

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102419\
 Data File : BG043330.D
 Acq On : 24 Oct 2019 14:20
 Operator : JU
 Sample : PB124233BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124233BL

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-BG102119.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	5.603	421	428	444	rBV	640682	1137259	29.82%	7.000%
2	7.201	691	700	712	rBV	675961	1230668	32.27%	7.575%
3	7.560	753	761	776	rBV	685831	1250556	32.79%	7.697%
4	8.024	832	840	850	rVB	146629	252063	6.61%	1.551%
5	8.347	886	895	909	rBV	559881	999465	26.21%	6.152%
6	9.187	1028	1038	1051	rBV	346514	724014	18.99%	4.456%
7	10.826	1308	1317	1328	rBV	126219	290198	7.61%	1.786%
8	13.276	1726	1734	1759	rBV	1096746	1959866	51.39%	12.063%
9	14.651	1961	1968	1984	rBV2	273338	487809	12.79%	3.002%
10	16.138	2214	2221	2246	rBV	941580	1711571	44.88%	10.534%
11	17.395	2429	2435	2451	rBV2	345929	642724	16.85%	3.956%
12	20.004	2873	2879	2899	rBV	2609424	3813458	100.00%	23.471%
13	21.661	3154	3161	3180	rBV2	412927	876076	22.97%	5.392%
14	24.863	3696	3706	3724	rBV2	274124	871621	22.86%	5.365%

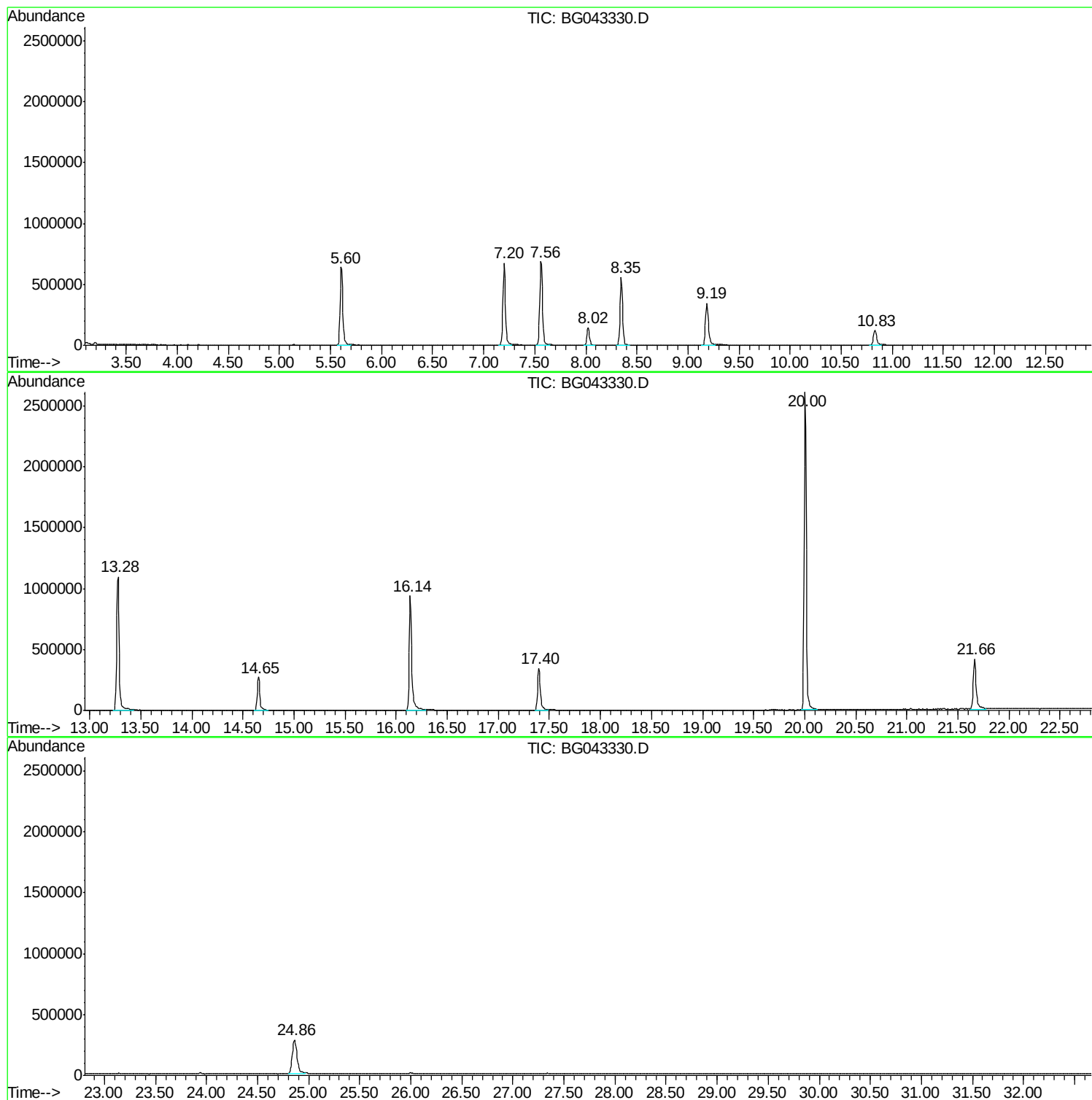
Sum of corrected areas: 16247348

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.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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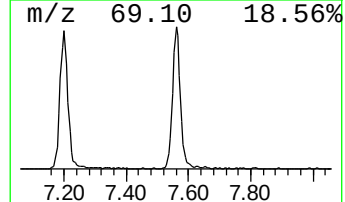
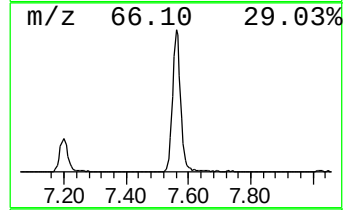
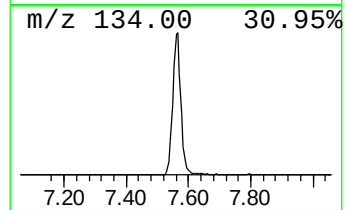
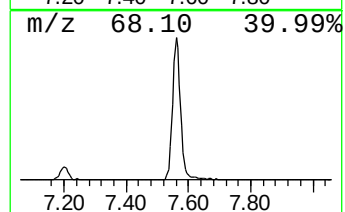
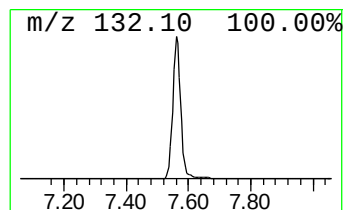
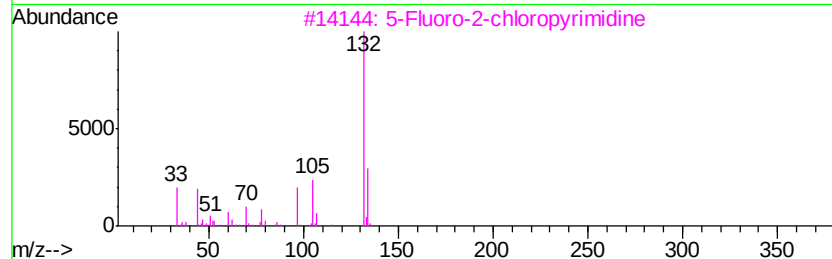
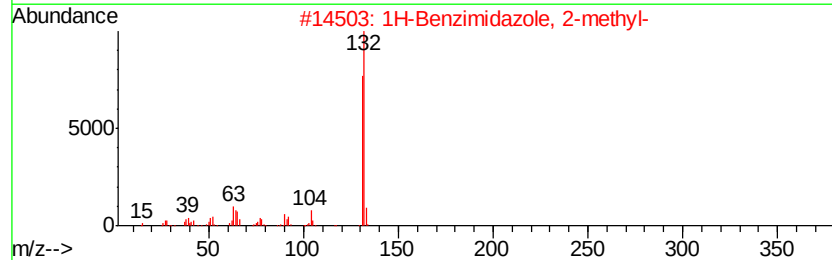
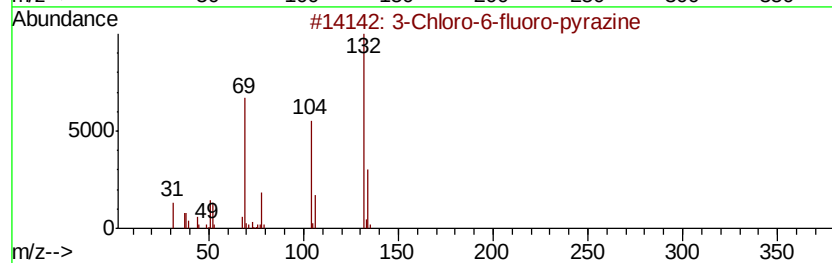
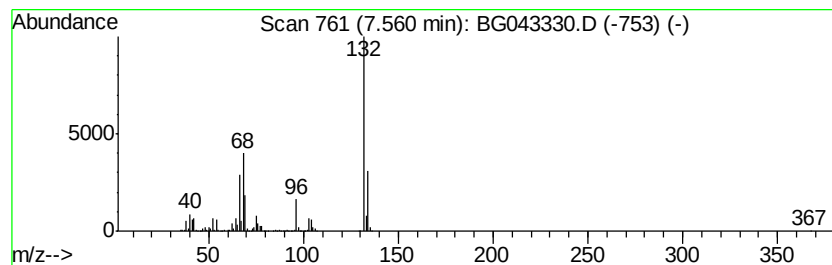
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Peak Number 1 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	99.23 ng	1250560	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
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
unknown7.56	7.56	99.2	ng	1250560	1	8.02	252063	20.0