

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092117\
 Data File : BG028871.D
 Acq On : 21 Sep 2017 23:43
 Operator : SJ/JU
 Sample : I5331-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HR-05-92017-B

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-BG091217.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	5.392	300	307	316	rBV	206146	323255	7.75%	1.900%
2	5.856	380	386	399	rBV	711683	1219870	29.25%	7.170%
3	7.443	647	656	670	rBV	661513	1199112	28.75%	7.048%
4	7.825	712	721	734	rBV	716942	1294207	31.03%	7.606%
5	8.289	792	800	807	rBV	121661	211735	5.08%	1.244%
6	8.612	846	855	864	rBV	568922	1000414	23.99%	5.880%
7	9.458	991	999	1013	rBV	452924	839482	20.13%	4.934%
8	11.109	1272	1280	1287	rBV	150406	289542	6.94%	1.702%
9	13.524	1684	1691	1703	rBV	1358169	2128762	51.05%	12.511%
10	14.905	1919	1926	1935	rBV2	264059	446232	10.70%	2.623%
11	16.397	2173	2180	2193	rBV	1072096	1805984	43.31%	10.614%
12	17.660	2387	2395	2405	rBV2	401051	657754	15.77%	3.866%
13	20.245	2829	2835	2848	rBV	3196044	4170221	100.00%	24.510%
14	21.985	3123	3131	3141	rBV2	401585	782448	18.76%	4.599%
15	25.457	3710	3722	3743	rBV2	162733	645489	15.48%	3.794%

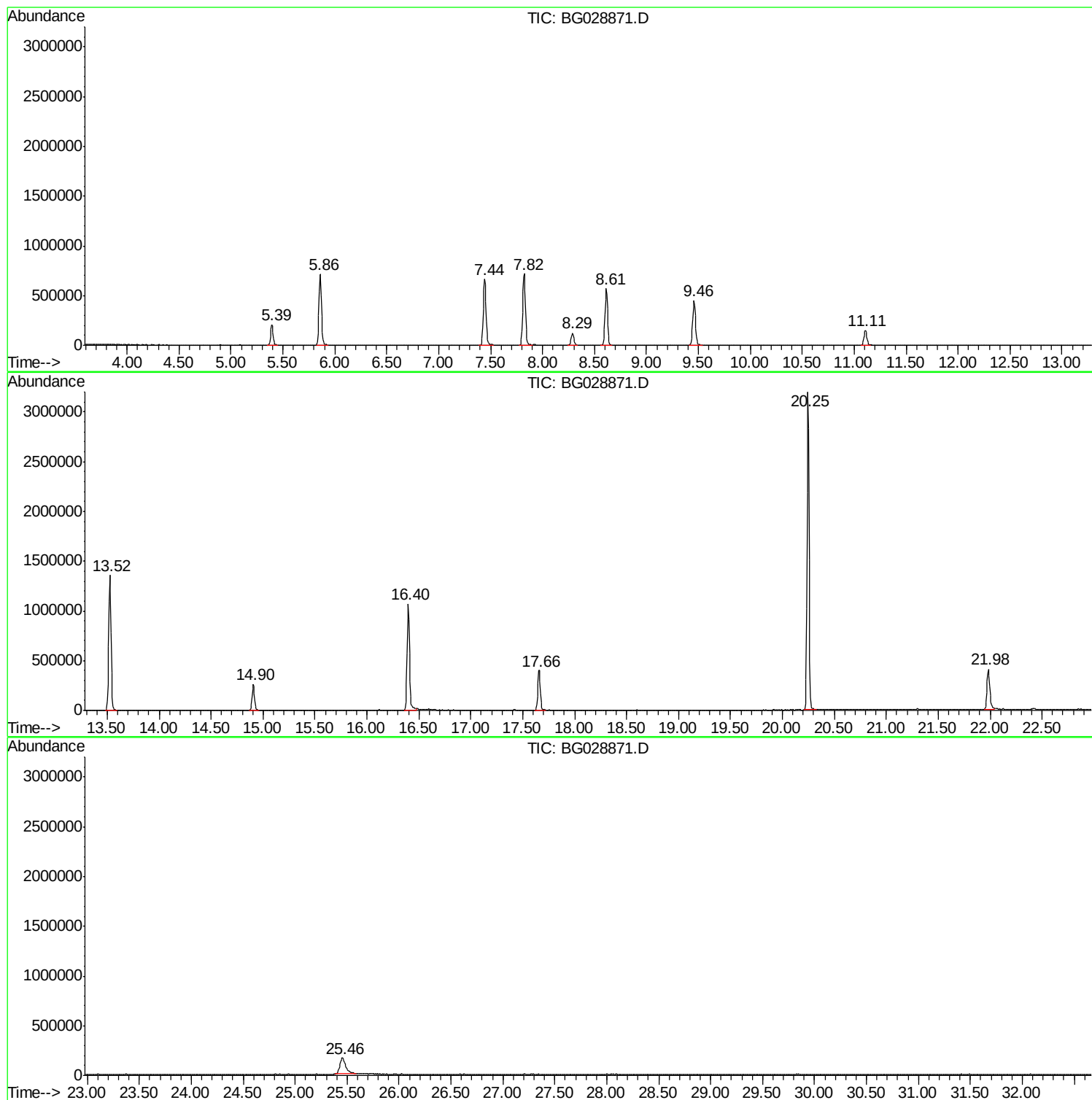
Sum of corrected areas: 17014507

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.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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ClientSampled :
HR-05-92017-B

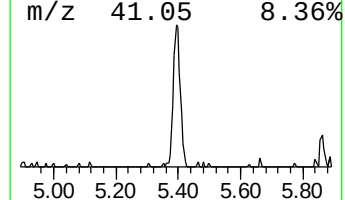
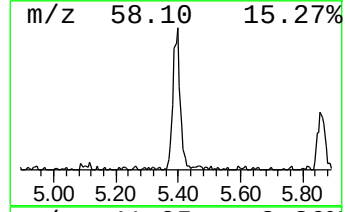
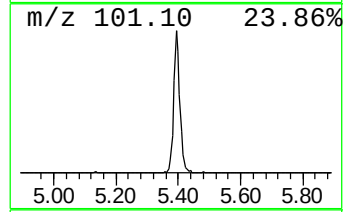
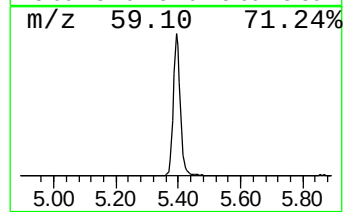
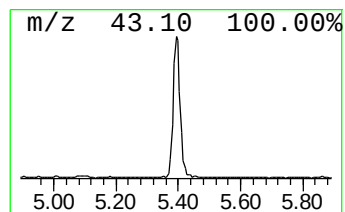
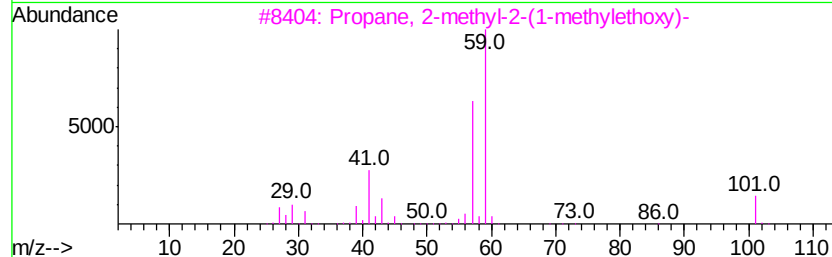
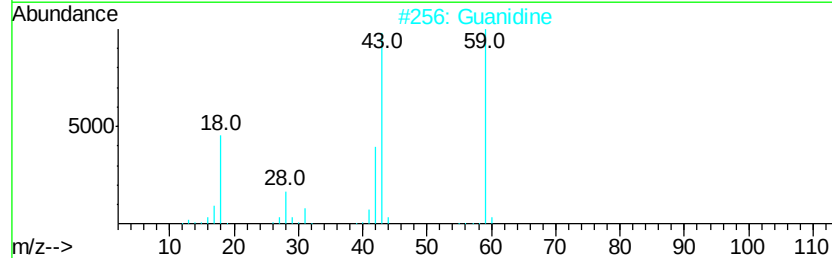
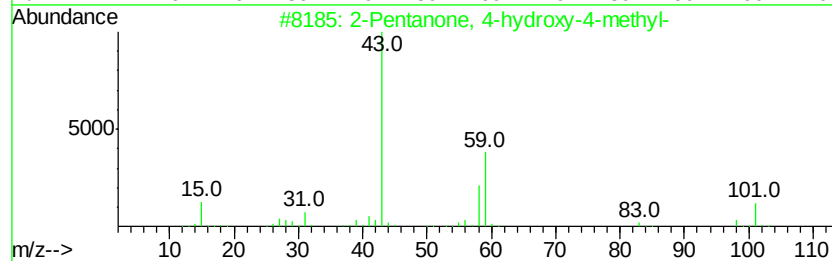
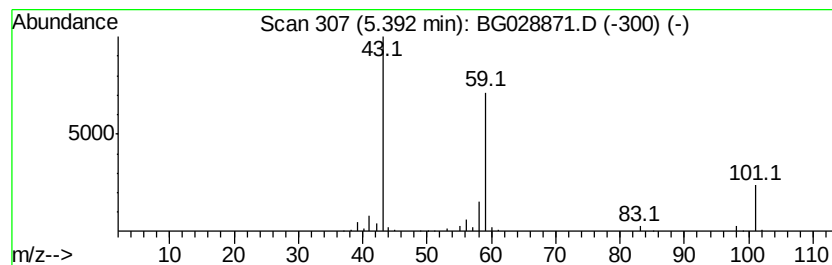
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.39	30.53 ng	323255	1,4-Dichlorobenzene-d4	8.29

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Guanidine	59	CH5N3	000113-00-8	43
3		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	42
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	32
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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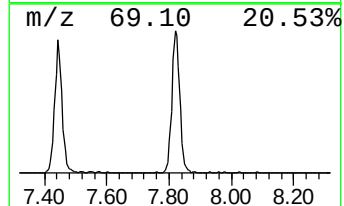
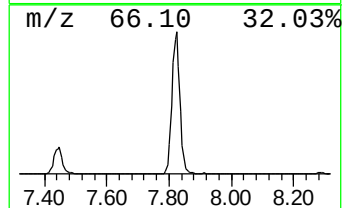
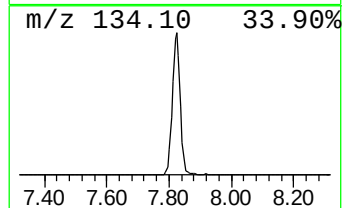
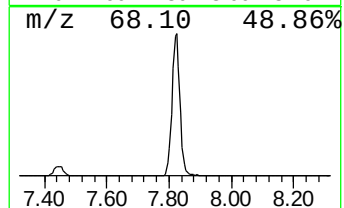
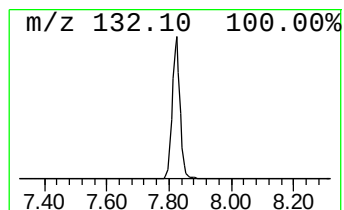
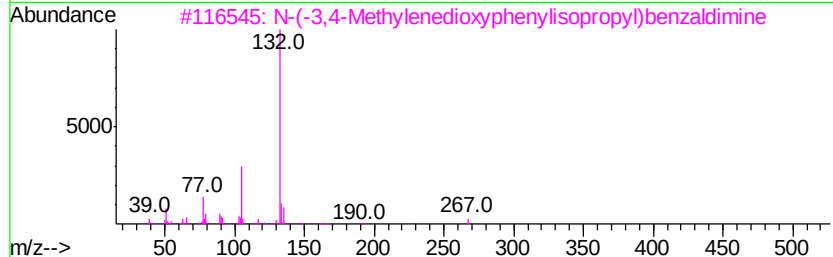
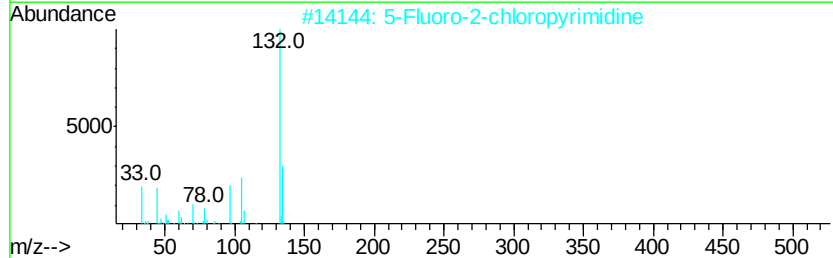
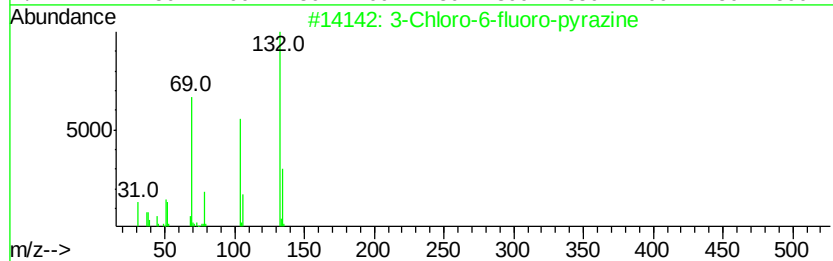
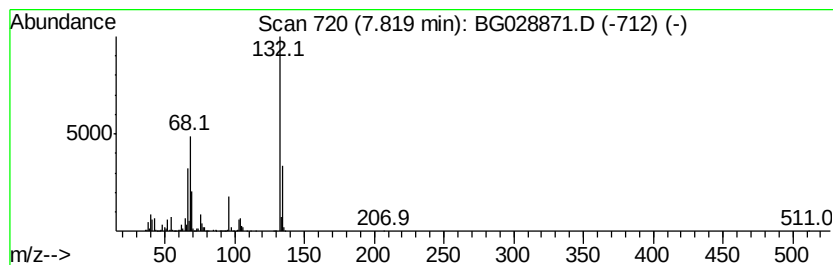
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	122.25 ng	1294210	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		N-(-3,4-Methylenedioxyphenylisopropyl)-	267	C17H17NO2	1000378-96-6	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.39	30.5	ng	323255	1	8.29	211735	20.0
unknown7.82	7.82	122.3	ng	1294210	1	8.29	211735	20.0