

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050115\
 Data File : BG016850.D
 Acq On : 1 May 2015 18:13
 Operator : TP/IZ
 Sample : G2032-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B3GW

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-BG042915.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	4.379	281	288	296	rBV2	28458	44485	1.70%	0.342%
2	4.867	366	371	391	rBV	282177	473138	18.09%	3.633%
3	6.482	640	646	670	rBV	143222	307113	11.74%	2.358%
4	6.800	694	700	723	rBV	646402	1100484	42.07%	8.450%
5	7.264	772	779	798	rBV	139552	272094	10.40%	2.089%
6	7.575	825	832	850	rBV	615619	1019735	38.98%	7.830%
7	8.421	969	976	1000	rBV	432589	818838	31.30%	6.288%
8	10.031	1244	1250	1267	rBV	200898	391613	14.97%	3.007%
9	12.552	1672	1679	1699	rBV	1279390	2048148	78.29%	15.727%
10	13.944	1909	1916	1932	rBV2	330892	545261	20.84%	4.187%
11	15.448	2166	2172	2191	rBV	879810	1350212	51.61%	10.368%
12	16.700	2379	2385	2399	rBV2	407242	647158	24.74%	4.969%
13	19.355	2832	2837	2849	rBV	2095385	2616131	100.00%	20.089%
14	20.907	3096	3101	3110	rBV	549122	707140	27.03%	5.430%
15	22.981	3448	3454	3465	rVB2	398775	681326	26.04%	5.232%

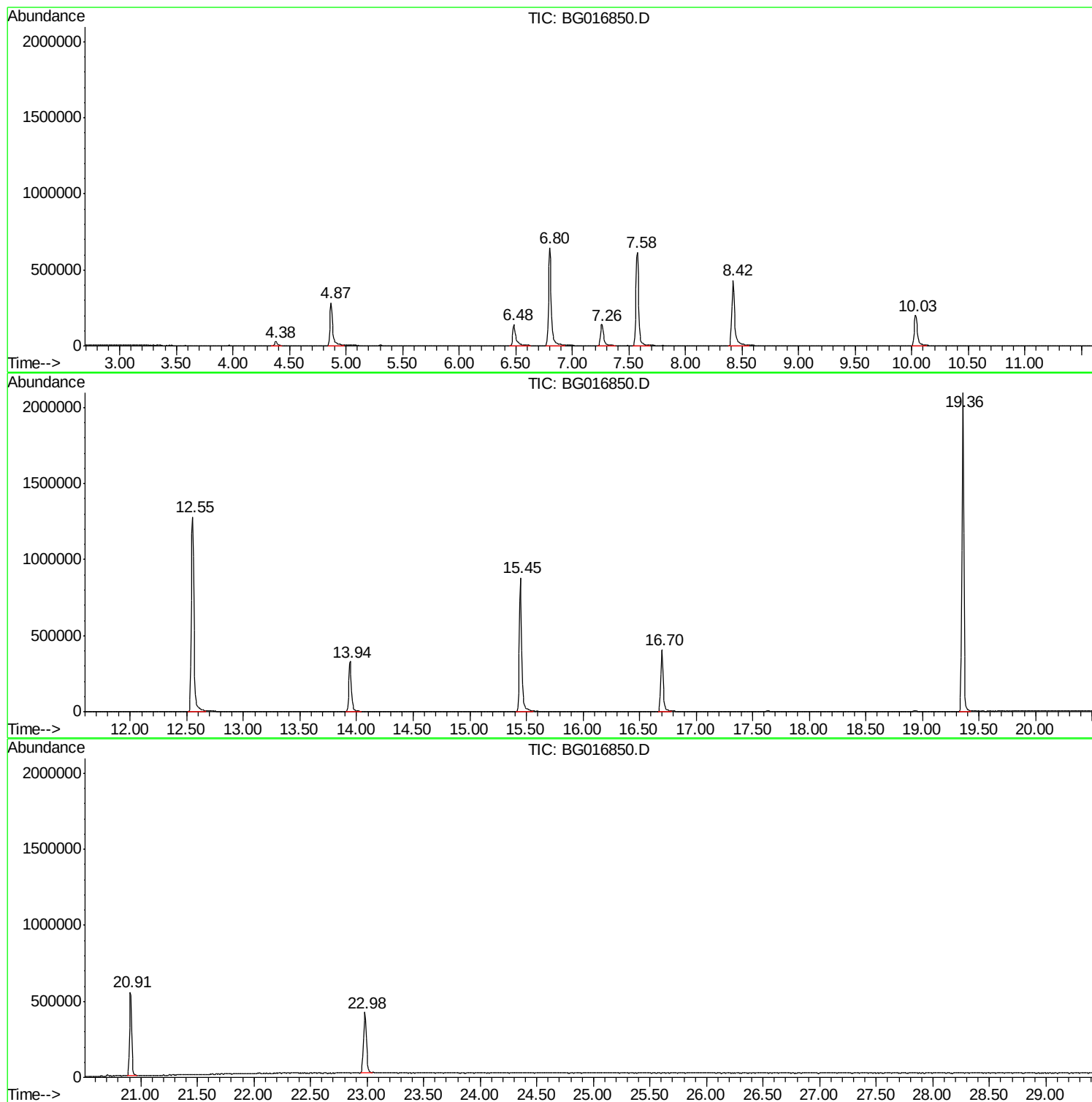
Sum of corrected areas: 13022876

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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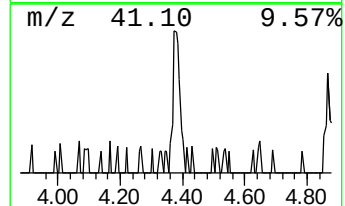
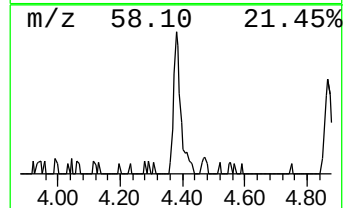
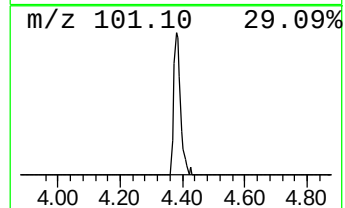
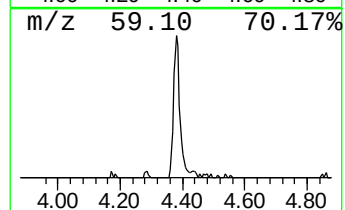
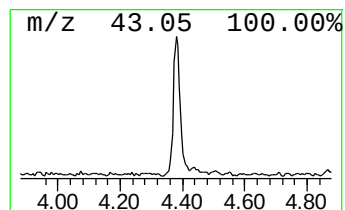
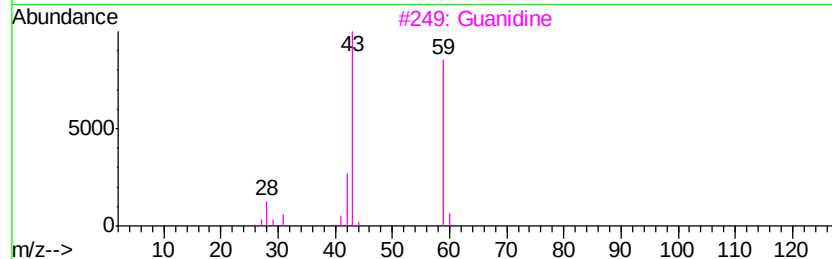
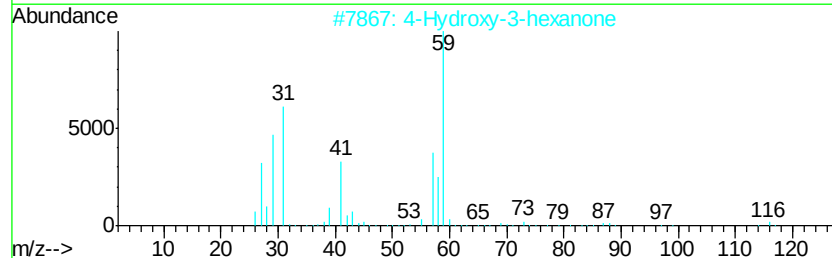
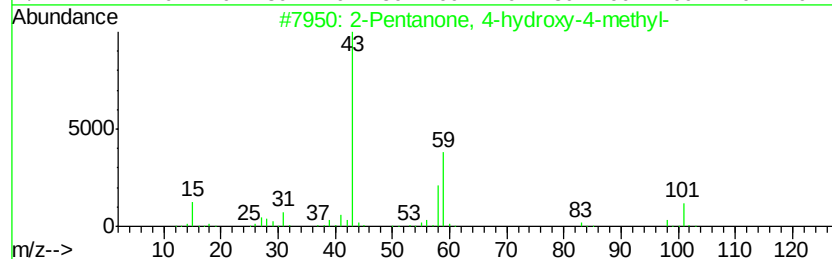
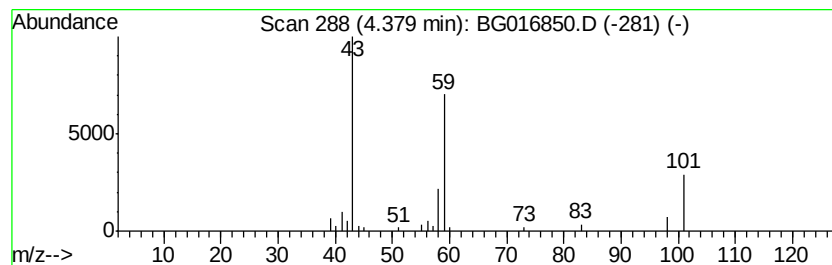
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	3.27 ng	44485	1,4-Dichlorobenzene-d4	7.26

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	25
3		Guanidine	59	CH5N3	000113-00-8	9
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		Hydrazine, (1-methylethyl)-	74	C3H10N2	002257-52-5	9



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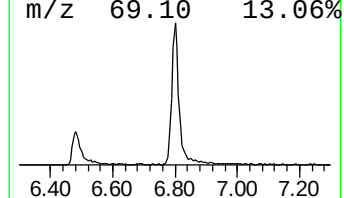
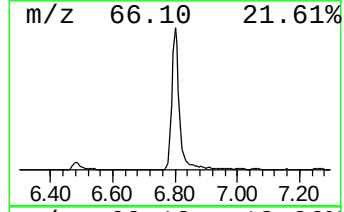
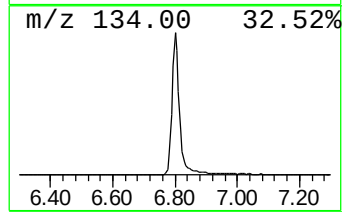
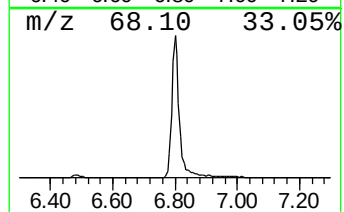
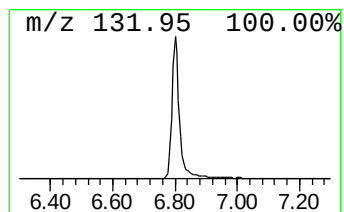
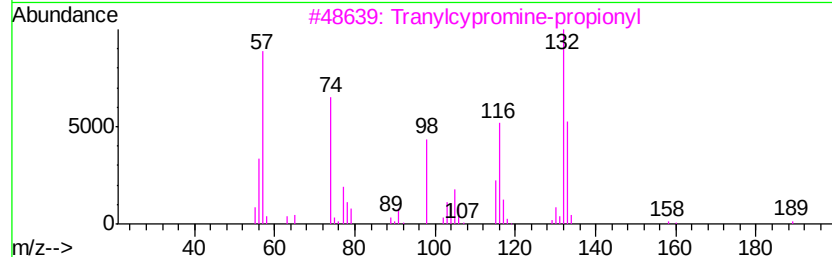
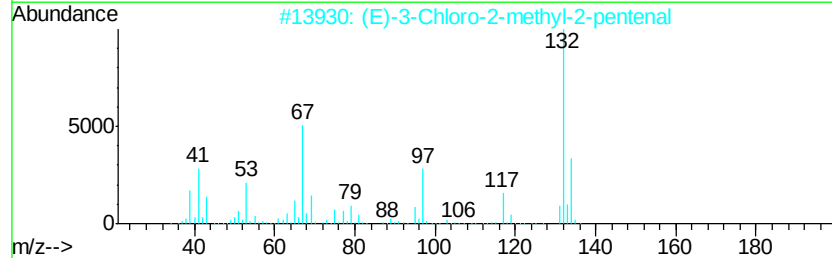
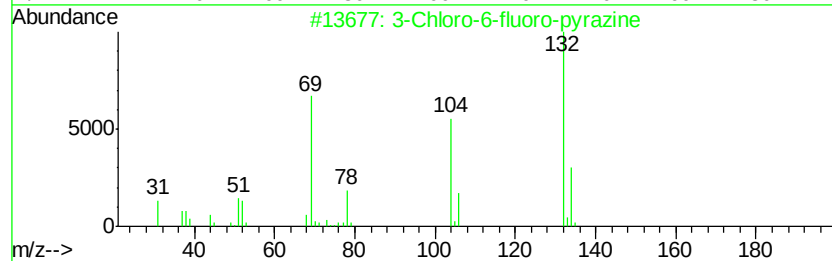
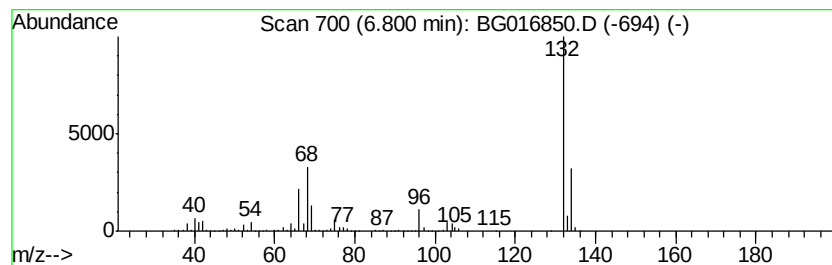
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Peak Number 2 unknown6.80 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	80.89 ng	1100480	1,4-Dichlorobenzene-d4	7.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.38	3.3	ng	44485	1	7.26	272094	20.0
unknown6.80	6.80	80.9	ng	1100480	1	7.26	272094	20.0