

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080218\
 Data File : BG036055.D
 Acq On : 2 Aug 2018 11:30
 Operator : JU/SJ
 Sample : PB111684BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111684BL

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	5.135	309	314	324	rBV	51507	79189	2.61%	0.668%
2	5.605	388	394	409	rBV	405722	688750	22.66%	5.806%
3	7.191	655	664	676	rBV	382476	726881	23.92%	6.127%
4	7.555	719	726	736	rBV	422216	719034	23.66%	6.061%
5	8.019	799	805	813	rBV	109337	191662	6.31%	1.616%
6	8.342	852	860	869	rBV	504107	874050	28.76%	7.368%
7	9.182	994	1003	1020	rBV	354309	690462	22.72%	5.820%
8	10.827	1276	1283	1297	rBV	107883	225458	7.42%	1.901%
9	13.265	1691	1698	1711	rBV	1014938	1705099	56.10%	14.373%
10	14.646	1926	1933	1947	rBV2	189630	356887	11.74%	3.008%
11	16.132	2178	2186	2198	rBV	435322	841304	27.68%	7.092%
12	17.389	2393	2400	2411	rBV	277301	514783	16.94%	4.339%
13	19.997	2837	2844	2859	rBV	2241455	3039297	100.00%	25.620%
14	21.654	3119	3126	3139	rBV2	337334	634457	20.88%	5.348%
15	24.849	3659	3670	3687	rBV2	174541	575748	18.94%	4.853%

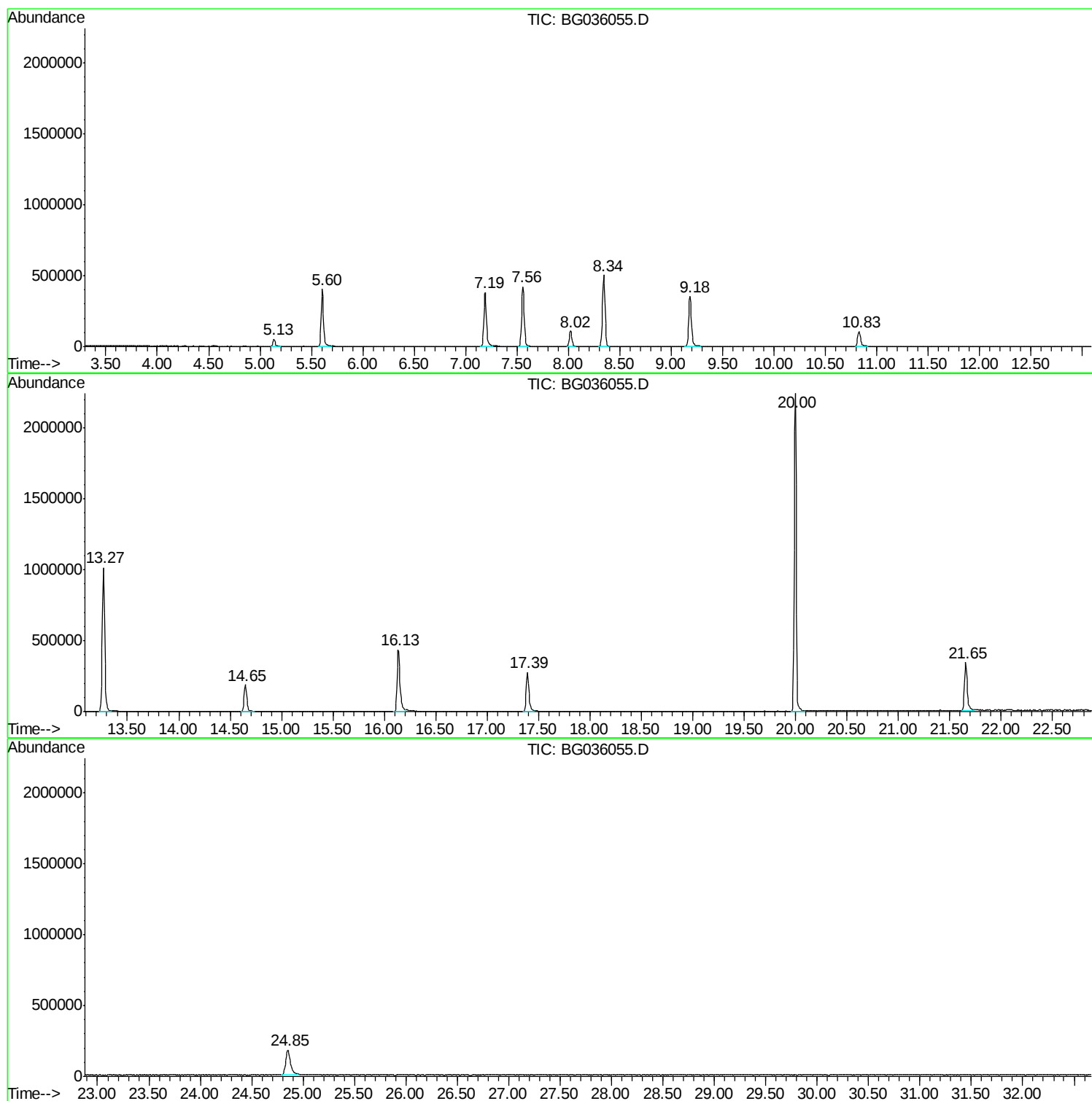
Sum of corrected areas: 11863061

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



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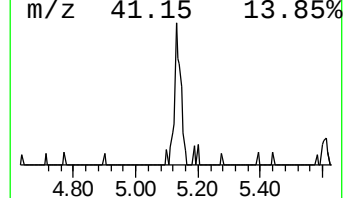
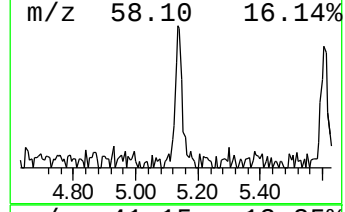
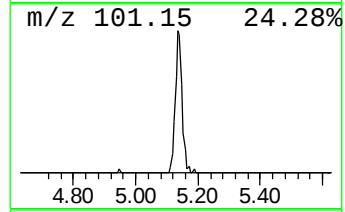
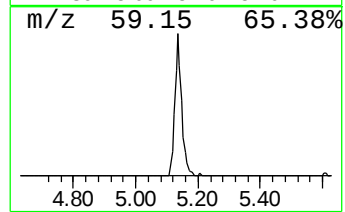
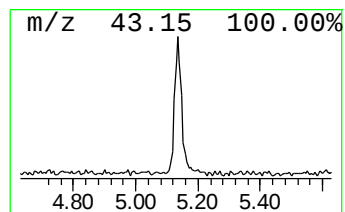
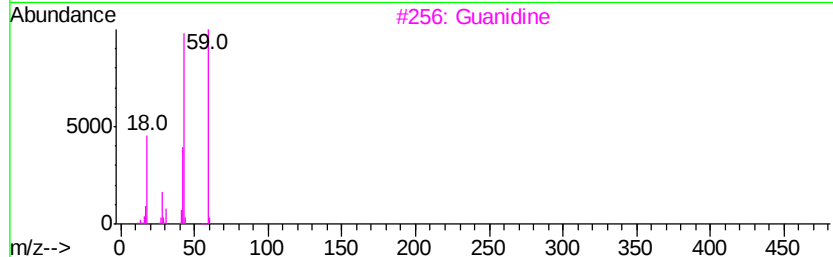
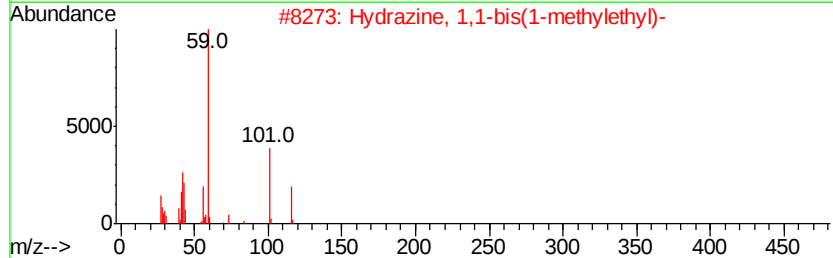
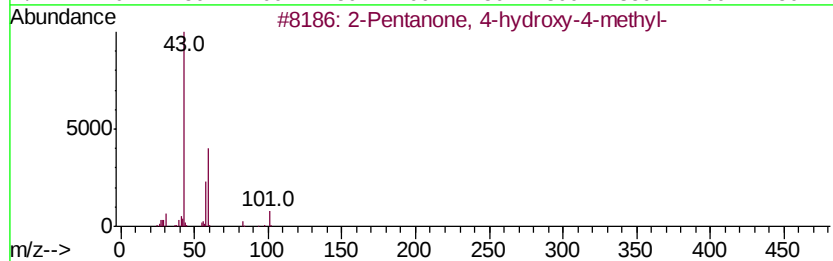
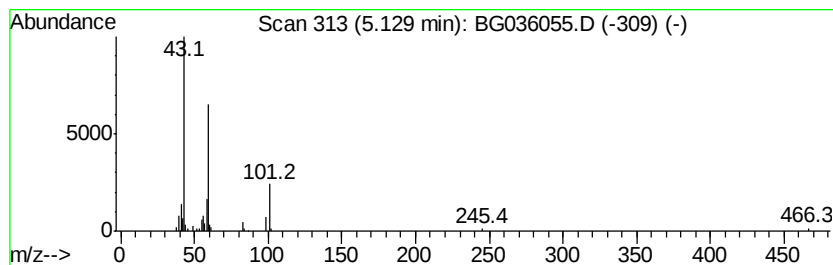
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	8.26 ng	79189	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	43
3			Guanidine	59	CH5N3	000113-00-8	43
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			3-Hexanol	102	C6H14O	000623-37-0	12



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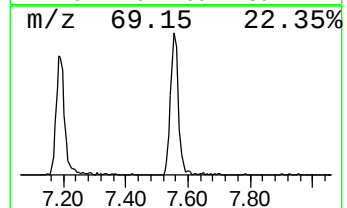
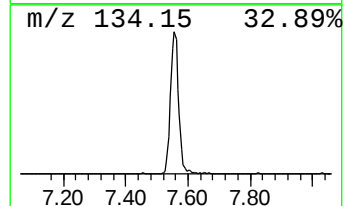
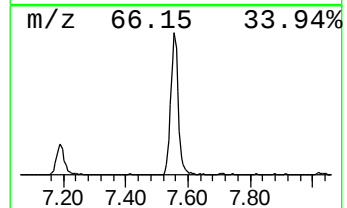
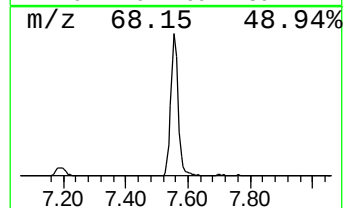
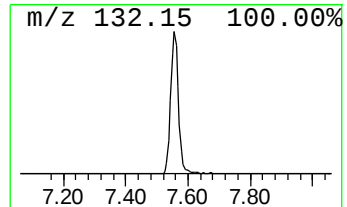
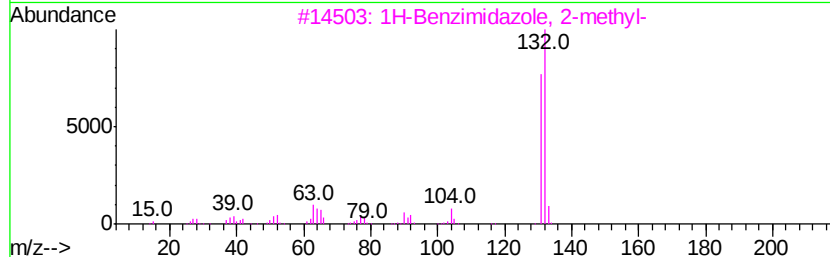
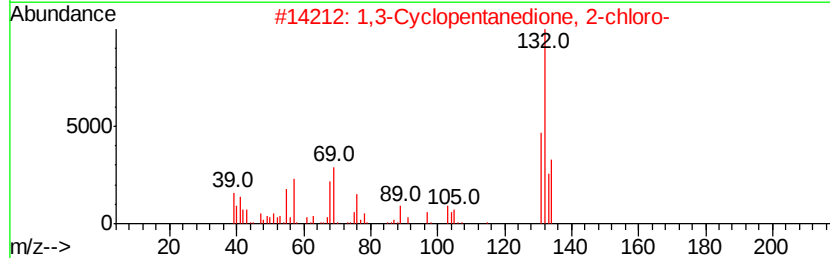
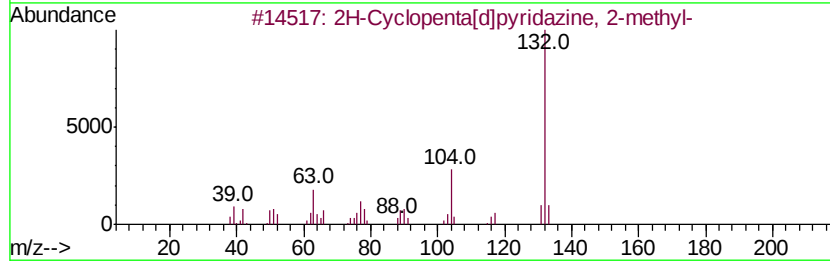
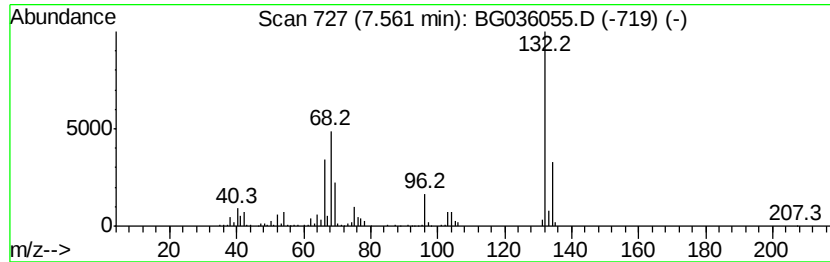
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Peak Number 2 unknown7.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	75.03 ng	719034	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		1H-Indazole, 3-methyl-	132	C8H8N2	1000316-00-2	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	8.3	ng	79189	1	8.02	191662	20.0
unknown7.56	7.56	75.0	ng	719034	1	8.02	191662	20.0