

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080818\
 Data File : BG036221.D
 Acq On : 9 Aug 2018 7:12
 Operator : JU/SJ
 Sample : J4350-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 080218-TIPC-E-U-M

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_G\METHODS\8270-BG071218.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	5.597	386	393	411	rBV	154101	304620	10.87%	3.108%
2	7.177	656	662	678	rBV	93231	212484	7.58%	2.168%
3	7.542	718	724	738	rBV	306117	557540	19.90%	5.688%
4	8.006	796	803	814	rVB3	89912	166227	5.93%	1.696%
5	8.329	851	858	872	rBV	401215	723166	25.81%	7.378%
6	9.169	993	1001	1023	rBV	315959	664549	23.72%	6.780%
7	10.814	1273	1281	1292	rBV	86070	194757	6.95%	1.987%
8	13.251	1689	1696	1713	rBV	906298	1562248	55.76%	15.939%
9	14.632	1925	1931	1946	rBV2	174442	323642	11.55%	3.302%
10	16.124	2179	2185	2201	rBV	455488	841089	30.02%	8.581%
11	17.381	2393	2399	2410	rBV2	247833	444083	15.85%	4.531%
12	19.984	2836	2842	2854	rBV	2083511	2801628	100.00%	28.583%
13	21.640	3116	3124	3140	rBV	268727	531284	18.96%	5.420%
14	24.824	3655	3666	3685	rBV2	139689	474337	16.93%	4.839%

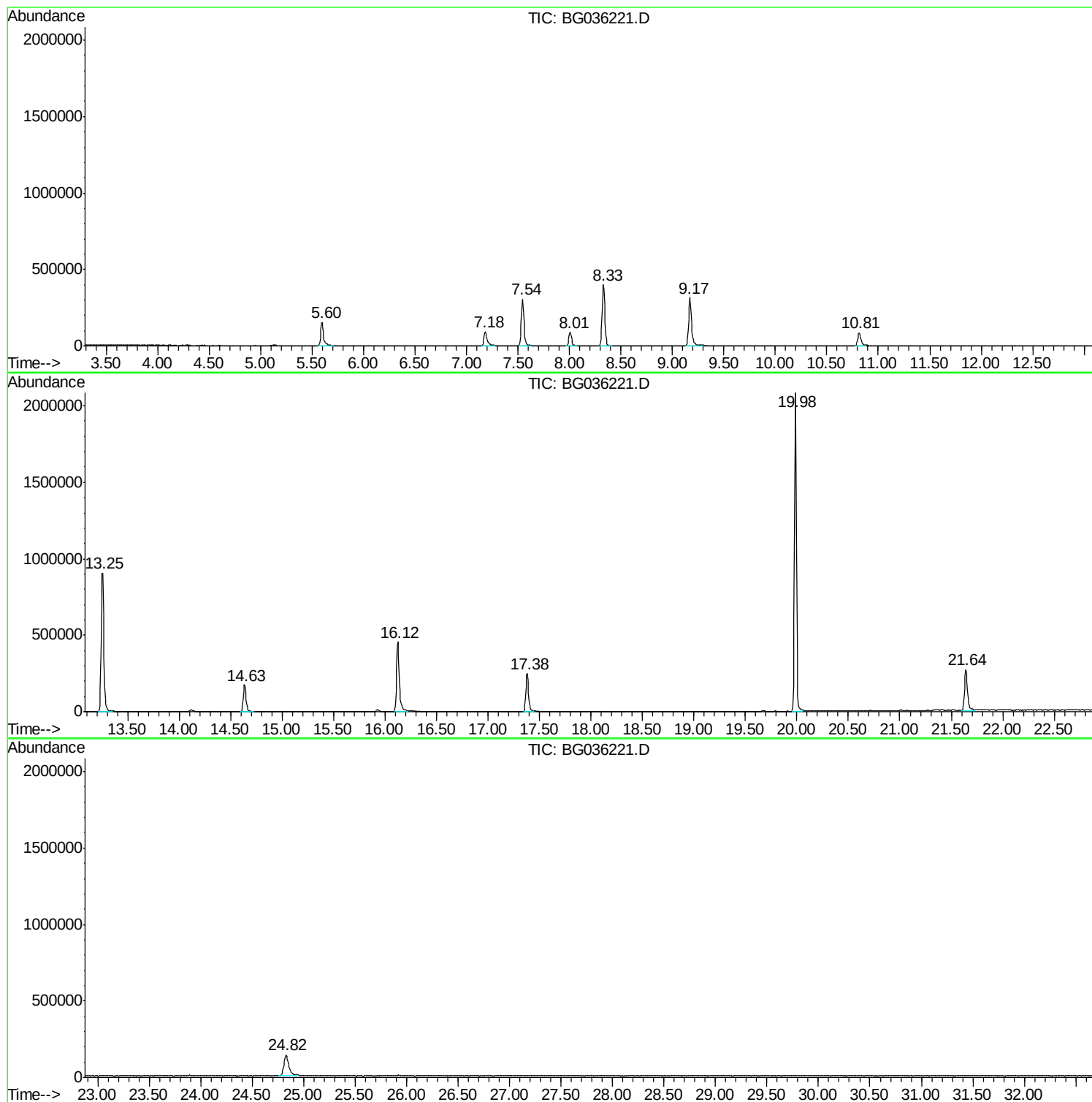
Sum of corrected areas: 9801654

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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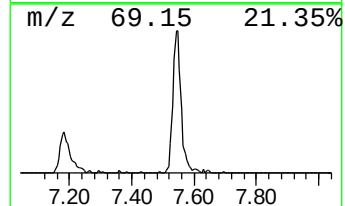
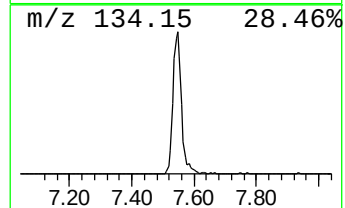
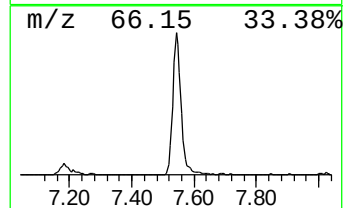
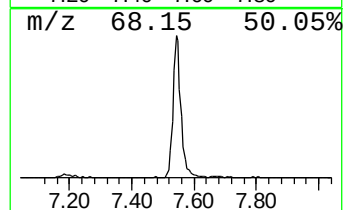
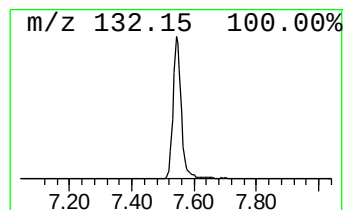
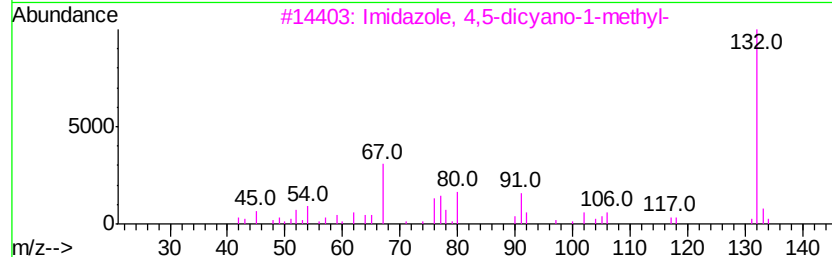
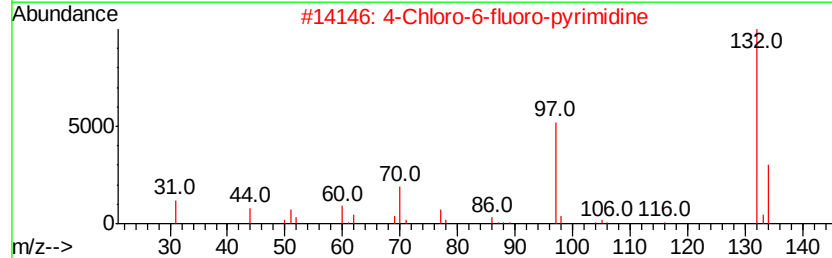
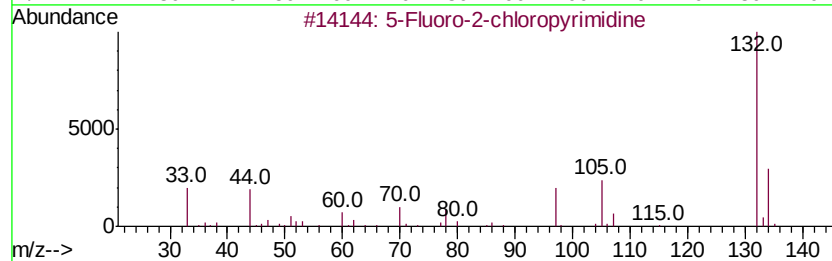
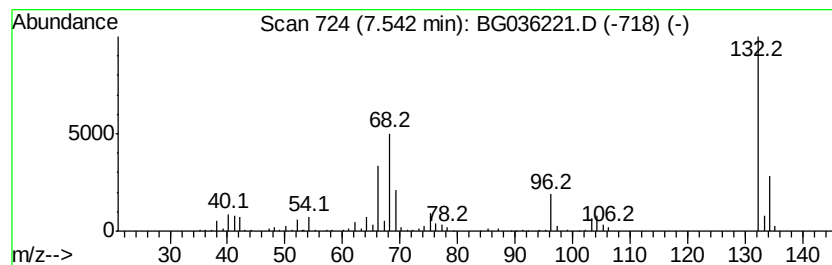
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Peak Number 1 unknown7.54 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	67.08 ng	557540	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	25
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
5		5-Aminoindole	132	C8H8N2	005192-03-0	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.54	7.54	67.1	ng	557540	1	8.01	166227	20.0