

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092718\
 Data File : BG037048.D
 Acq On : 28 Sep 2018 13:26
 Operator : JU/SJ
 Sample : PB113326BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113326BL

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	5.145	312	316	326	rVB	26692	42158	1.46%	0.299%
2	5.615	390	396	422	rBV	580478	1078304	37.33%	7.654%
3	7.201	659	666	685	rBV	574250	1148732	39.77%	8.154%
4	7.572	722	729	749	rBV	569806	1103657	38.21%	7.834%
5	8.042	802	809	821	rBV2	118401	242930	8.41%	1.724%
6	8.359	856	863	885	rBV	442644	869772	30.11%	6.174%
7	9.199	999	1006	1027	rBV	324066	724545	25.08%	5.143%
8	10.844	1279	1286	1301	rBV	157154	331582	11.48%	2.354%
9	13.288	1695	1702	1722	rBV	929444	1626219	56.30%	11.543%
10	14.663	1930	1936	1952	rBV3	260706	480662	16.64%	3.412%
11	16.150	2182	2189	2212	rBV	736045	1326735	45.93%	9.417%
12	17.413	2397	2404	2417	rBV	381288	661705	22.91%	4.697%
13	20.016	2841	2847	2863	rBV	2101968	2888508	100.00%	20.503%
14	21.679	3123	3130	3145	rBV	425979	803344	27.81%	5.702%
15	24.881	3665	3675	3695	rVB	247073	759408	26.29%	5.390%

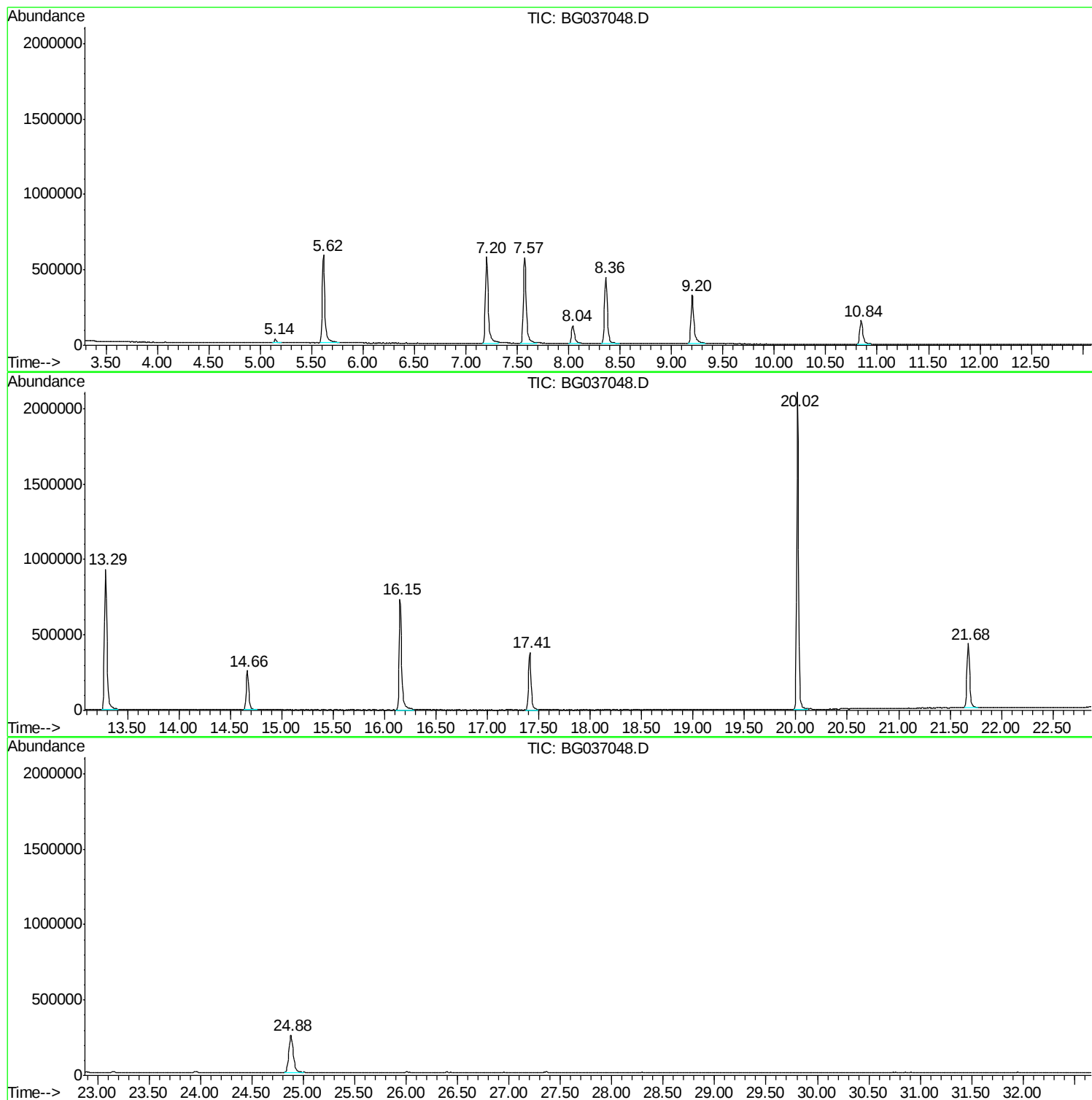
Sum of corrected areas: 14088261

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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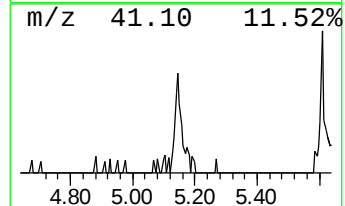
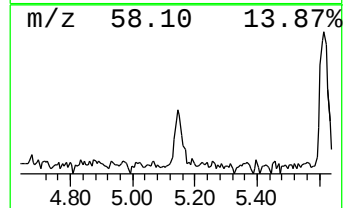
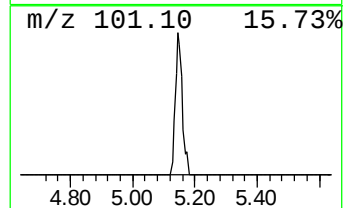
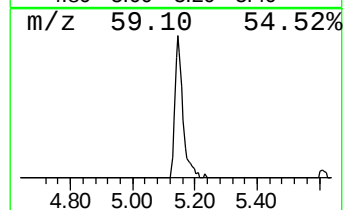
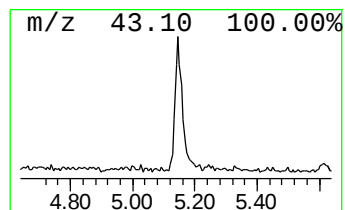
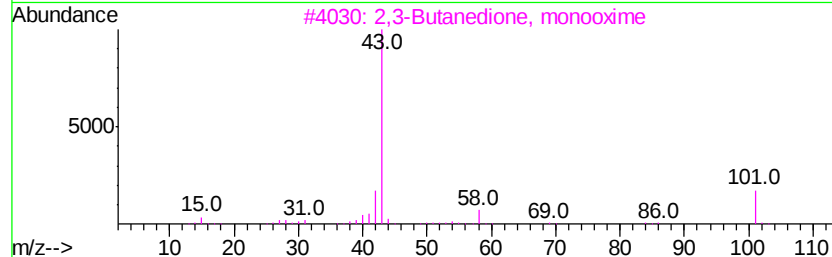
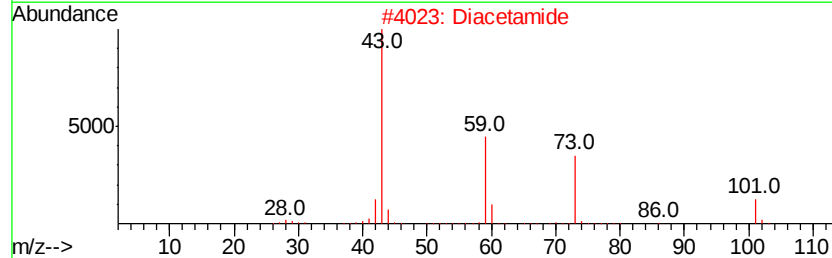
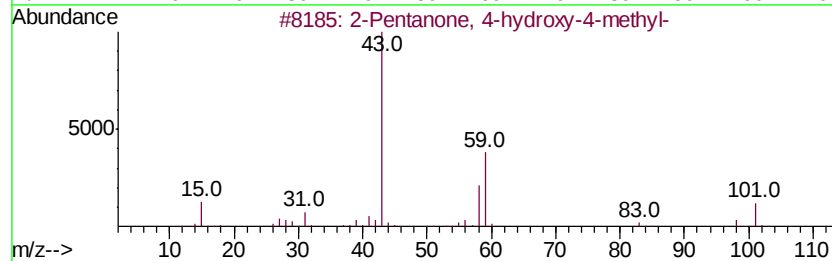
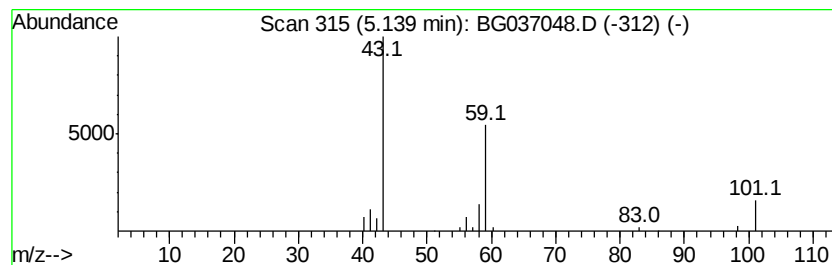
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	3.47 ng	42158	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		Diacetamide	101	C4H7NO2	000625-77-4	9
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		3,3-Dimethylbutylamine	101	C6H15N	015673-00-4	7
5		Propylamine	59	C3H9N	000107-10-8	7



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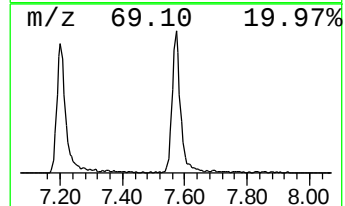
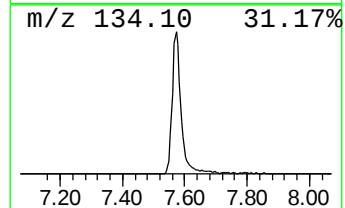
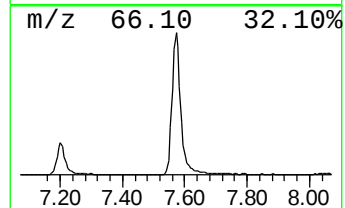
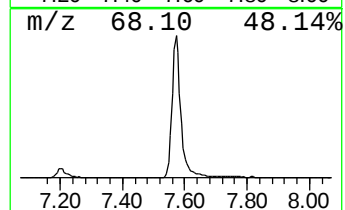
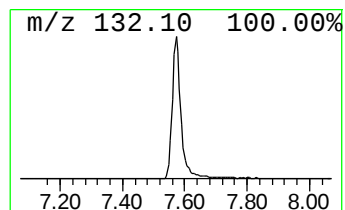
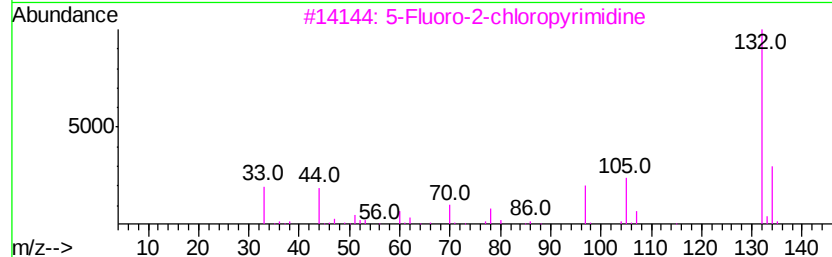
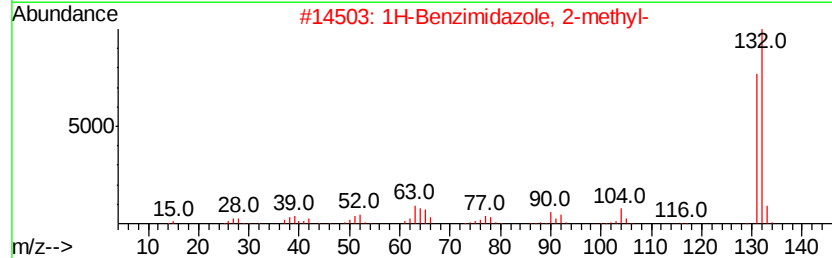
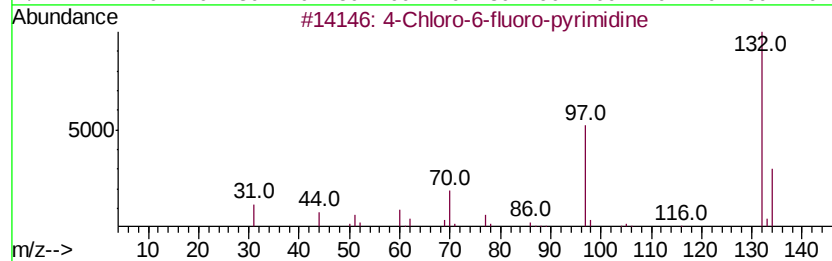
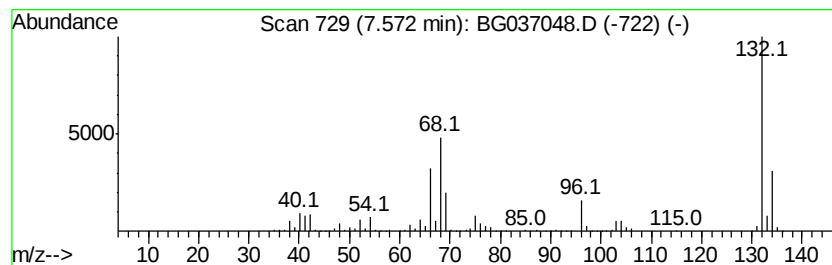
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	90.86 ng	1103660	1,4-Dichlorobenzene-d4	8.04

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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	3.5	ng	42158	1	8.04	242930	20.0
unknown7.57	7.57	90.9	ng	1103660	1	8.04	242930	20.0