

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090419\
 Data File : BG042714.D
 Acq On : 4 Sep 2019 19:15
 Operator : HP/JU
 Sample : PB122793BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB122793BL

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-BG082219.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	5.555	413	420	434	rBV	394452	632320	25.43%	6.102%
2	7.165	686	694	712	rBV	377948	669659	26.93%	6.462%
3	7.529	749	756	770	rBV	415919	722140	29.04%	6.968%
4	7.999	829	836	843	rBV	93974	166089	6.68%	1.603%
5	8.323	884	891	902	rBV	361110	616819	24.81%	5.952%
6	9.169	1026	1035	1051	rBV	210567	407992	16.41%	3.937%
7	10.820	1309	1316	1335	rBV	101163	217092	8.73%	2.095%
8	13.276	1727	1734	1749	rBV	804153	1300124	52.29%	12.546%
9	14.662	1963	1970	1983	rBV	193223	350578	14.10%	3.383%
10	16.149	2217	2223	2241	rBV	554414	1025749	41.25%	9.898%
11	17.412	2431	2438	2453	rBV	273461	479786	19.29%	4.630%
12	20.021	2876	2882	2905	rVB	1840372	2486607	100.00%	23.995%
13	21.689	3159	3166	3181	rBV	334666	629994	25.34%	6.079%
14	24.938	3706	3719	3741	rBV	185699	658113	26.47%	6.351%

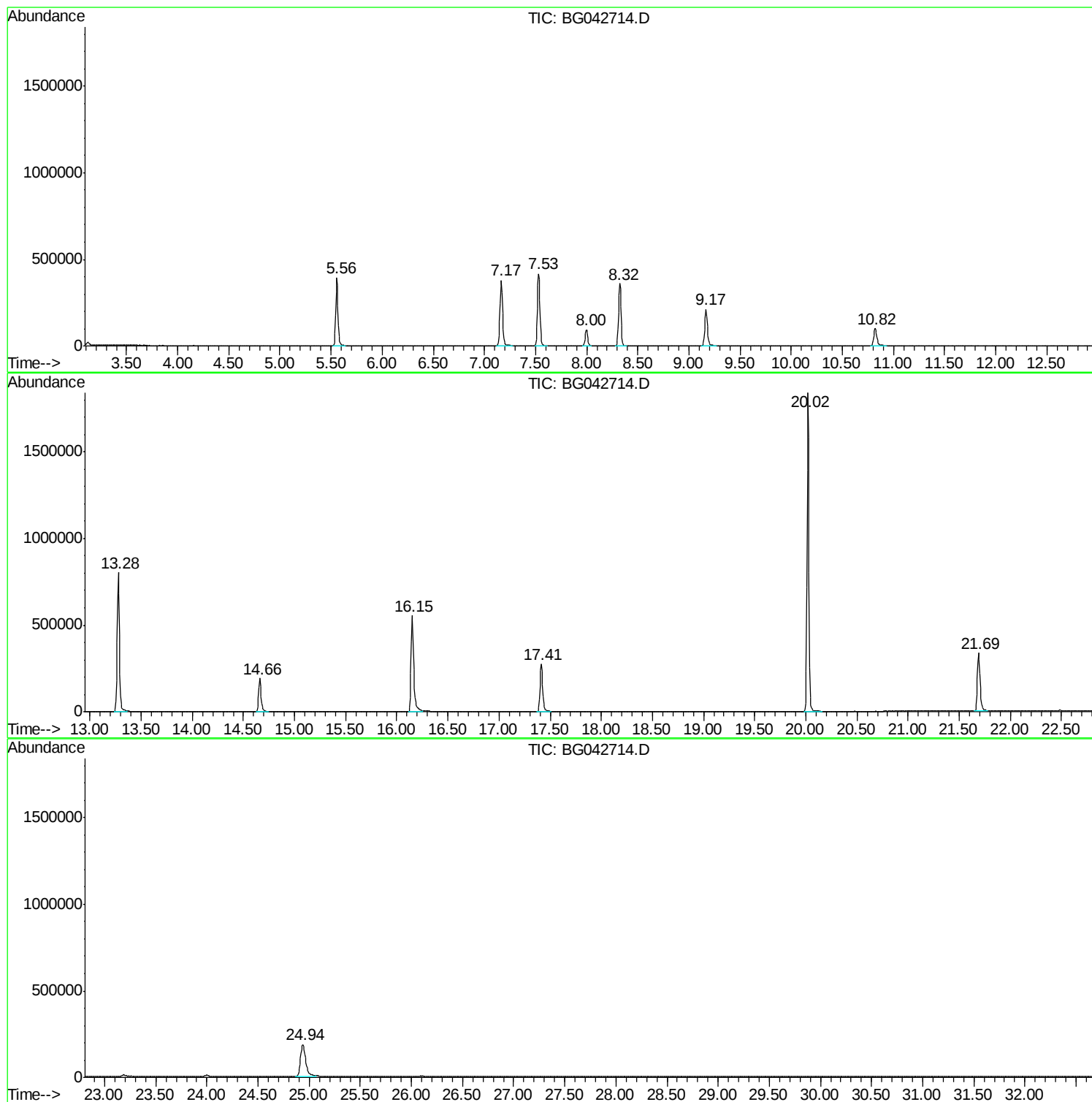
Sum of corrected areas: 10363062

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.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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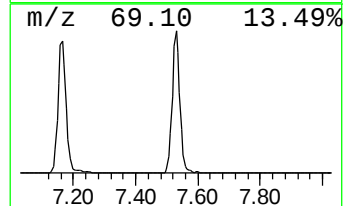
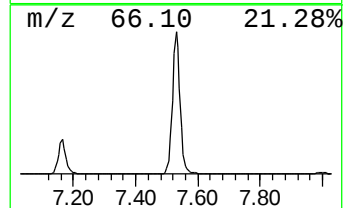
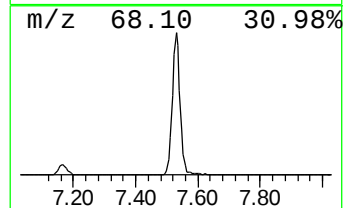
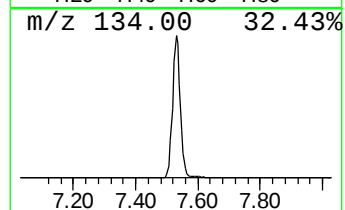
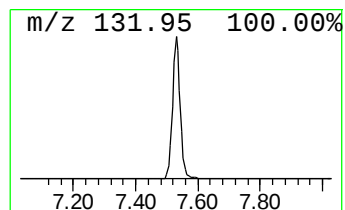
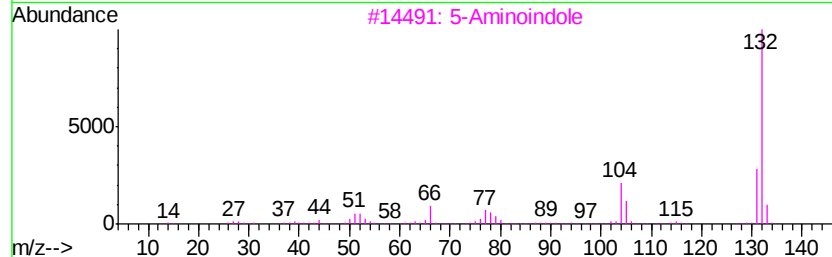
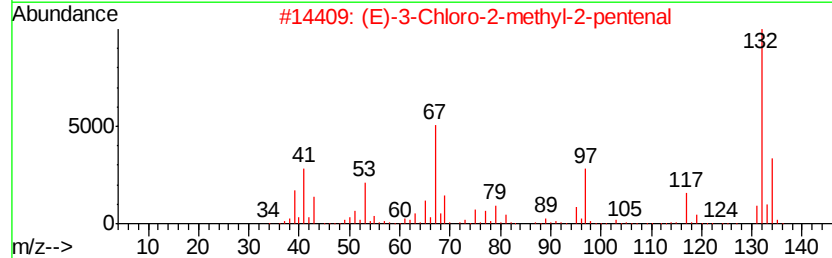
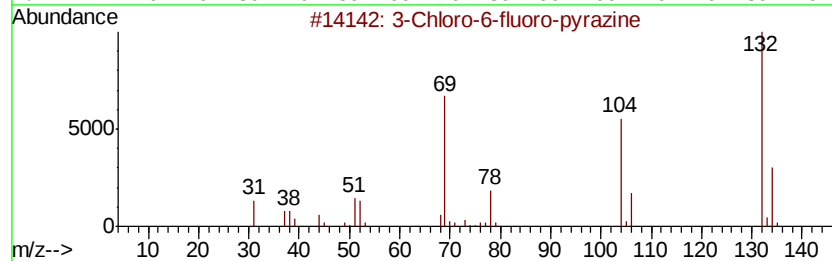
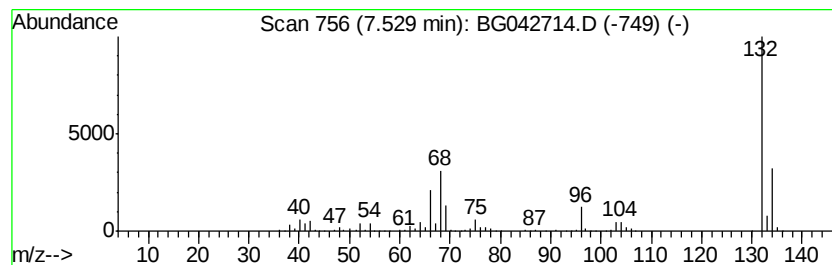
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Peak Number 1 unknown7.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	86.96 ng	722140	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.53	7.53	87.0	ng	722140	1	8.00	166089	20.0