

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100118\
Data File : BG037101.D
Acq On : 2 Oct 2018 12:11
Operator : JU/SJ
Sample : PB113413BL
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB113413BL

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-BG092018.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	5.147	311	316	326	rVB	41614	72087	1.65%	0.311%
2	5.617	389	396	418	rBV	1037875	1934905	44.30%	8.359%
3	7.203	659	666	688	rBV	1028217	2052118	46.98%	8.866%
4	7.574	722	729	748	rBV	1025856	1983393	45.41%	8.569%
5	8.038	802	808	823	rBV	192425	396175	9.07%	1.712%
6	8.361	855	863	881	rBV	801112	1558293	35.68%	6.732%
7	9.201	998	1006	1029	rBV	608401	1327176	30.39%	5.734%
8	10.840	1276	1285	1306	rBV	260644	579254	13.26%	2.503%
9	13.284	1694	1701	1724	rBV	1695809	2967886	67.95%	12.822%
10	14.665	1926	1936	1949	rBV2	431323	793728	18.17%	3.429%
11	16.152	2182	2189	2215	rBV	1048178	1938682	44.39%	8.376%
12	17.409	2397	2403	2424	rBV2	564501	988555	22.63%	4.271%
13	20.018	2840	2847	2862	rBV	3043709	4367621	100.00%	18.870%
14	21.675	3122	3129	3148	rVB2	578605	1118078	25.60%	4.830%
15	24.877	3663	3674	3698	rBV2	312940	1068301	24.46%	4.615%

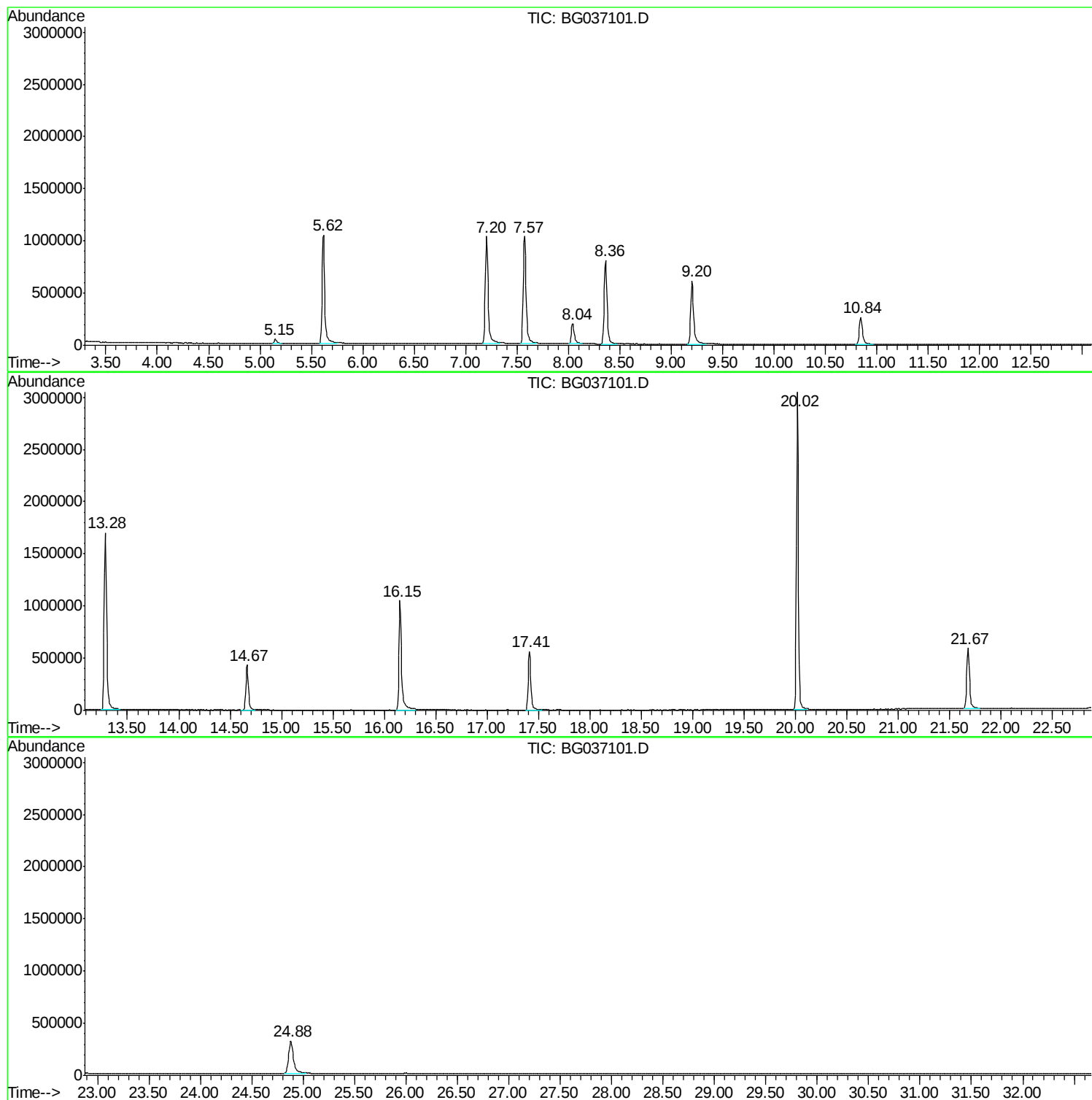
Sum of corrected areas: 23146252

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



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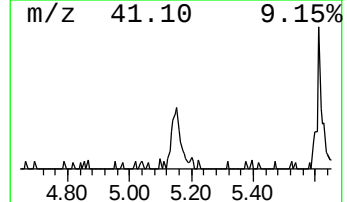
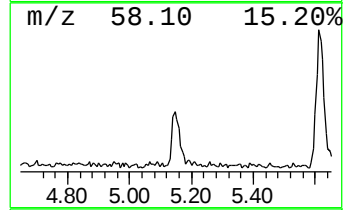
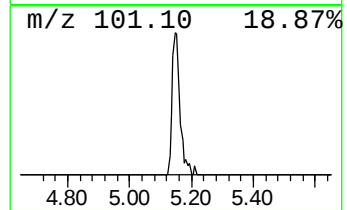
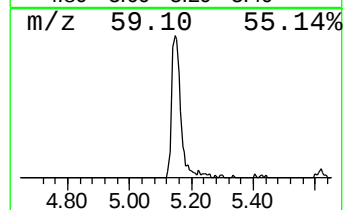
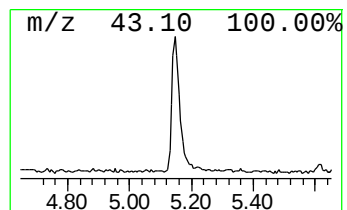
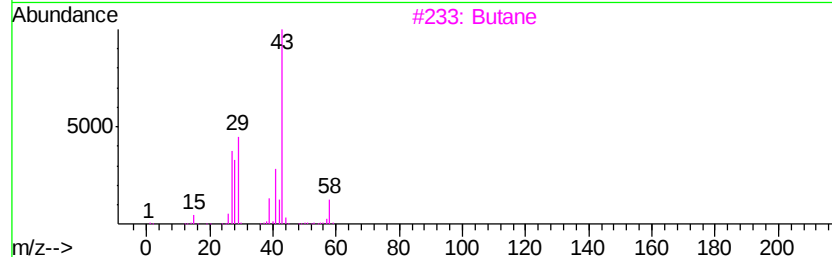
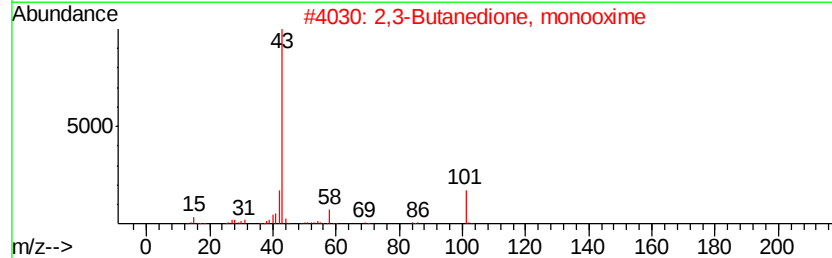
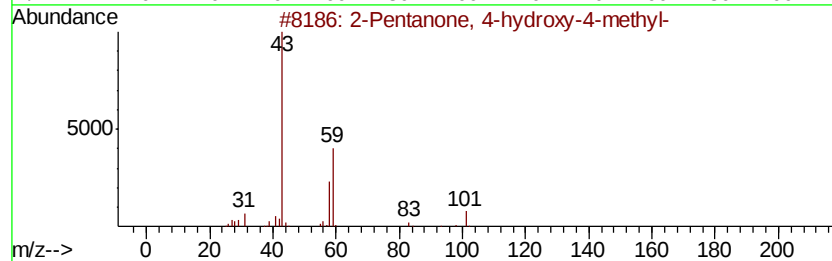
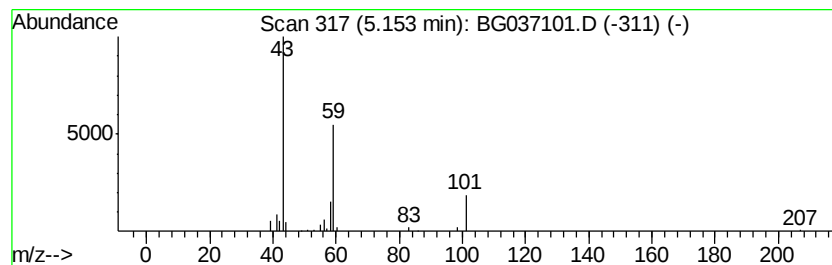
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.64 ng	72087	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Butane	58	C4H10	000106-97-8	9
4			Di-n-propyl ether	102	C6H14O	000111-43-3	9
5			Cycloserine	102	C3H6N2O2	000068-41-7	9



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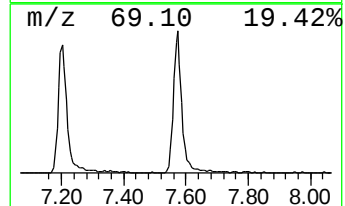
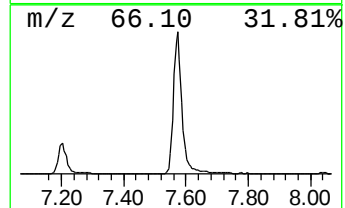
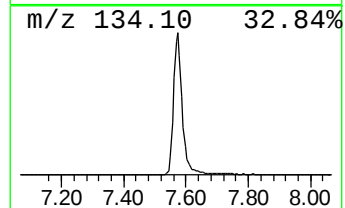
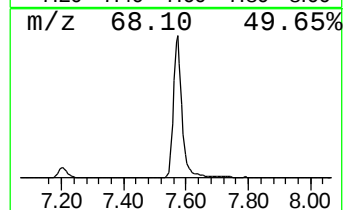
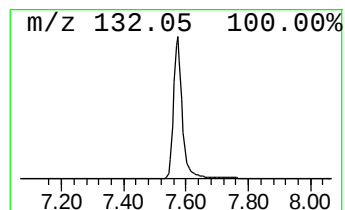
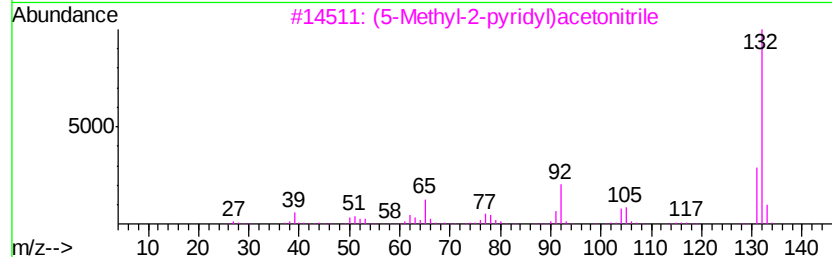
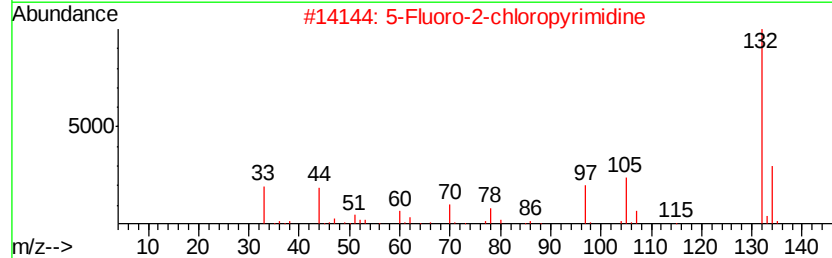
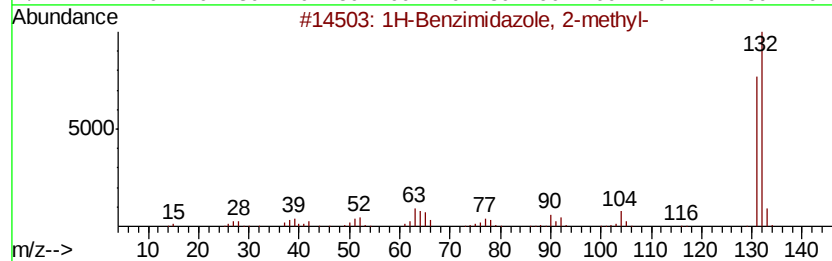
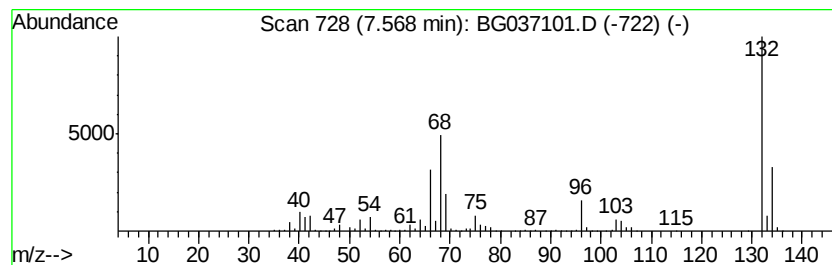
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	100.13 ng	1983390	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	3.6	ng	72087	1	8.04	396175	20.0
unknown7.57	7.57	100.1	ng	1983390	1	8.04	396175	20.0