

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092419\
 Data File : VX012640.D
 Acq On : 24 Sep 2019 18:04
 Operator : JC/SP
 Sample : K5019-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 QNWP8-MW30S-GW

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_X\METHOD\82X092319W.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	1.070	26	30	42	rBV	1220049	1435862	100.00%	23.569%
2	1.265	59	62	70	rBV	24970	31729	2.21%	0.521%
3	2.594	273	280	295	rBV	38674	74257	5.17%	1.219%
4	5.490	744	755	768	rBV	97984	281549	19.61%	4.621%
5	5.655	768	782	799	rVV	187801	524927	36.56%	8.616%
6	6.057	835	848	861	rBV	120458	323776	22.55%	5.315%
7	6.850	967	978	994	rBV	287366	678401	47.25%	11.136%
8	8.709	1276	1283	1301	rBV	563613	893249	62.21%	14.662%
9	10.111	1506	1513	1527	rBV	517621	720880	50.21%	11.833%
10	11.135	1675	1681	1694	rBV	369004	485051	33.78%	7.962%
11	12.074	1829	1835	1844	rBV	430948	526941	36.70%	8.649%
12	12.940	1972	1977	1985	rVB	12056	16331	1.14%	0.268%
13	13.806	2115	2119	2122	rBV2	12889	16604	1.16%	0.273%
14	16.409	2540	2546	2556	rVB2	44757	82601	5.75%	1.356%

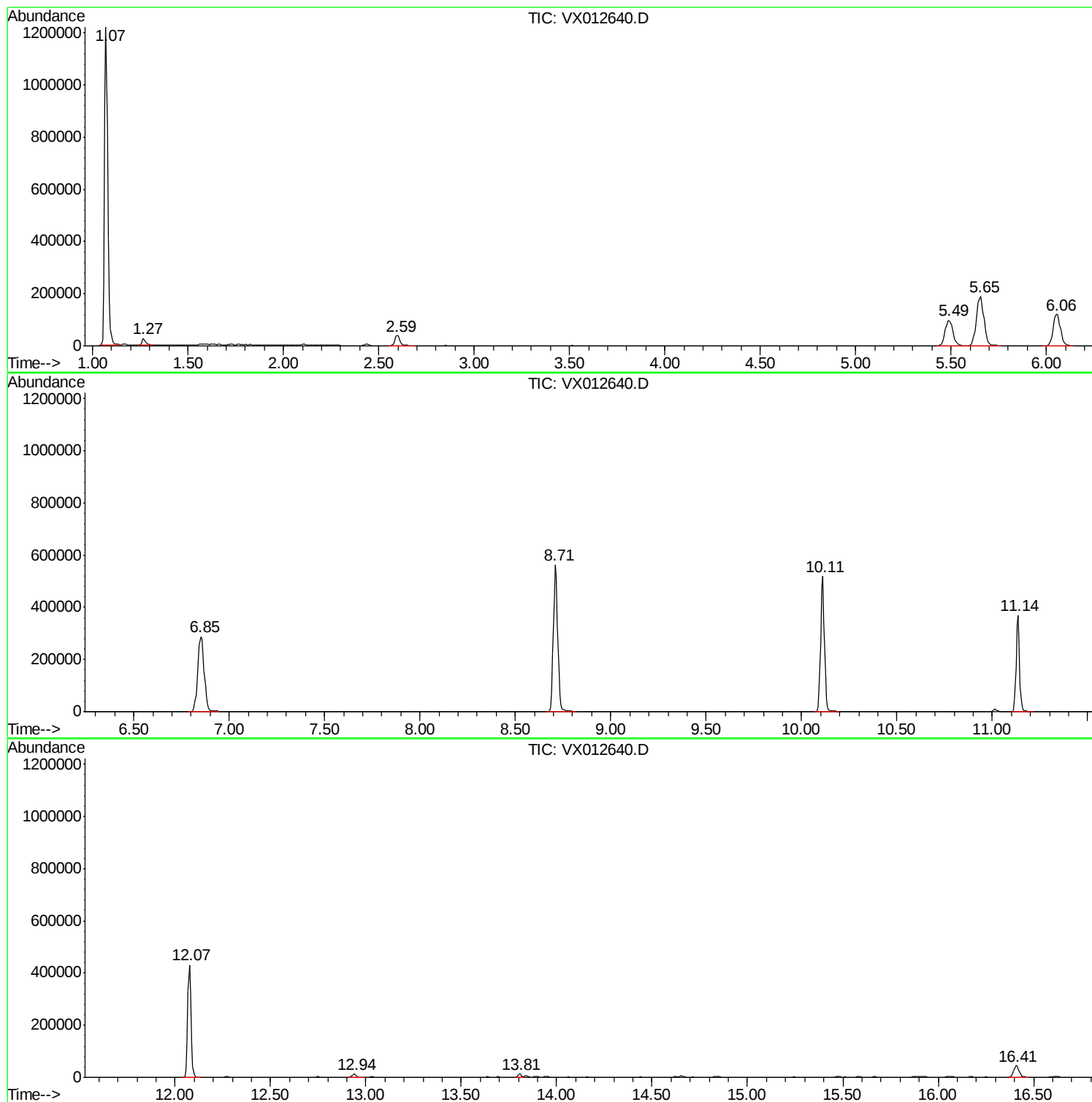
Sum of corrected areas: 6092158

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



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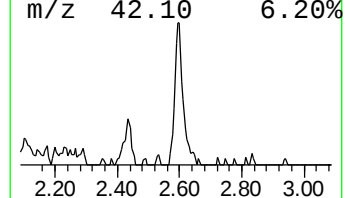
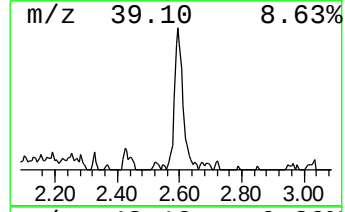
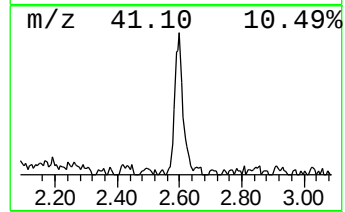
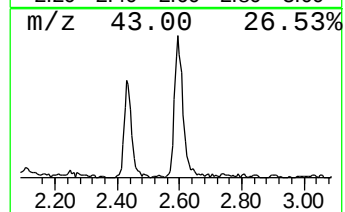
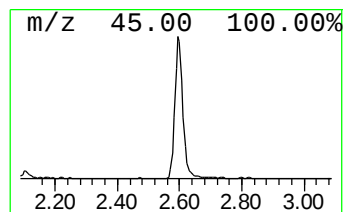
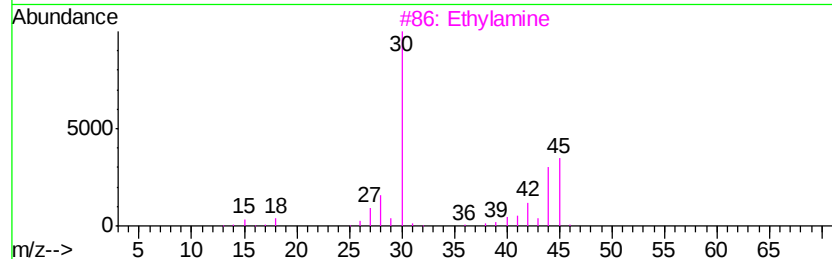
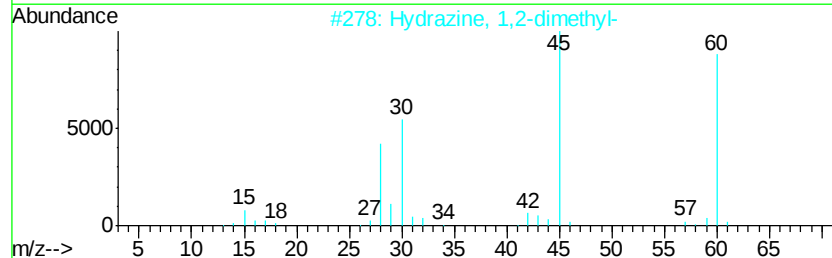
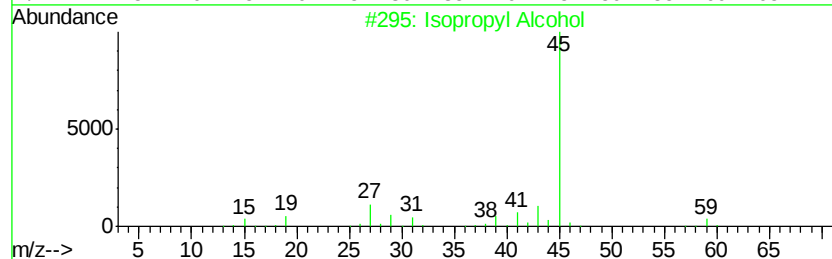
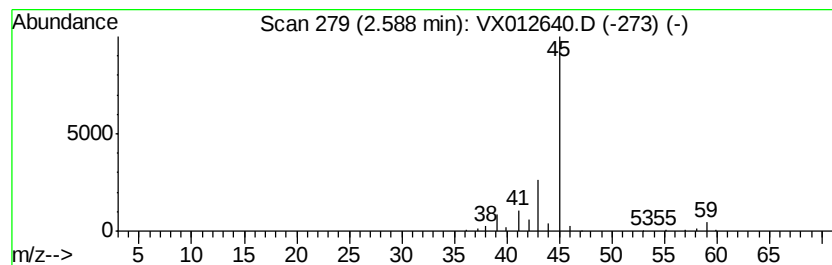
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Peak Number 2 Isopropyl Alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.59	7.07 ug/l	74257	Pentafluorobenzene	5.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	83
2			Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3			Ethylamine	45	C2H7N	000075-04-7	5
4			2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	4
5			Formamide	45	CH3NO	000075-12-7	4



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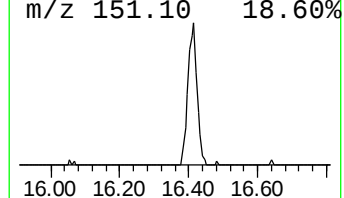
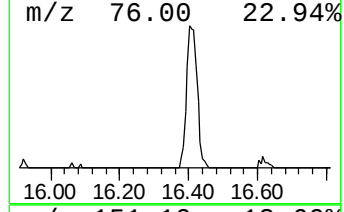
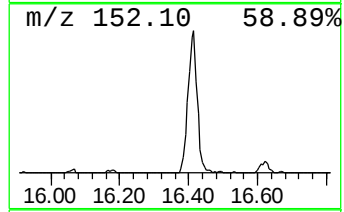
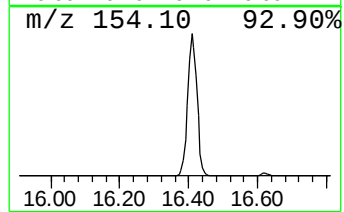
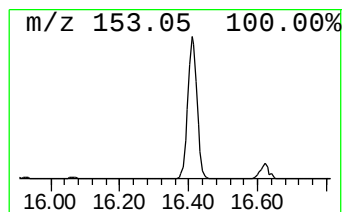
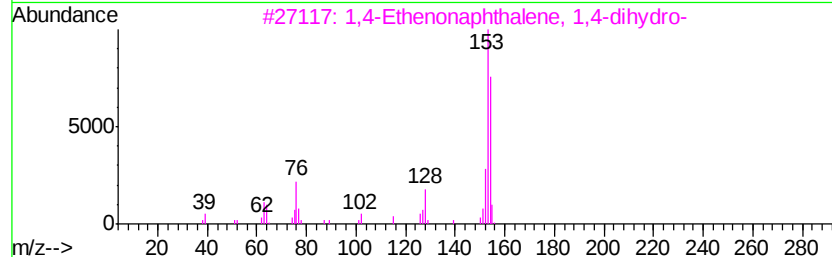
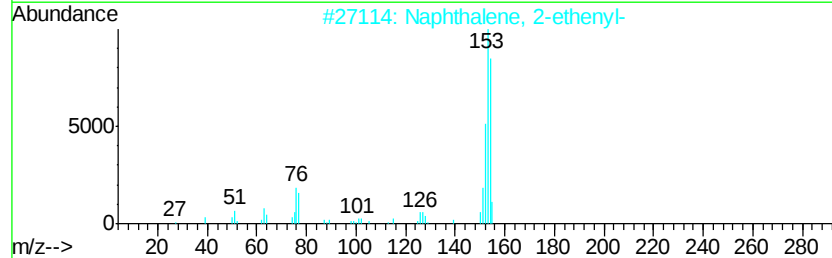
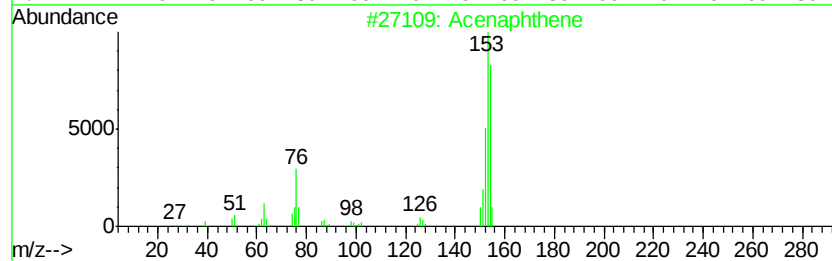
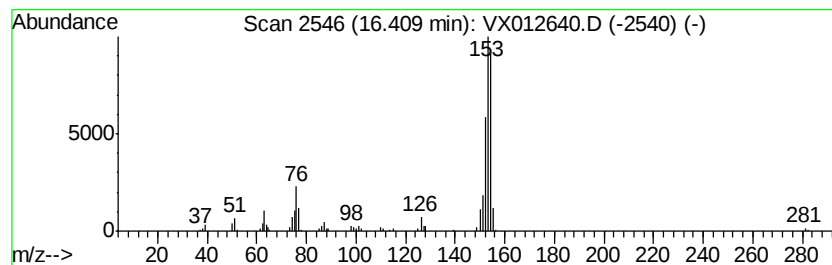
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Peak Number 3 Acenaphthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.41	7.84 ug/l	82601	1,4-Dichlorobenzene-d4	12.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	95
2			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	76
3			1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	59
4			.beta.-(1-Naphthyl)acrylic acid	198	C13H10O2	013026-12-5	50
5			Acenaphthylene, 5-bromo-1,2-dihy...	232	C12H9Br	002051-98-1	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl Alcohol	2.59	7.1	ug/l	74257	1	5.65	524927	50.0
Acenaphthene	16.41	7.8	ug/l	82601	4	12.07	526941	50.0