

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062117\
 Data File : BG027583.D
 Acq On : 22 Jun 2017 6:26
 Operator : SJ/JU
 Sample : PB99997BL
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99997BL

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-BG062117.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	5.639	292	298	309	rVB	63043	102857	4.63%	1.114%
2	6.086	367	374	387	rBV	398359	662058	29.83%	7.169%
3	7.637	630	638	653	rBV	378564	690111	31.09%	7.473%
4	8.037	698	706	717	rVB	369753	674364	30.38%	7.302%
5	8.501	778	785	794	rVB	77047	137268	6.18%	1.486%
6	8.830	833	841	850	rBV	257747	473571	21.33%	5.128%
7	9.664	975	983	994	rBV	211911	399045	17.98%	4.321%
8	11.321	1256	1265	1272	rBV	103244	198000	8.92%	2.144%
9	13.712	1665	1672	1679	rBV	607742	958555	43.18%	10.380%
10	15.087	1899	1906	1914	rBV2	166299	275735	12.42%	2.986%
11	16.574	2150	2159	2172	rBV	595654	951578	42.87%	10.304%
12	17.837	2367	2374	2382	rBV2	277732	437475	19.71%	4.737%
13	20.416	2806	2813	2824	rVB	1691672	2219740	100.00%	24.036%
14	22.202	3110	3117	3127	rVB	296424	551037	24.82%	5.967%
15	25.839	3725	3736	3748	rBV2	158940	503517	22.68%	5.452%

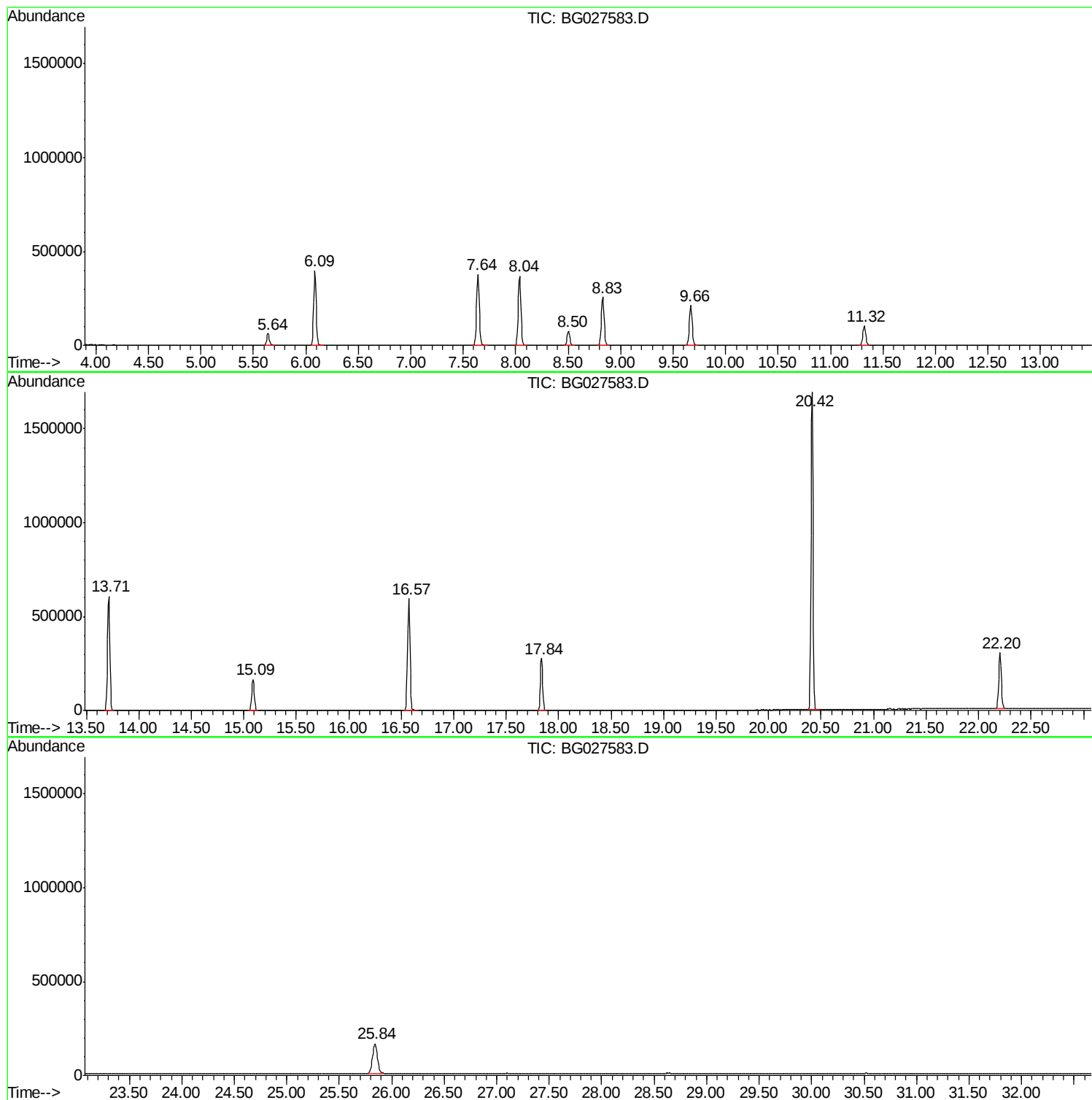
Sum of corrected areas: 9234911

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062117.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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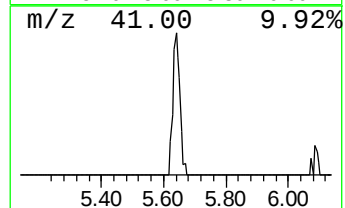
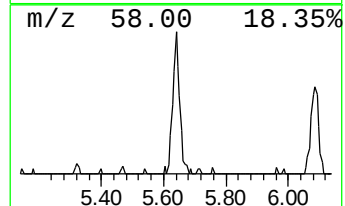
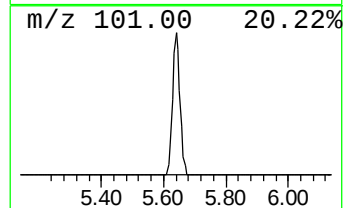
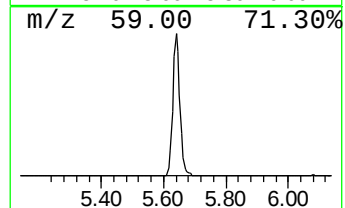
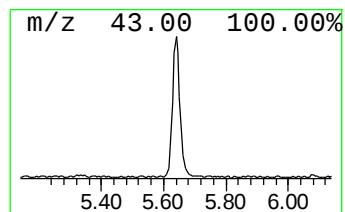
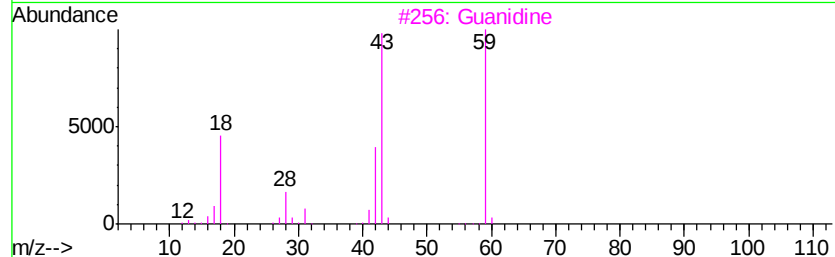
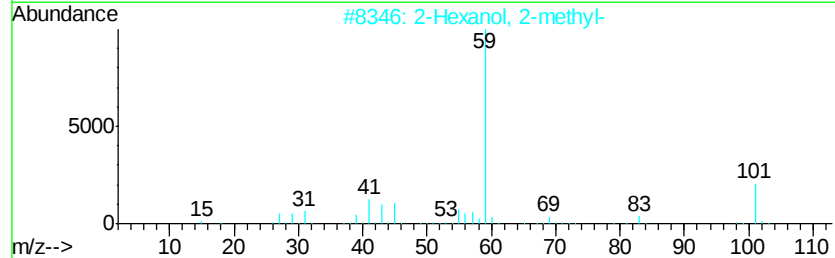
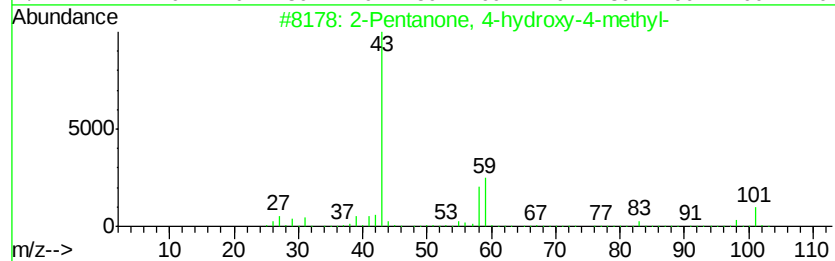
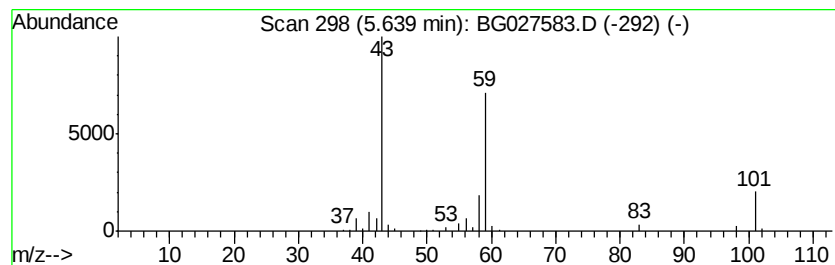
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	14.99 ng	102857	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	53
3		Guanidine	59	CH5N3	000113-00-8	43
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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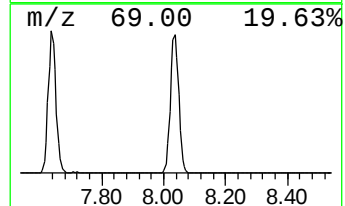
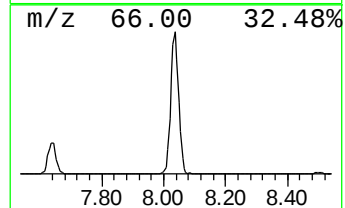
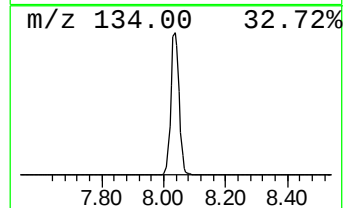
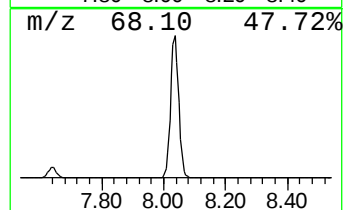
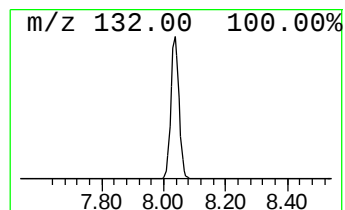
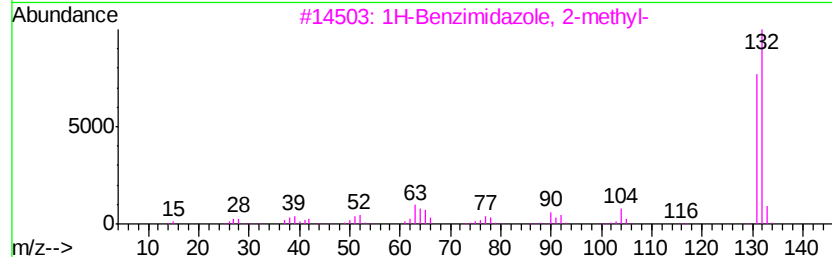
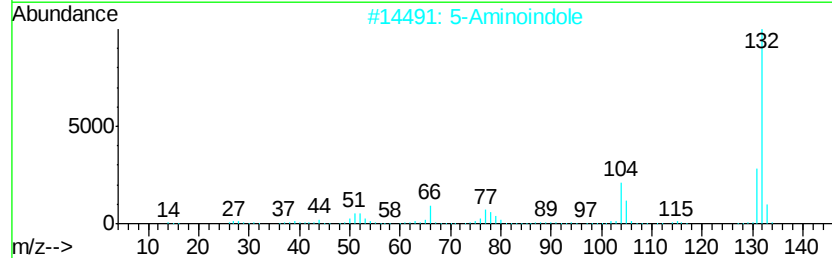
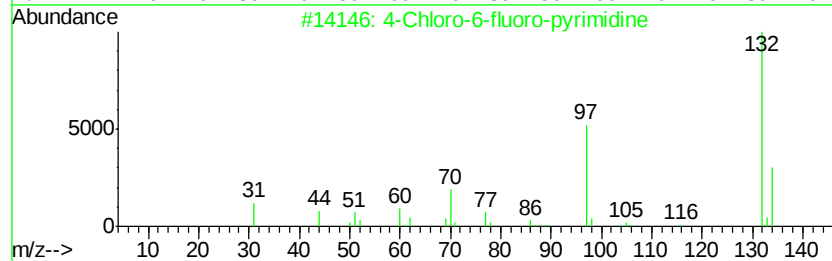
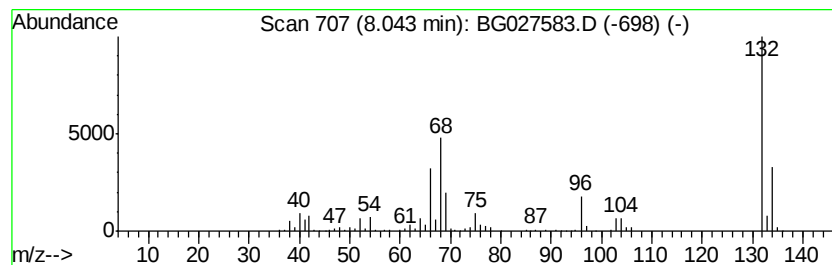
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Peak Number 2 unknown8.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	98.26 ng	674364	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
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...	5.64	15.0	ng	102857	1	8.50	137268	20.0
unknown8.04	8.04	98.3	ng	674364	1	8.50	137268	20.0