

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041718\
 Data File : BG033858.D
 Acq On : 17 Apr 2018 23:30
 Operator : SJ/JU
 Sample : PB108313BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB108313BL

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-BG031918.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	6.266	500	507	533	rBV	248379	913602	25.58%	6.249%
2	7.964	788	796	823	rBV2	197989	1006357	28.17%	6.883%
3	8.346	854	861	880	rBV	292400	959997	26.87%	6.566%
4	8.839	936	945	961	rBV3	40202	166296	4.66%	1.137%
5	9.156	991	999	1021	rBV	310260	844067	23.63%	5.773%
6	10.044	1142	1150	1170	rBV2	127661	587100	16.44%	4.016%
7	11.706	1426	1433	1449	rBV	68138	239766	6.71%	1.640%
8	14.074	1829	1836	1863	rBV	759800	1909843	53.46%	13.063%
9	15.443	2063	2069	2086	rBV2	198432	466074	13.05%	3.188%
10	16.924	2313	2321	2340	rBV	698869	1700121	47.59%	11.629%
11	18.181	2528	2535	2557	rBV2	292134	709545	19.86%	4.853%
12	20.702	2958	2964	2981	rBV	2320923	3572246	100.00%	24.434%
13	22.523	3266	3274	3287	rBV2	352941	797615	22.33%	5.456%
14	26.266	3900	3911	3929	rBV3	202575	747465	20.92%	5.113%

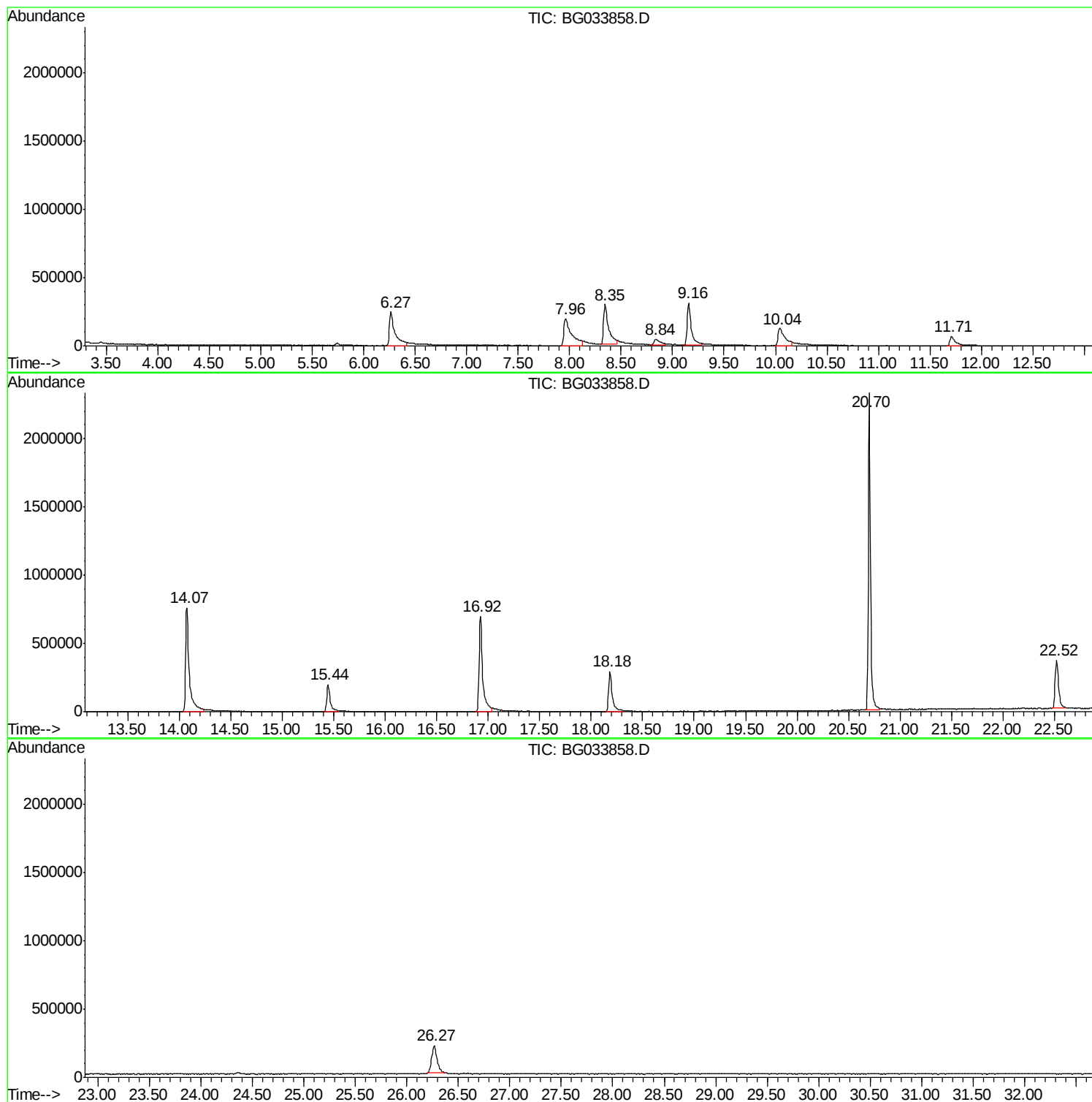
Sum of corrected areas: 14620094

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.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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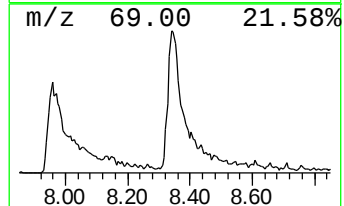
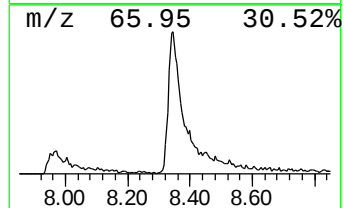
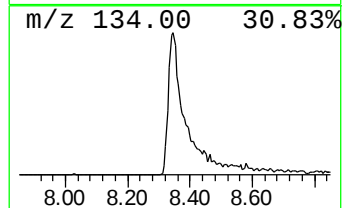
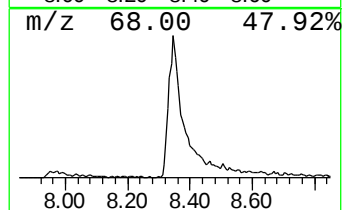
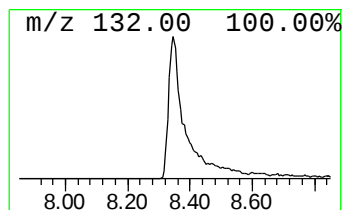
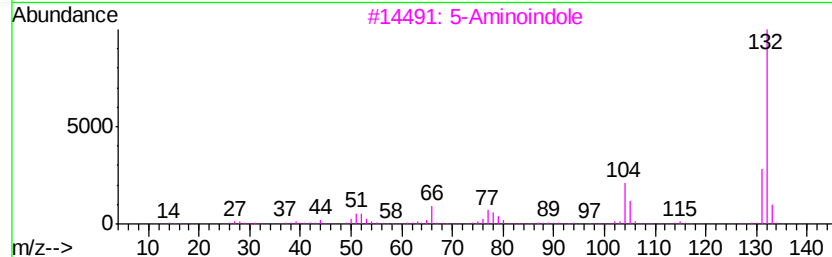
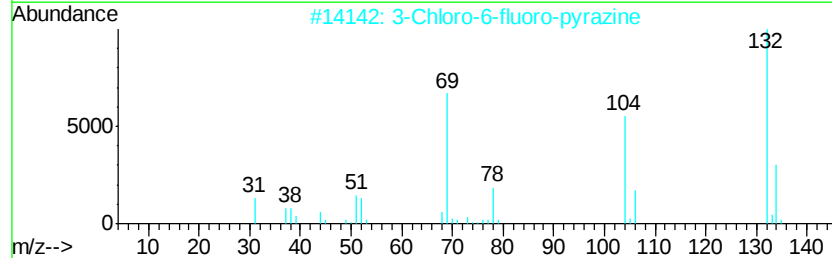
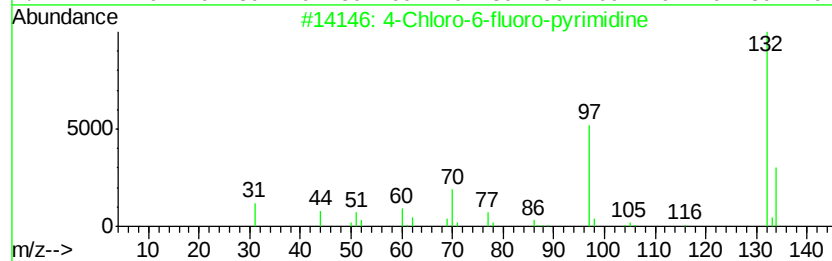
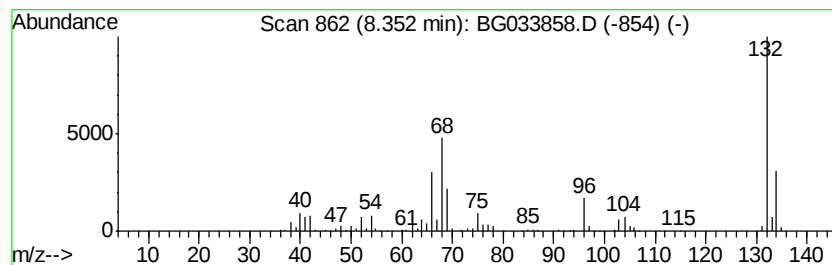
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Peak Number 1 unknown8.35 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.35	115.46 ng	959997	1,4-Dichlorobenzene-d4	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown8.35	8.35	115.5	ng	959997	1	8.84	166296	20.0