

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062618\
 Data File : BG035434.D
 Acq On : 27 Jun 2018 7:15
 Operator : JU/SJ
 Sample : PB110599BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110599BL

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-BG060718.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	5.162	313	319	326	rBV	59712	94109	2.60%	0.543%
2	5.626	391	398	409	rBV	725720	1174987	32.44%	6.776%
3	7.212	660	668	679	rBV	768589	1318598	36.41%	7.605%
4	7.582	724	731	741	rBV	710483	1236965	34.15%	7.134%
5	8.052	802	811	818	rVB	178734	308080	8.51%	1.777%
6	8.370	859	865	873	rBV	519440	926903	25.59%	5.346%
7	9.210	1000	1008	1018	rBV2	433195	777571	21.47%	4.484%
8	10.855	1278	1288	1297	rBV	235258	418920	11.57%	2.416%
9	13.298	1696	1704	1715	rBV	1255400	1980619	54.68%	11.423%
10	14.673	1931	1938	1951	rVB2	418657	676987	18.69%	3.904%
11	16.159	2184	2191	2201	rBV	1099720	1693068	46.74%	9.764%
12	17.416	2398	2405	2412	rBV2	615046	950578	26.24%	5.482%
13	20.019	2842	2848	2854	rBV	2781372	3621941	100.00%	20.889%
14	21.681	3124	3131	3141	rBV2	704987	1147151	31.67%	6.616%
15	24.894	3666	3678	3691	rBV2	329371	1012911	27.97%	5.842%

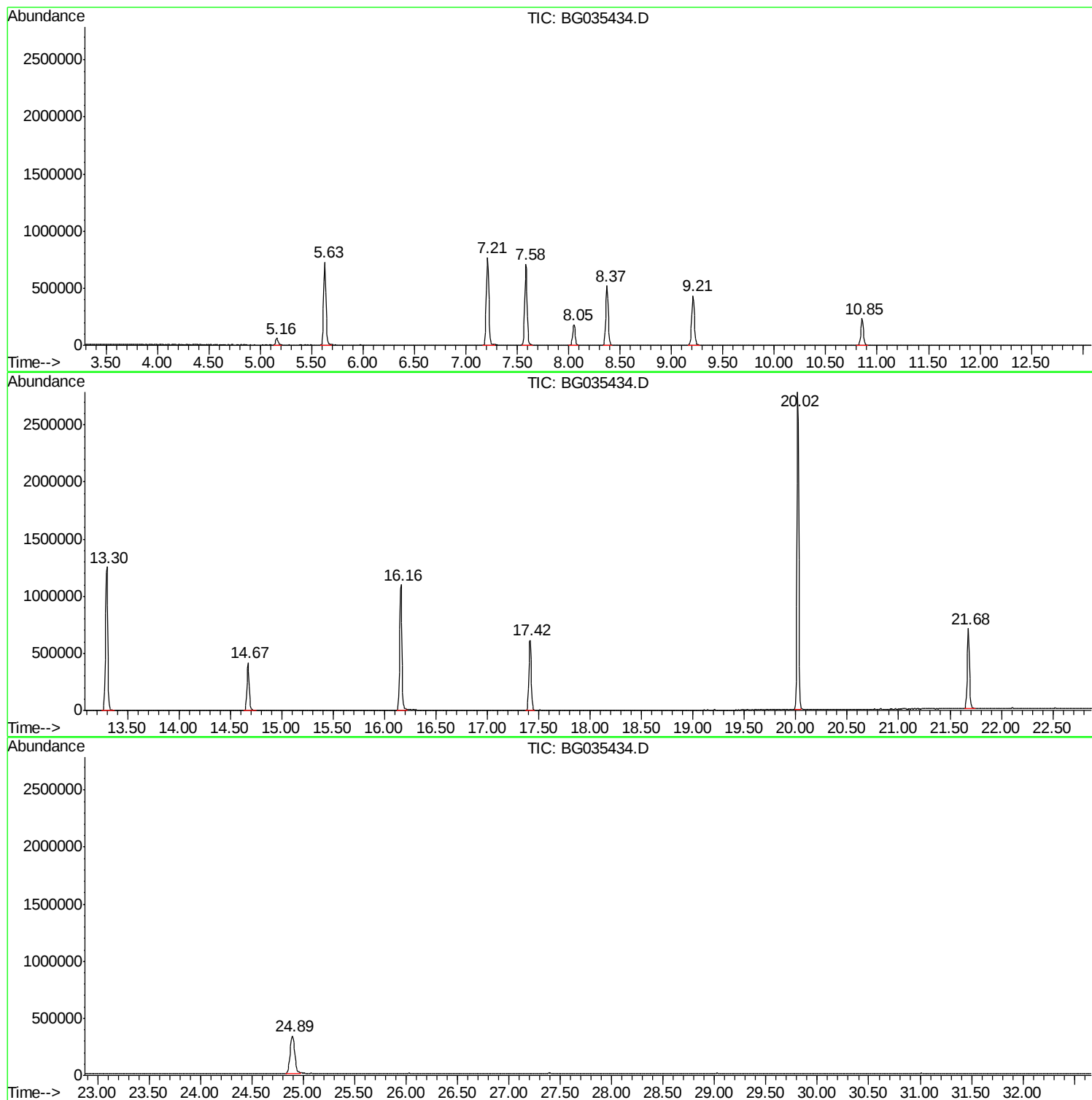
Sum of corrected areas: 17339388

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



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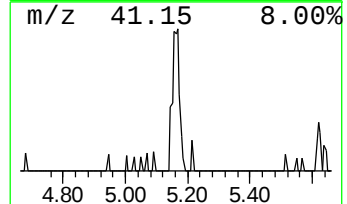
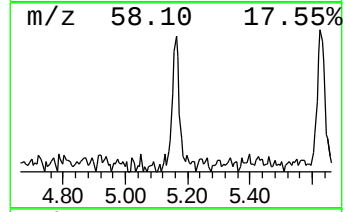
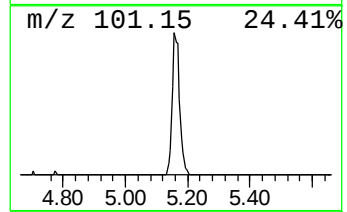
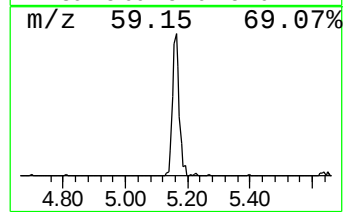
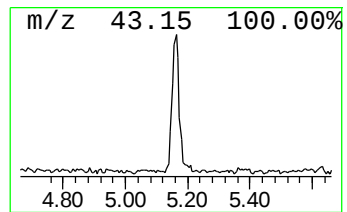
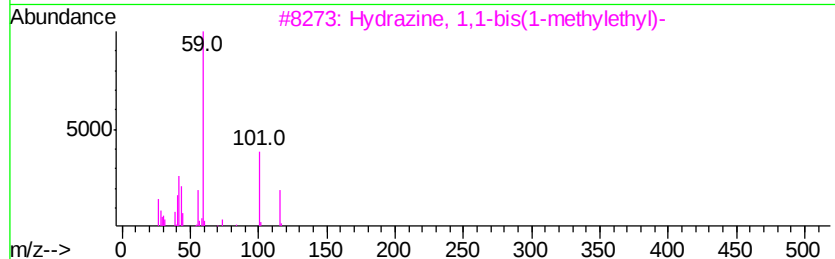
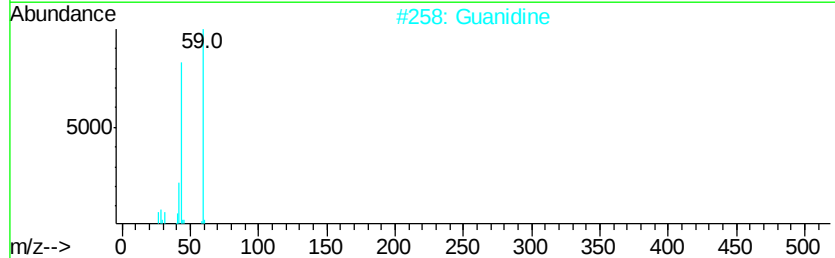
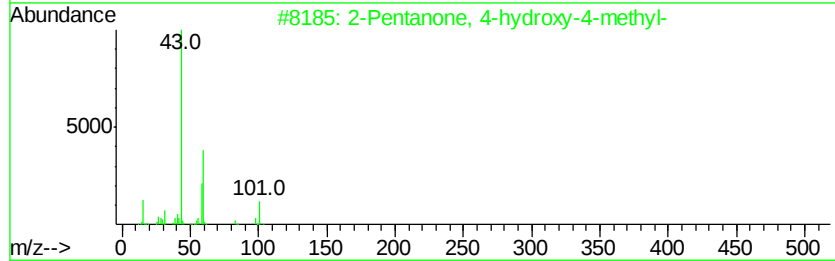
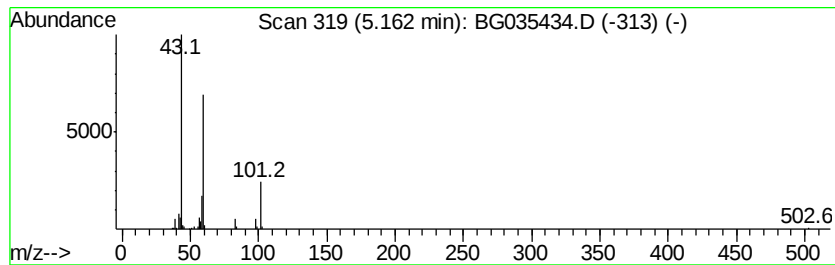
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	6.11 ng	94109	1,4-Dichlorobenzene-d4	8.05

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2		Guanidine	59	CH5N3	000113-00-8	37
3		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
4		2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	28
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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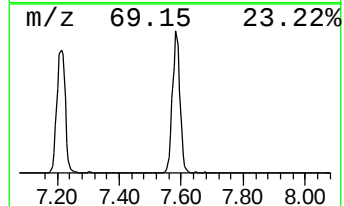
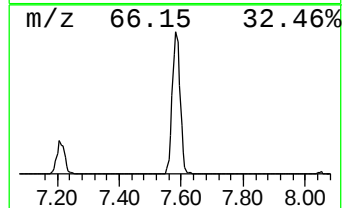
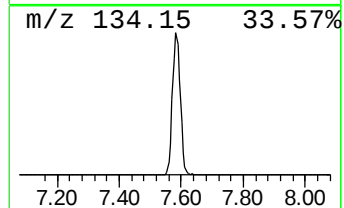
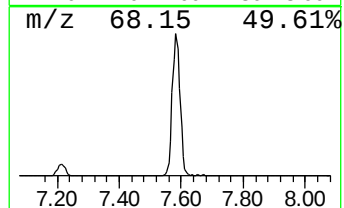
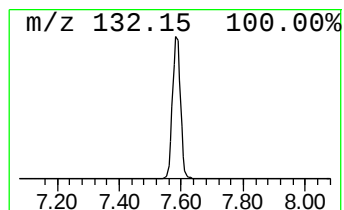
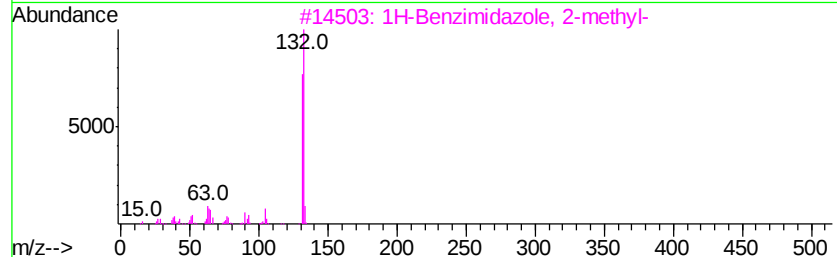
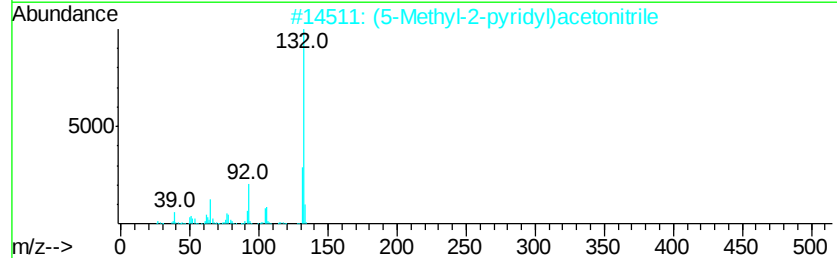
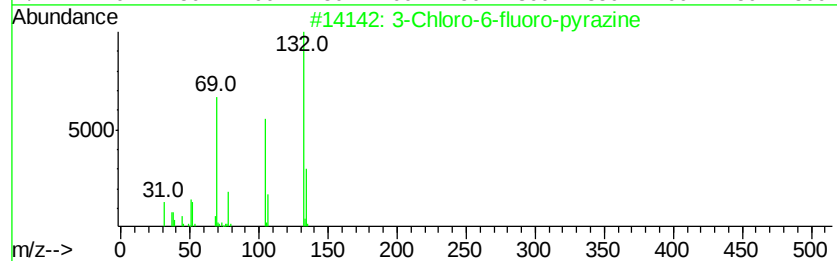
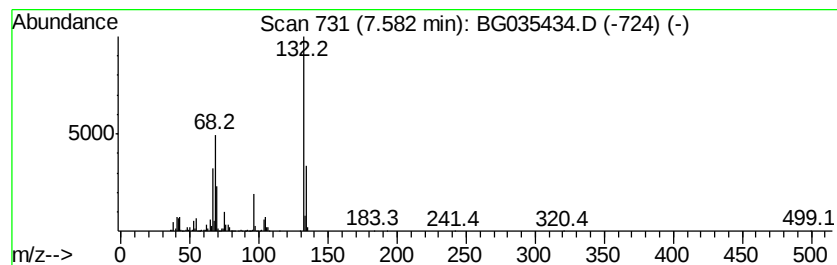
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	80.30 ng	1236970	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	6.1	ng	94109	1	8.05	308080	20.0
unknown7.58	7.58	80.3	ng	1236970	1	8.05	308080	20.0