

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092117\
 Data File : BG028870.D
 Acq On : 21 Sep 2017 23:03
 Operator : SJ/JU
 Sample : I5331-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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

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	317	rBV	190424	297117	7.12%	1.762%
2	5.862	379	387	397	rBV	716731	1212112	29.06%	7.190%
3	7.442	649	656	674	rBV	647095	1194470	28.64%	7.086%
4	7.824	712	721	733	rBV	720815	1281608	30.73%	7.602%
5	8.288	790	800	808	rBV	119844	219228	5.26%	1.300%
6	8.612	847	855	869	rBV	547707	992100	23.79%	5.885%
7	9.458	991	999	1015	rBV	441190	835304	20.03%	4.955%
8	11.109	1272	1280	1287	rBV	144267	284070	6.81%	1.685%
9	13.523	1684	1691	1701	rBV	1325633	2111318	50.62%	12.524%
10	14.904	1920	1926	1936	rBV2	244801	428942	10.28%	2.544%
11	16.397	2173	2180	2194	rBV	1021616	1774285	42.54%	10.525%
12	17.660	2387	2395	2408	rBV	376816	626917	15.03%	3.719%
13	20.245	2829	2835	2848	rBV	3073839	4171042	100.00%	24.743%
14	21.984	3124	3131	3143	rBV2	389199	760245	18.23%	4.510%
15	25.451	3709	3721	3749	rBV2	165048	669010	16.04%	3.969%

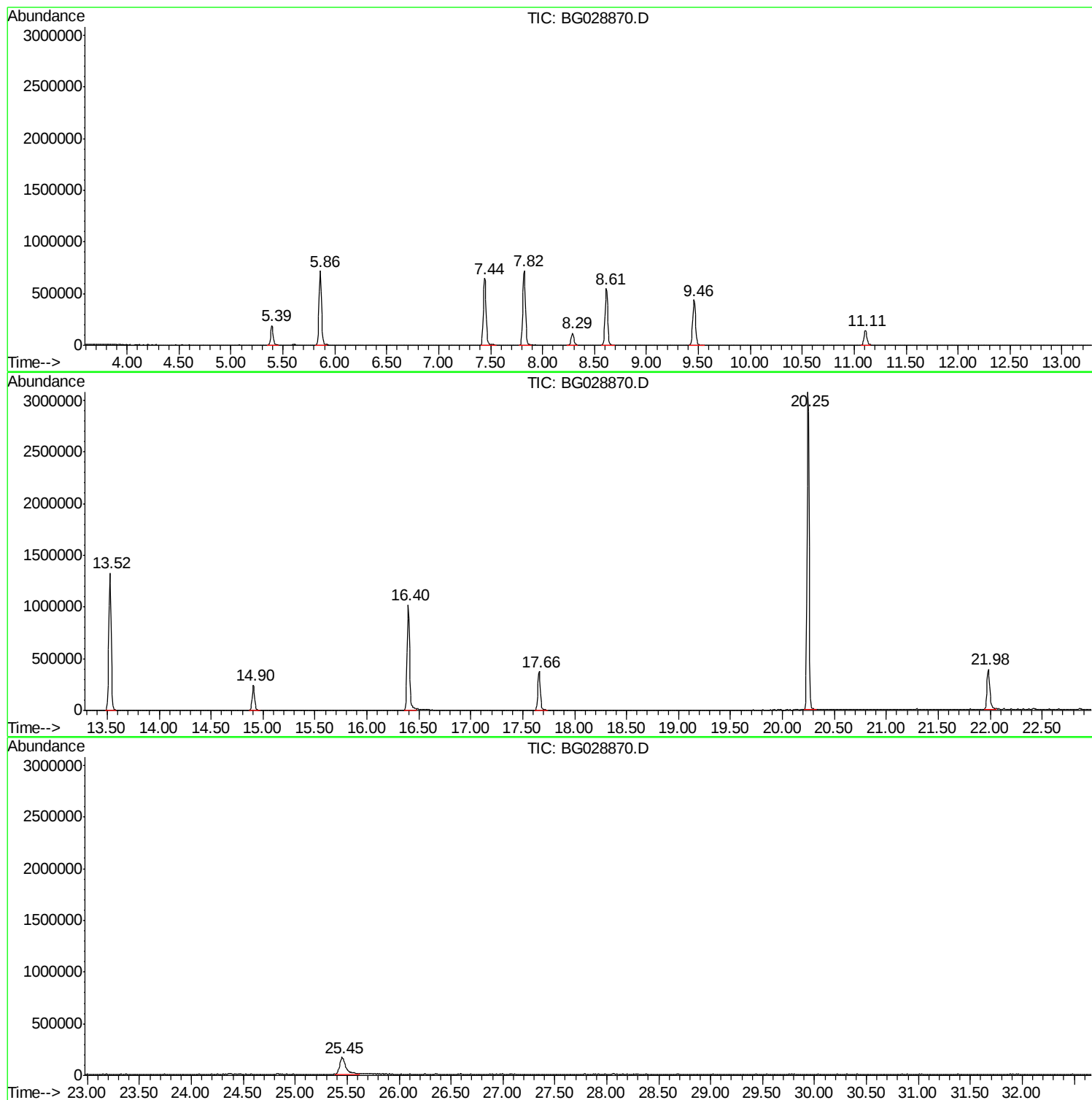
Sum of corrected areas: 16857768

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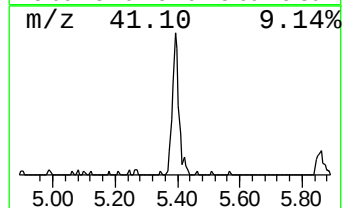
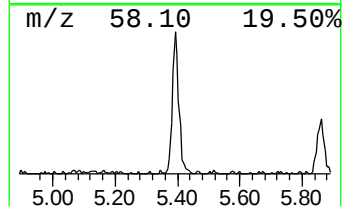
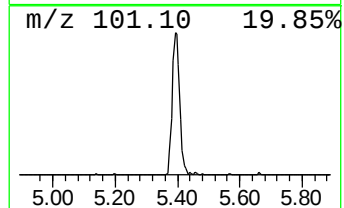
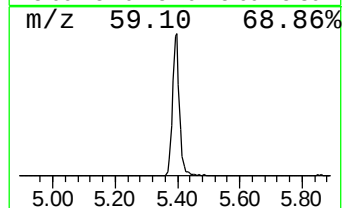
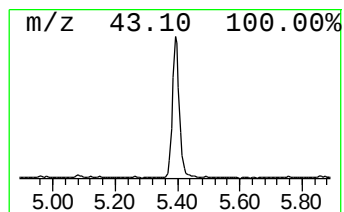
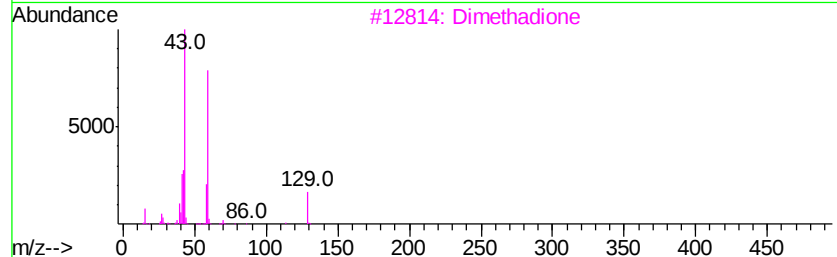
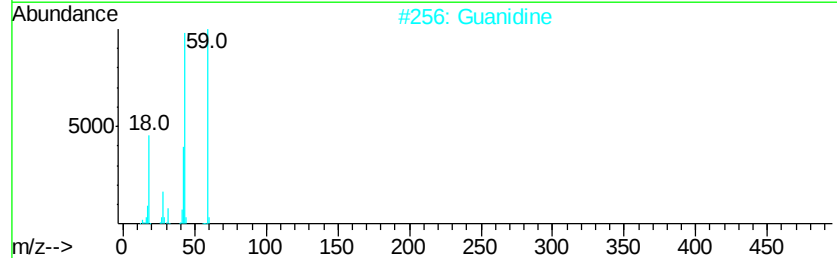
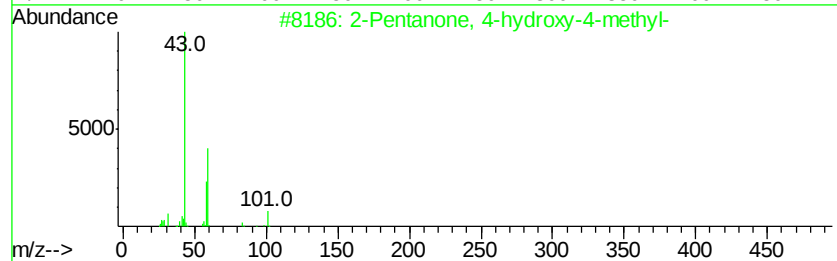
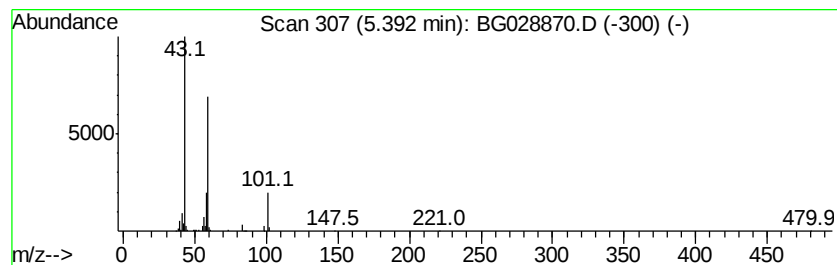
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.39	27.11 ng	297117	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	45
2	Guanidine	59	CH5N3	000113-00-8	38
3	Dimethadione	129	C5H7NO3	000695-53-4	33
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28



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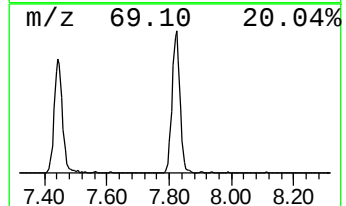
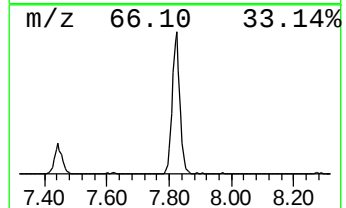
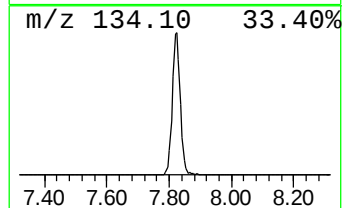
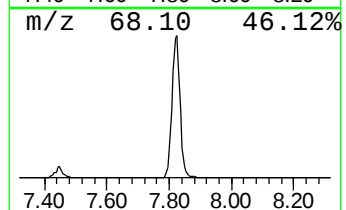
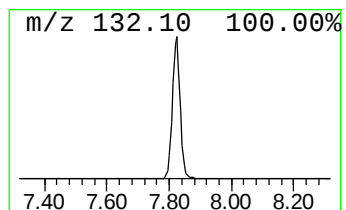
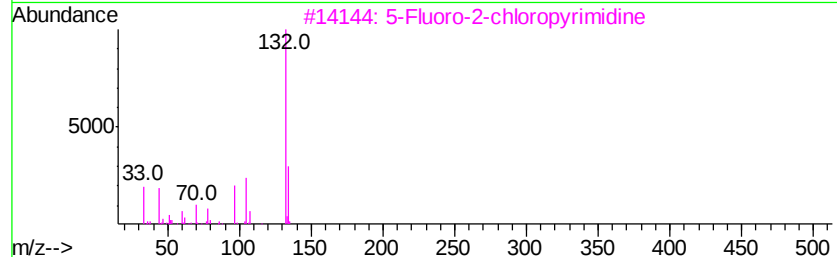
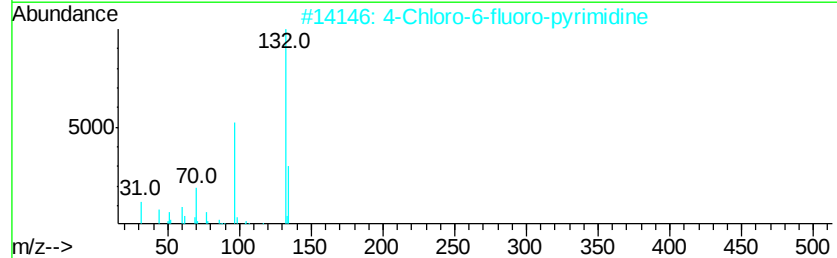
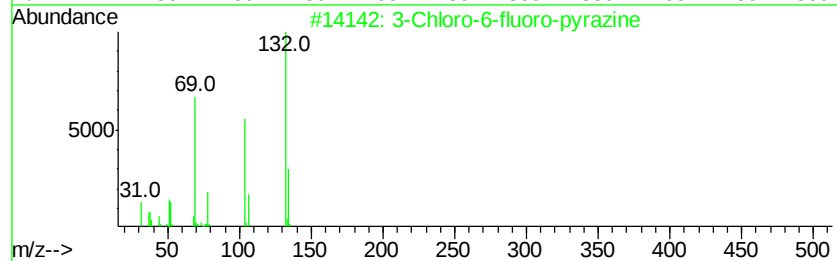
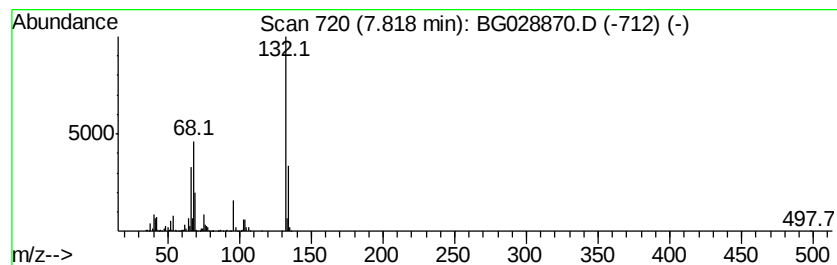
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	116.92 ng	1281610	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		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.39	27.1	ng	297117	1	8.29	219228	20.0
unknown7.82	7.82	116.9	ng	1281610	1	8.29	219228	20.0