

Data Path : Z:\HPCHEM1\BNA G\DATA\BG063017\
 Data File : BG027760.D
 Acq On : 30 Jun 2017 19:08
 Operator : SJ/JU
 Sample : PB100317BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100317BL

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-BG062817.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	5.604	286	292	303	rBV	85494	142427	4.21%	0.944%
2	6.056	362	369	385	rBV	647602	1101749	32.59%	7.304%
3	7.619	627	635	651	rBV	652464	1177895	34.85%	7.809%
4	8.012	692	702	710	rBV	634161	1141609	33.77%	7.568%
5	8.477	774	781	789	rBV	114943	202671	6.00%	1.344%
6	8.806	828	837	848	rVB	447712	824417	24.39%	5.465%
7	9.640	971	979	989	rBV	367265	699335	20.69%	4.636%
8	11.297	1253	1261	1272	rVB	174730	329156	9.74%	2.182%
9	13.694	1661	1669	1679	rBV	1132949	1831577	54.19%	12.142%
10	15.069	1896	1903	1913	rBV3	290502	485253	14.36%	3.217%
11	16.555	2148	2156	2173	rBV	1008796	1603858	47.45%	10.633%
12	17.819	2364	2371	2382	rBV	431511	678816	20.08%	4.500%
13	20.398	2803	2810	2816	rBV	2554648	3380189	100.00%	22.409%
14	22.178	3105	3113	3124	rBV	402961	756932	22.39%	5.018%
15	25.798	3715	3729	3744	rBV	228574	728275	21.55%	4.828%

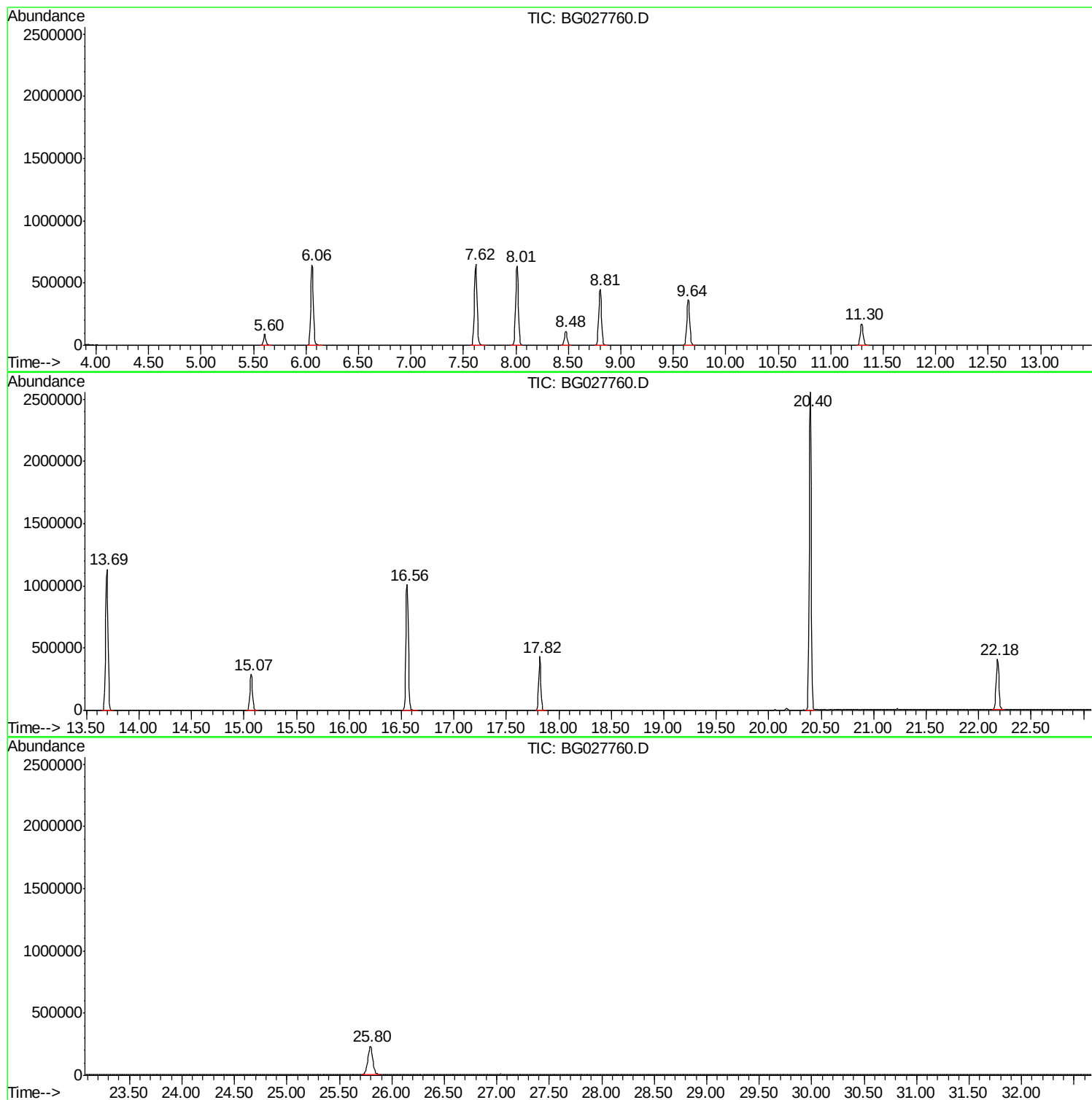
Sum of corrected areas: 15084159

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062817.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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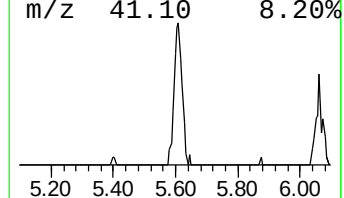
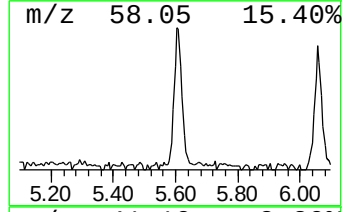
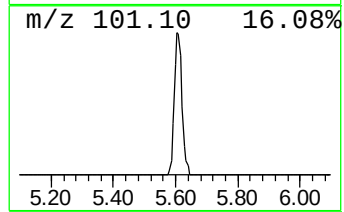
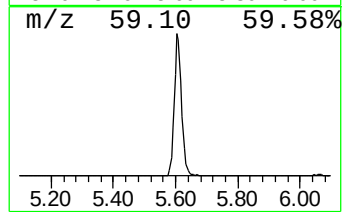
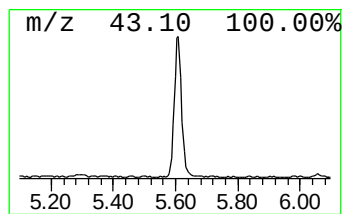
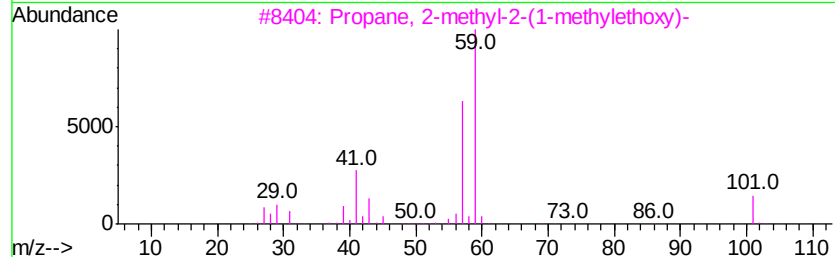
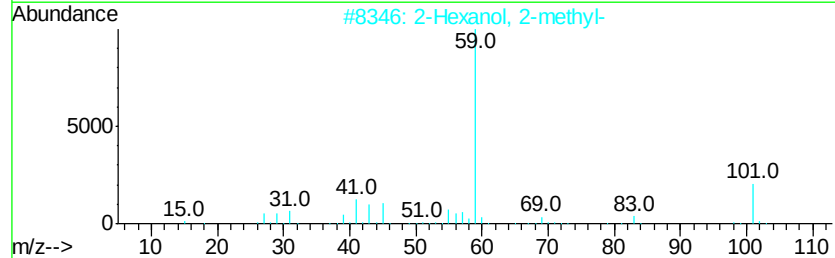
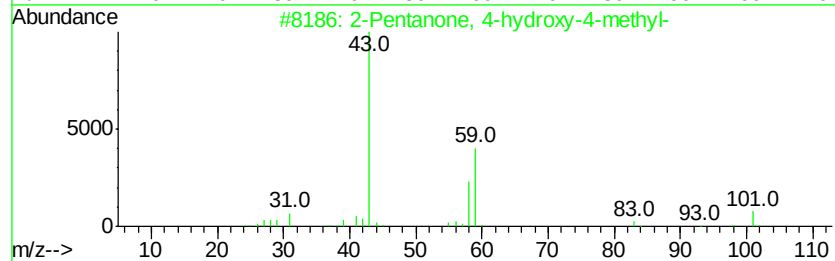
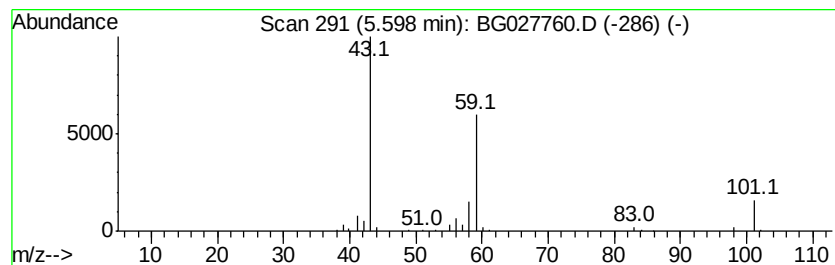
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.60	14.06 ng	142427	1,4-Dichlorobenzene-d4	8.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			Guanidine	59	CH5N3	000113-00-8	9



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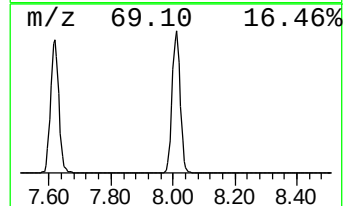
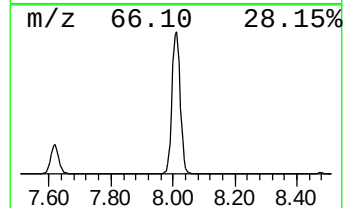
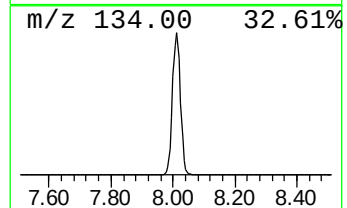
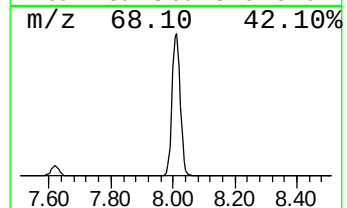
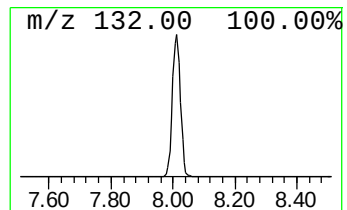
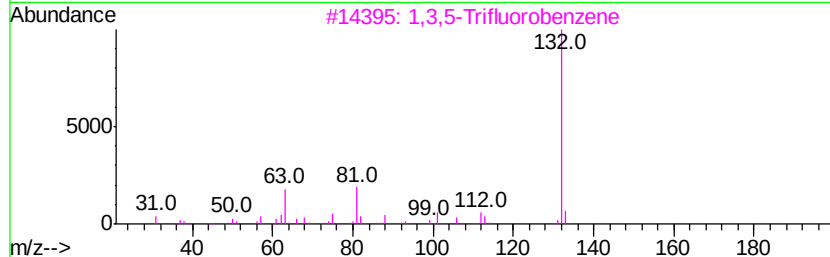
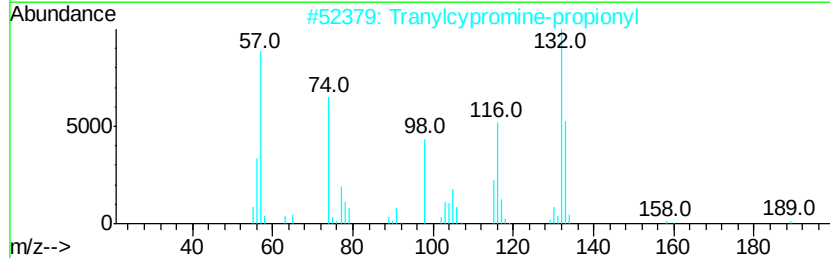
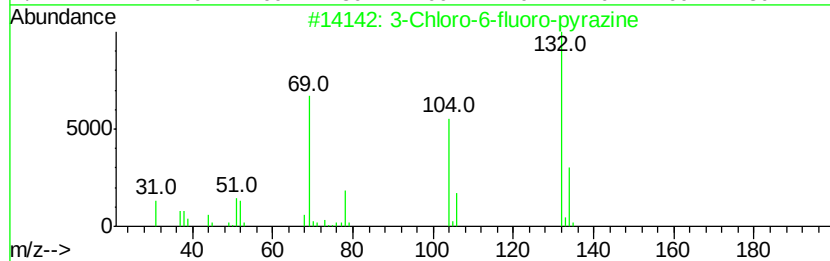
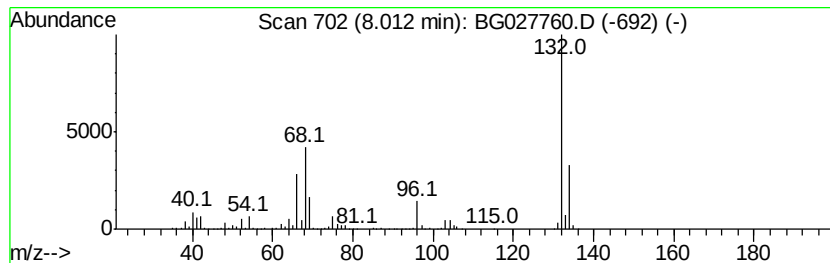
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Peak Number 2 unknown8.01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.01	112.66 ng	1141610	1,4-Dichlorobenzene-d4	8.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.60	14.1	ng	142427	1	8.48	202671	20.0
unknown8.01	8.01	112.7	ng	1141610	1	8.48	202671	20.0