

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG072618\
 Data File : BG035915.D
 Acq On : 26 Jul 2018 19:12
 Operator : JU/SJ
 Sample : PB111449BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111449BL

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	325	rBV	81723	129738	3.31%	0.819%
2	5.609	388	395	413	rBV	621677	1046995	26.71%	6.613%
3	7.195	658	665	682	rBV	650769	1164853	29.72%	7.357%
4	7.559	720	727	739	rBV	635369	1112775	28.39%	7.028%
5	8.023	800	806	816	rBV	127036	213152	5.44%	1.346%
6	8.346	851	861	873	rBV	484831	855202	21.82%	5.401%
7	9.186	996	1004	1020	rBV	362858	693013	17.68%	4.377%
8	10.831	1274	1284	1298	rBV	132062	278262	7.10%	1.757%
9	13.275	1692	1700	1711	rBV	1102752	1854017	47.30%	11.710%
10	14.649	1928	1934	1948	rBV2	263688	473883	12.09%	2.993%
11	16.135	2181	2187	2201	rBV	971803	1652819	42.16%	10.439%
12	17.393	2393	2401	2414	rBV	416424	732542	18.69%	4.627%
13	20.001	2838	2845	2859	rBV	2934039	3920048	100.00%	24.759%
14	21.657	3120	3127	3143	rBV2	473591	887639	22.64%	5.606%
15	24.853	3660	3671	3686	rBV3	258395	818140	20.87%	5.167%

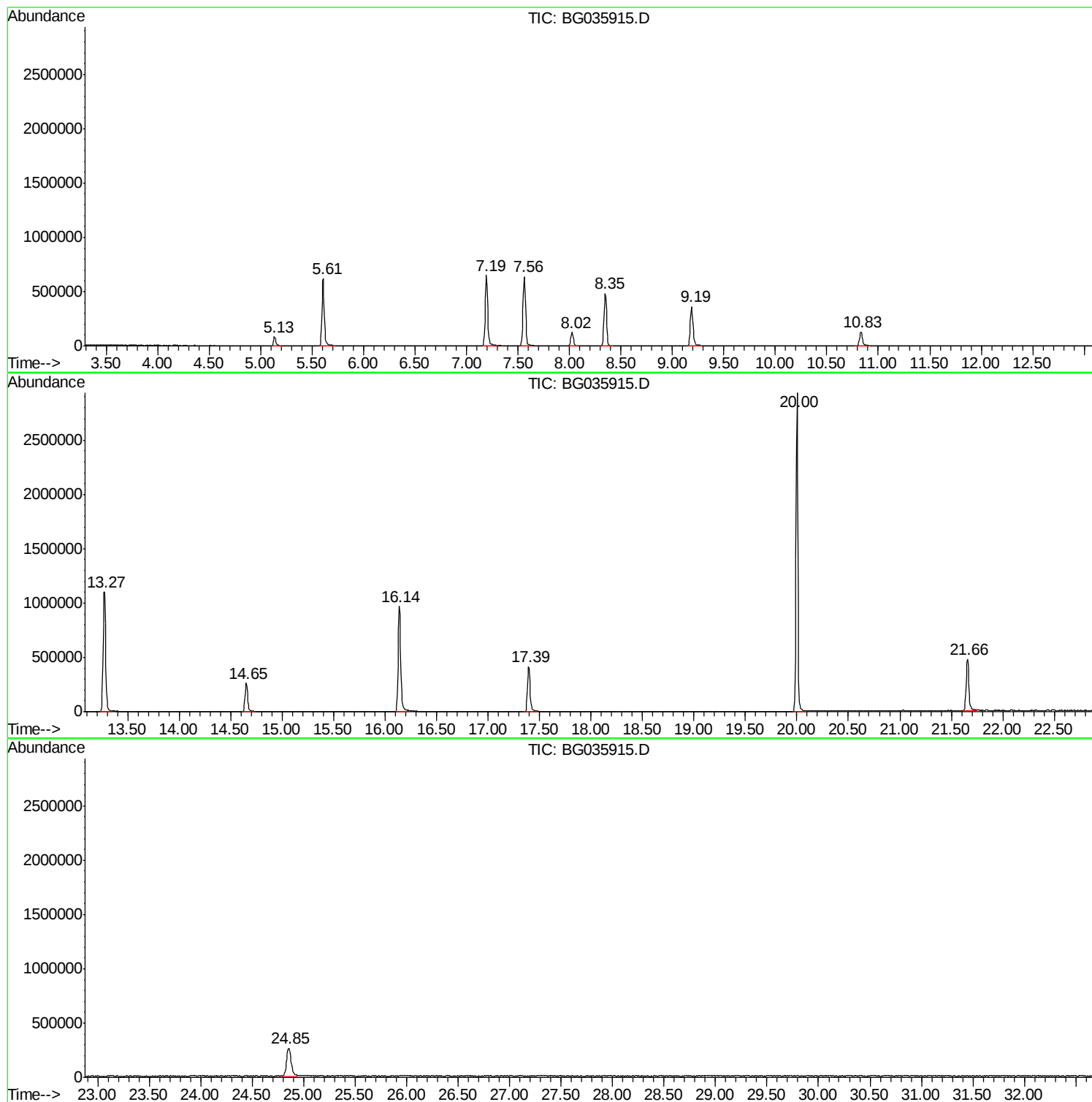
Sum of corrected areas: 15833078

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



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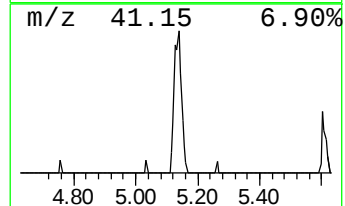
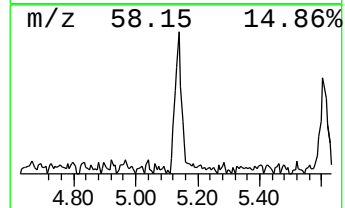
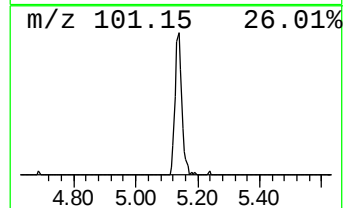
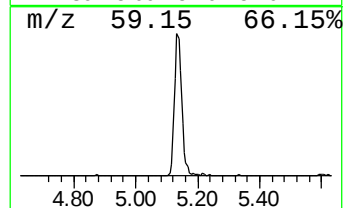
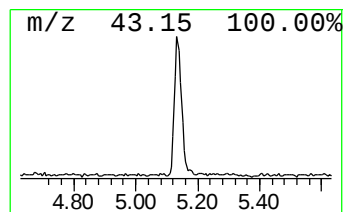
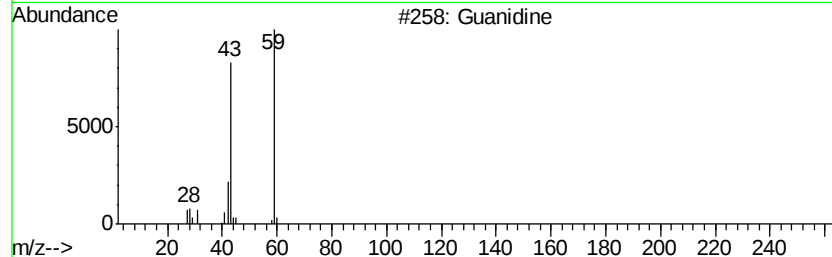
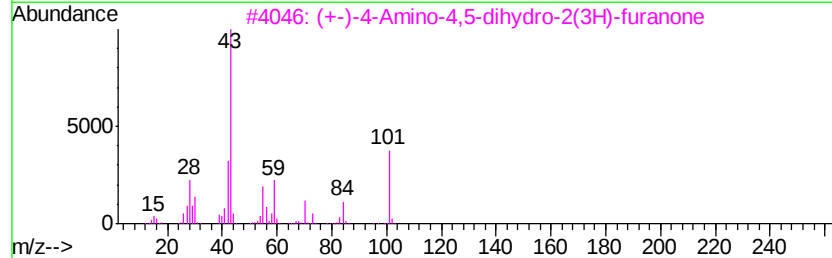
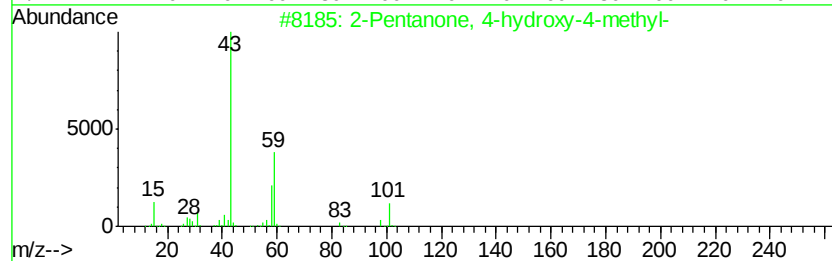
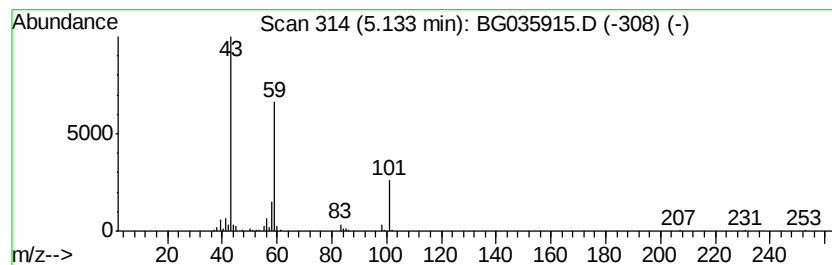
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	12.17 ng	129738	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	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3		Guanidine	59	CH5N3	000113-00-8	47
4		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	38
5		3-Hexanol	102	C6H14O	000623-37-0	37



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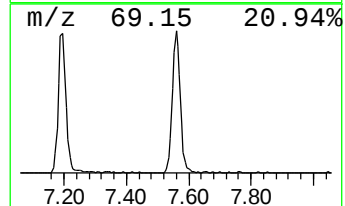
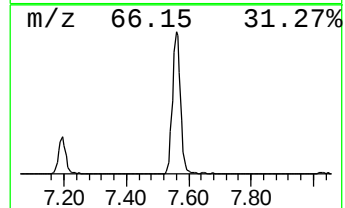
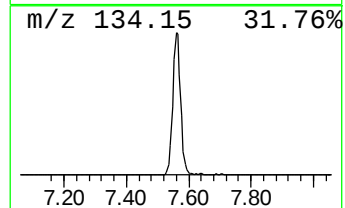
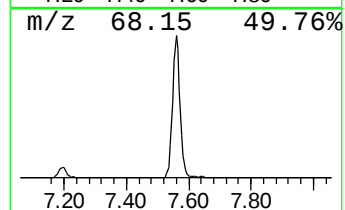
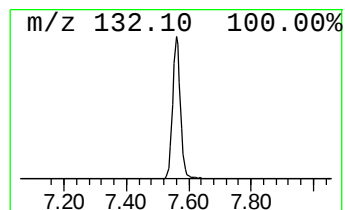
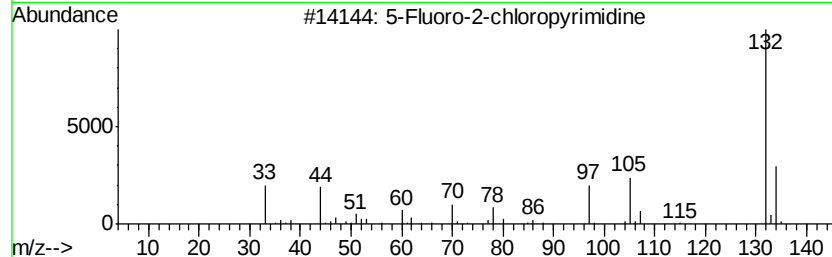
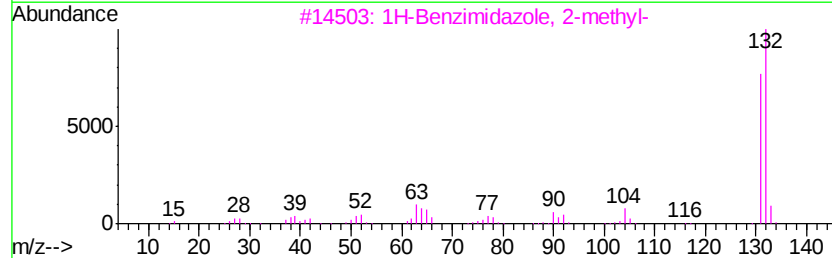
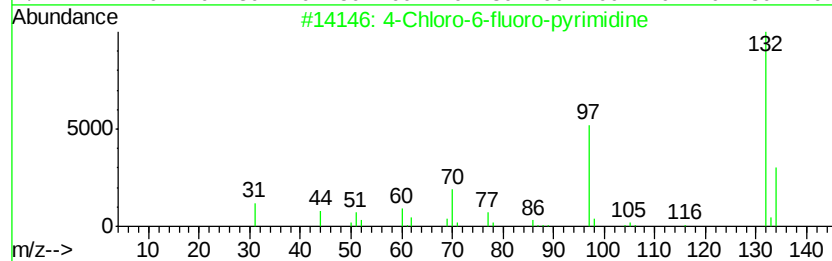
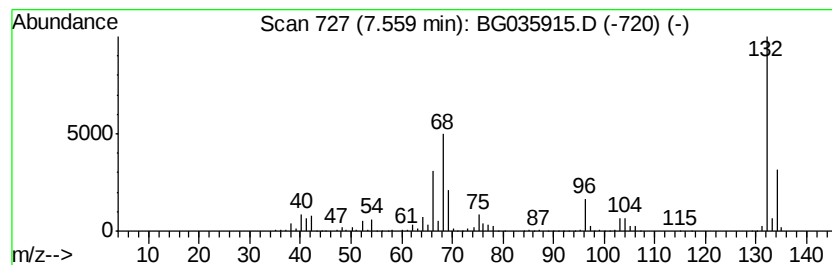
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	104.41 ng	1112780	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		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	14



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	12.2	ng	129738	1	8.02	213152	20.0
unknown7.56	7.56	104.4	ng	1112780	1	8.02	213152	20.0