

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062617\
 Data File : BG027680.D
 Acq On : 26 Jun 2017 18:31
 Operator : SJ/JU
 Sample : I3876-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CAN-FB062217

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	5.627	287	296	303	rBV	15037	23167	1.06%	0.327%
2	6.074	365	372	382	rVB	130762	226966	10.36%	3.208%
3	7.625	630	636	646	rBB	76579	140379	6.41%	1.984%
4	8.018	696	703	714	rBV	243274	448724	20.48%	6.342%
5	8.488	775	783	790	rBV2	47458	83614	3.82%	1.182%
6	8.817	831	839	849	rBB	215527	403933	18.43%	5.709%
7	9.652	973	981	992	rBV	187963	344784	15.73%	4.873%
8	11.303	1253	1262	1270	rBV	63598	120244	5.49%	1.699%
9	13.700	1662	1670	1677	rBV	551377	884448	40.36%	12.501%
10	15.075	1898	1904	1912	rVB2	110176	185835	8.48%	2.627%
11	16.561	2150	2157	2166	rBV	604632	956483	43.64%	13.519%
12	17.825	2365	2372	2382	rBV	205731	316921	14.46%	4.479%
13	20.404	2804	2811	2820	rBV	1609177	2191553	100.00%	30.975%
14	22.184	3106	3114	3125	rVB	208735	390946	17.84%	5.526%
15	25.809	3718	3731	3744	rBV2	112589	357263	16.30%	5.049%

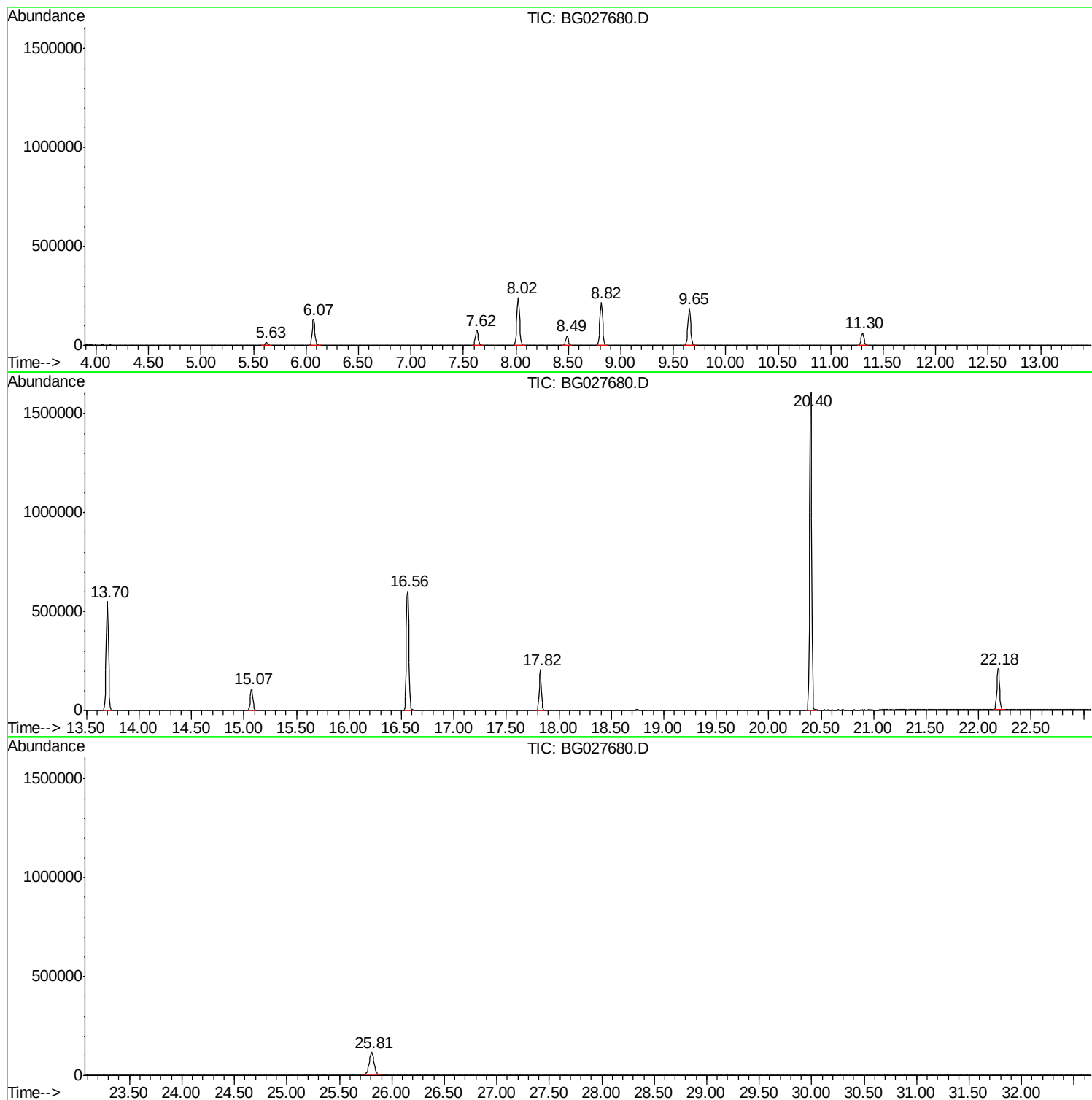
Sum of corrected areas: 7075260

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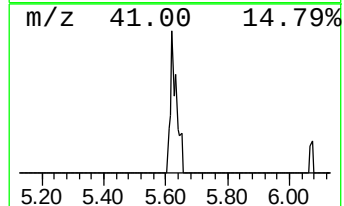
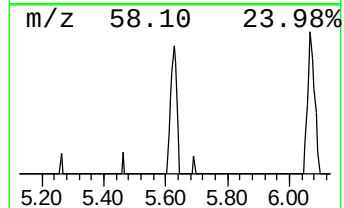
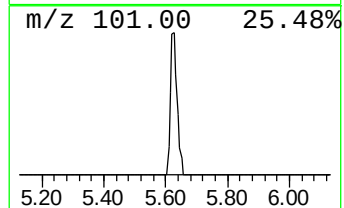
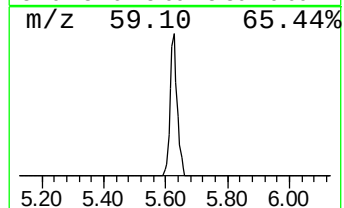
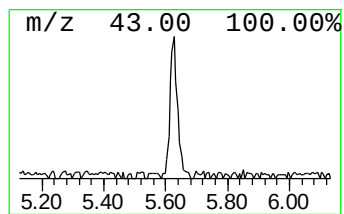
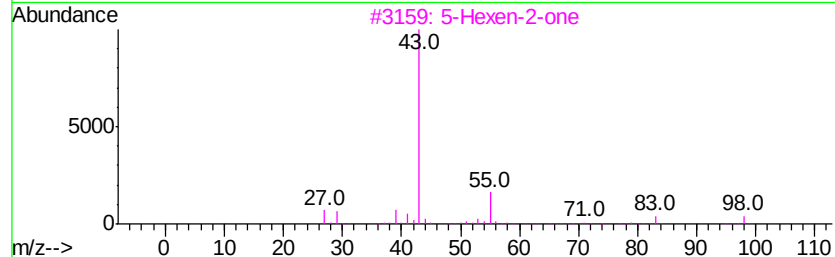
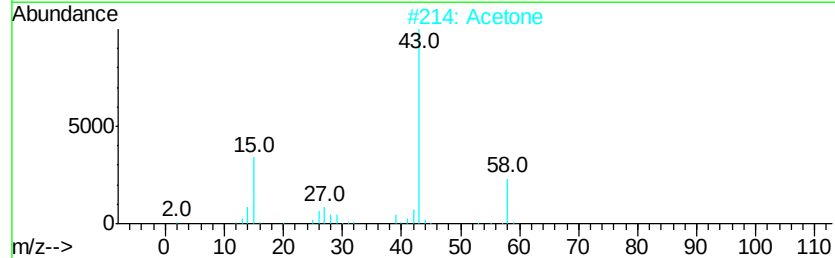
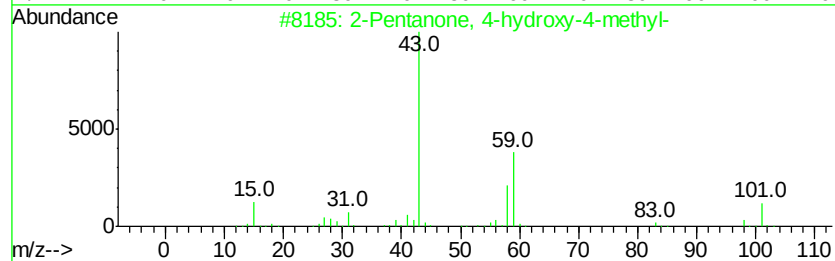
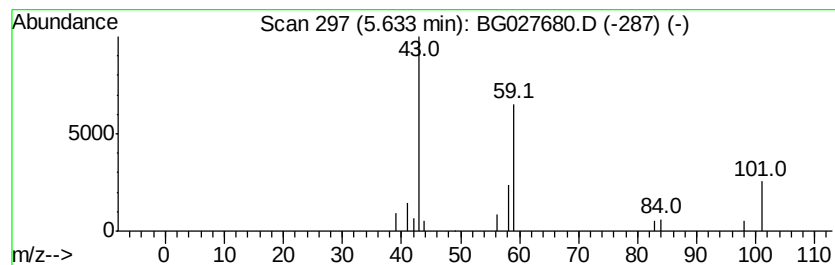
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	5.54 ng	23167	1,4-Dichlorobenzene-d4	8.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2	Acetone	58	C3H6O	000067-64-1	9
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9
4	Guanidine	59	CH5N3	000113-00-8	9
5	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9



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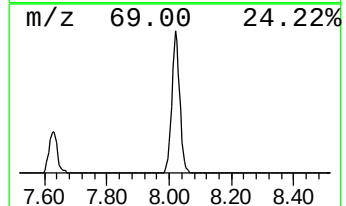
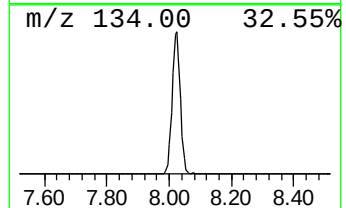
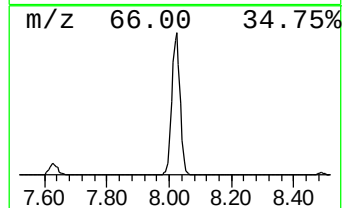
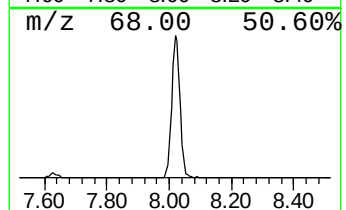
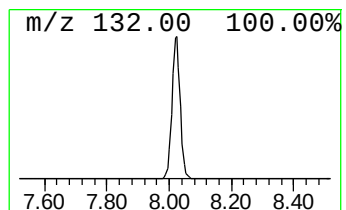
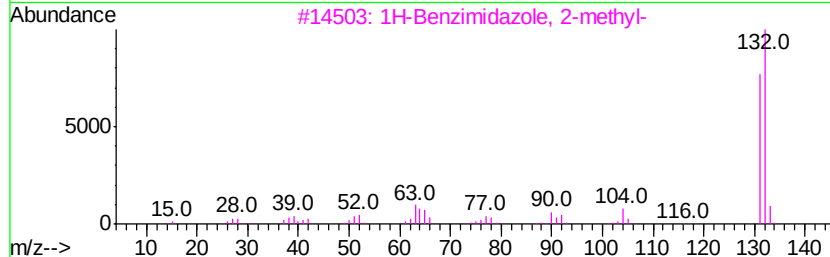
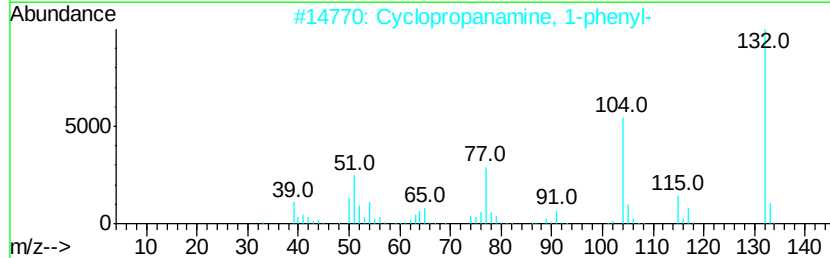
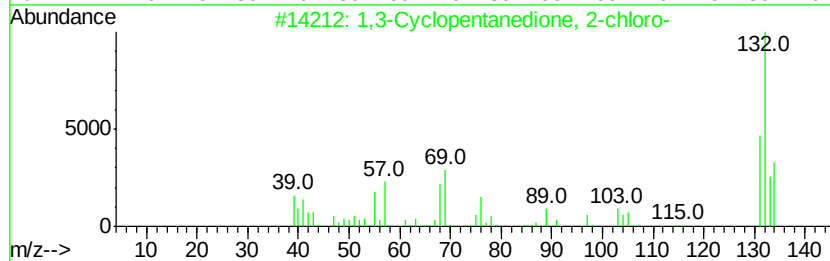
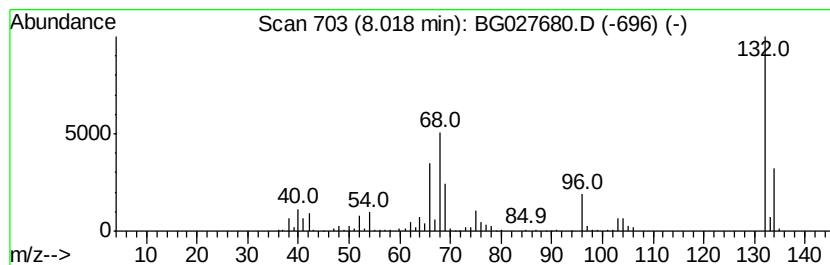
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Peak Number 2 unknown8.02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	107.33 ng	448724	1,4-Dichlorobenzene-d4	8.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	14
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.63	5.5	ng	23167	1	8.49	83614	20.0
unknown8.02	8.02	107.3	ng	448724	1	8.49	83614	20.0