

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD081618\
Data File : VD059813.D
Acq On : 16 Aug 2018 16:22
Operator : VA/AP
Sample : J4460-03
Misc : 5.02g/5ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
RA-081018-SW-003

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_D\METHOD\82D081518S.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	1.211	96	100	103	rBV	138887	317349	41.41%	6.576%
2	1.270	103	106	124	rVB	92797	298326	38.93%	6.182%
3	3.050	283	287	292	rVB2	10256	21127	2.76%	0.438%
4	4.821	461	467	474	rBV2	7523	24206	3.16%	0.502%
5	5.529	533	539	543	rBV	171974	460822	60.13%	9.549%
6	5.618	543	548	562	rVB	283275	766383	100.00%	15.881%
7	5.992	580	586	598	rBB	142367	369713	48.24%	7.661%
8	6.671	649	655	664	rBV	309903	756563	98.72%	15.678%
9	8.687	855	860	872	rBB	351736	760110	99.18%	15.751%
10	9.464	934	939	944	rBV	16031	42703	5.57%	0.885%
11	10.674	1058	1062	1071	rVB	377515	648289	84.59%	13.434%
12	13.872	1381	1387	1397	rBB	129485	336809	43.95%	6.979%
13	15.091	1500	1511	1516	rBV3	4797	23347	3.05%	0.484%

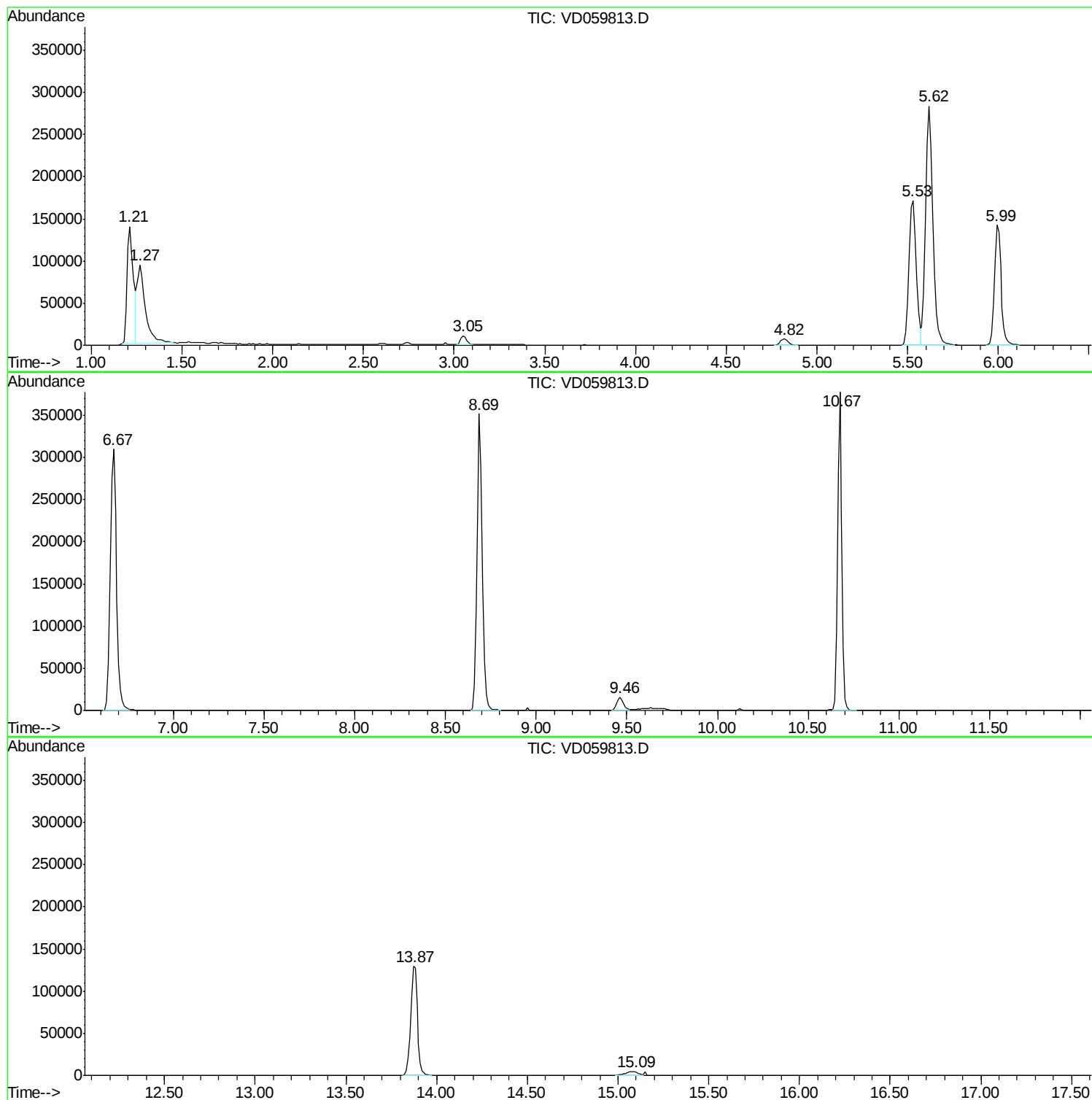
Sum of corrected areas: 4825747

Data Path : Z:\V0ASRV\HPCHEM1\MSV0A D\DATA\VD081618\
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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D081518S.M
Quant Title : SW846 8260

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



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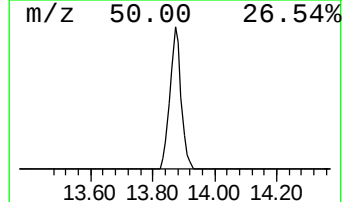
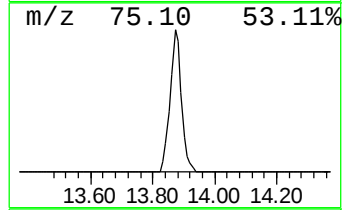
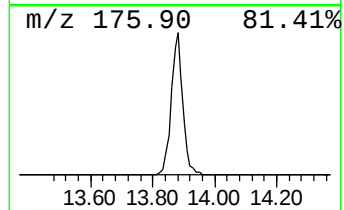
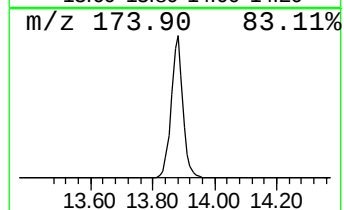
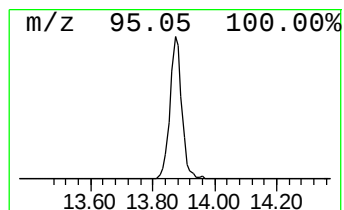
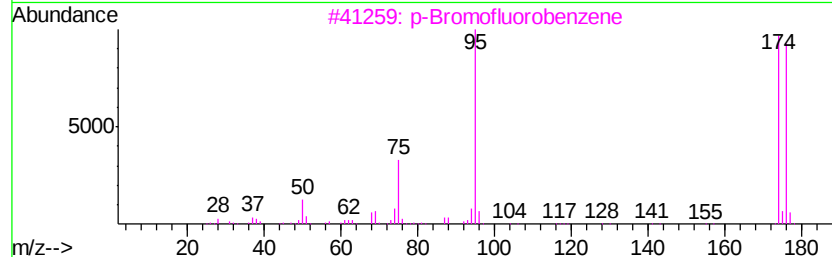
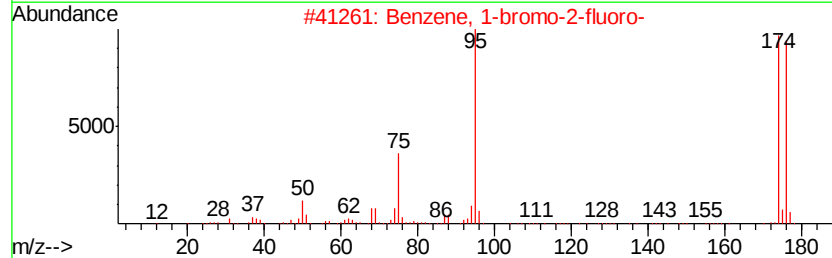
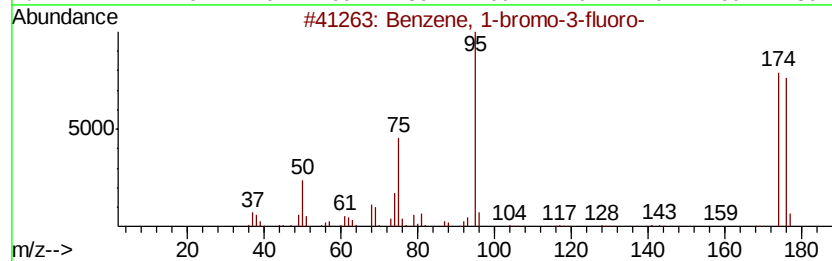
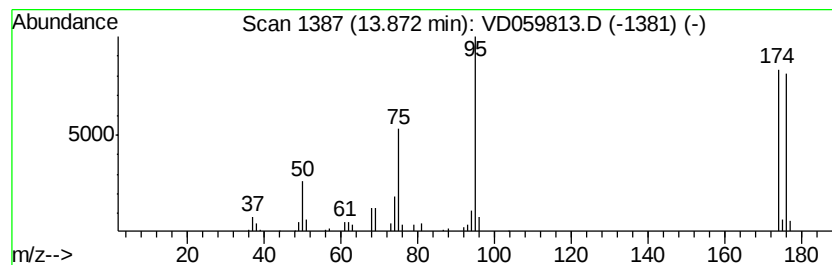
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	25.98 ug/l	336809	Chlorobenzene-d5	10.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-fluoro-	174	C6H4BrF	001073-06-9	98
2		Benzene, 1-bromo-2-fluoro-	174	C6H4BrF	001072-85-1	94
3		p-Bromofluorobenzene	174	C6H4BrF	000460-00-4	91
4		4(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	019808-30-1	23
5		2-Naphthalenol, 3-methoxy-	174	C11H10O2	018515-11-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-bromo-...	13.87	26.0	ug/l	336809	3	10.67	648289	50.0