

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061517\
 Data File : BG027453.D
 Acq On : 15 Jun 2017 22:23
 Operator : SJ/MA
 Sample : PB99795BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99795BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061417.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	5.658	329	335	345	rBV	87619	141698	3.09%	0.777%
2	6.105	403	411	425	rBV	779929	1340380	29.20%	7.355%
3	7.656	667	675	692	rBV	762772	1411425	30.74%	7.744%
4	8.050	734	742	756	rBV	711792	1329291	28.95%	7.294%
5	8.520	813	822	829	rBV	122394	214057	4.66%	1.175%
6	8.849	869	878	889	rVB	503251	939021	20.45%	5.152%
7	9.683	1011	1020	1028	rBV	418072	801738	17.46%	4.399%
8	11.340	1293	1302	1311	rVB	174407	327048	7.12%	1.795%
9	13.731	1701	1709	1716	rBV	1208534	1985337	43.24%	10.894%
10	15.106	1936	1943	1951	rBV2	280174	467295	10.18%	2.564%
11	16.593	2189	2196	2212	rBV	1276477	2015856	43.91%	11.061%
12	17.856	2404	2411	2419	rBV2	455882	748095	16.29%	4.105%
13	20.435	2843	2850	2858	rBV	3389194	4590954	100.00%	25.191%
14	22.233	3148	3156	3164	rVB	525761	961065	20.93%	5.273%
15	25.888	3767	3778	3795	rVB2	287279	951652	20.73%	5.222%

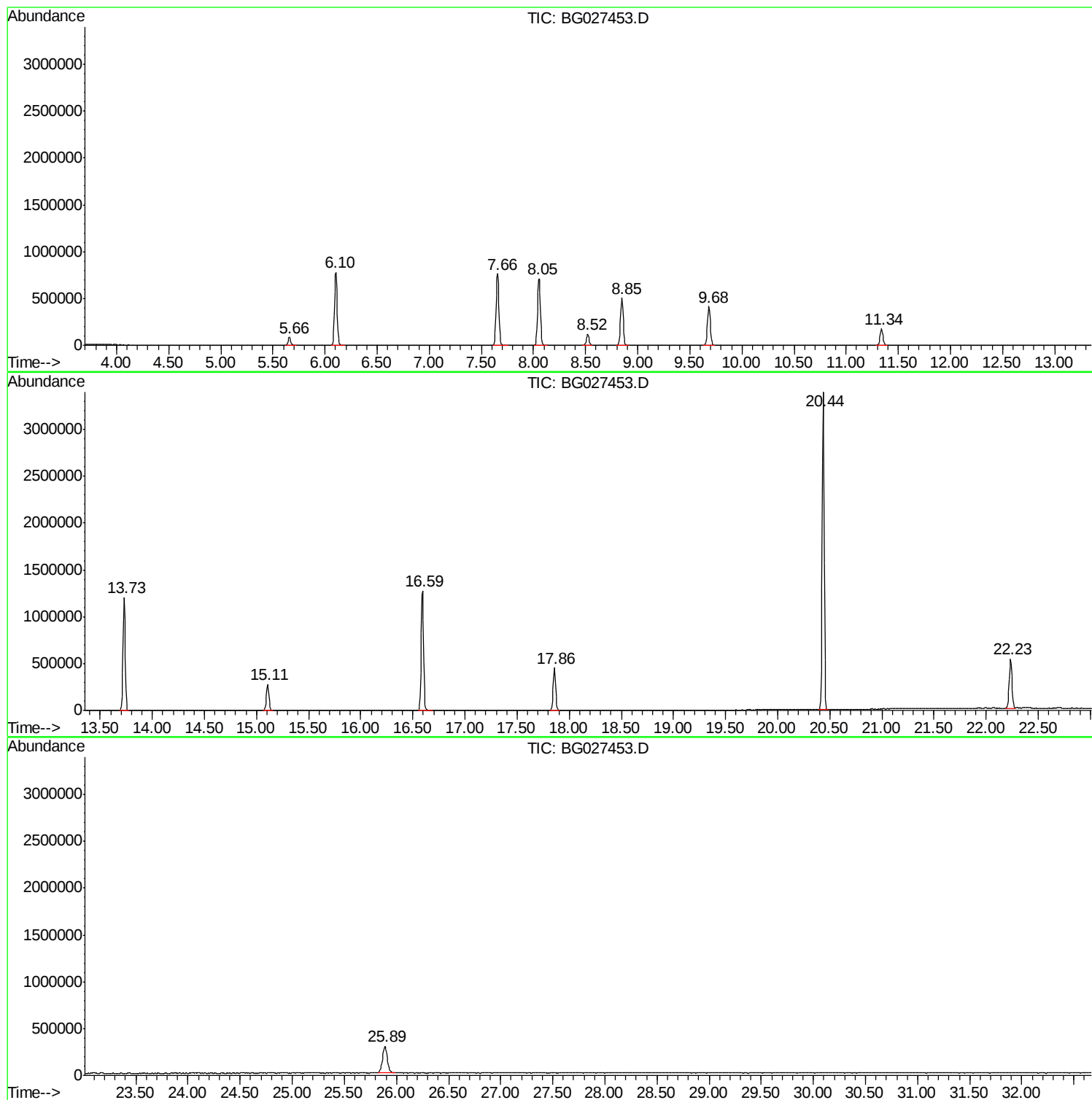
Sum of corrected areas: 18224912

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061417.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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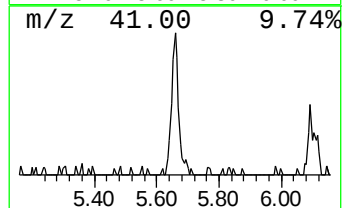
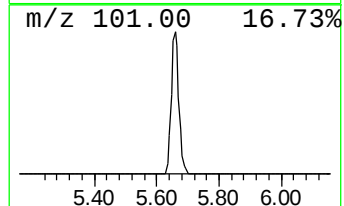
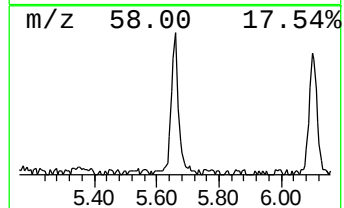
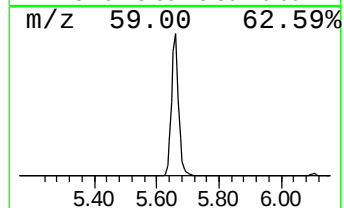
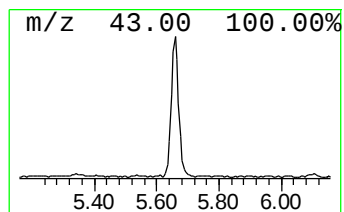
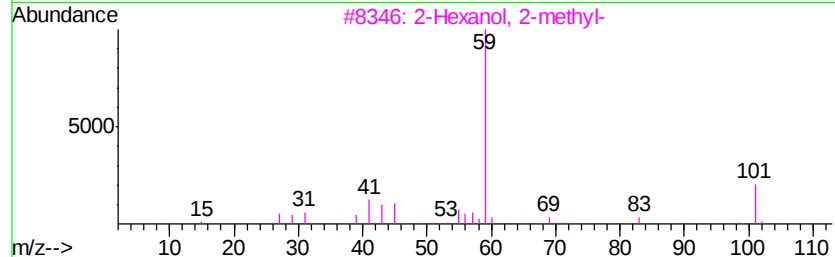
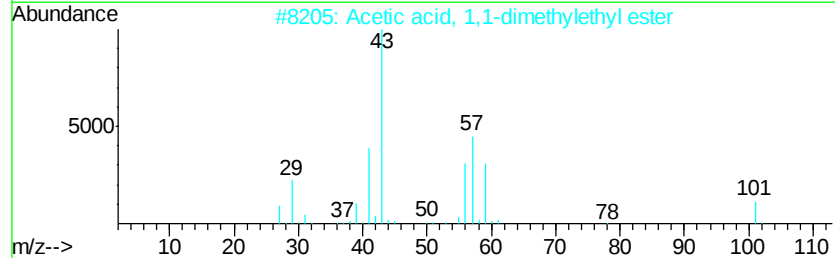
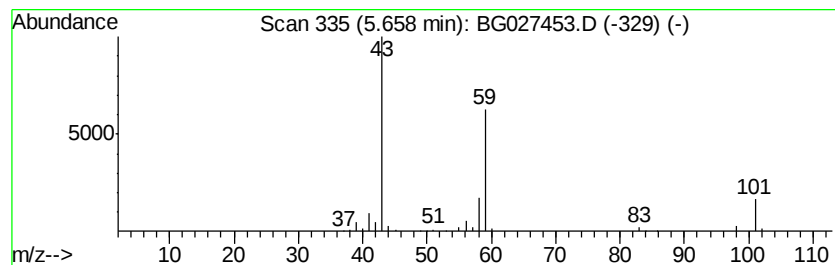
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.66	13.24 ng	141698	1,4-Dichlorobenzene-d4	8.52

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28	
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16	



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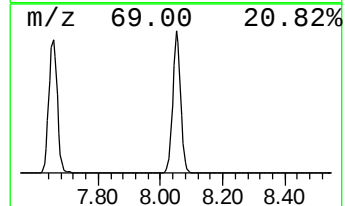
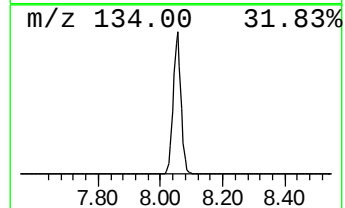
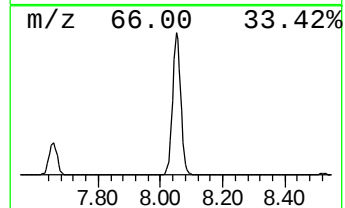
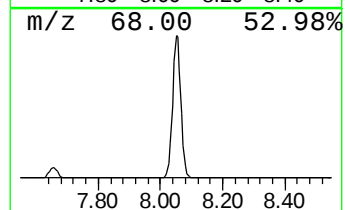
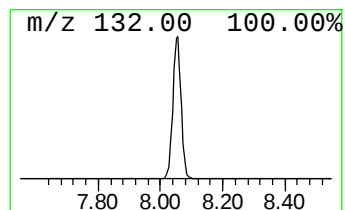
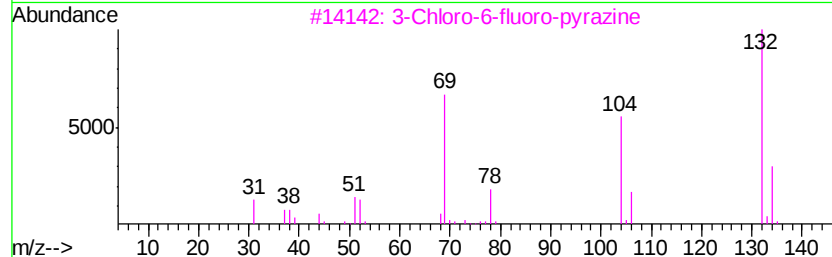
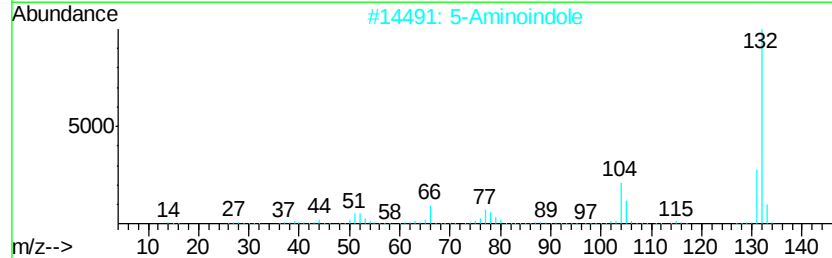
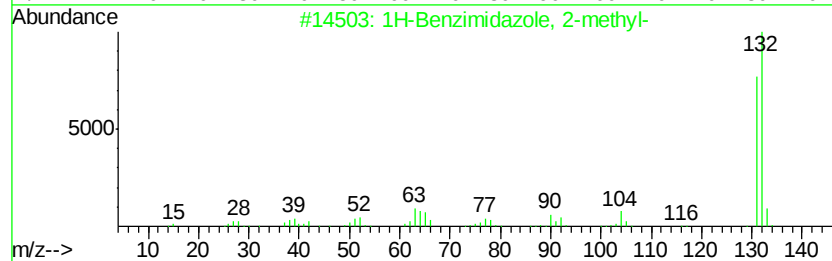
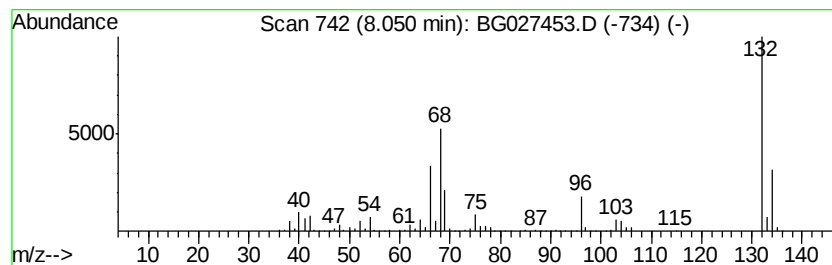
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Peak Number 2 unknown8.05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.05	124.20 ng	1329290	1,4-Dichlorobenzene-d4	8.52	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2	5-Aminoindole	132	C8H8N2	005192-03-0	22
3	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.66	13.2	ng	141698	1	8.52	214057	20.0
unknown8.05	8.05	124.2	ng	1329290	1	8.52	214057	20.0