

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020918\
 Data File : BG032643.D
 Acq On : 9 Feb 2018 22:43
 Operator : SJ/JU
 Sample : PB106437BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB106437BL

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:\HPCHEM1\BNA_G\METHODS\8270-BG012918.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	6.122	475	482	495	rBV	365970	710019	22.34%	5.471%
2	7.802	761	768	790	rBV	378432	823407	25.91%	6.345%
3	8.196	828	835	854	rBV	396937	825074	25.97%	6.358%
4	8.684	910	918	925	rBV2	68297	147012	4.63%	1.133%
5	9.013	967	974	988	rBV	318354	646494	20.35%	4.982%
6	9.882	1114	1122	1142	rBV	205158	484650	15.25%	3.735%
7	11.557	1399	1407	1420	rBV	106439	240122	7.56%	1.850%
8	13.942	1806	1813	1831	rBV	953064	1698568	53.45%	13.089%
9	15.323	2041	2048	2063	rBV2	251932	463991	14.60%	3.575%
10	16.798	2290	2299	2317	rBV	848451	1476601	46.47%	11.378%
11	18.061	2506	2514	2528	rBV2	356822	641621	20.19%	4.944%
12	20.599	2940	2946	2961	rBV	2128689	3177585	100.00%	24.485%
13	22.391	3244	3251	3263	rBV2	410725	824653	25.95%	6.355%
14	26.057	3863	3875	3894	rBV2	234678	817629	25.73%	6.300%

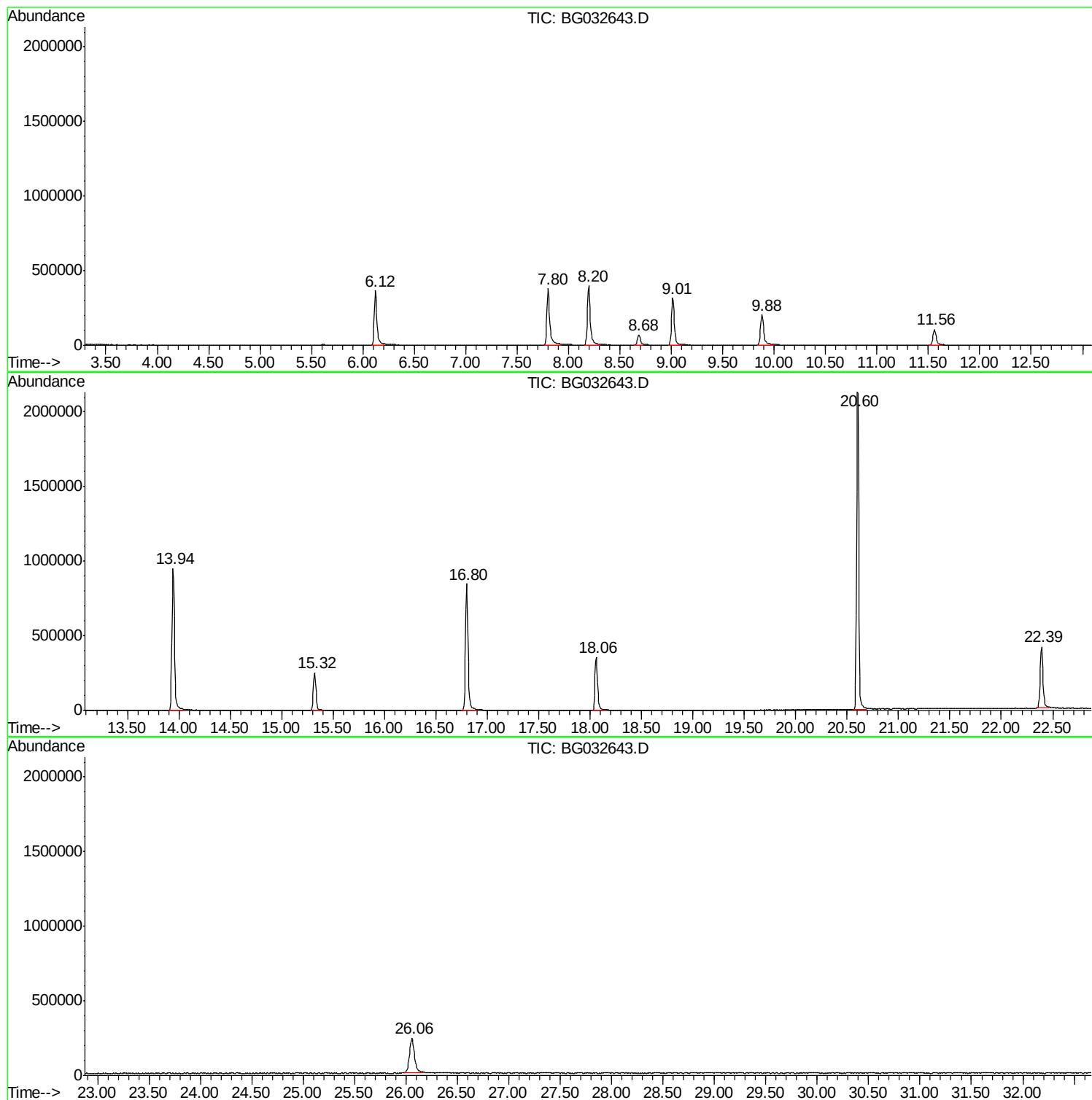
Sum of corrected areas: 12977426

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.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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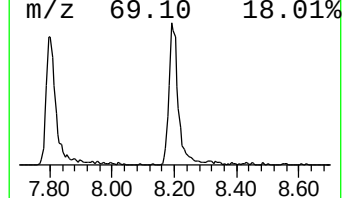
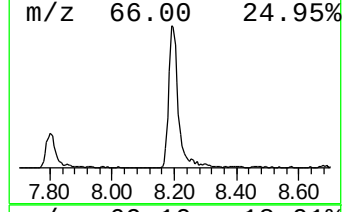
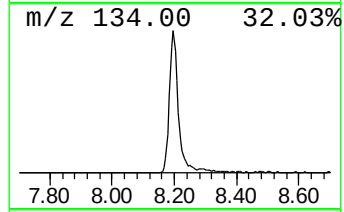
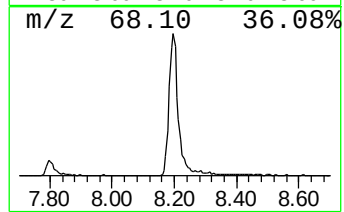
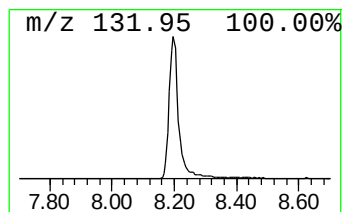
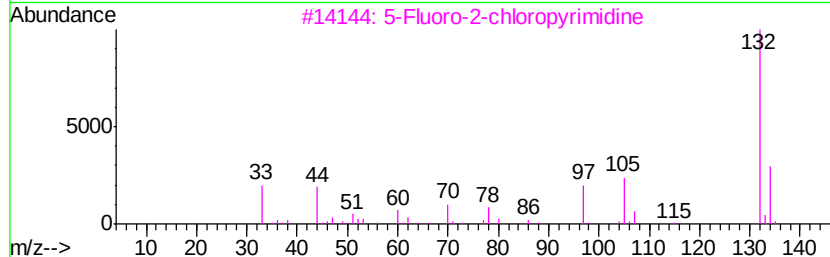
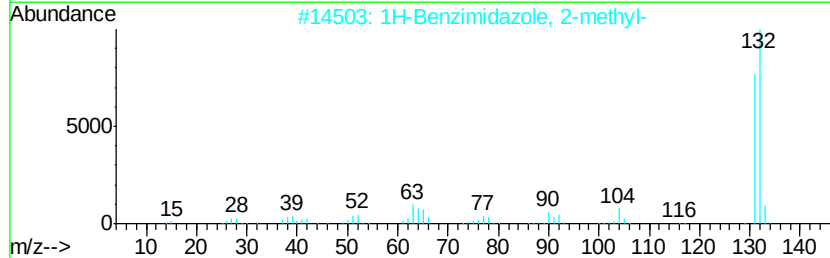
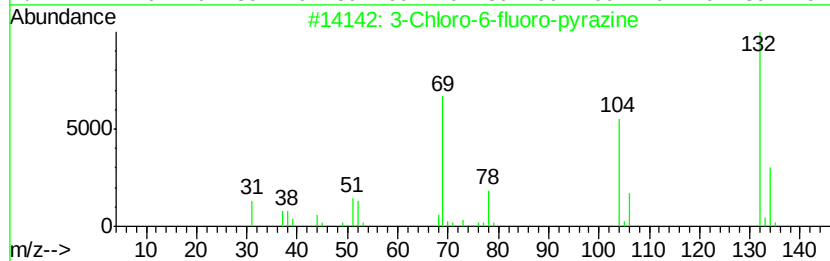
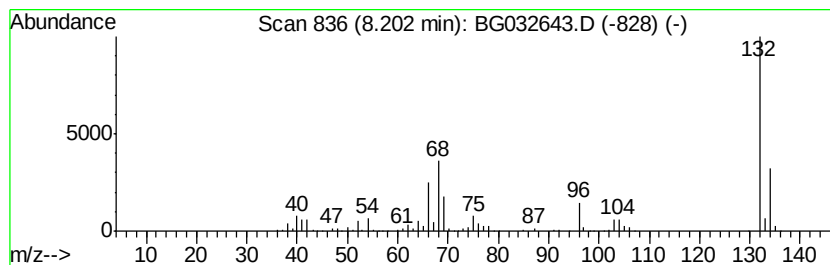
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Peak Number 1 unknown8.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	112.25 ng	825074	1,4-Dichlorobenzene-d4	8.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
5		5-Aminoindole	132	C8H8N2	005192-03-0	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown8.20	8.20	112.3	ng	825074	1	8.68	147012	20.0