

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092017\
 Data File : BG028837.D
 Acq On : 20 Sep 2017 18:33
 Operator : SJ/JU
 Sample : PB102470TB
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB102470TB

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	302	307	317	rBV	90578	145329	4.79%	1.057%
2	5.856	379	386	400	rBV	585870	997513	32.88%	7.257%
3	7.443	647	656	671	rBV	586319	1054123	34.75%	7.669%
4	7.825	711	721	731	rBV	579733	1016091	33.50%	7.392%
5	8.289	793	800	811	rVB	122202	207437	6.84%	1.509%
6	8.618	847	856	864	rVB	414817	756963	24.95%	5.507%
7	9.458	991	999	1011	rBV	329628	628967	20.73%	4.576%
8	11.109	1271	1280	1292	rBV	153860	283103	9.33%	2.060%
9	13.530	1684	1692	1702	rBV	1046757	1686793	55.61%	12.272%
10	14.910	1919	1927	1935	rBV2	263140	444124	14.64%	3.231%
11	16.397	2173	2180	2198	rBV	876610	1475959	48.66%	10.738%
12	17.660	2388	2395	2404	rBV	396311	626813	20.66%	4.560%
13	20.245	2829	2835	2849	rBV	2253306	3033516	100.00%	22.069%
14	21.984	3124	3131	3147	rBV2	396735	734030	24.20%	5.340%
15	25.451	3711	3721	3735	rBV2	191744	654686	21.58%	4.763%

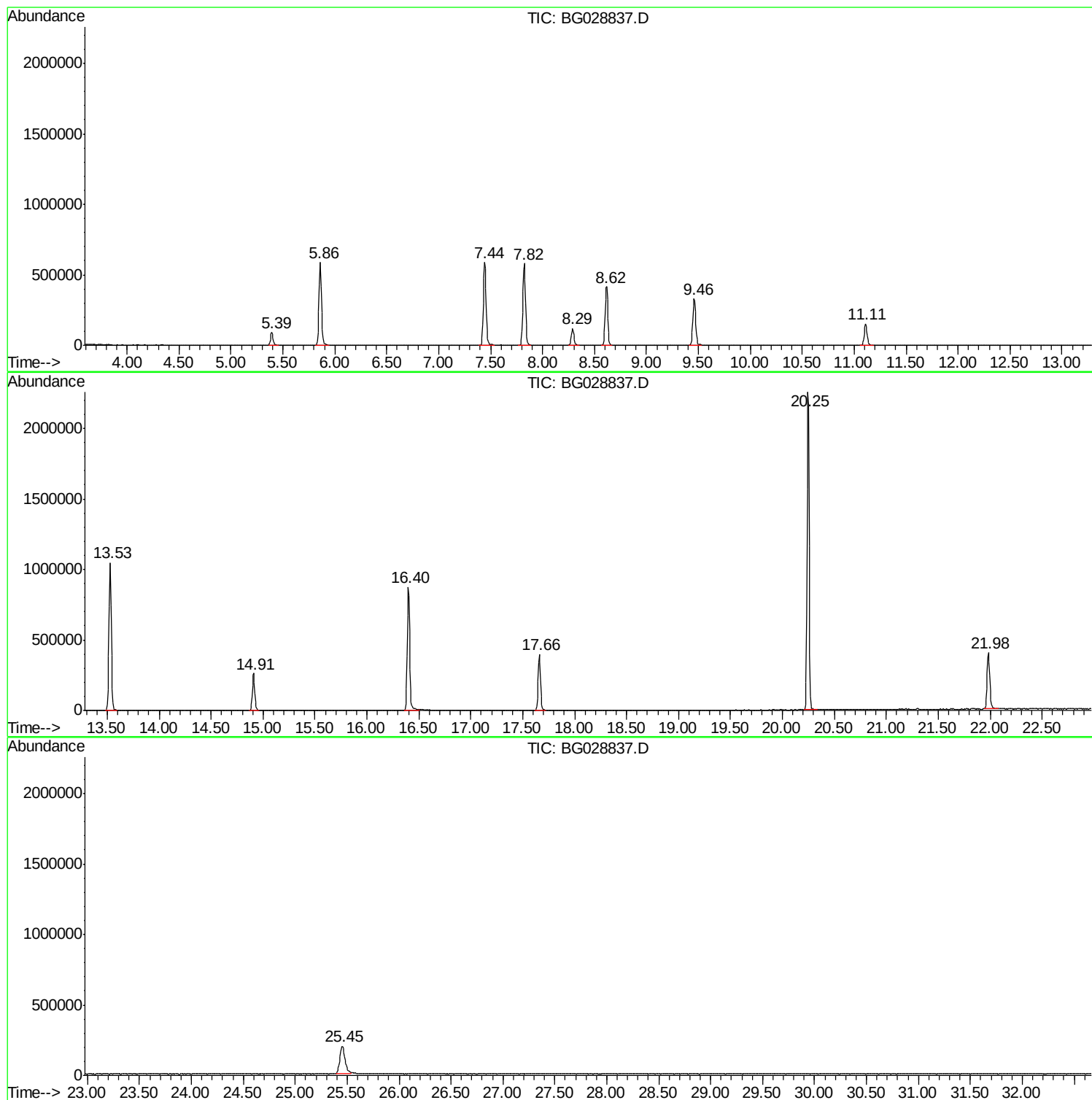
Sum of corrected areas: 13745447

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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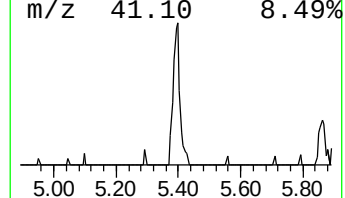
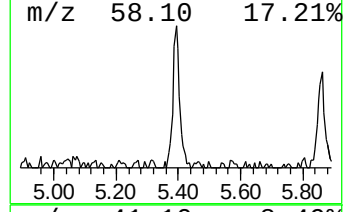
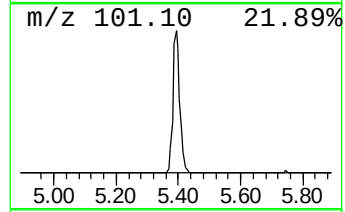
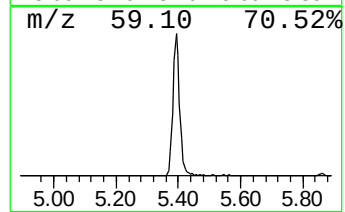
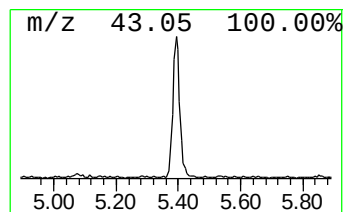
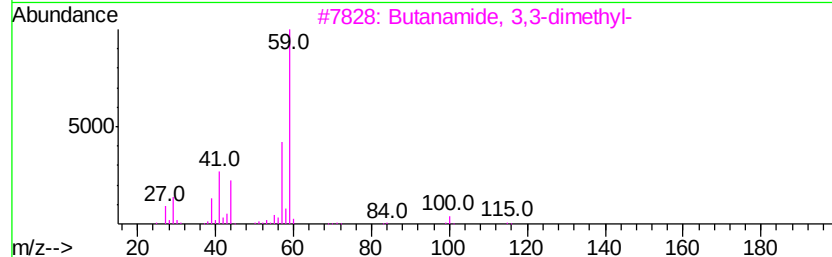
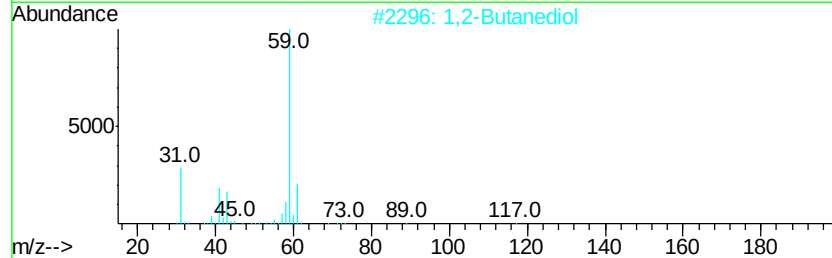
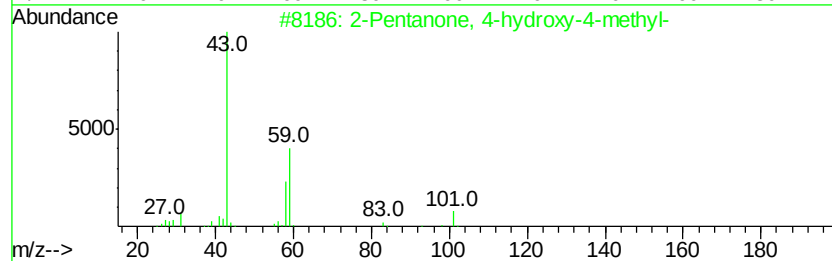
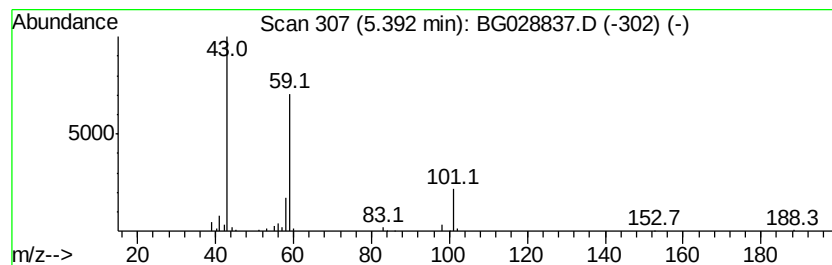
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.39	14.01 ng	145329	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		1,2-Butanediol	90	C4H10O2	000584-03-2	17
3		Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	17
4		Oxirane, tetramethyl-	100	C6H12O	005076-20-0	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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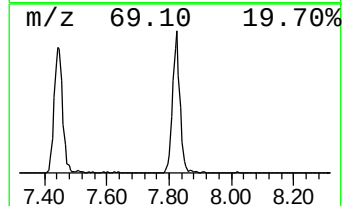
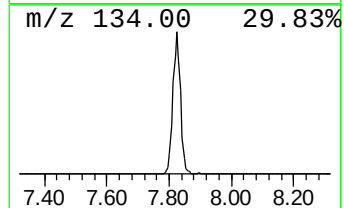
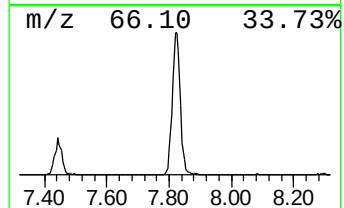
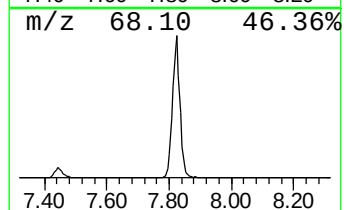
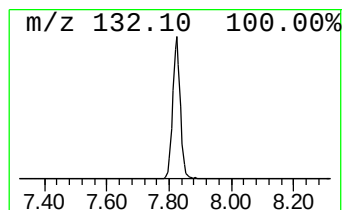
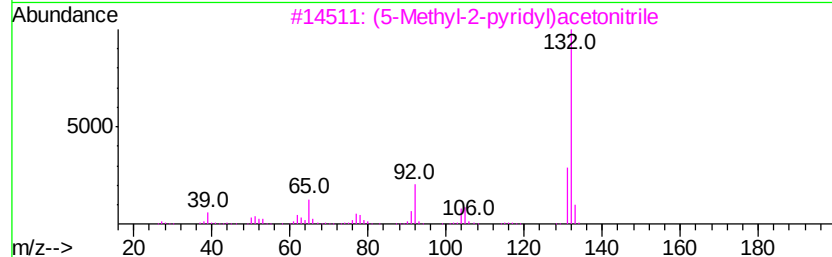
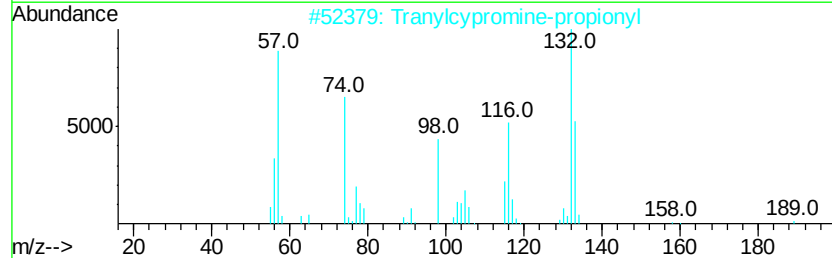
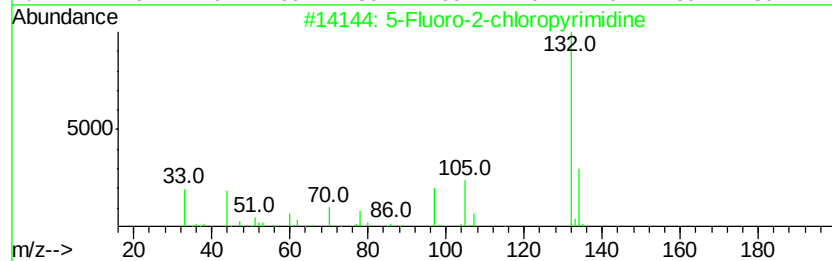
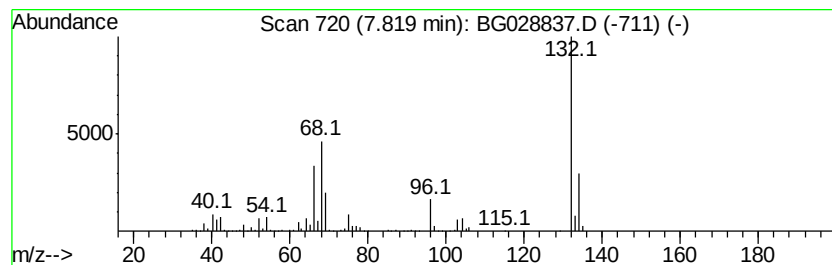
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	97.97 ng	1016090	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	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	14.0	ng	145329	1	8.29	207437	20.0
unknown7.82	7.82	98.0	ng	1016090	1	8.29	207437	20.0