

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080218\
 Data File : BG036054.D
 Acq On : 2 Aug 2018 10:51
 Operator : JU/SJ
 Sample : PB111656BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111656BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.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	5.133	308	314	321	rBV	84442	133302	4.04%	0.977%
2	5.603	388	394	413	rBV	581696	985595	29.89%	7.224%
3	7.189	657	664	676	rBV	568054	1019473	30.91%	7.472%
4	7.559	719	727	735	rBV	571423	1020982	30.96%	7.483%
5	8.023	798	806	815	rVB	119108	208849	6.33%	1.531%
6	8.346	851	861	871	rBV	452283	804807	24.41%	5.899%
7	9.186	996	1004	1019	rBV	312891	627504	19.03%	4.599%
8	10.825	1277	1283	1292	rBV2	126172	251178	7.62%	1.841%
9	13.269	1692	1699	1719	rBV	948897	1566654	47.51%	11.482%
10	14.649	1927	1934	1949	rBV2	220980	399129	12.10%	2.925%
11	16.135	2180	2187	2214	rBV	701852	1307345	39.64%	9.582%
12	17.392	2393	2401	2417	rBV2	322385	597740	18.13%	4.381%
13	19.995	2838	2844	2859	rBV	2319475	3297665	100.00%	24.169%
14	21.651	3120	3126	3141	rBV2	398575	749626	22.73%	5.494%
15	24.853	3660	3671	3688	rBV2	219666	674157	20.44%	4.941%

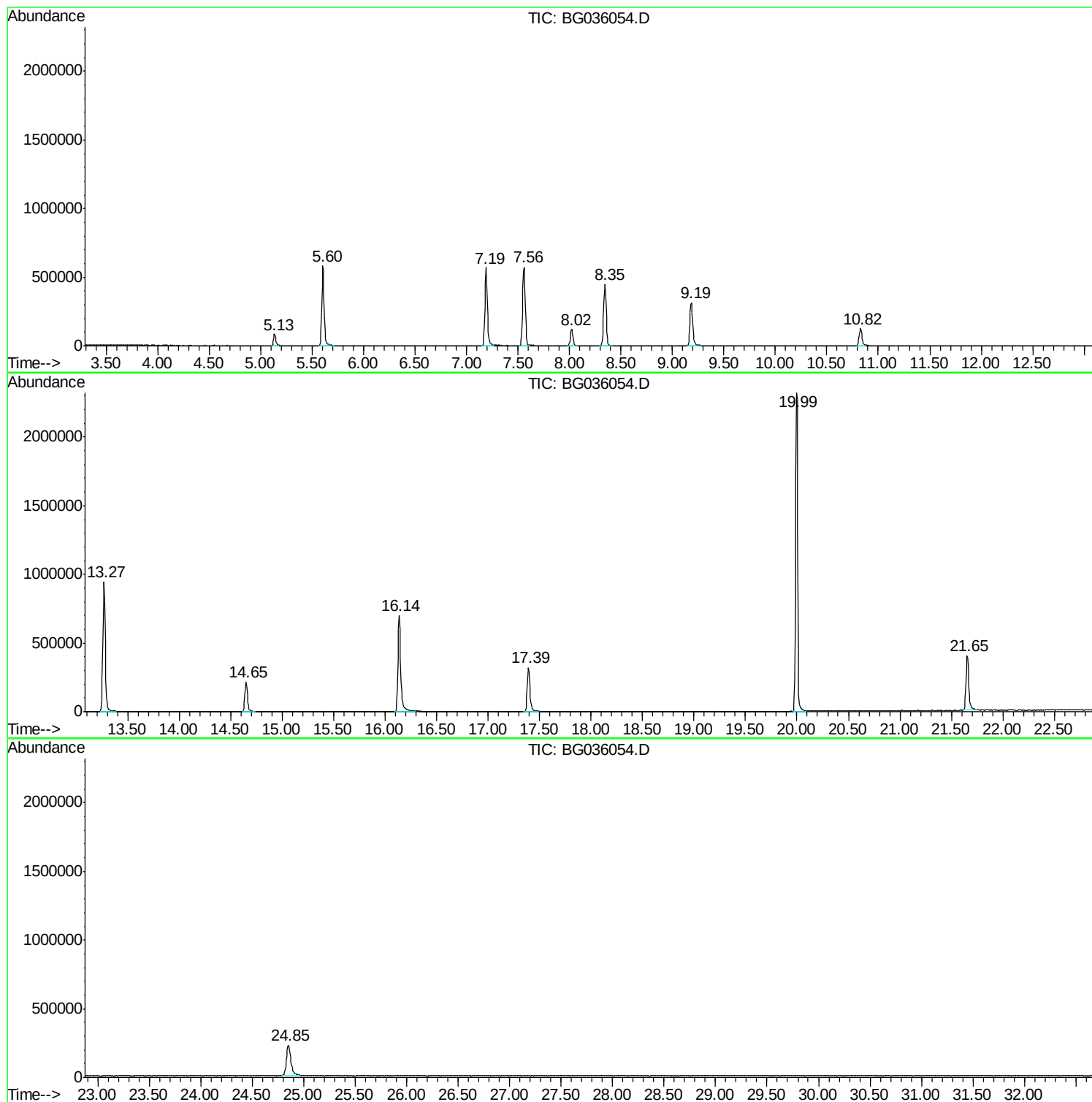
Sum of corrected areas: 13644006

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



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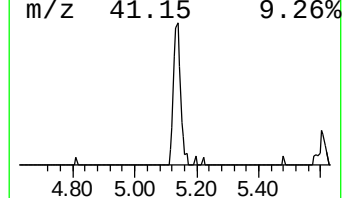
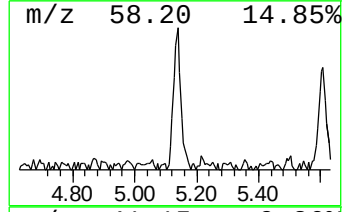
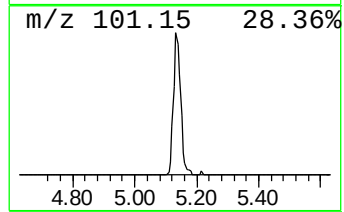
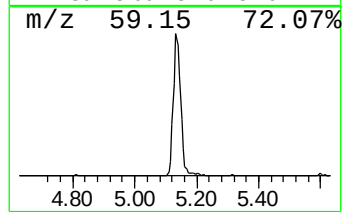
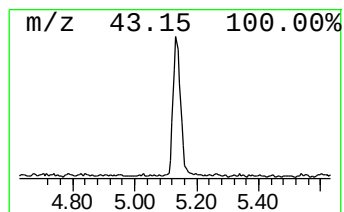
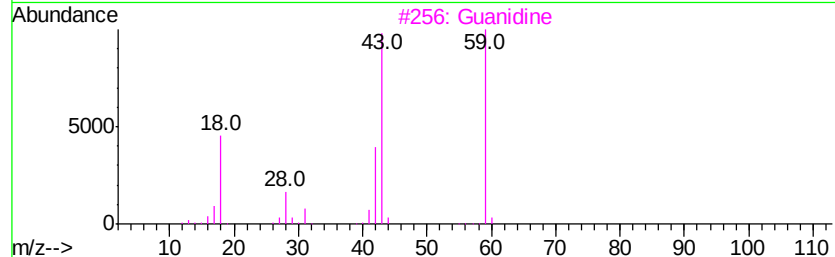
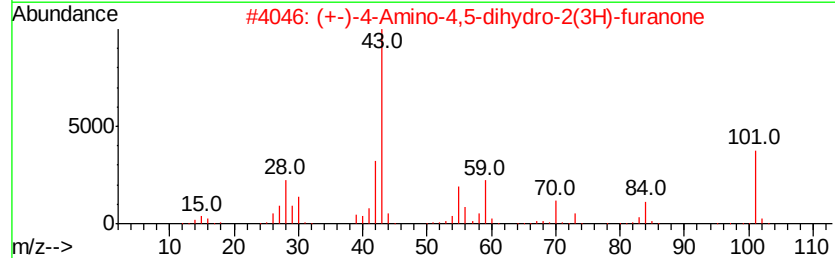
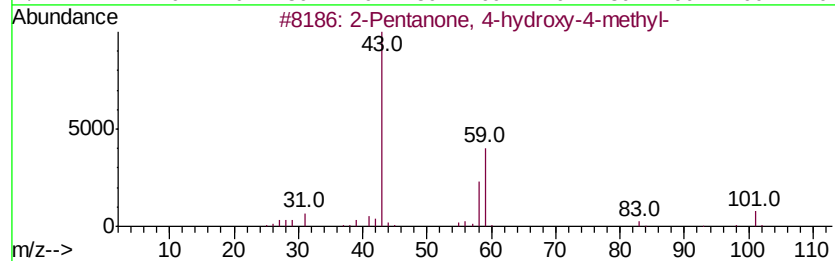
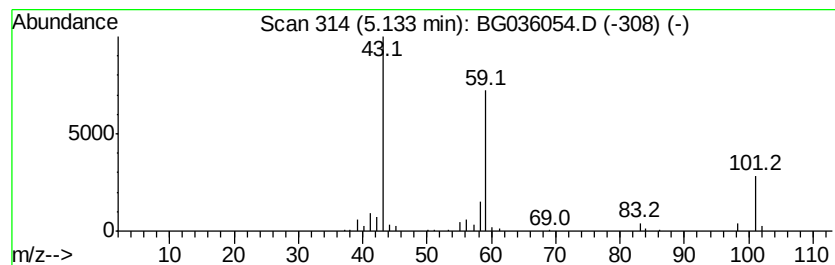
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	12.77 ng	133302	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	47
3		Guanidine	59	CH5N3	000113-00-8	43
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33



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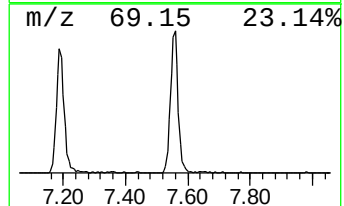
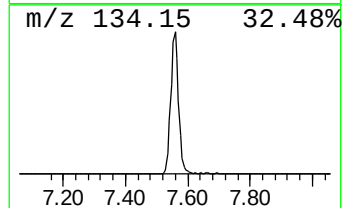
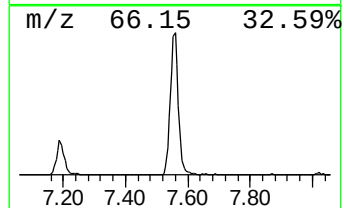
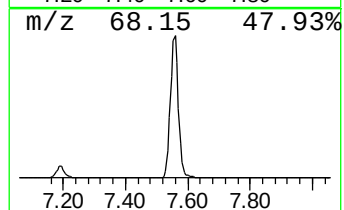
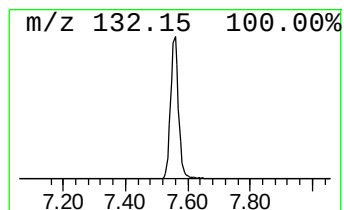
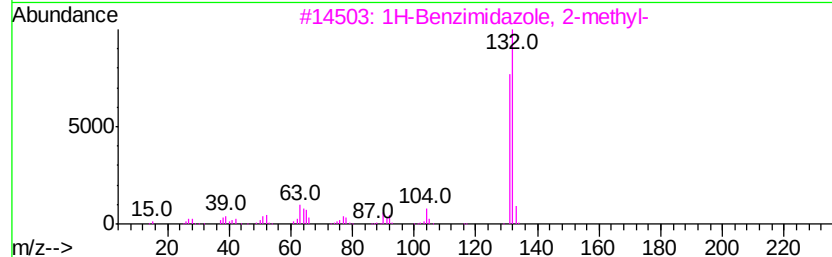
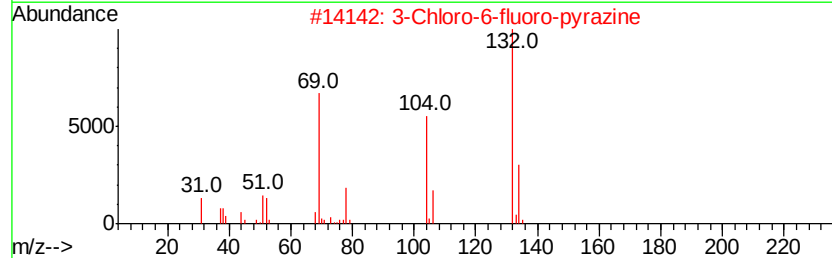
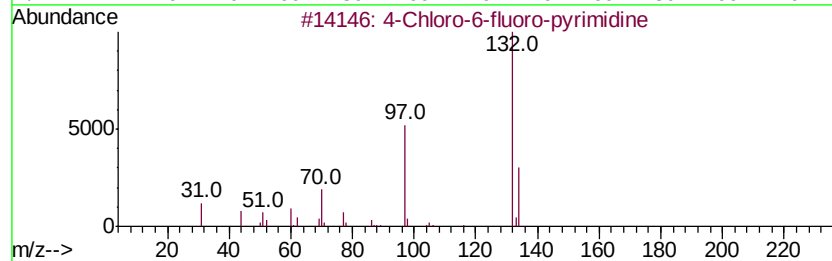
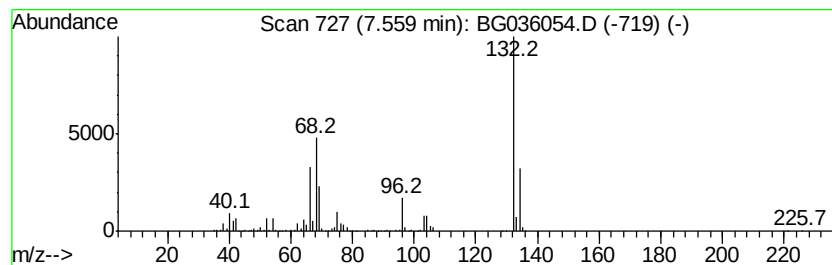
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	97.77 ng	1020980	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	12.8	ng	133302	1	8.02	208849	20.0
unknown7.56	7.56	97.8	ng	1020980	1	8.02	208849	20.0