

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092217\
 Data File : BG028879.D
 Acq On : 22 Sep 2017 18:01
 Operator : SJ/JU
 Sample : PB102487BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB102487BL

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	5.379	299	305	313	rBV	65692	103504	3.63%	0.947%
2	5.844	377	384	397	rBV	409565	711756	24.97%	6.509%
3	7.430	648	654	667	rBV	382892	723909	25.39%	6.620%
4	7.812	711	719	731	rBV	400586	710735	24.93%	6.500%
5	8.276	790	798	807	rVB2	75688	137090	4.81%	1.254%
6	8.605	846	854	864	rBV	297759	545318	19.13%	4.987%
7	9.445	990	997	1012	rBV	225603	440761	15.46%	4.031%
8	11.096	1268	1278	1288	rBV	92882	183303	6.43%	1.676%
9	13.511	1682	1689	1700	rBV	711909	1127882	39.57%	10.315%
10	14.898	1917	1925	1933	rBV2	159634	281802	9.89%	2.577%
11	16.384	2170	2178	2193	rBV	775685	1267348	44.46%	11.590%
12	17.642	2386	2392	2400	rVB2	318499	500629	17.56%	4.578%
13	20.227	2826	2832	2846	rBV	2080902	2850651	100.00%	26.070%
14	20.996	2955	2963	2964	rBV7	18160	35620	1.25%	0.326%
15	21.960	3120	3127	3136	rBV	370847	663317	23.27%	6.066%
16	25.409	3703	3714	3730	rBV2	206582	650815	22.83%	5.952%

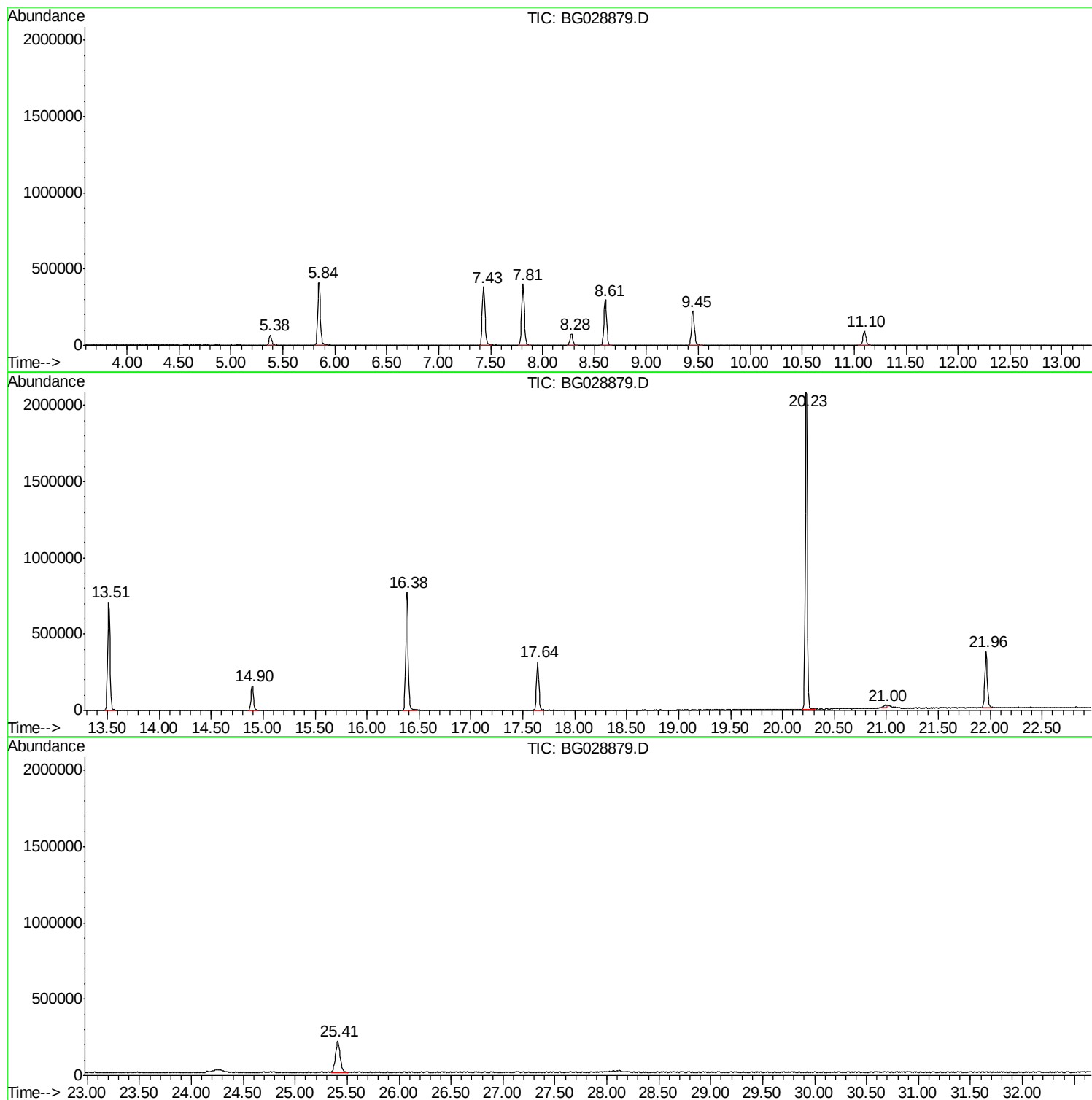
Sum of corrected areas: 10934440

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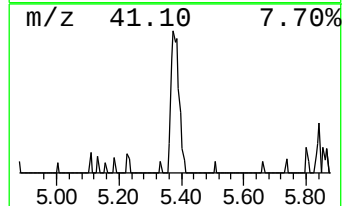
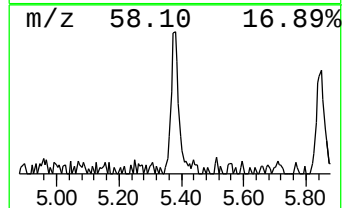
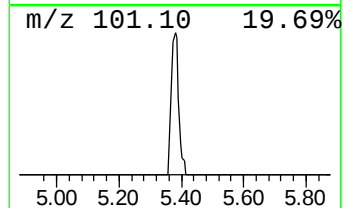
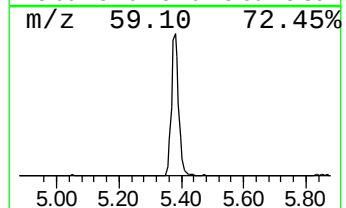
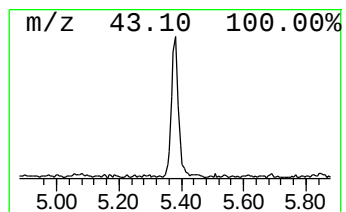
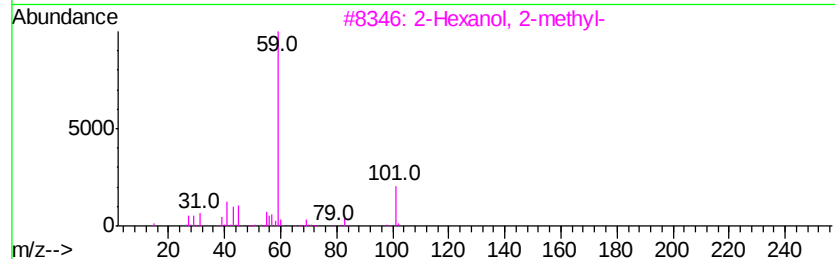
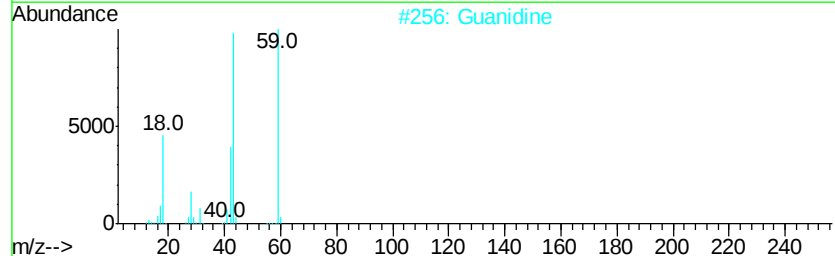
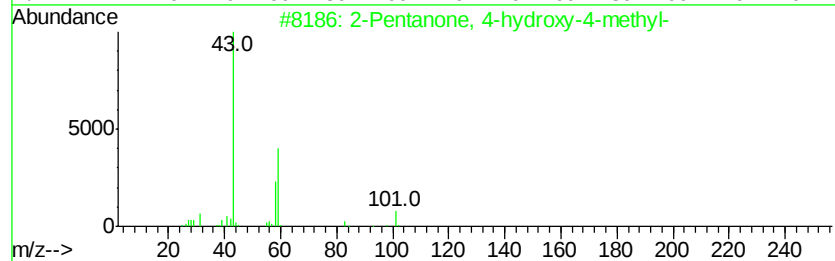
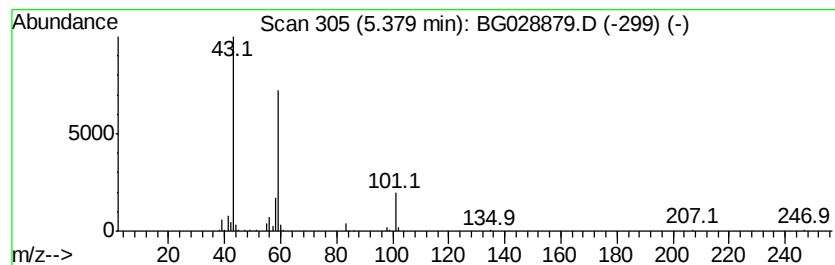
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.38	15.10 ng	103504	1,4-Dichlorobenzene-d4	8.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Guanidine	59	CH5N3	000113-00-8	43
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	37
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25



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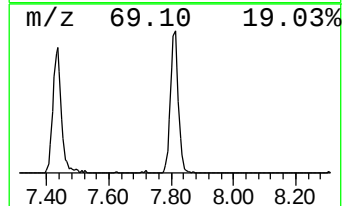
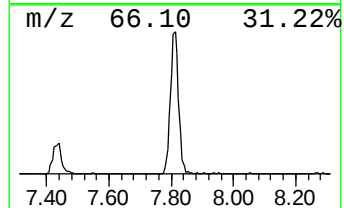
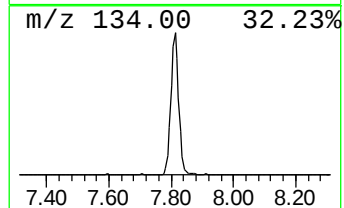
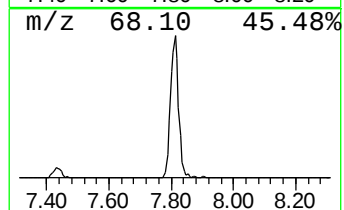
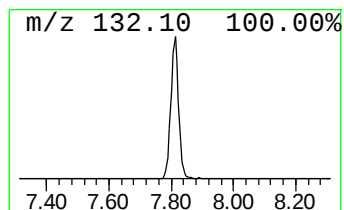
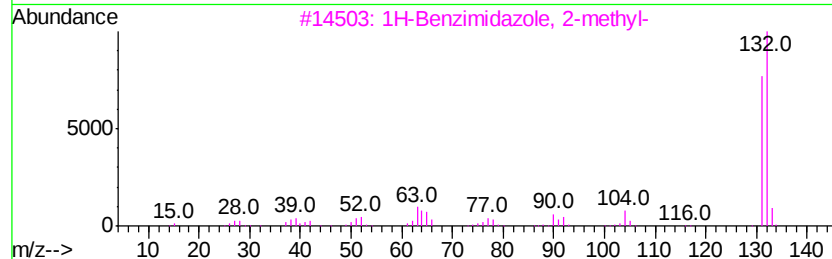
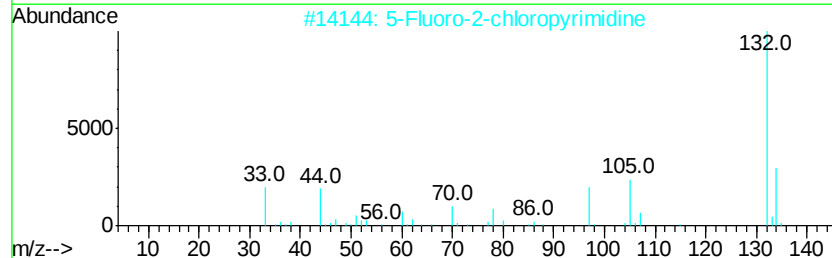
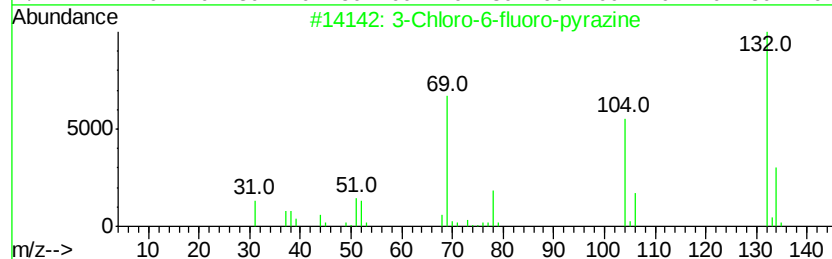
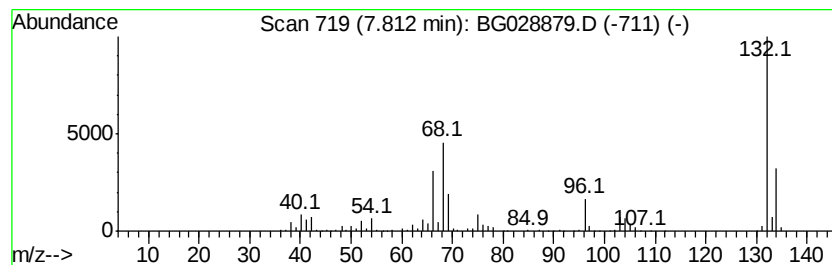
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Peak Number 2 unknown7.81 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	103.69 ng	710735	1,4-Dichlorobenzene-d4	8.28

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		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		5-Aminoindole	132	C8H8N2	005192-03-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.38	15.1	ng	103504	1	8.28	137090	20.0
unknown7.81	7.81	103.7	ng	710735	1	8.28	137090	20.0