

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073118\
 Data File : VX003781.D
 Acq On : 01 Aug 2018 05:59
 Operator : JC/SP
 Sample : J4143-06
 Misc : 7.05G/5.0mL/MSVOA X/SOIL
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GW-01(10-12)

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\82X073018S.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	1.069	75	79	92	rBV	1685891	1979397	100.00%	48.338%
2	1.166	92	95	101	rVB	34945	33933	1.71%	0.829%
3	1.386	127	131	139	rVB3	13294	23319	1.18%	0.569%
4	2.446	298	305	315	rVB	109916	188337	9.51%	4.599%
5	2.568	318	325	331	rBV	22063	38407	1.94%	0.938%
6	4.702	665	675	695	rBV2	18731	76817	3.88%	1.876%
7	5.501	796	806	818	rBV2	46546	124128	6.27%	3.031%
8	5.659	822	832	851	rVB	111606	310678	15.70%	7.587%
9	6.068	890	899	909	rBV	43330	113083	5.71%	2.762%
10	6.860	1019	1029	1042	rBV	143009	331725	16.76%	8.101%
11	8.720	1327	1334	1341	rVB	246219	375750	18.98%	9.176%
12	10.116	1557	1563	1570	rBV	184157	241675	12.21%	5.902%
13	11.140	1726	1731	1736	rVB	94368	115703	5.85%	2.826%
14	12.079	1880	1885	1891	rBV	100852	120731	6.10%	2.948%
15	16.420	2591	2597	2605	rVB2	11578	21202	1.07%	0.518%

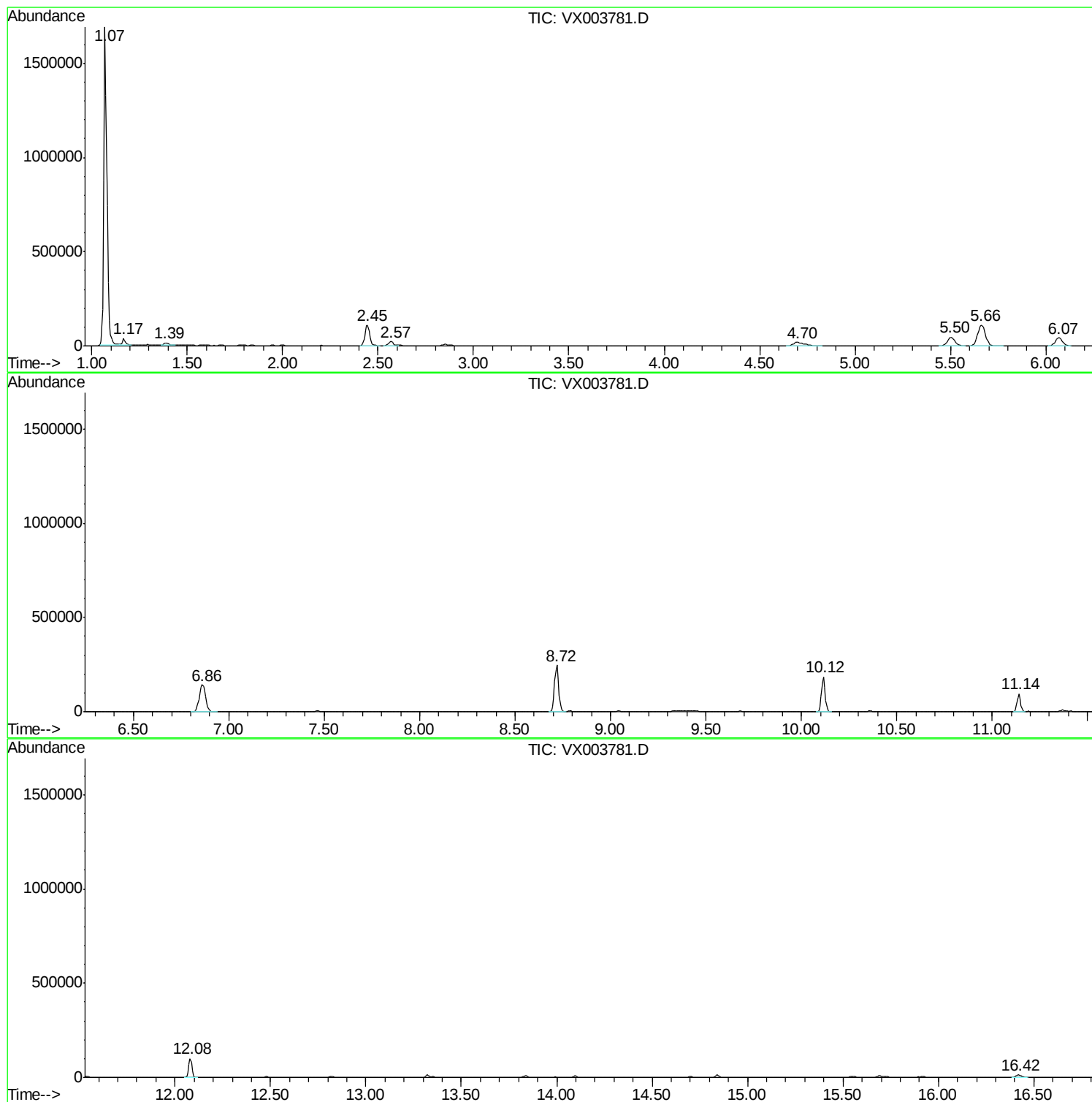
Sum of corrected areas: 4094885

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



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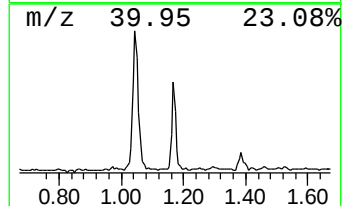
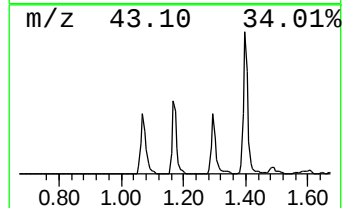
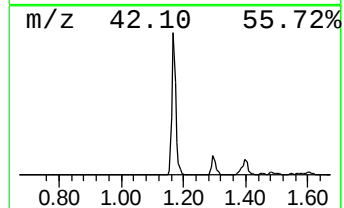
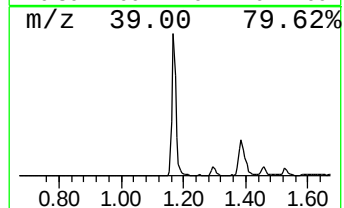
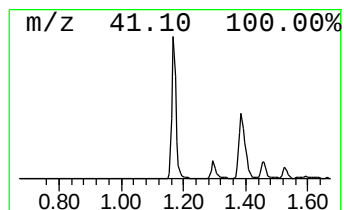
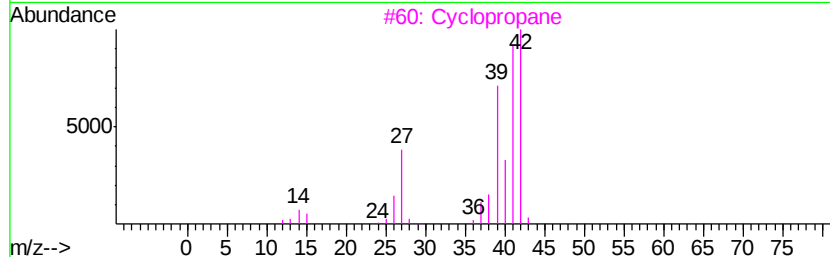
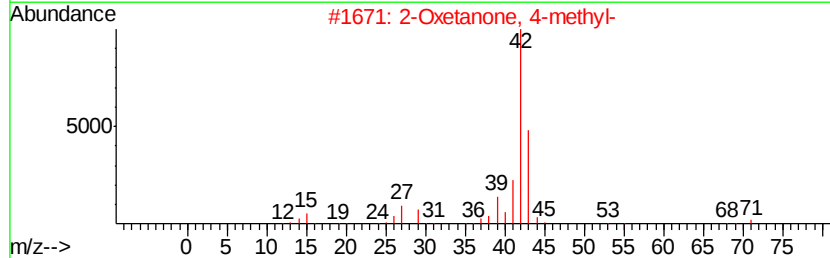
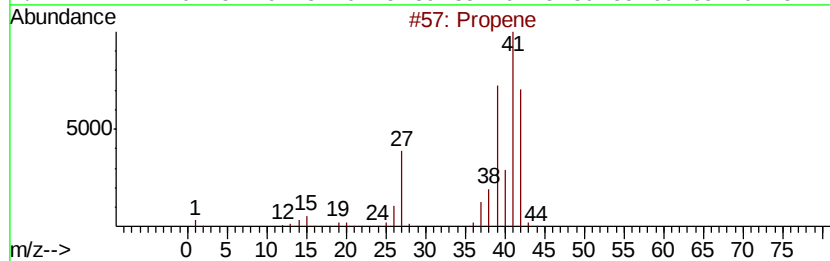
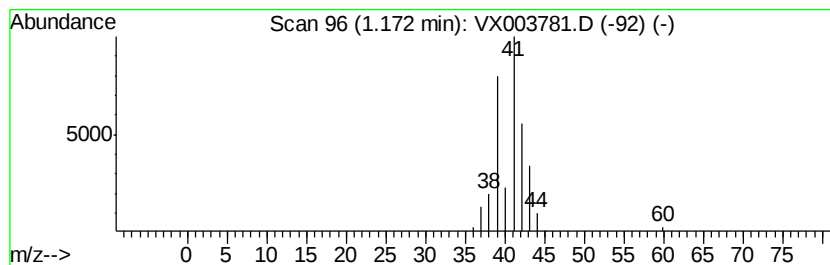
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Peak Number 2 Propene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.17	5.46 ug/l	33933	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propene	42	C3H6	000115-07-1	90
2		2-Oxetanone, 4-methyl-	86	C4H6O2	003068-88-0	9
3		Cyclopropane	42	C3H6	000075-19-4	9
4		Cyclopropene	40	C3H4	002781-85-3	9
5		4-Amino-1-butanol	89	C4H11NO	013325-10-5	2



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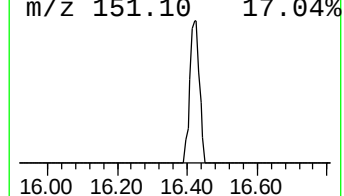
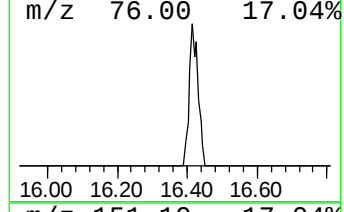
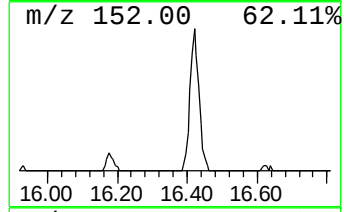
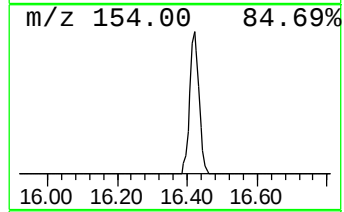
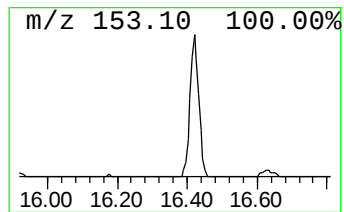
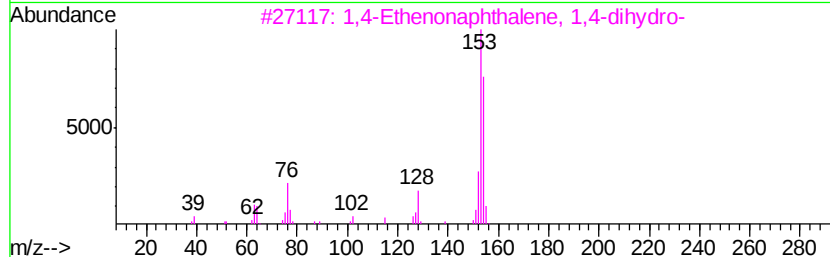
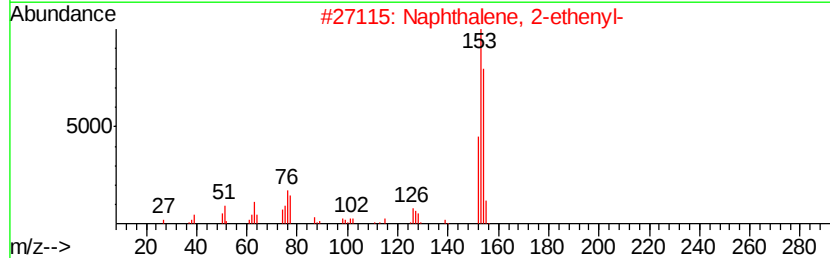
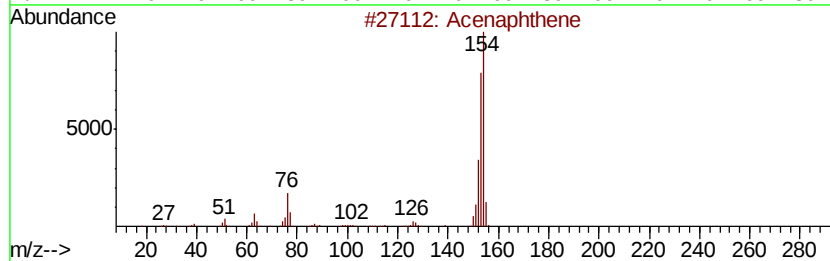
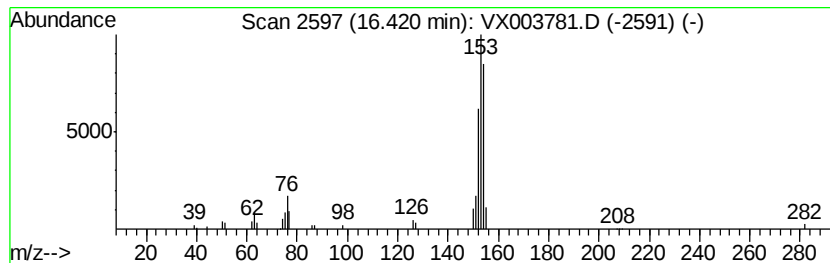
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Peak Number 3 Acenaphthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.42	8.78 ug/l	21202	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acenaphthene	154	C12H10	000083-32-9	81
2		Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	76
3		1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	59
4		3-Fluoro-para-anisaldehyde	154	C8H7FO2	000351-54-2	53
5		Biphenyl	154	C12H10	000092-52-4	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.17	5.5	ug/l	33933	1	5.66	310678	50.0
Acenaphthene	16.42	8.8	ug/l	21202	4	12.08	120731	50.0