

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111718\
 Data File : VX006067.D
 Acq On : 17 Nov 2018 14:39
 Operator : JC/MD
 Sample : J5984-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 STORAGE-BLANK-SAMPLE-MANA

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_X\METHOD\82X110718W.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	1.282	18	21	27	rVB	10801	11377	1.92%	0.292%
2	1.629	71	78	82	rVB5	3922	8425	1.42%	0.216%
3	5.494	699	712	727	rBV	64043	180900	30.49%	4.649%
4	5.659	727	739	753	rBV	110584	307604	51.85%	7.905%
5	6.061	794	805	819	rBV2	70291	184956	31.17%	4.753%
6	6.854	921	935	949	rBV	159418	380428	64.12%	9.776%
7	8.713	1233	1240	1257	rBV	328537	515518	86.89%	13.247%
8	9.414	1344	1355	1368	rBV	186192	456546	76.95%	11.732%
9	10.109	1463	1469	1479	rBV	335740	456129	76.88%	11.721%
10	11.134	1632	1637	1652	rBV	253253	339615	57.24%	8.727%
11	11.371	1669	1676	1688	rVB	384382	593301	100.00%	15.246%
12	11.676	1709	1726	1737	rVB4	7970	33276	5.61%	0.855%
13	12.079	1786	1792	1805	rBV	345985	423425	71.37%	10.881%

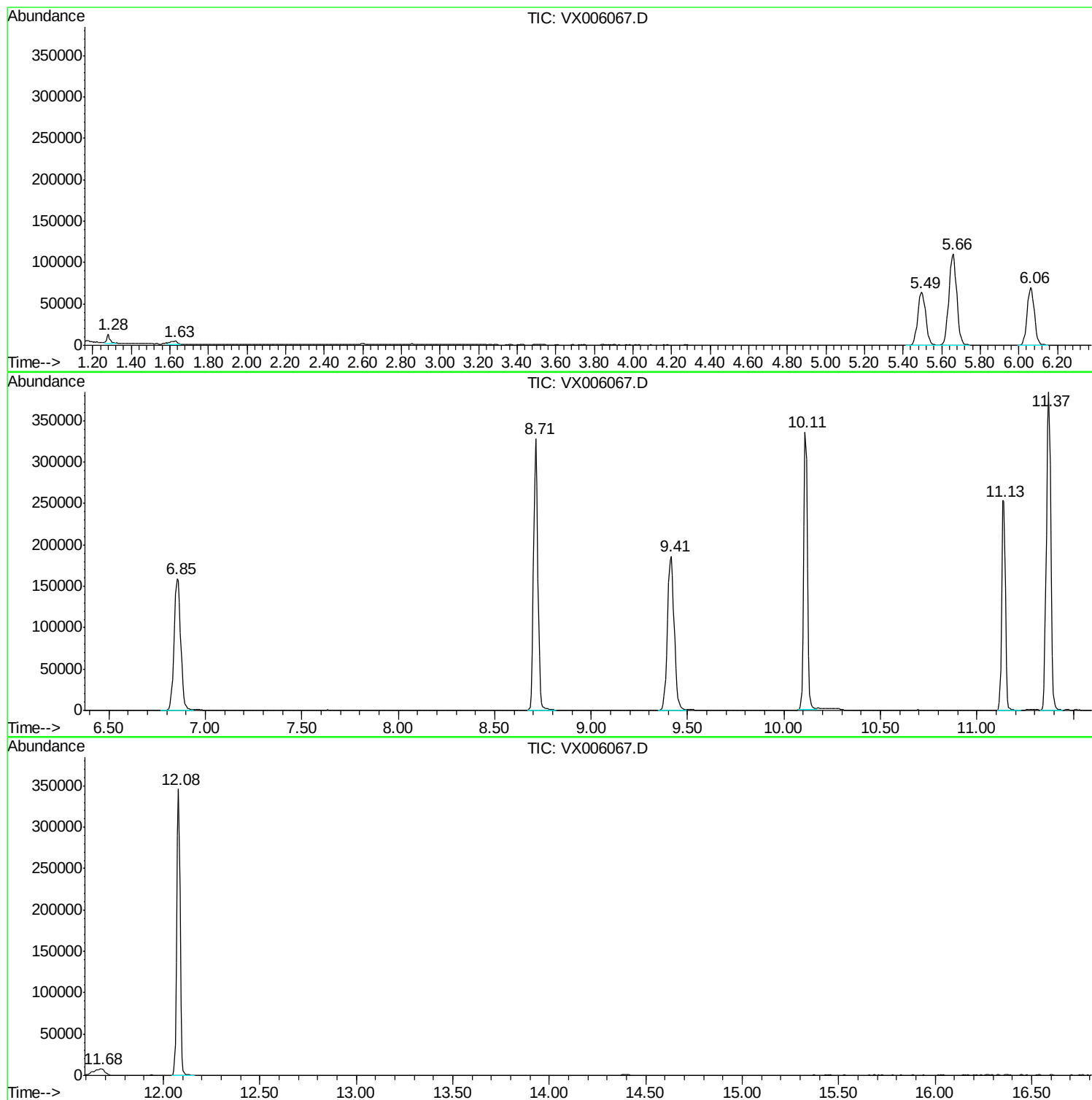
Sum of corrected areas: 3891500

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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



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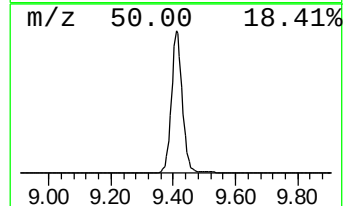
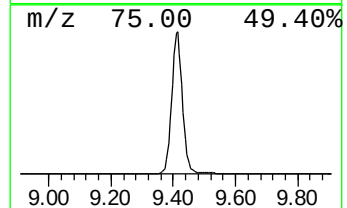
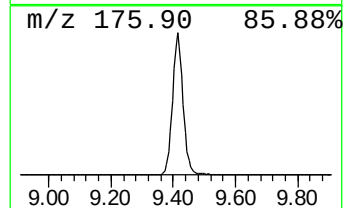
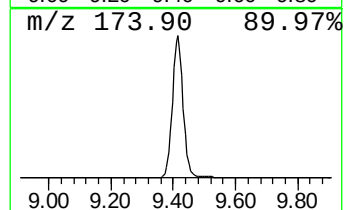
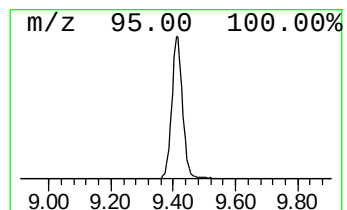
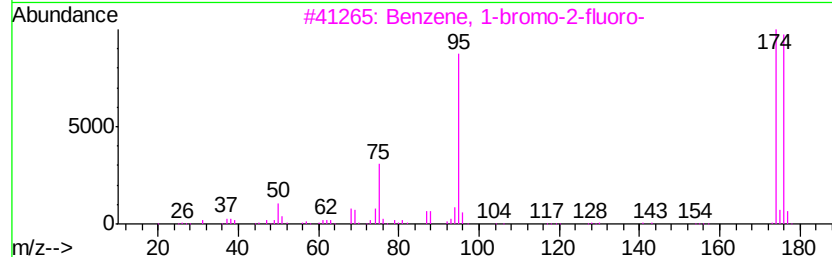
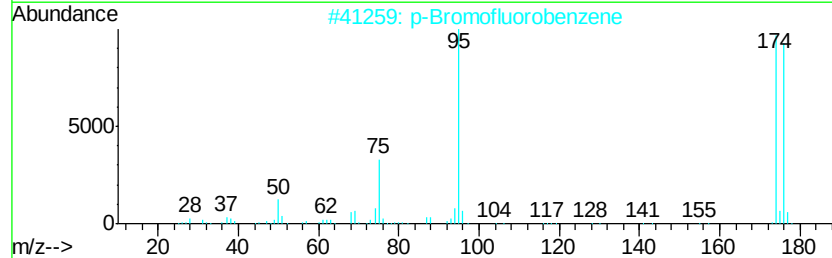
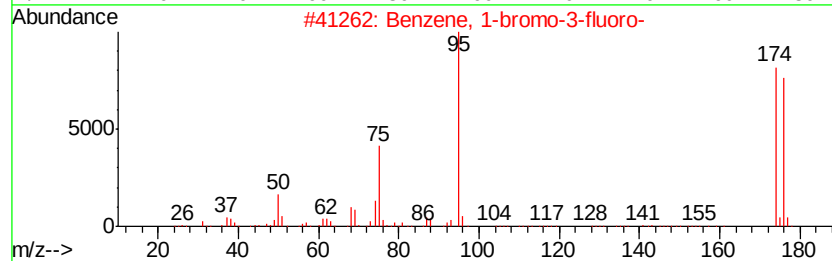
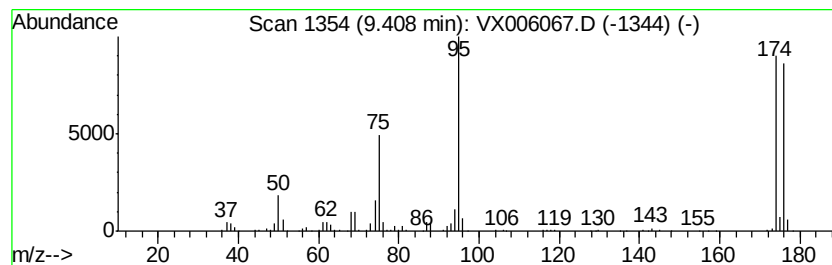
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-bromo-3-fluoro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.41	50.05 ug/l	456546	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-fluoro-	174	C6H4BrF	001073-06-9	98
2		p-Bromofluorobenzene	174	C6H4BrF	000460-00-4	94
3		Benzene, 1-bromo-2-fluoro-	174	C6H4BrF	001072-85-1	94
4		Pyrazole, 3-(2-bromoethyl)-	174	C5H7BrN2	211738-72-6	45
5		4(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	019808-30-1	37



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-bromo-...	9.41	50.0	ug/l	456546	3	10.11	456129	50.0