

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD123019\
 Data File : VD064680.D
 Acq On : 30 Dec 2019 16:35
 Operator : VA/SY
 Sample : K6456-03
 Misc : 6.49G/5.00ml/MSVOA D/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SB-04

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\82D122419S.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	1.956	3	6	12	rVB3	23085	35681	1.42%	0.191%
2	4.967	505	518	528	rBV6	21185	66444	2.64%	0.356%
3	7.714	976	985	994	rBV3	13444	35018	1.39%	0.188%
4	7.920	1008	1020	1024	rBV	345434	820029	32.61%	4.392%
5	7.979	1024	1030	1043	rVB	635259	1433304	57.00%	7.677%
6	8.332	1079	1090	1101	rBV	296489	670361	26.66%	3.590%
7	8.867	1172	1181	1192	rBV	912308	1814764	72.17%	9.720%
8	9.614	1296	1308	1323	rBV	718017	2118971	84.27%	11.349%
9	10.344	1424	1432	1445	rBV	1386244	2514503	100.00%	13.467%
10	11.649	1646	1654	1668	rBV	1242021	2094596	83.30%	11.218%
11	11.802	1671	1680	1692	rVB	804668	1566236	62.29%	8.389%
12	12.238	1747	1754	1760	rBV	24917	42404	1.69%	0.227%
13	12.638	1815	1822	1833	rVB	936213	1564839	62.23%	8.381%
14	13.196	1909	1917	1928	rBV	1108685	1942208	77.24%	10.402%
15	13.585	1976	1983	1992	rVB	1210574	1951839	77.62%	10.454%

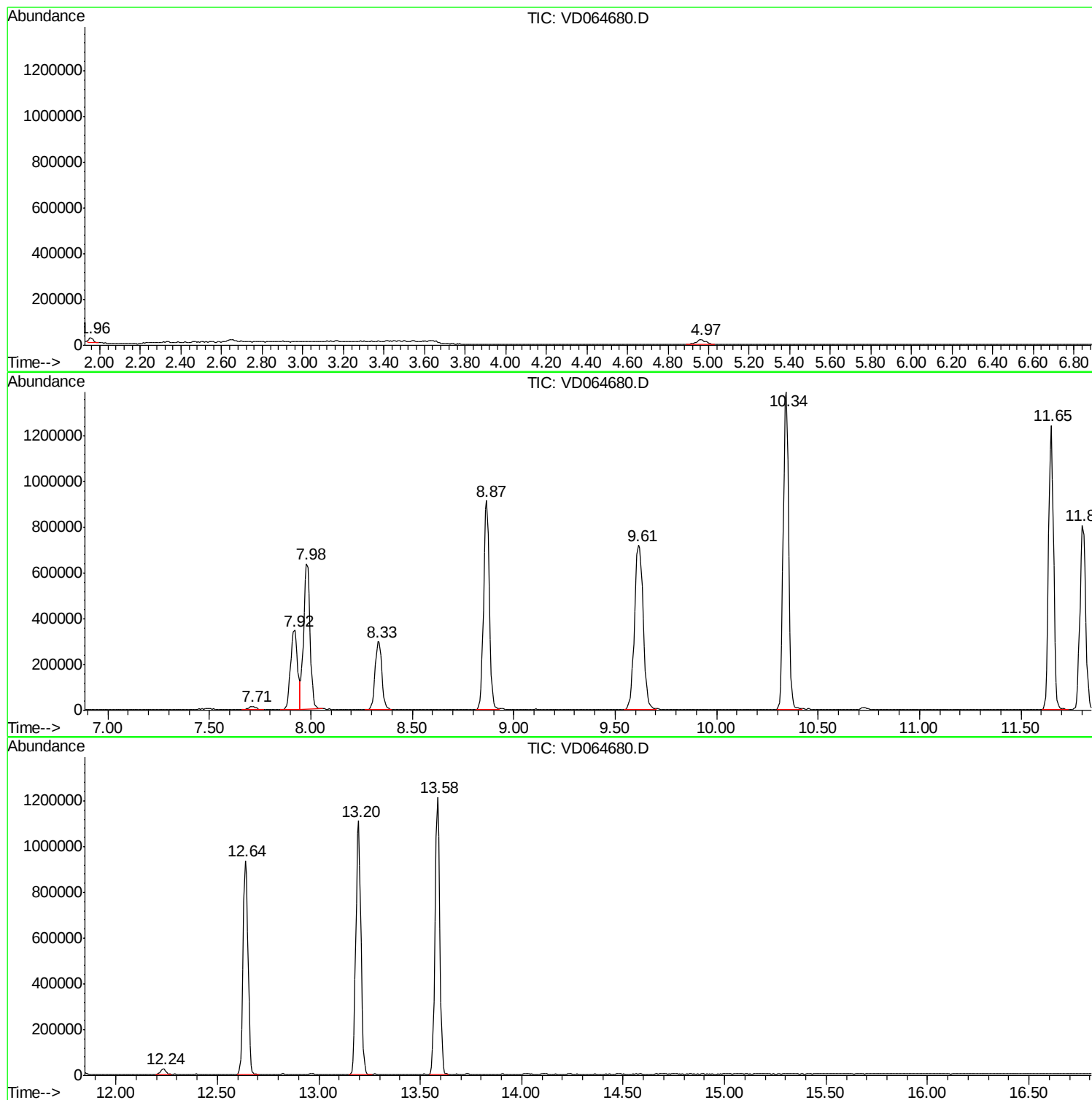
Sum of corrected areas: 18671197

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122419S.M
Quant Title : SW846 8260

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



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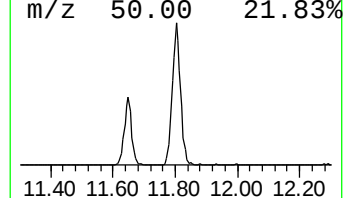
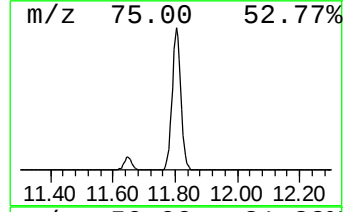
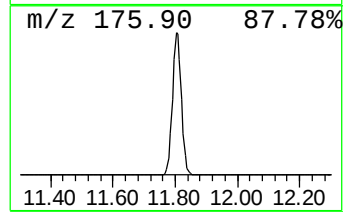
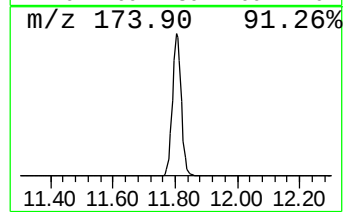
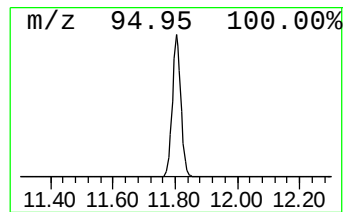
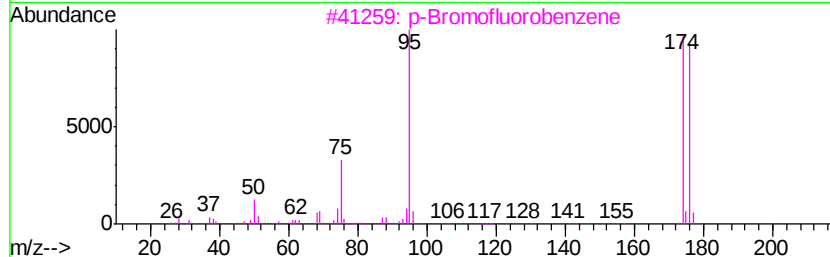
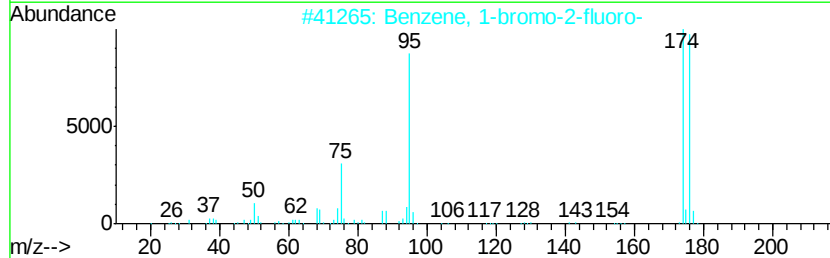
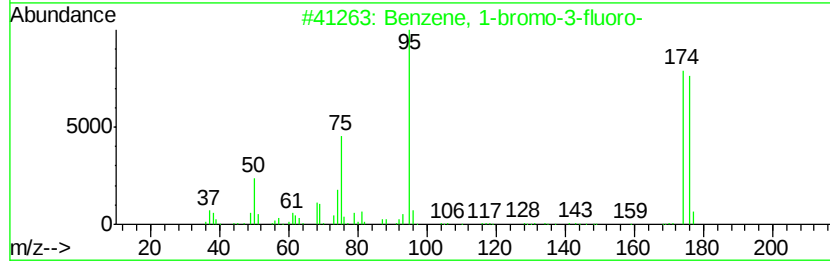
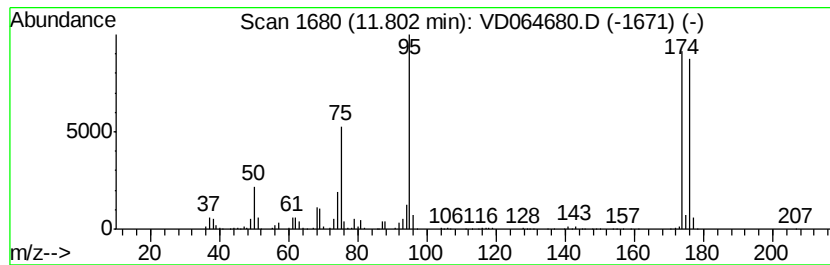
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.80	37.39 ug/l	1566240	Chlorobenzene-d5	11.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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	93
4		4(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	019808-30-1	37
5		3-Chloro-6-methoxy-4-methylpyrid...	174	C6H7ClN2O2	050450-89-0	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-bromo-...	11.80	37.4	ug/l	1566240	3	11.65	2094600	50.0