

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG072618\
 Data File : BG035914.D
 Acq On : 26 Jul 2018 18:33
 Operator : JU/SJ
 Sample : PB111441BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111441BL

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	5.137	309	315	323	rBV	77395	120534	3.24%	0.773%
2	5.607	388	395	412	rBV	606427	1016258	27.30%	6.519%
3	7.193	657	665	676	rBV	642898	1117631	30.02%	7.169%
4	7.558	720	727	740	rBV	613970	1084939	29.15%	6.959%
5	8.027	801	807	819	rVB	129811	228338	6.13%	1.465%
6	8.345	854	861	873	rVB	457856	820945	22.05%	5.266%
7	9.185	994	1004	1018	rBV	344907	683102	18.35%	4.382%
8	10.830	1276	1284	1295	rBV	147904	303279	8.15%	1.945%
9	13.273	1691	1700	1713	rBV	1105698	1817840	48.83%	11.660%
10	14.648	1927	1934	1943	rBV2	279177	504027	13.54%	3.233%
11	16.140	2181	2188	2204	rBV	899652	1602549	43.05%	10.279%
12	17.391	2395	2401	2412	rBV	467133	788095	21.17%	5.055%
13	20.000	2839	2845	2858	rBV	2806813	3722511	100.00%	23.878%
14	21.656	3120	3127	3144	rBV	523465	926735	24.90%	5.944%
15	24.858	3661	3672	3685	rBV2	281802	853171	22.92%	5.473%

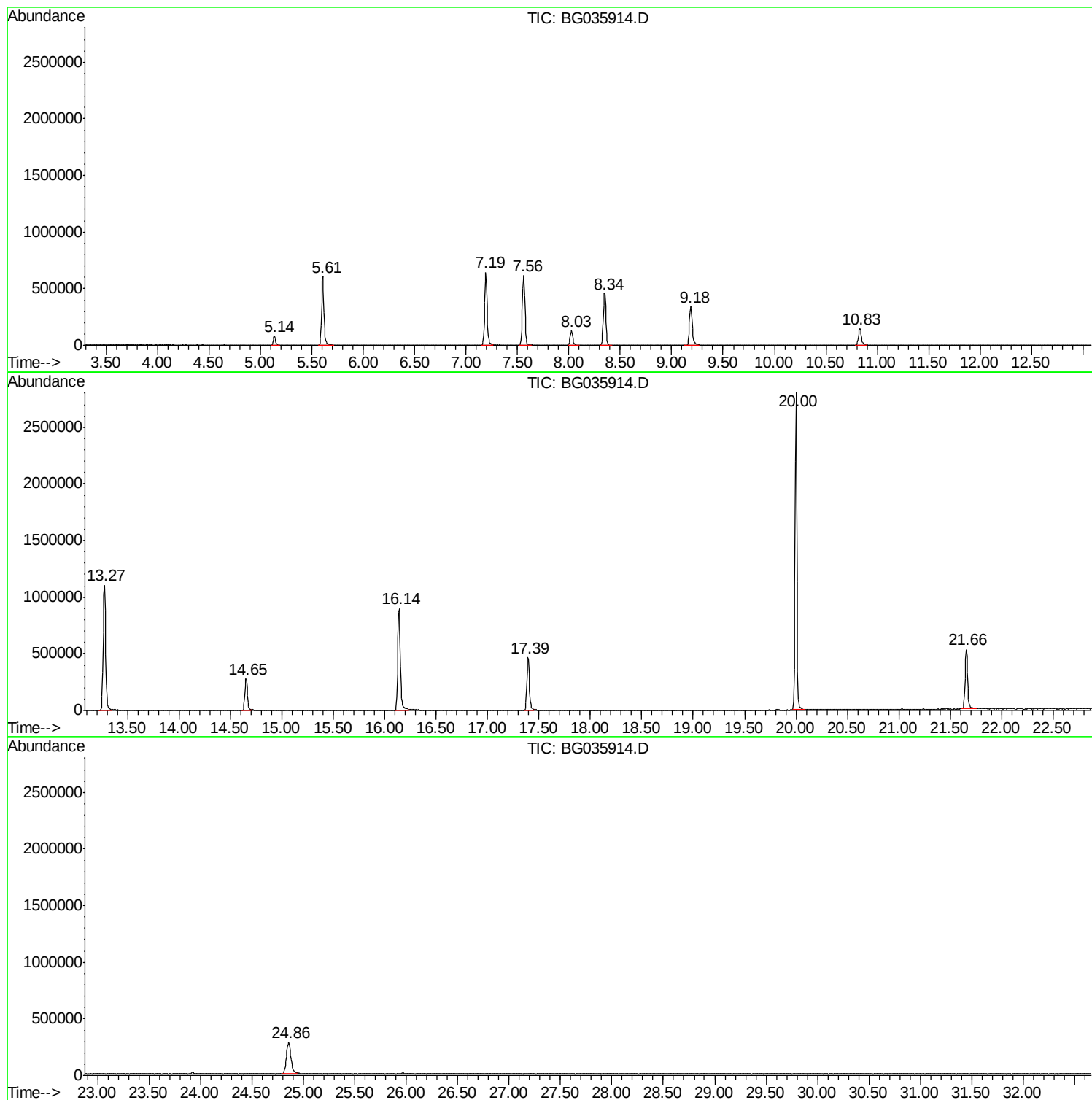
Sum of corrected areas: 15589954

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



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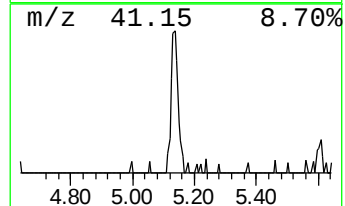
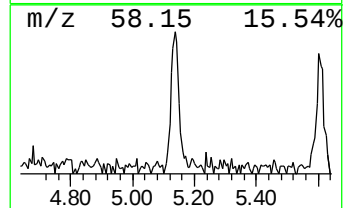
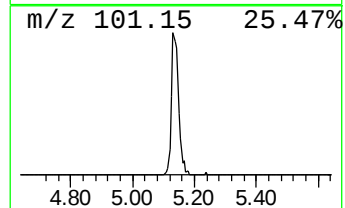
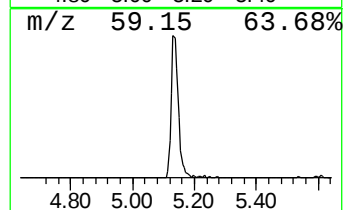
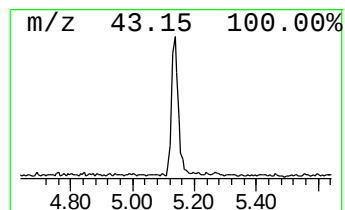
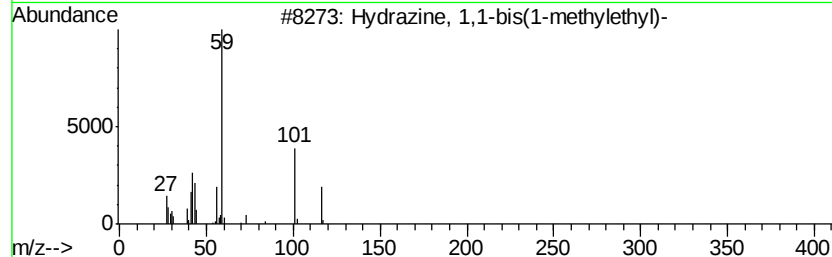
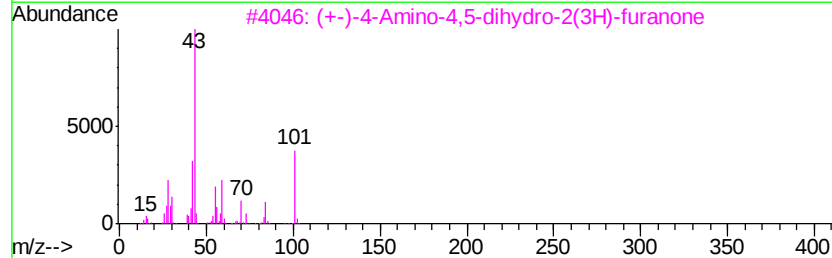
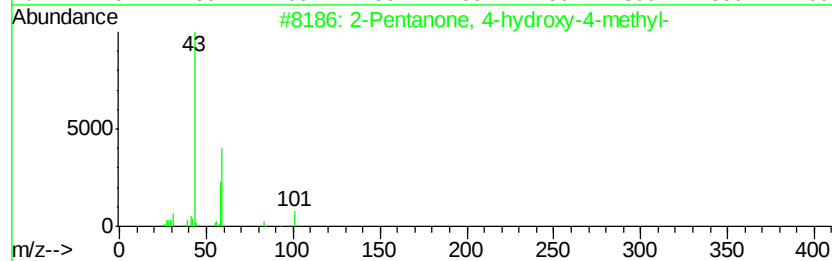
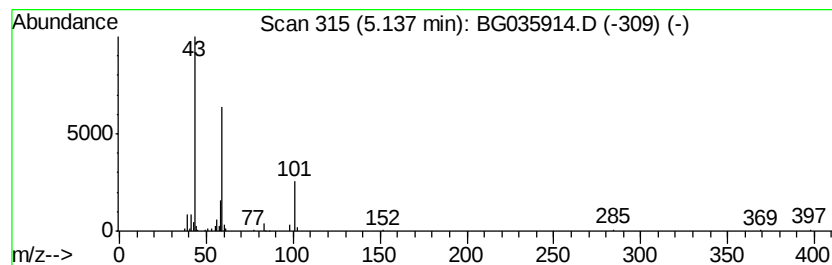
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	10.56 ng	120534	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			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	43
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			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	27



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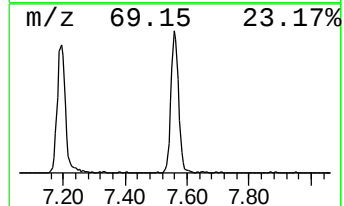
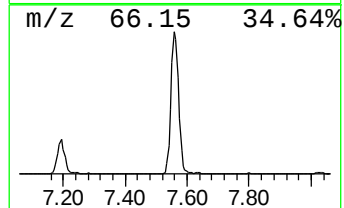
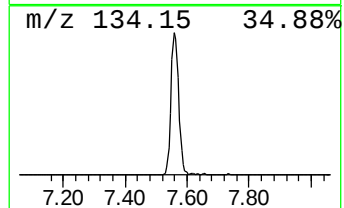
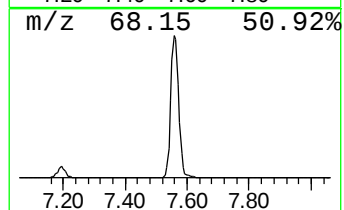
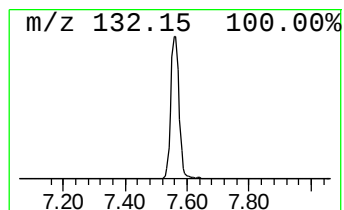
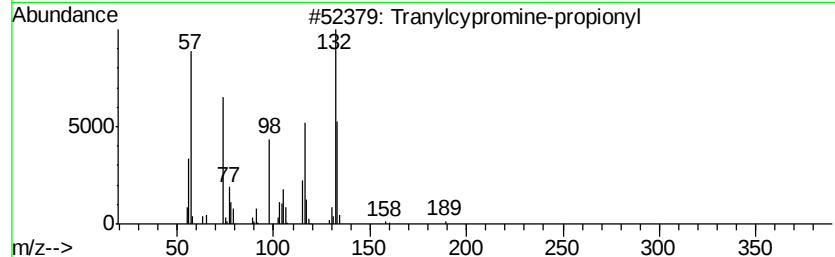
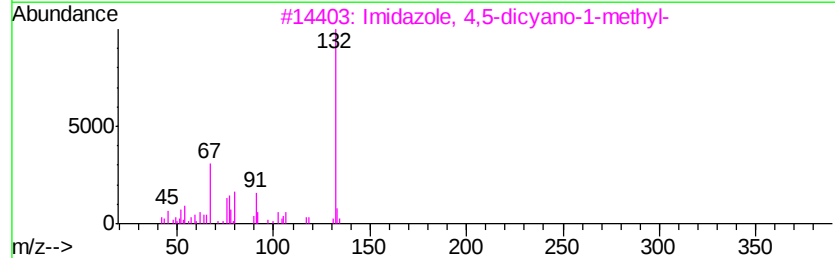
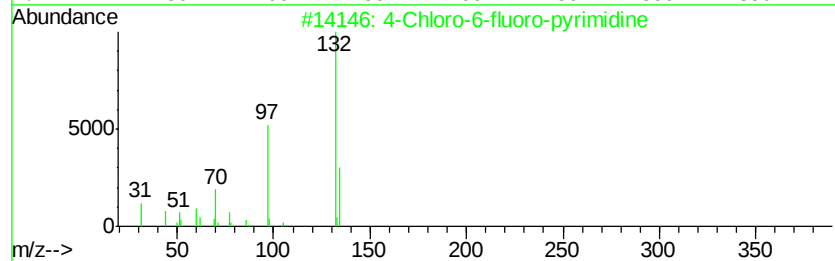
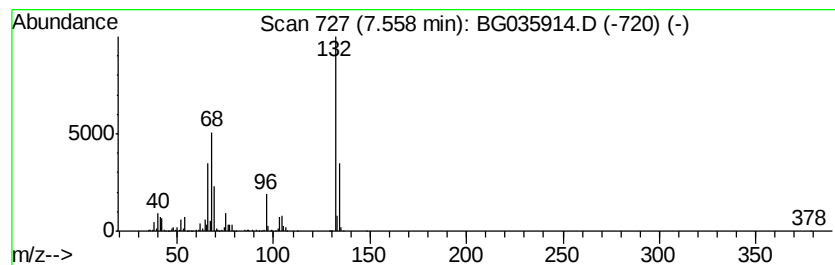
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	95.03 ng	1084940	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		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	25
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	10.6	ng	120534	1	8.02	228338	20.0
unknown7.56	7.56	95.0	ng	1084940	1	8.02	228338	20.0