

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081018\
 Data File : BG036272.D
 Acq On : 10 Aug 2018 18:04
 Operator : JU/SJ
 Sample : PB111936BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111936BL

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	5.126	307	313	323	rBV	80191	128176	4.41%	1.053%
2	5.596	386	393	411	rBV	521836	923734	31.81%	7.591%
3	7.182	656	663	677	rBV	538547	987512	34.00%	8.115%
4	7.546	713	725	743	rBV	557185	983646	33.87%	8.083%
5	8.004	796	803	811	rBV2	93620	162705	5.60%	1.337%
6	8.327	850	858	870	rBV	421044	760930	26.20%	6.253%
7	9.167	993	1001	1013	rBV	290825	602915	20.76%	4.954%
8	10.812	1273	1281	1290	rBV	88659	187409	6.45%	1.540%
9	13.256	1689	1697	1710	rBV	885643	1523636	52.46%	12.520%
10	14.636	1923	1932	1947	rBV2	160269	313647	10.80%	2.577%
11	16.123	2178	2185	2210	rBV	701385	1302739	44.86%	10.705%
12	17.380	2393	2399	2413	rBV	240018	441147	15.19%	3.625%
13	19.982	2836	2842	2861	rBV	2191558	2904331	100.00%	23.866%
14	21.639	3118	3124	3135	rBV2	242712	505653	17.41%	4.155%
15	24.828	3657	3667	3681	rBV2	125321	441203	15.19%	3.626%

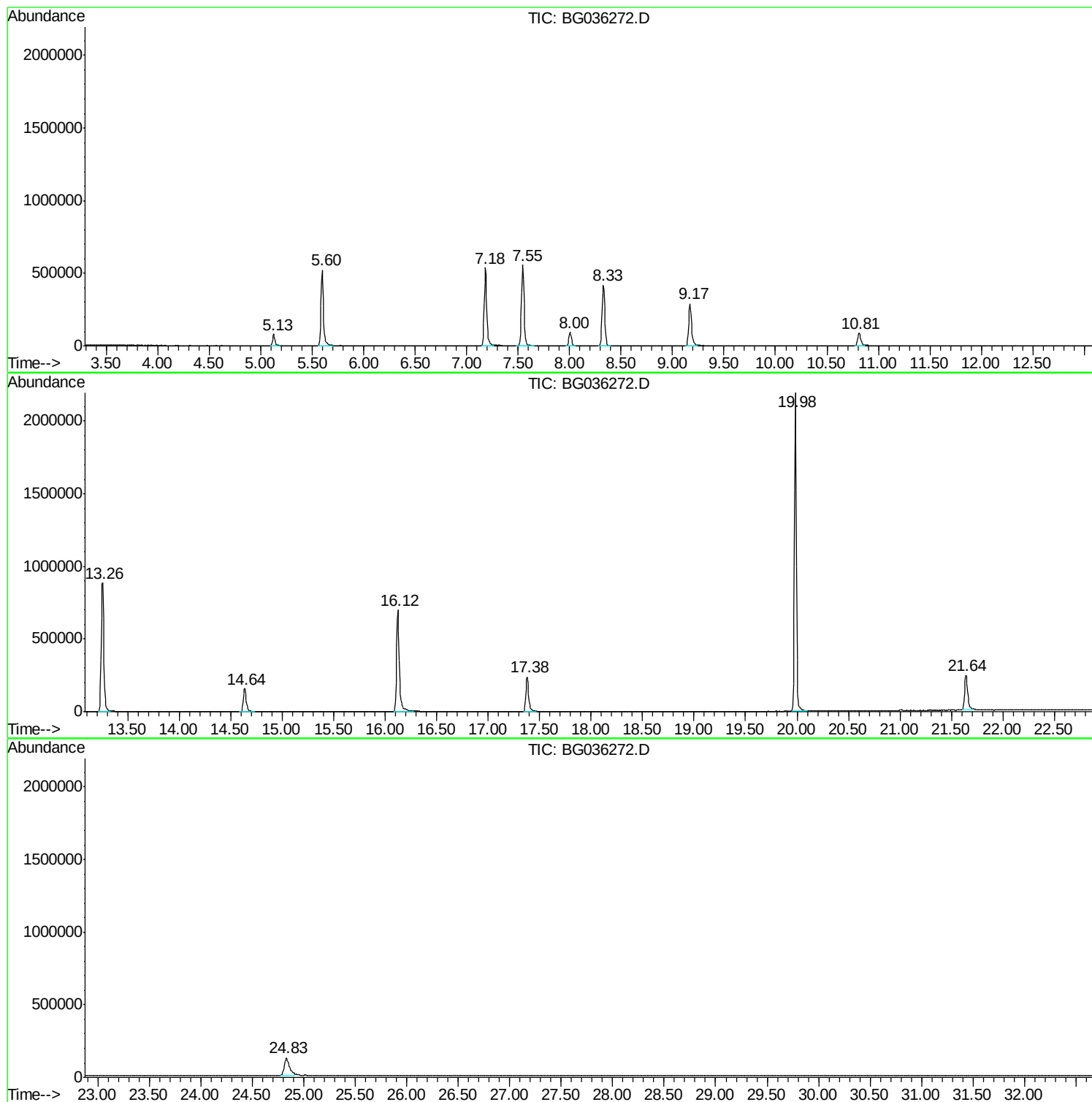
Sum of corrected areas: 12169383

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



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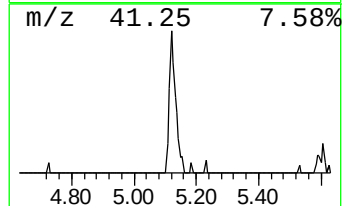
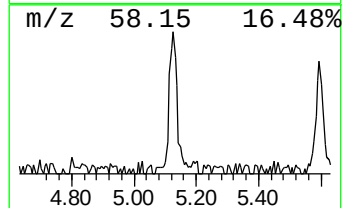
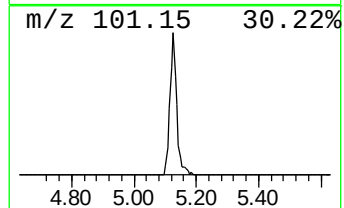
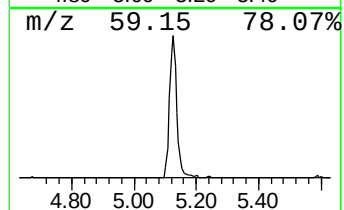
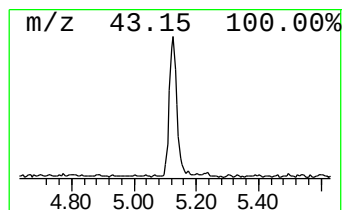
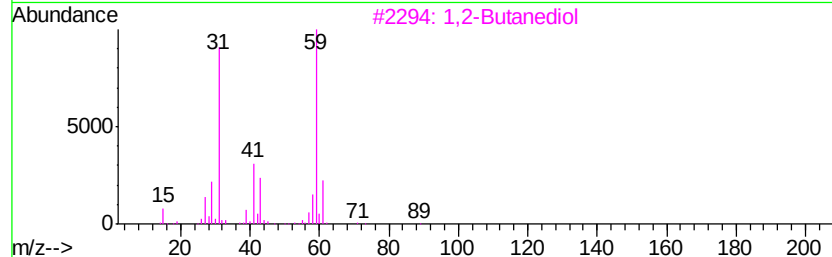
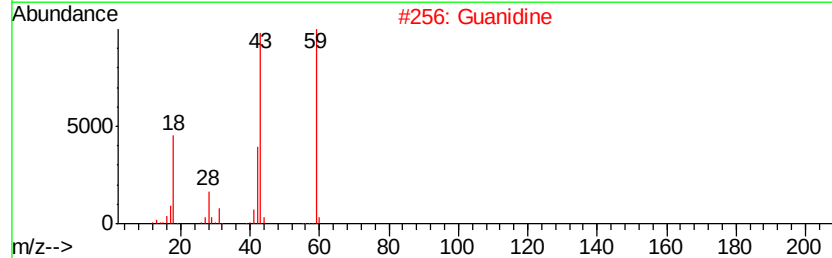
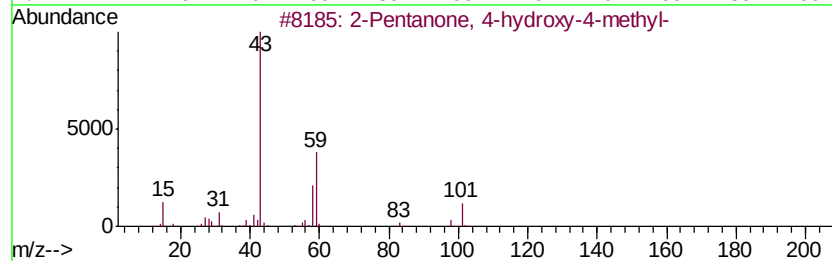
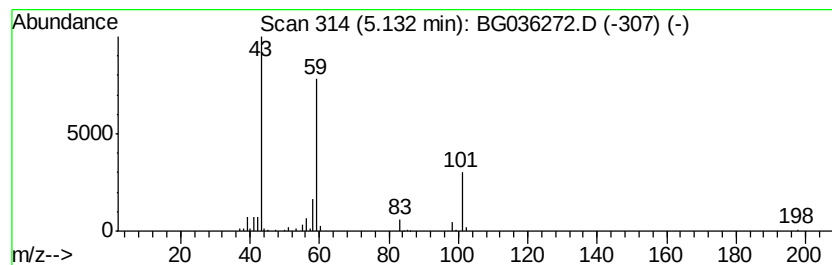
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	15.76 ng	128176	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Guanidine	59	CH5N3	000113-00-8	50
3		1,2-Butanediol	90	C4H10O2	000584-03-2	38
4		Dimethadione	129	C5H7NO3	000695-53-4	38
5		Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	36



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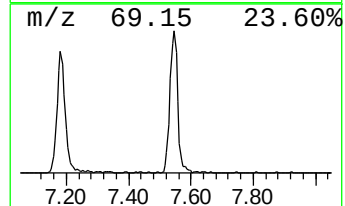
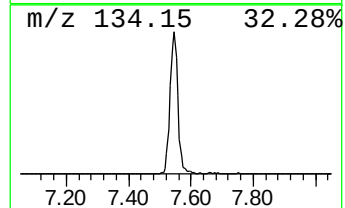
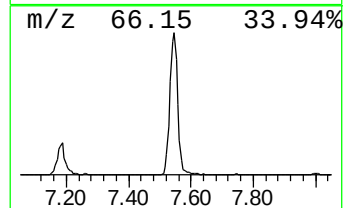
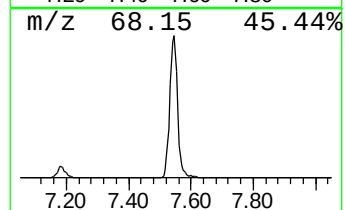
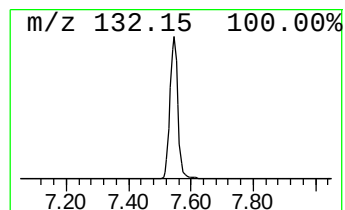
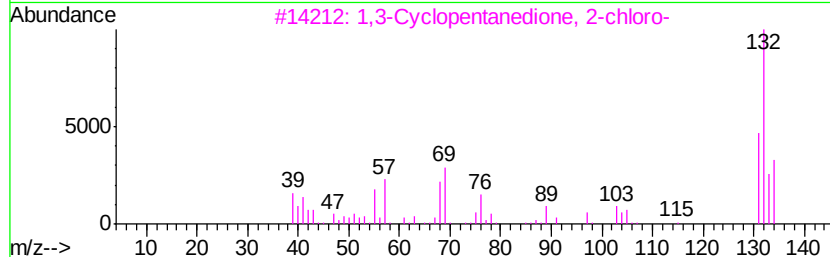
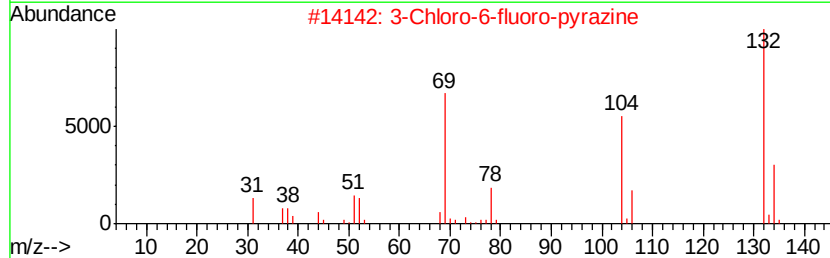
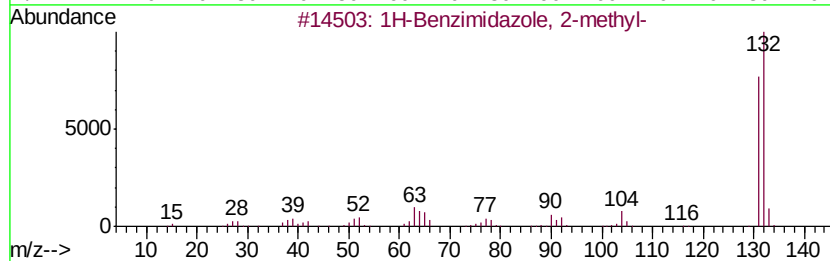
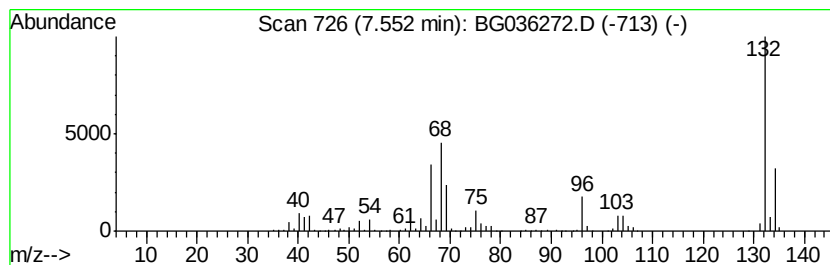
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Peak Number 2 unknown7.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	120.91 ng	983646	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	15.8	ng	128176	1	8.00	162705	20.0
unknown7.55	7.55	120.9	ng	983646	1	8.00	162705	20.0