

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090419\
Data File : BG042704.D
Acq On : 4 Sep 2019 12:52
Operator : HP/JU
Sample : PB122757BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB122757BL

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	5.556	414	420	434	rBV	321483	520858	18.05%	4.751%
2	7.166	686	694	713	rBV	299589	550206	19.06%	5.019%
3	7.530	749	756	771	rBV	344108	593902	20.58%	5.417%
4	7.995	828	835	843	rBV	112392	190291	6.59%	1.736%
5	8.324	883	891	900	rBV	433277	742071	25.71%	6.769%
6	9.164	1026	1034	1047	rBV	249707	487834	16.90%	4.450%
7	10.821	1309	1316	1327	rBV	122831	241686	8.37%	2.205%
8	13.277	1727	1734	1748	rBV	933547	1525845	52.87%	13.918%
9	14.663	1962	1970	1985	rBV	217764	400361	13.87%	3.652%
10	16.150	2217	2223	2245	rBV	422461	802601	27.81%	7.321%
11	17.413	2431	2438	2453	rBV2	314434	543552	18.83%	4.958%
12	20.022	2876	2882	2894	rBV	2134106	2886088	100.00%	26.325%
13	21.690	3158	3166	3183	rBV	390326	709291	24.58%	6.470%
14	24.934	3706	3718	3741	rBV	219515	768703	26.63%	7.012%

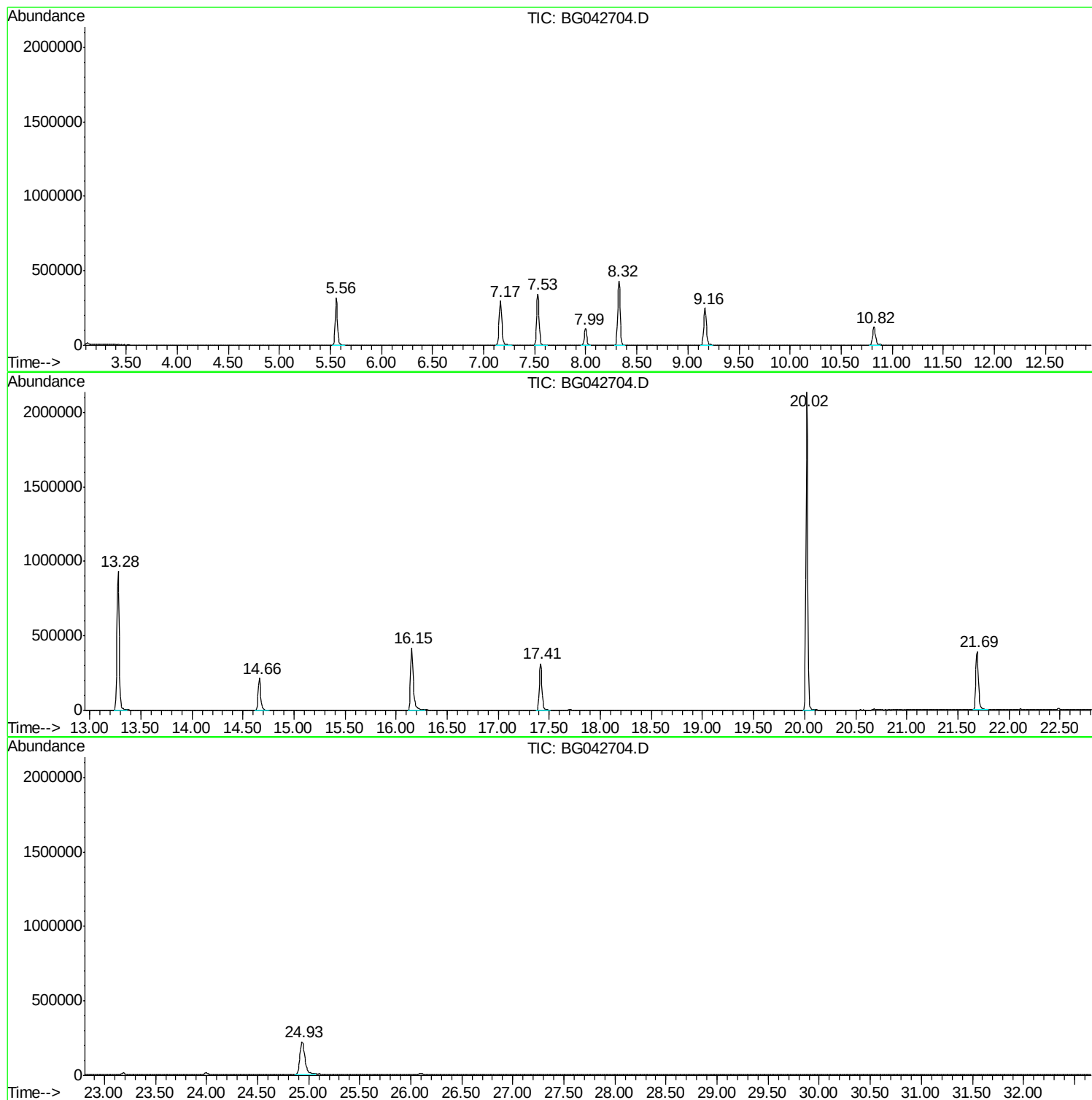
Sum of corrected areas: 10963289

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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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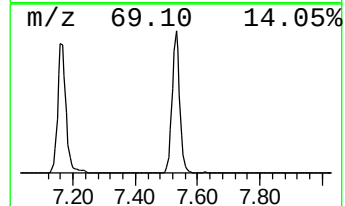
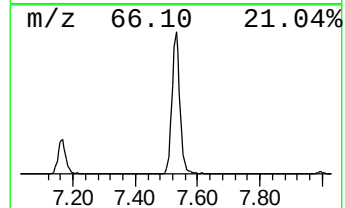
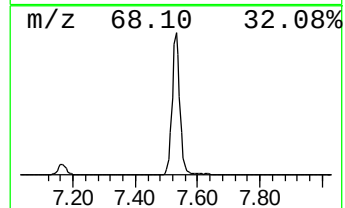
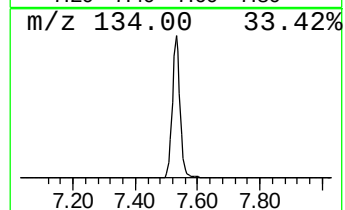
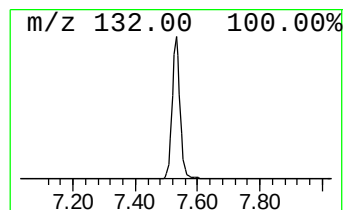
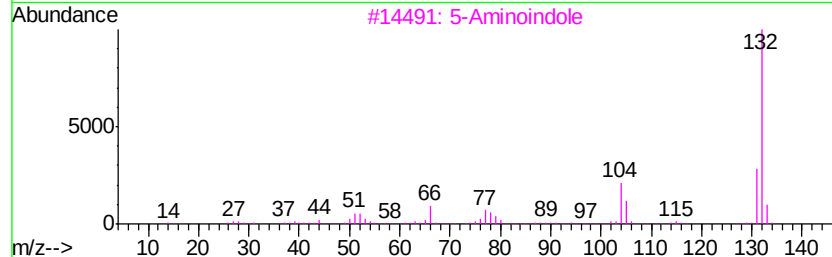
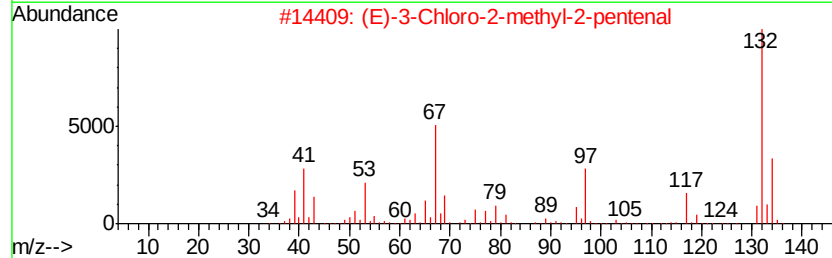
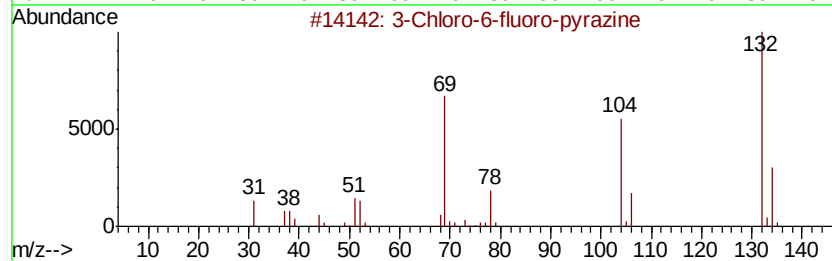
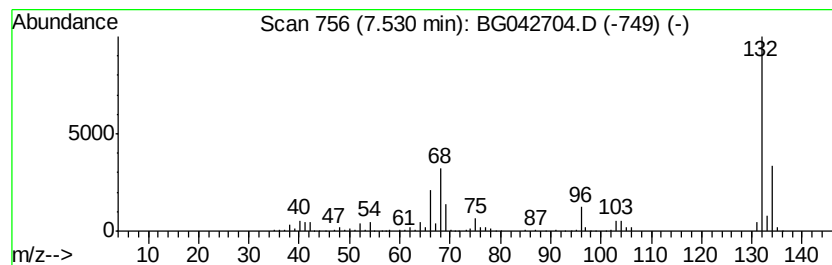
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Peak Number 1 unknown7.53 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	62.42 ng	593902	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	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	16
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	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	62.4	ng	593902	1	8.00	190291	20.0