

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX121418\
 Data File : VX006521.D
 Acq On : 14 Dec 2018 16:58
 Operator : JC/MD
 Sample : J6396-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-15B_R1

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\82X112918W.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	1.617	62	76	79	rBV	45560	109156	3.37%	0.456%
2	2.452	179	213	214	rBV3	180341	1107456	34.23%	4.631%
3	3.050	301	311	318	rBV2	58265	146483	4.53%	0.612%
4	3.202	327	336	345	rBV	121424	295261	9.13%	1.235%
5	5.501	702	713	727	rBV2	375590	1068712	33.03%	4.469%
6	5.659	727	739	753	rVV	668862	1878534	58.06%	7.855%
7	6.068	794	806	824	rVB	397908	1049181	32.42%	4.387%
8	6.860	925	936	948	rBV	1005768	2349449	72.61%	9.824%
9	8.713	1233	1240	1255	rBV	2158976	3235717	100.00%	13.530%
10	9.829	1414	1423	1433	rBV	1133618	2187171	67.59%	9.145%
11	10.109	1463	1469	1484	rVB	2053273	2825350	87.32%	11.814%
12	11.134	1631	1637	1646	rBV	1689088	2233325	69.02%	9.338%
13	11.512	1693	1699	1712	rVB	1847490	2680316	82.84%	11.207%
14	12.079	1786	1792	1806	rBV	2255632	2749534	84.97%	11.497%

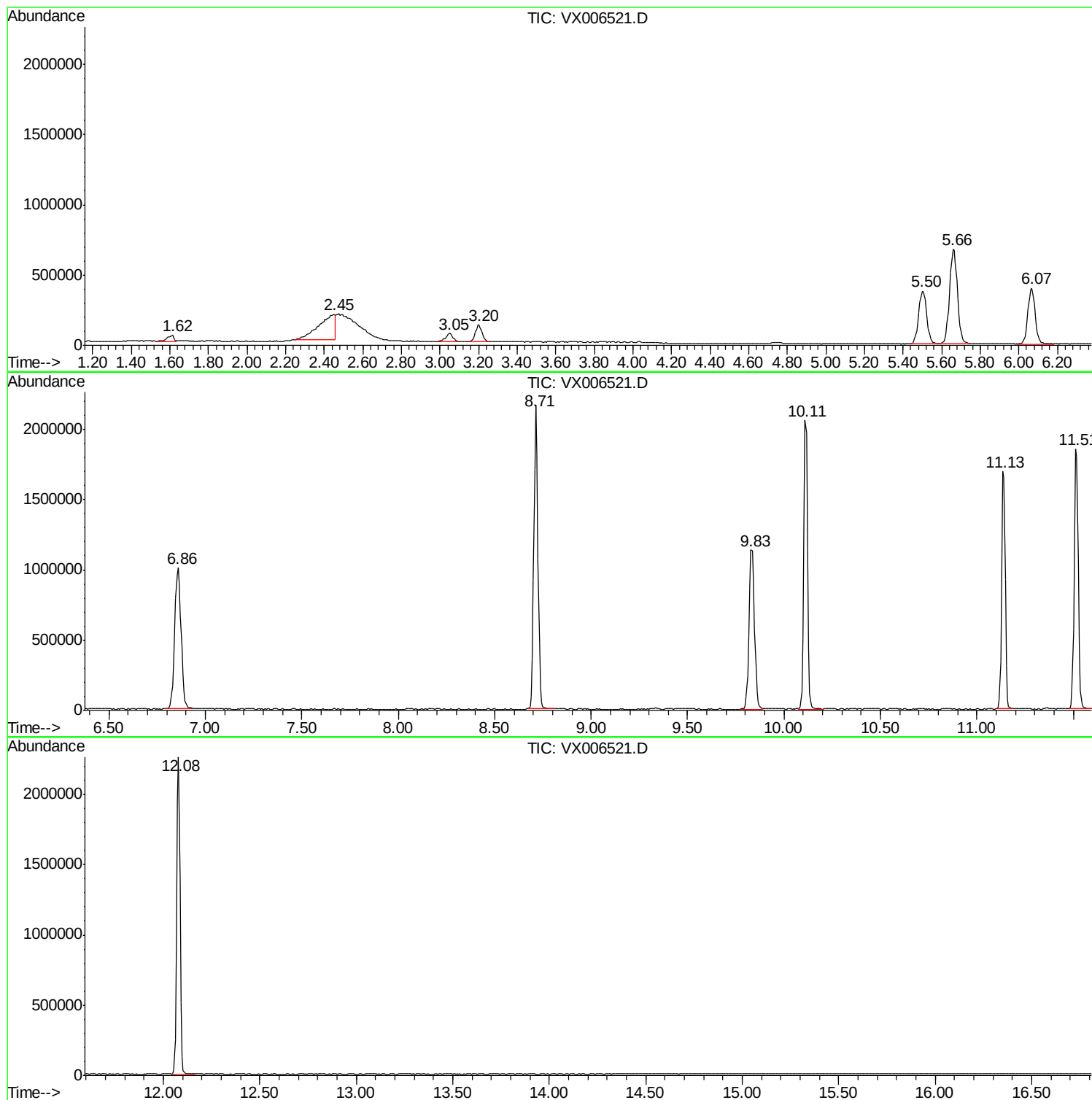
Sum of corrected areas: 23915645

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112918W.M
Quant Title : SW846 8260

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



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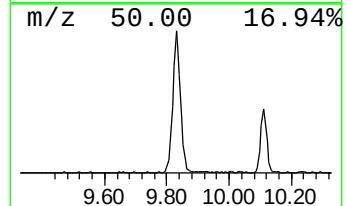
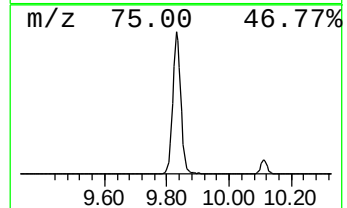
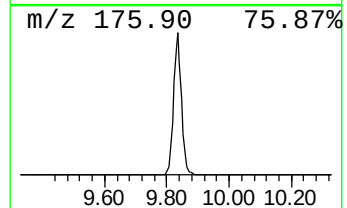
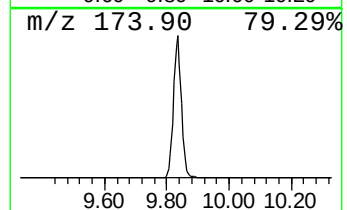
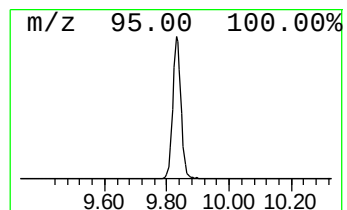
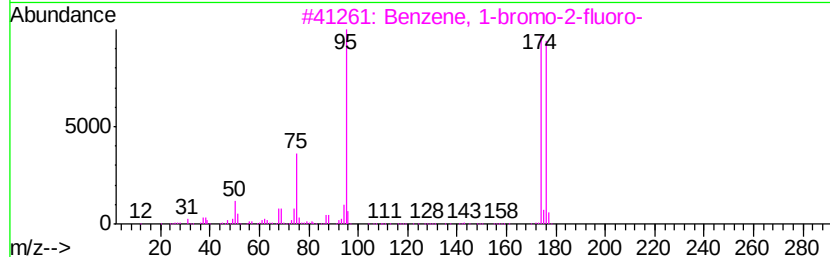
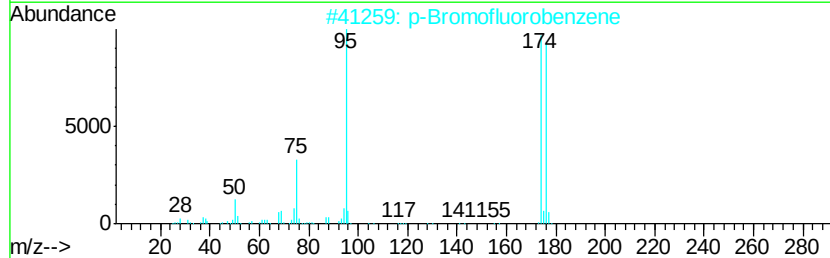
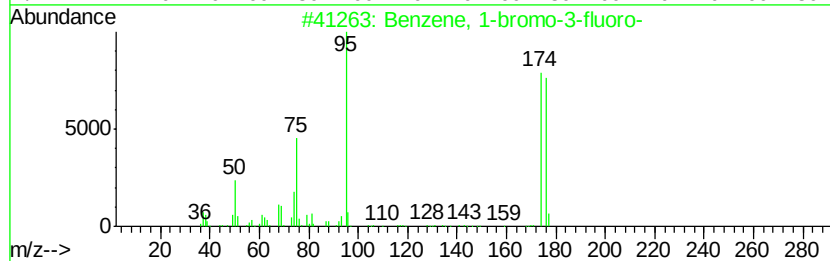
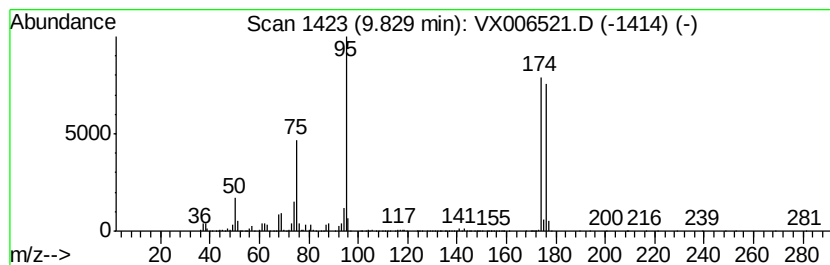
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Quant Title : SW846 8260

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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.
9.83	38.71 ug/l	2187170	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-fluoro-	174	C6H4BrF	001073-06-9	95
2		p-Bromofluorobenzene	174	C6H4BrF	000460-00-4	94
3		Benzene, 1-bromo-2-fluoro-	174	C6H4BrF	001072-85-1	93
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-...	9.83	38.7	ug/l	2187170	3	10.12	2825350	50.0