

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG032218\
 Data File : BG033307.D
 Acq On : 22 Mar 2018 22:58
 Operator : SJ/JU
 Sample : PB107545BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107545BL

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	5.793	420	426	437	rBV2	31192	56613	1.28%	0.309%
2	6.305	507	513	531	rBV	671554	1358350	30.80%	7.412%
3	7.991	792	800	815	rBV	694380	1510267	34.24%	8.241%
4	8.402	862	870	893	rBV	648872	1418866	32.17%	7.742%
5	8.890	946	953	967	rBV	84589	211923	4.80%	1.156%
6	9.225	1002	1010	1023	rBV	503962	1013470	22.98%	5.530%
7	10.088	1150	1157	1176	rBV	386029	914096	20.72%	4.988%
8	11.775	1437	1444	1456	rBV	132198	303497	6.88%	1.656%
9	14.137	1839	1846	1867	rBV	1383839	2400838	54.43%	13.101%
10	15.506	2072	2079	2093	rBV2	288538	534934	12.13%	2.919%
11	16.980	2323	2330	2347	rBV	972452	1700341	38.55%	9.278%
12	18.238	2537	2544	2562	rBV	450918	786967	17.84%	4.294%
13	20.758	2967	2973	2989	rBV	3253678	4410786	100.00%	24.068%
14	22.591	3278	3285	3298	rVB	469467	878805	19.92%	4.795%
15	26.381	3919	3930	3948	rBV2	236443	826416	18.74%	4.509%

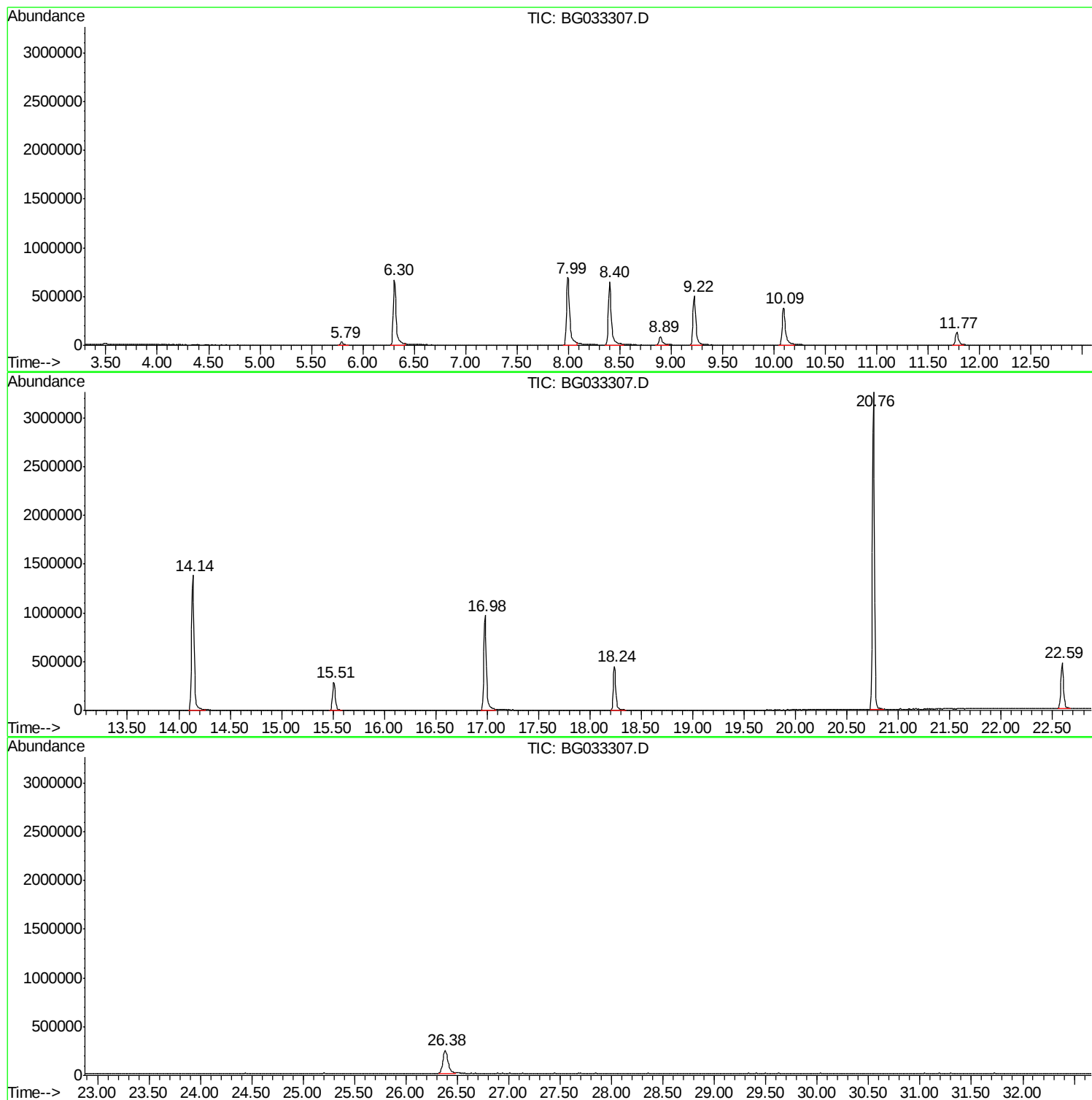
Sum of corrected areas: 18326169

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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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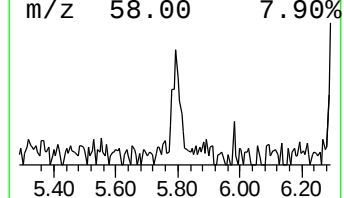
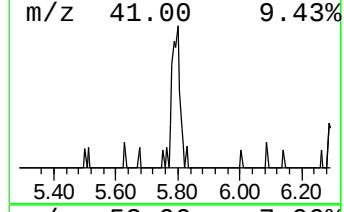
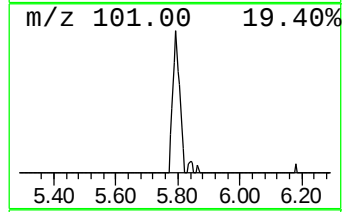
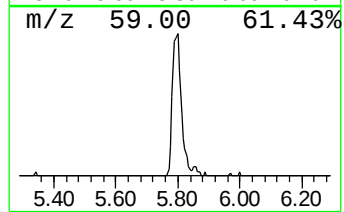
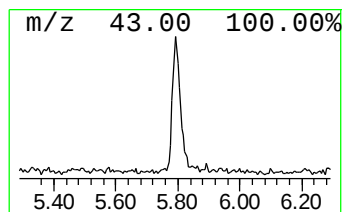
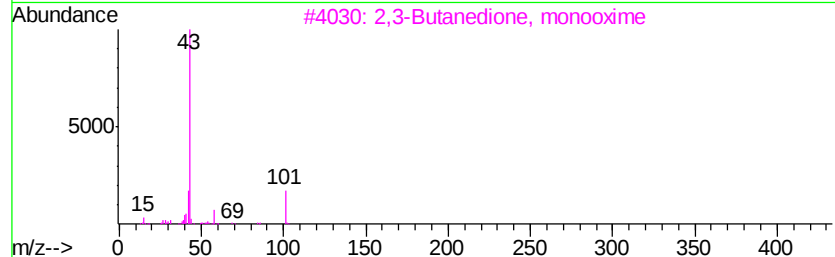
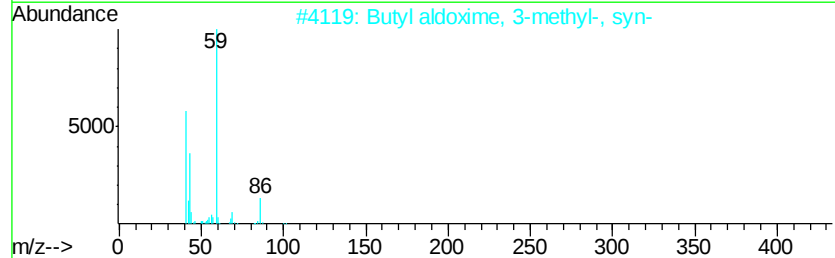
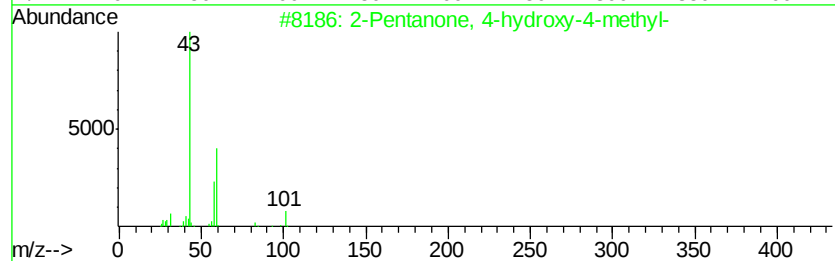
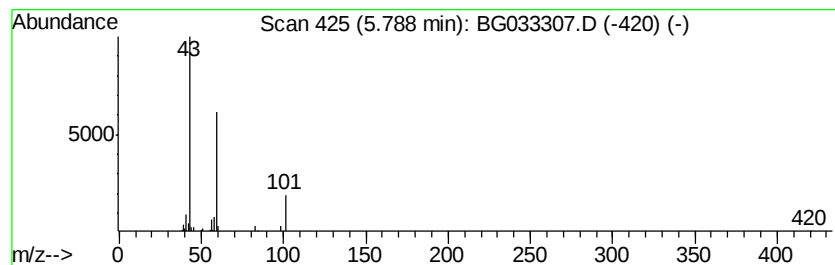
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.79	5.34 ng	56613	1,4-Dichlorobenzene-d4	8.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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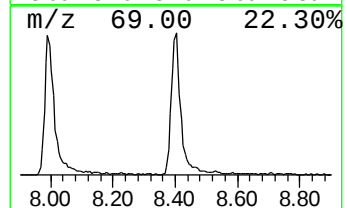
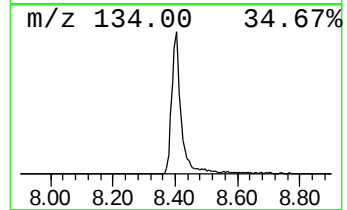
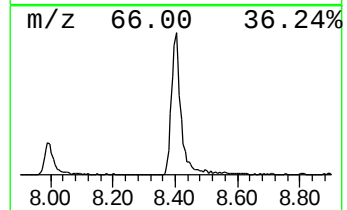
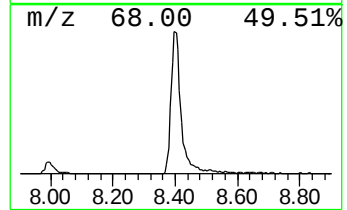
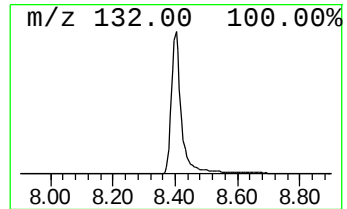
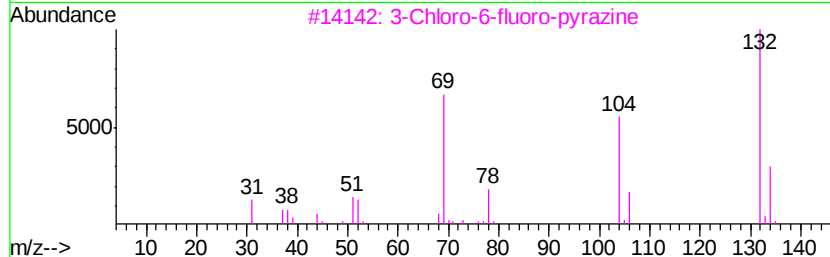
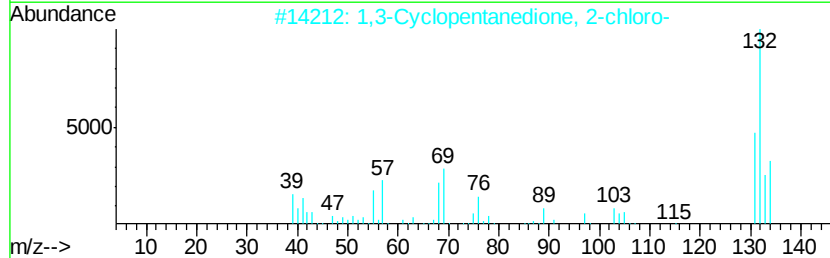
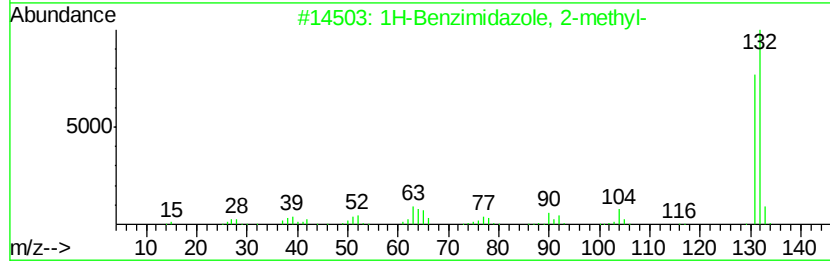
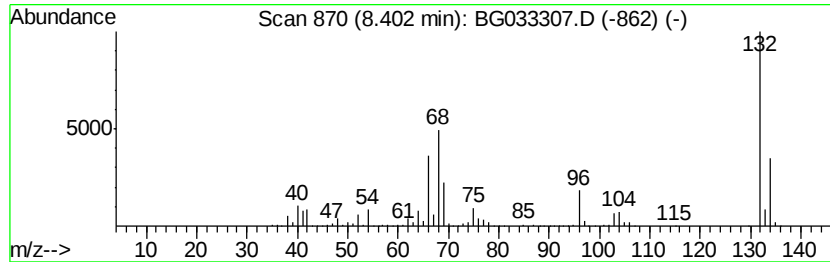
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Peak Number 2 unknown8.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	133.90 ng	1418870	1,4-Dichlorobenzene-d4	8.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.79	5.3	ng	56613	1	8.90	211923	20.0
unknown8.40	8.40	133.9	ng	1418870	1	8.90	211923	20.0