

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032718\
 Data File : BG033364.D
 Acq On : 27 Mar 2018 16:27
 Operator : SJ/JU
 Sample : PB107684BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107684BL

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	5.773	418	423	430	rBV	33263	57370	1.33%	0.305%
2	6.284	503	510	528	rBV	660576	1311930	30.50%	6.980%
3	7.970	789	797	822	rBV	672618	1475175	34.30%	7.849%
4	8.376	857	866	895	rBV	648844	1392725	32.38%	7.410%
5	8.869	941	950	960	rBV	99362	231866	5.39%	1.234%
6	9.198	995	1006	1020	rBV	507093	1035218	24.07%	5.508%
7	10.068	1146	1154	1172	rBV	364974	860828	20.01%	4.580%
8	11.748	1431	1440	1455	rBV	151837	345126	8.02%	1.836%
9	14.110	1835	1842	1861	rBV	1324194	2238723	52.05%	11.912%
10	15.485	2068	2076	2089	rBV2	320547	564649	13.13%	3.004%
11	16.954	2318	2326	2350	rBV	1214615	2129960	49.52%	11.333%
12	18.211	2533	2540	2560	rBV2	487463	849610	19.75%	4.521%
13	20.732	2963	2969	2981	rBV	3054156	4301178	100.00%	22.885%
14	22.559	3273	3280	3293	rBV2	526383	1017670	23.66%	5.415%
15	26.331	3909	3922	3936	rBV	273981	982519	22.84%	5.228%

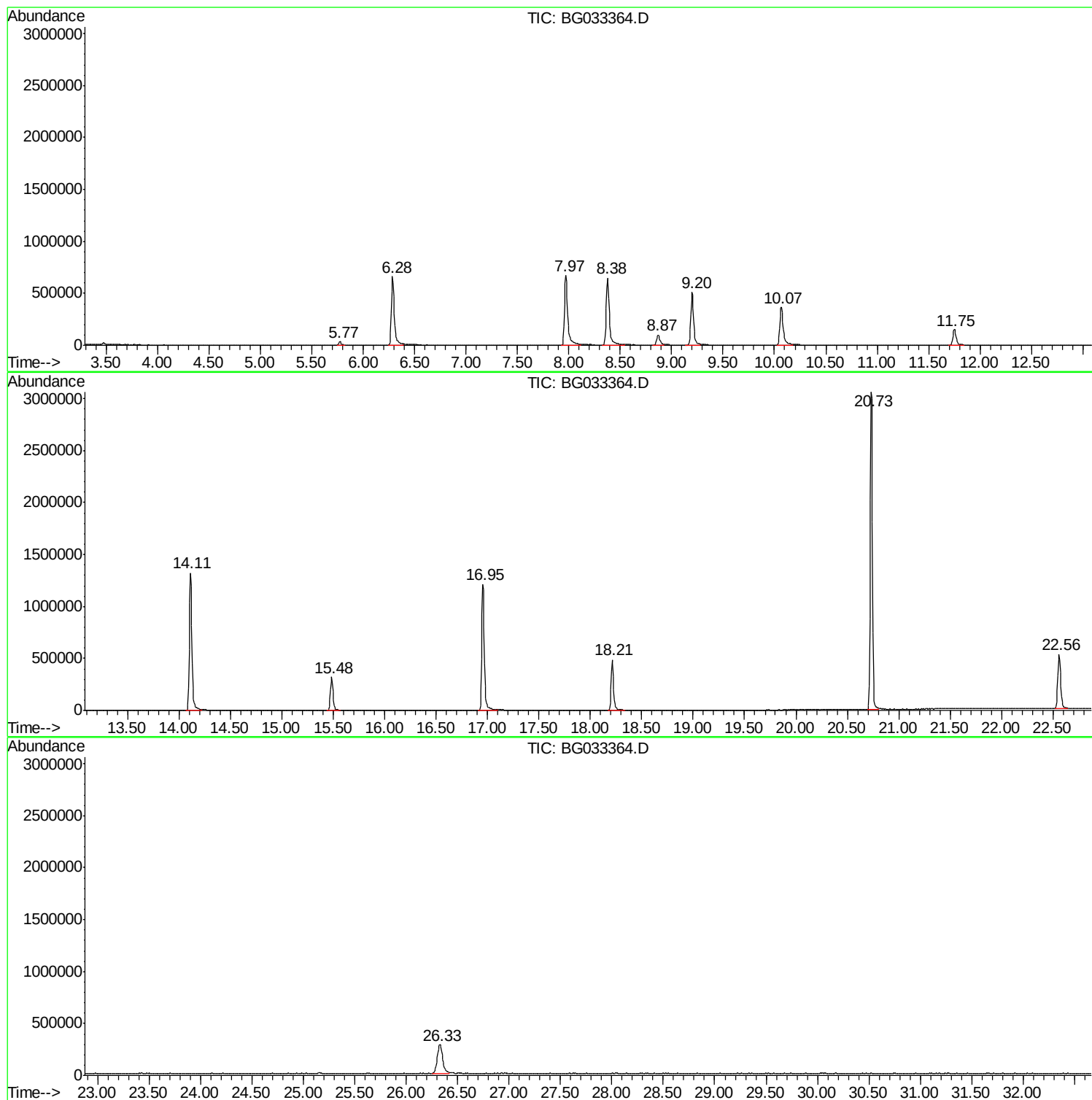
Sum of corrected areas: 18794547

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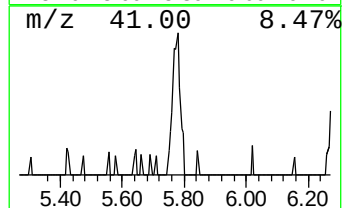
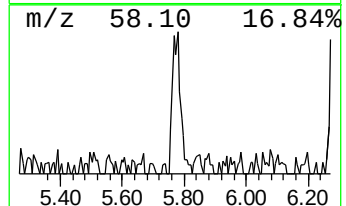
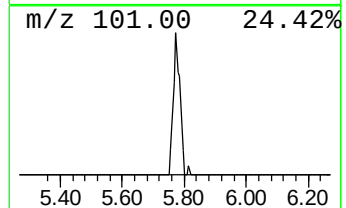
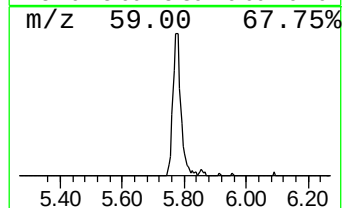
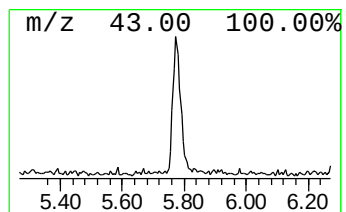
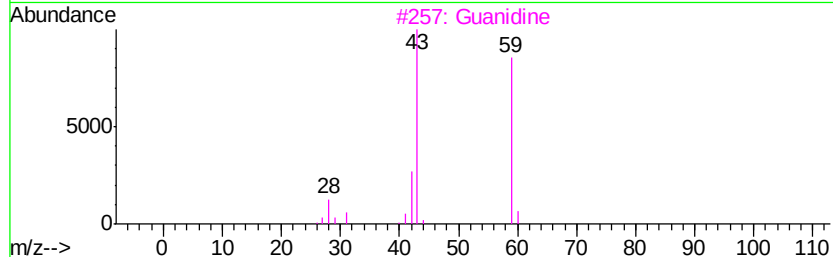
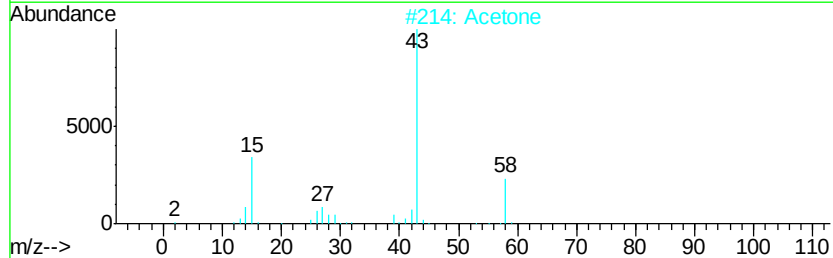
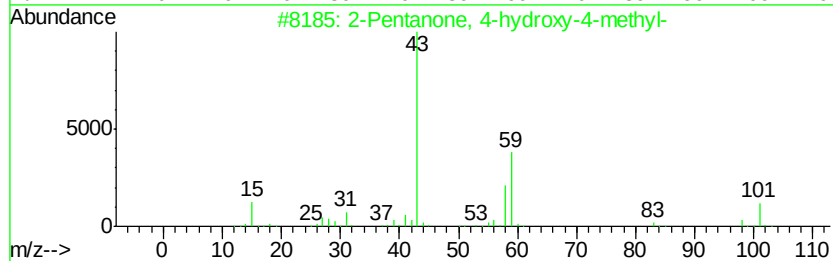
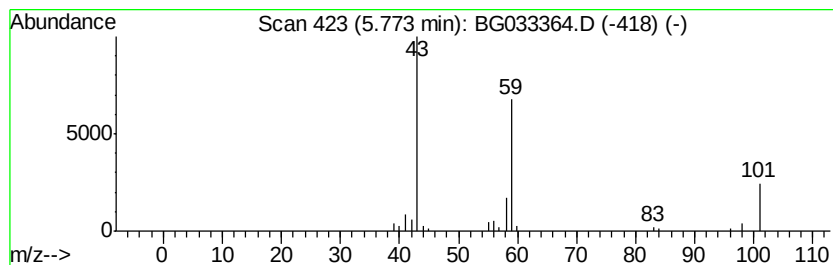
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.77	4.95 ng	57370	1,4-Dichlorobenzene-d4	8.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
2			Acetone	58	C3H6O	000067-64-1	9
3			Guanidine	59	CH5N3	000113-00-8	9
4			Acetamide	59	C2H5NO	000060-35-5	9
5			3-Hexanol	102	C6H14O	000623-37-0	9



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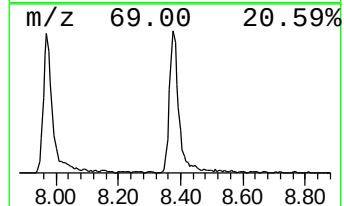
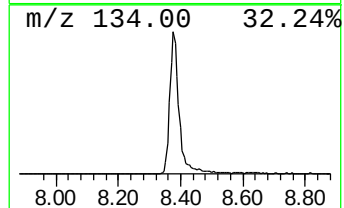
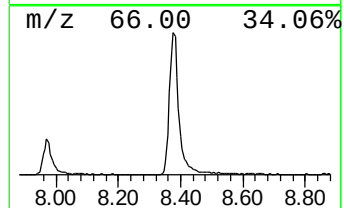
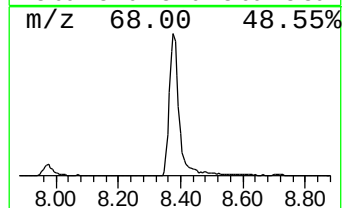
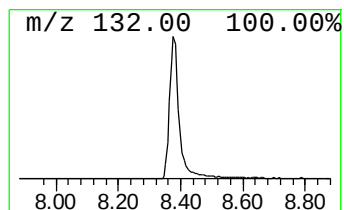
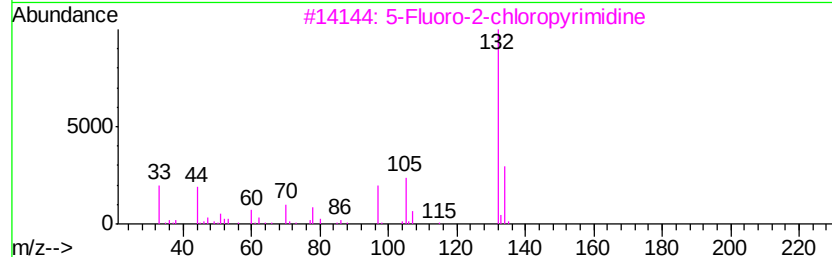
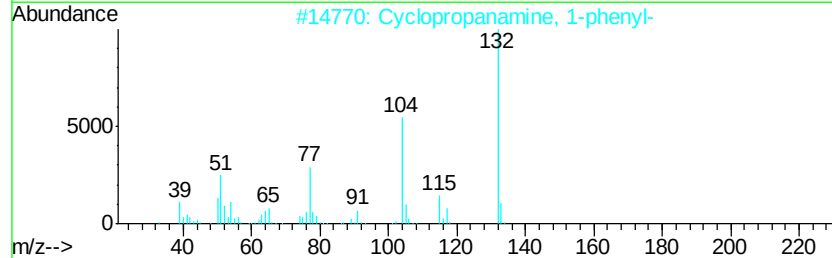
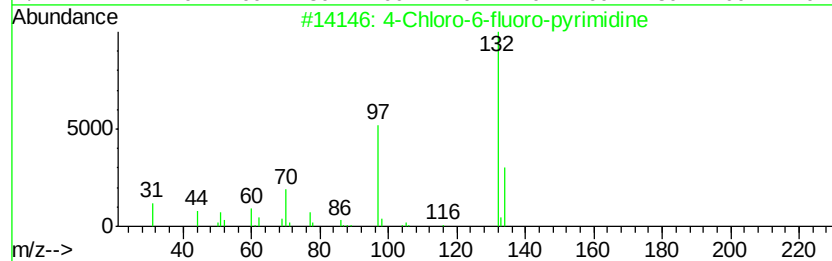
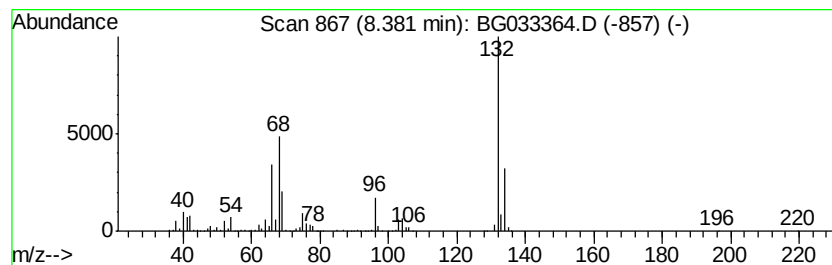
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Peak Number 2 unknown8.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	120.13 ng	1392730	1,4-Dichlorobenzene-d4	8.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	12
5		N-(-3,4-Methylenedioxyphenyl)isop...	267	C17H17NO2	1000378-96-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.77	5.0	ng	57370	1	8.87	231866	20.0
unknown8.38	8.38	120.1	ng	1392730	1	8.87	231866	20.0