

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071318\
 Data File : BG035641.D
 Acq On : 13 Jul 2018 22:40
 Operator : JU/SJ
 Sample : PB111092BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111092BL

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	5.162	314	319	328	rBV	116610	180596	3.82%	0.882%
2	5.632	390	399	410	rBV	933549	1535883	32.52%	7.499%
3	7.218	660	669	683	rBV	974765	1717203	36.36%	8.384%
4	7.588	724	732	741	rBV	910260	1595996	33.79%	7.792%
5	8.058	805	812	819	rBV	147526	260236	5.51%	1.271%
6	8.381	857	867	877	rBV	642042	1142010	24.18%	5.576%
7	9.216	1001	1009	1022	rBV	554772	1017873	21.55%	4.969%
8	10.860	1281	1289	1297	rBV	201723	364329	7.71%	1.779%
9	13.298	1697	1704	1713	rBV	1569228	2532512	53.62%	12.364%
10	14.673	1932	1938	1948	rBV2	339595	575747	12.19%	2.811%
11	16.165	2184	2192	2207	rBV	1306680	2127898	45.06%	10.389%
12	17.416	2399	2405	2413	rBV	530981	830116	17.58%	4.053%
13	20.024	2842	2849	2855	rBV	3612804	4722677	100.00%	23.057%
14	21.681	3123	3131	3139	rBV2	580768	994317	21.05%	4.854%
15	24.900	3668	3679	3691	rBV2	305736	885043	18.74%	4.321%

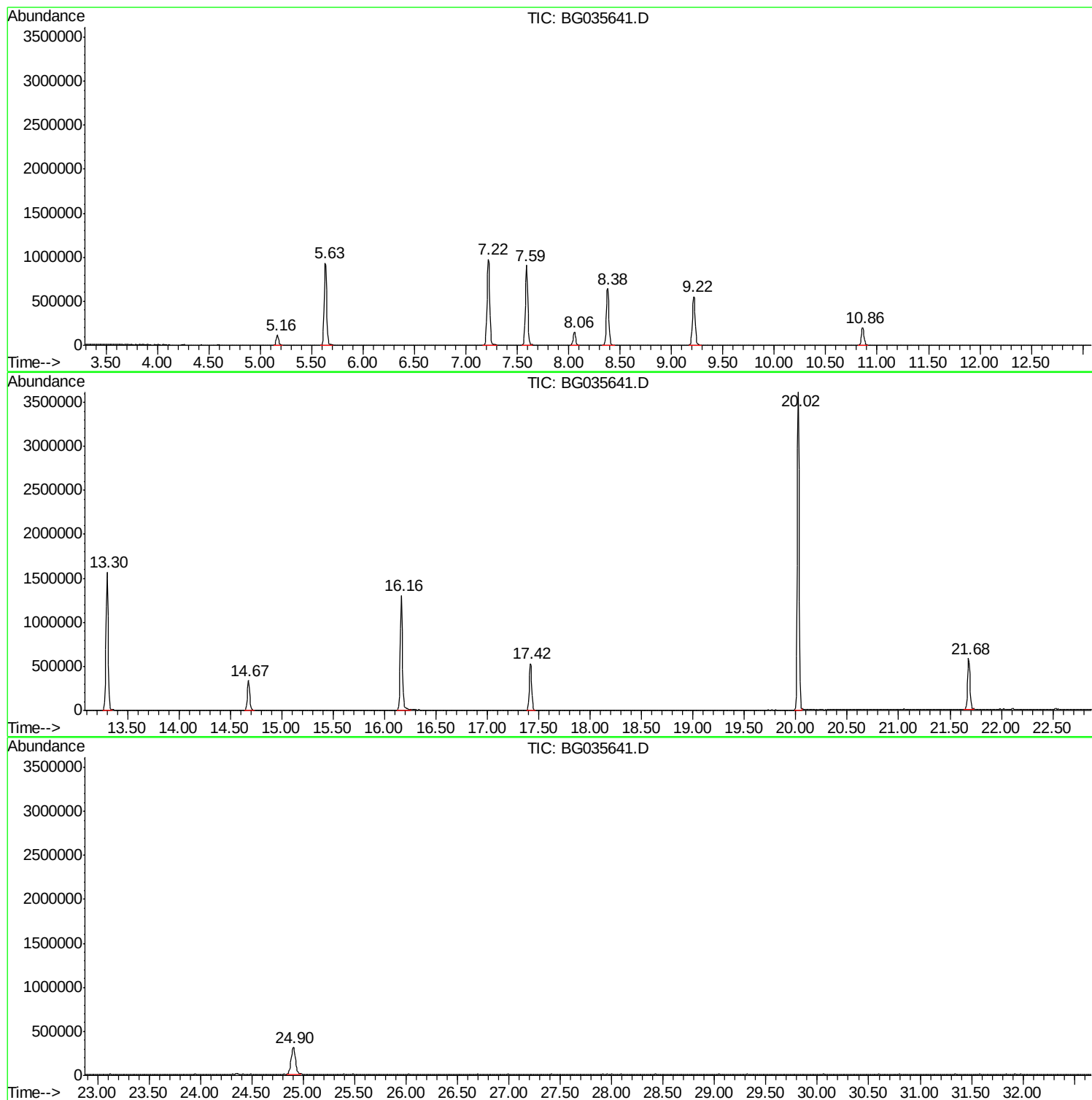
Sum of corrected areas: 20482436

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



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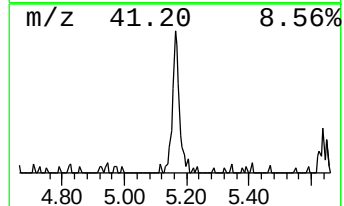
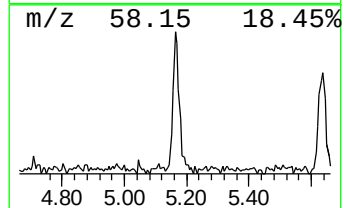
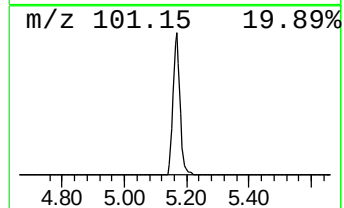
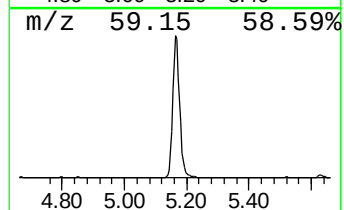
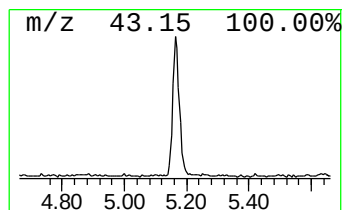
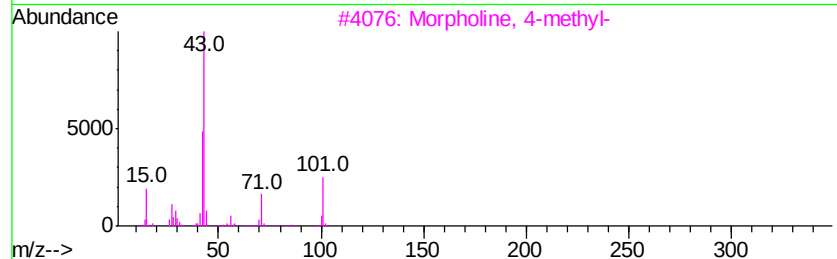
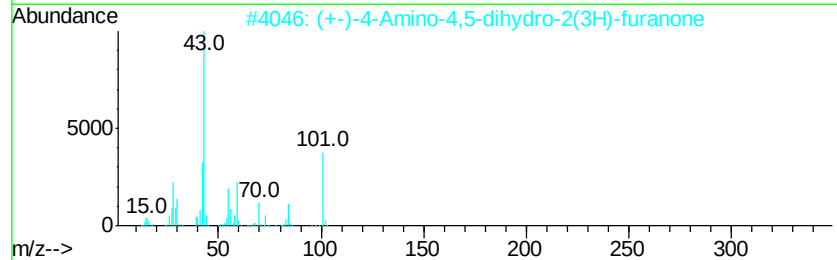
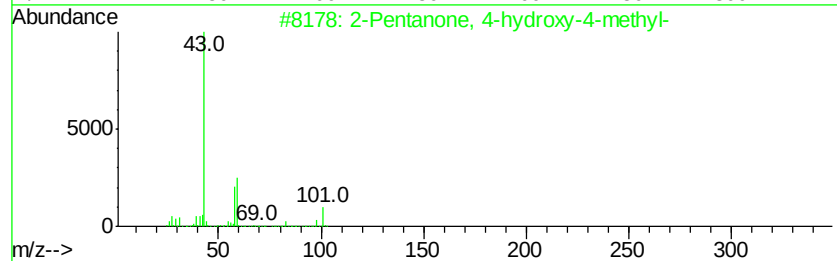
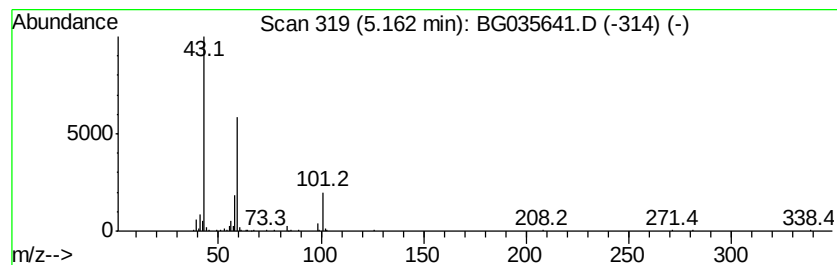
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	13.88 ng	180596	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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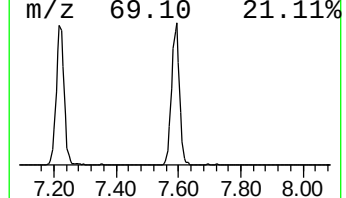
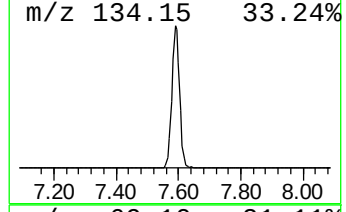
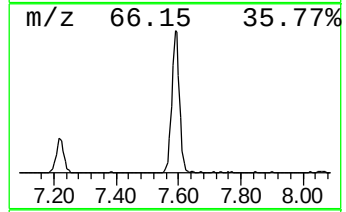
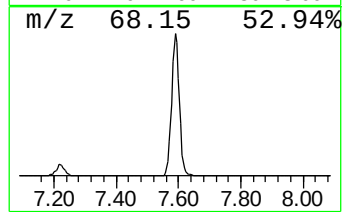
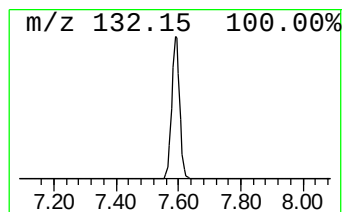
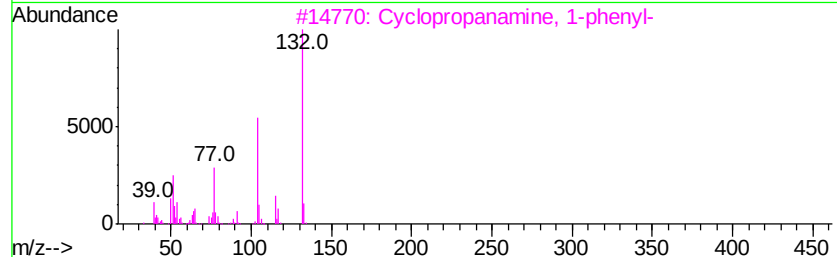
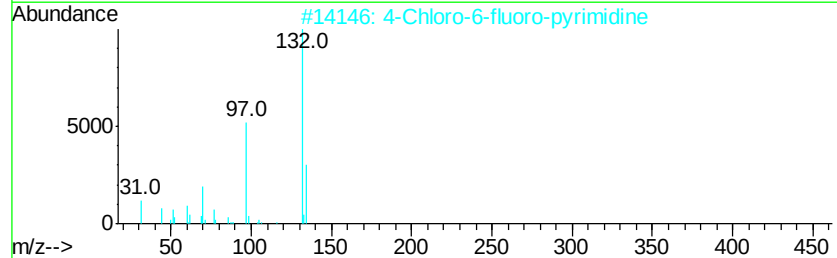
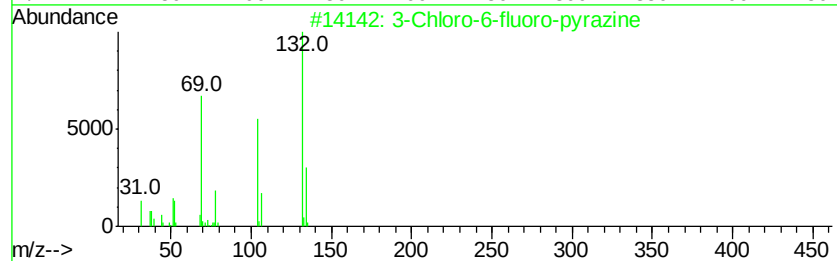
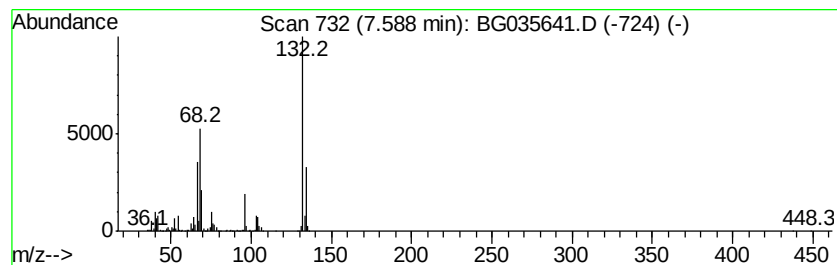
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	122.66 ng	1596000	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	13.9	ng	180596	1	8.06	260236	20.0
unknown7.59	7.59	122.7	ng	1596000	1	8.06	260236	20.0