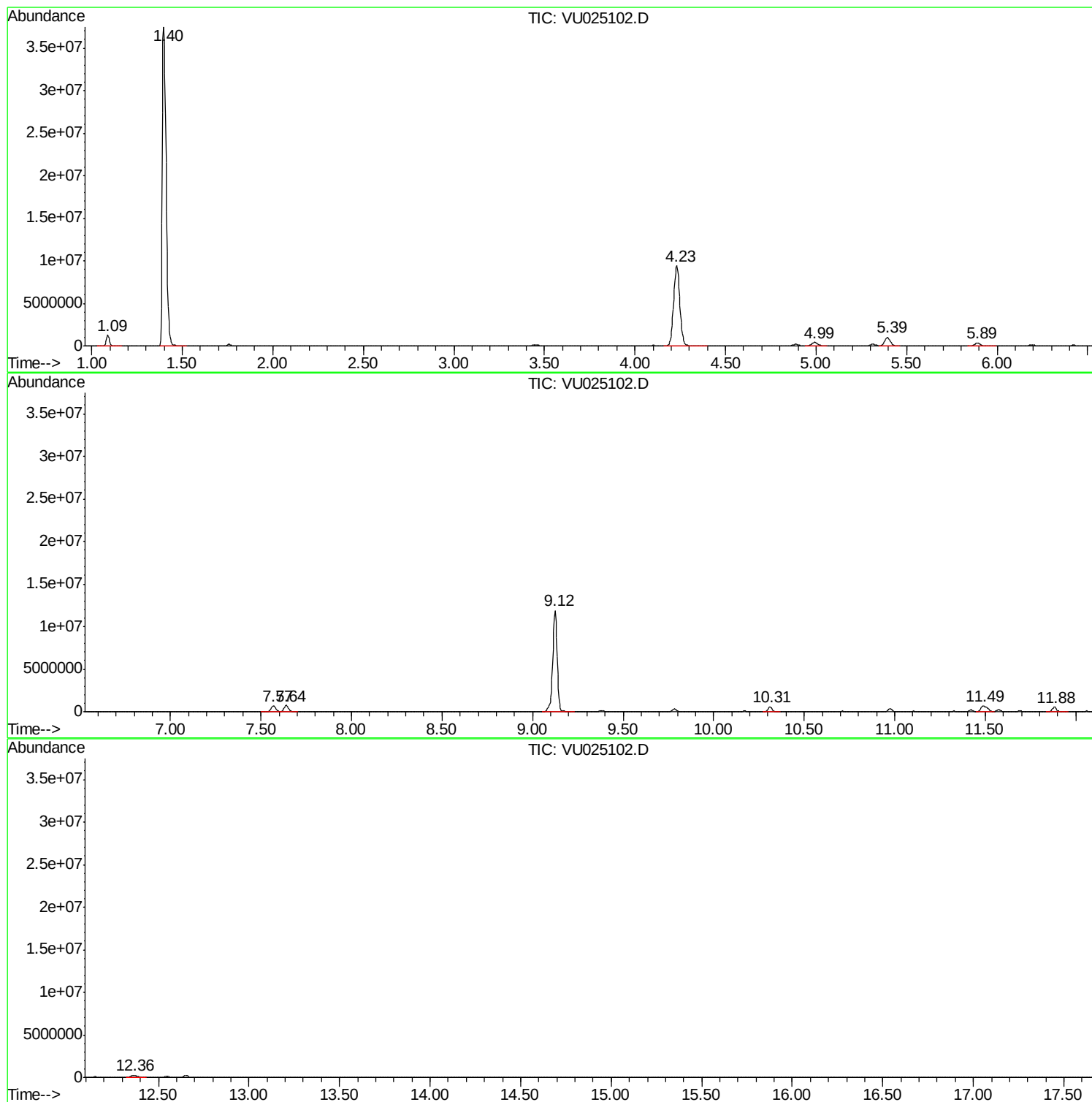
Sum of corrected areas: 109688590

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



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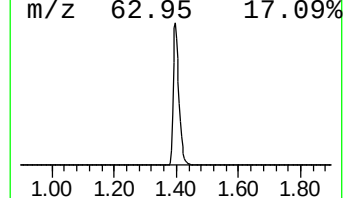
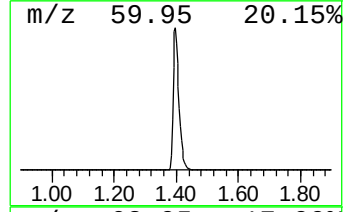
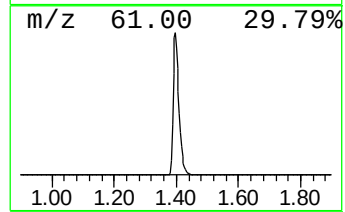
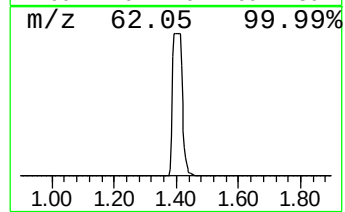
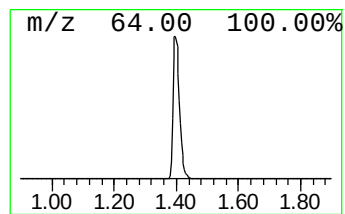
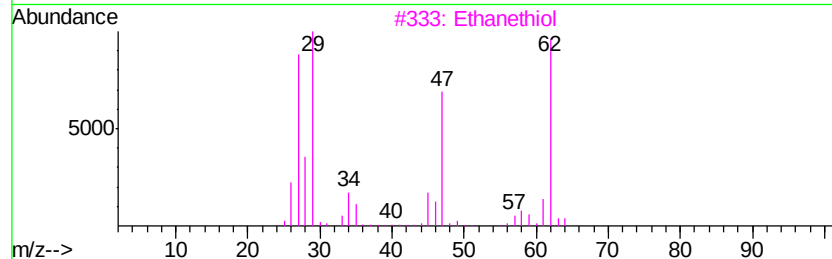
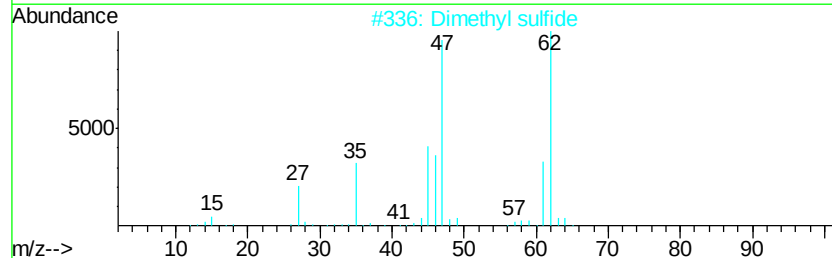
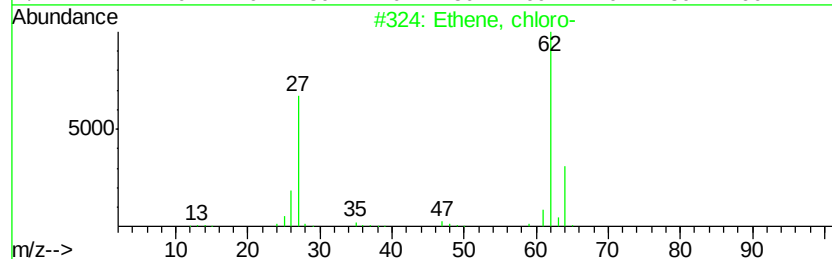
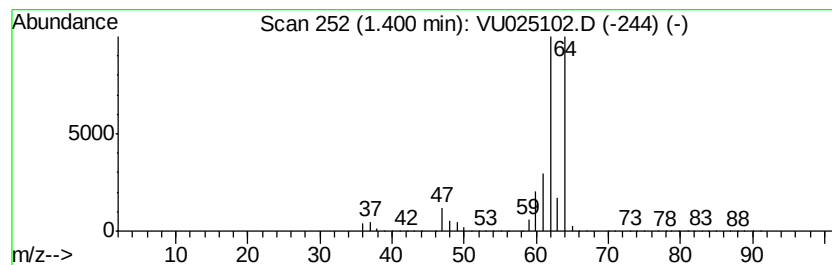
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Peak Number 2 unknown1.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.40	2786.44 ug/l	54607200	Pentafluorobenzene	4.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethene, chloro-	62	C2H3Cl	000075-01-4	23
2		Dimethyl sulfide	62	C2H6S	000075-18-3	9
3		Ethanethiol	62	C2H6S	000075-08-1	7
4		Phosphine, ethyl-	62	C2H7P	000593-68-0	7
5		Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	7



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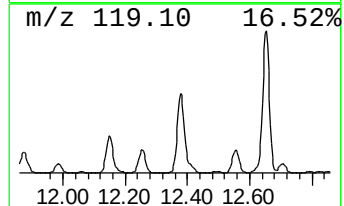
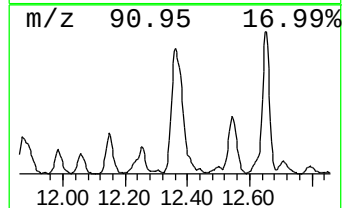
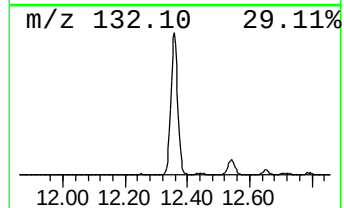
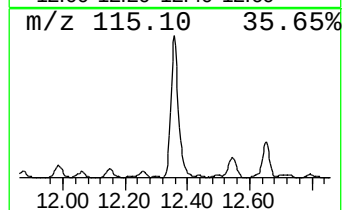
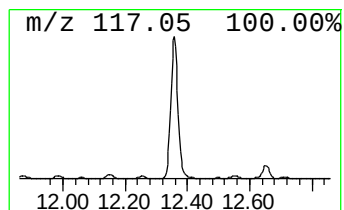
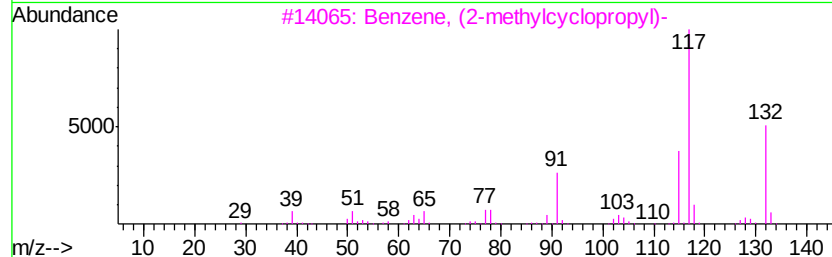
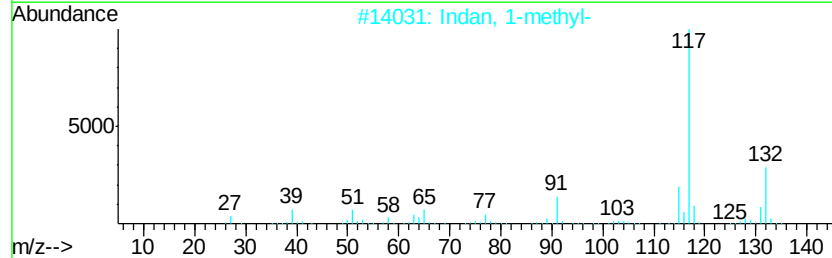
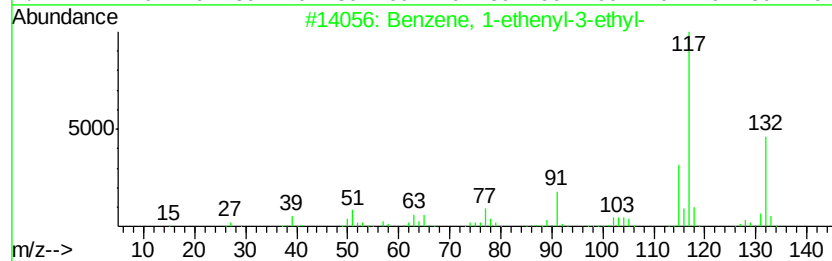
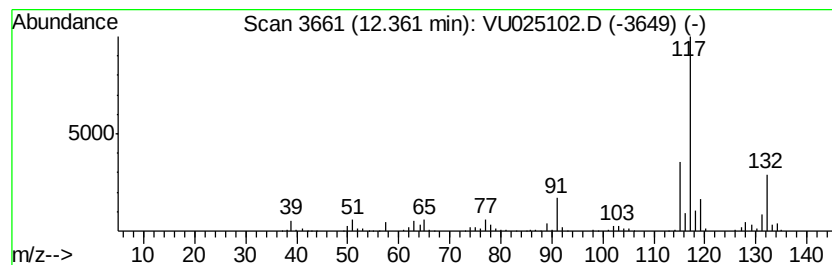
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Peak Number 3 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	17.98 ug/l	599634	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 90
2		Indan, 1-methyl-	132 C10H12	000767-58-8 90
3		Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0 87
4		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 86
5		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 83



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.40	1.40	2786.4	ug/l	54607200	1	4.99	979874	50.0
Benzene, 1-etheny...	12.36	18.0	ug/l	599634	4	11.49	1667780	50.0