

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
 Data File : BM018288.D
 Acq On : 19 Dec 2018 01:24
 Operator : JU/SJ
 Sample : J6398-17
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FB-01_121118

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.122	372	380	391	rBV	1247074	1793748	24.93%	4.761%
2	6.687	639	646	658	rBV	664681	1094372	15.21%	2.905%
3	7.028	696	704	716	rBV	2472354	3767395	52.36%	9.999%
4	7.487	776	782	790	rBV	563796	868791	12.08%	2.306%
5	7.798	827	835	847	rBV	2246109	3455630	48.03%	9.171%
6	8.645	972	979	994	rBV	1700155	2793324	38.82%	7.414%
7	10.251	1245	1252	1274	rBV	601888	1085963	15.09%	2.882%
8	12.751	1670	1677	1693	rBV	3967847	5845267	81.24%	15.514%
9	14.145	1907	1914	1926	rBV2	756083	1214555	16.88%	3.224%
10	15.657	2164	2171	2191	rBV	2662650	4075398	56.64%	10.816%
11	16.910	2378	2384	2401	rBV	841214	1294774	18.00%	3.436%
12	19.562	2829	2835	2845	rBV	5601023	7194848	100.00%	19.096%
13	21.127	3096	3101	3109	rBV	1229791	1573099	21.86%	4.175%
14	23.286	3462	3468	3477	rVB	871028	1620882	22.53%	4.302%

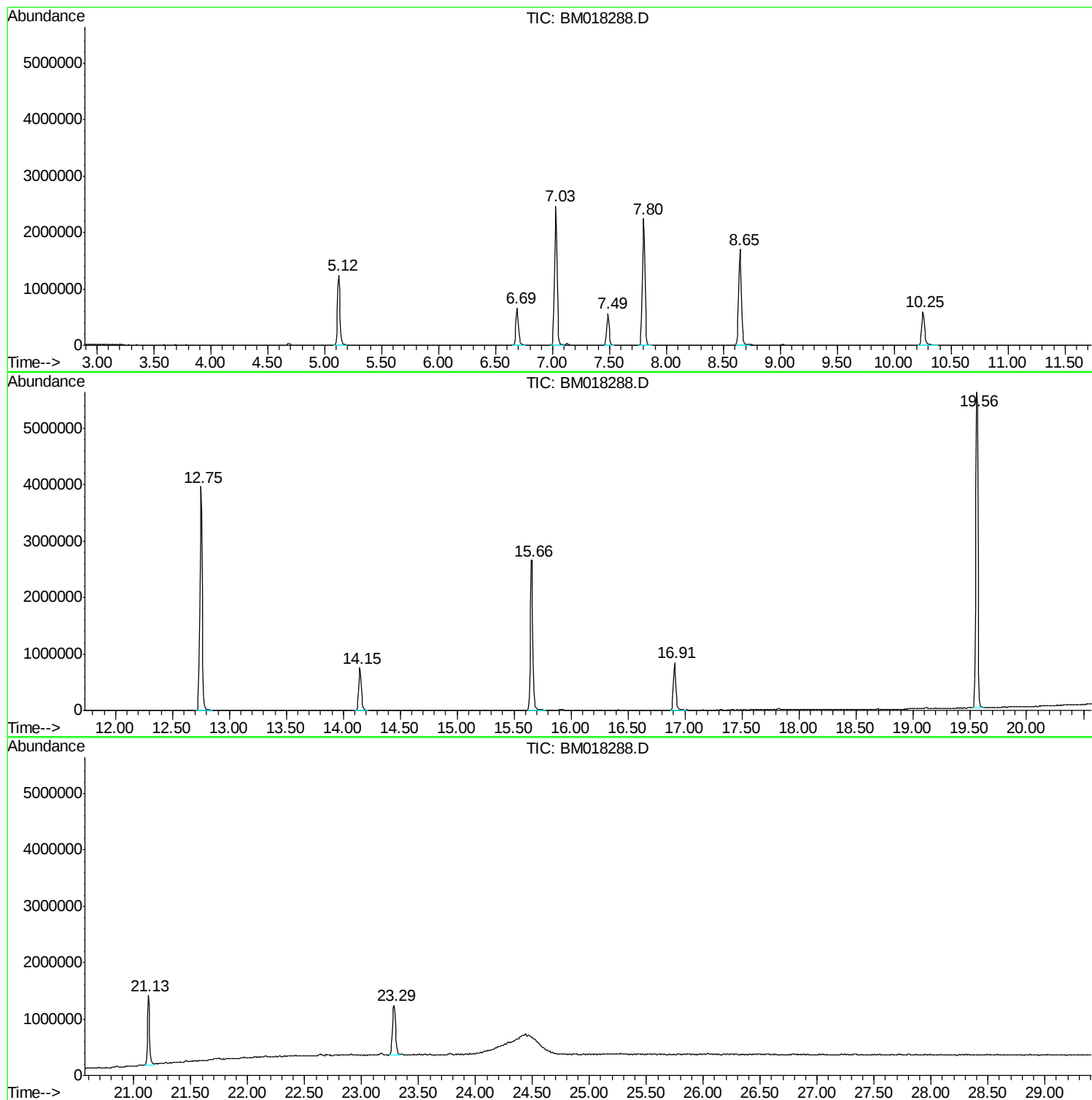
Sum of corrected areas: 37678046

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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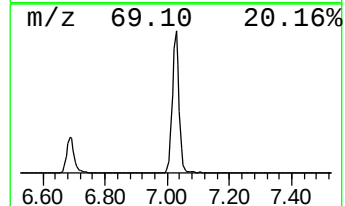
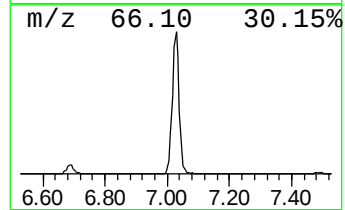
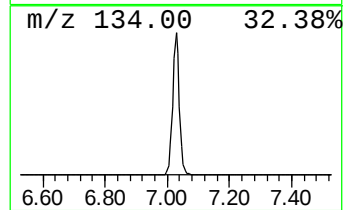
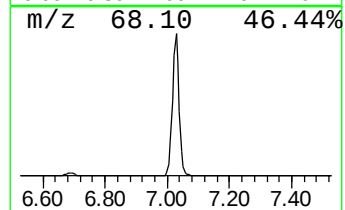
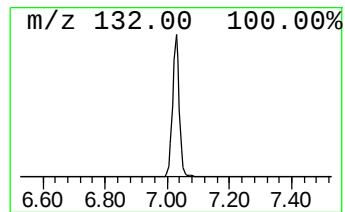
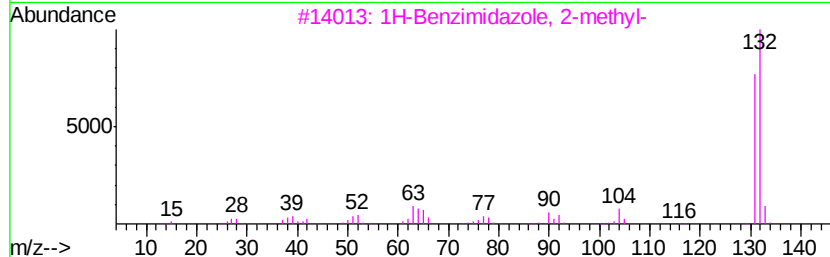
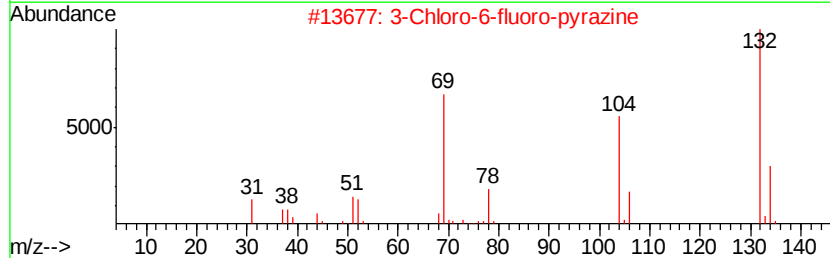
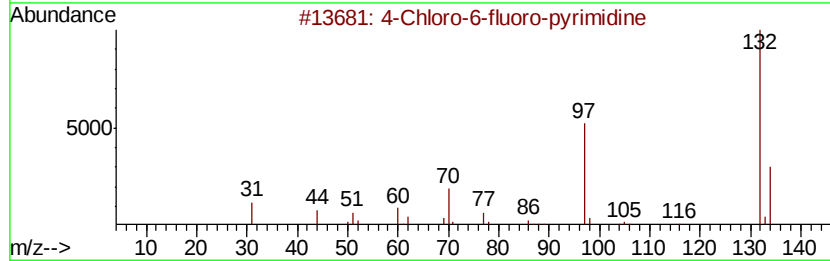
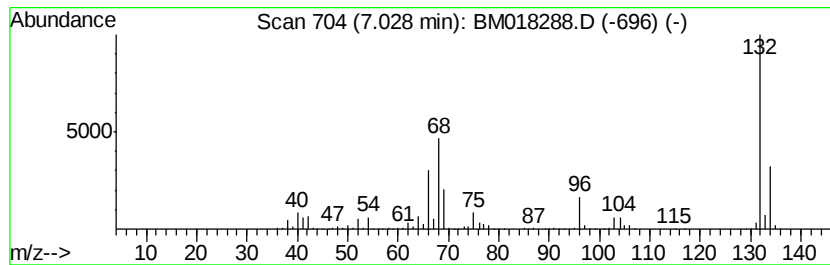
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	86.73 ng	3767400	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.03	7.03	86.7	ng	3767400	1	7.49	868791	20.0