

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091718\
 Data File : BG036750.D
 Acq On : 17 Sep 2018 12:09
 Operator : JU/SJ
 Sample : PB112885BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112885BL

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-BG082318.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	5.158	313	318	327	rVB	47923	74000	1.79%	0.376%
2	5.622	390	397	407	rBV	964987	1508515	36.55%	7.664%
3	7.209	658	667	679	rBV	960187	1621466	39.29%	8.238%
4	7.579	722	730	738	rBV	872627	1513161	36.67%	7.687%
5	8.049	803	810	817	rBV	195477	327143	7.93%	1.662%
6	8.372	857	865	874	rBV	690122	1182404	28.65%	6.007%
7	9.212	999	1008	1015	rBV	545136	982902	23.82%	4.993%
8	10.857	1280	1288	1296	rBV	250799	454273	11.01%	2.308%
9	13.296	1696	1703	1715	rBV	1553310	2368614	57.39%	12.033%
10	14.676	1930	1938	1948	rBV2	401030	658959	15.97%	3.348%
11	16.157	2184	2190	2204	rBV	1137368	1838803	44.56%	9.342%
12	17.420	2398	2405	2413	rBV	594053	920410	22.30%	4.676%
13	20.023	2842	2848	2861	rBV	3152567	4126958	100.00%	20.966%
14	21.680	3123	3130	3141	rBV2	662943	1095311	26.54%	5.565%
15	24.888	3666	3676	3688	rBV2	356008	1010720	24.49%	5.135%

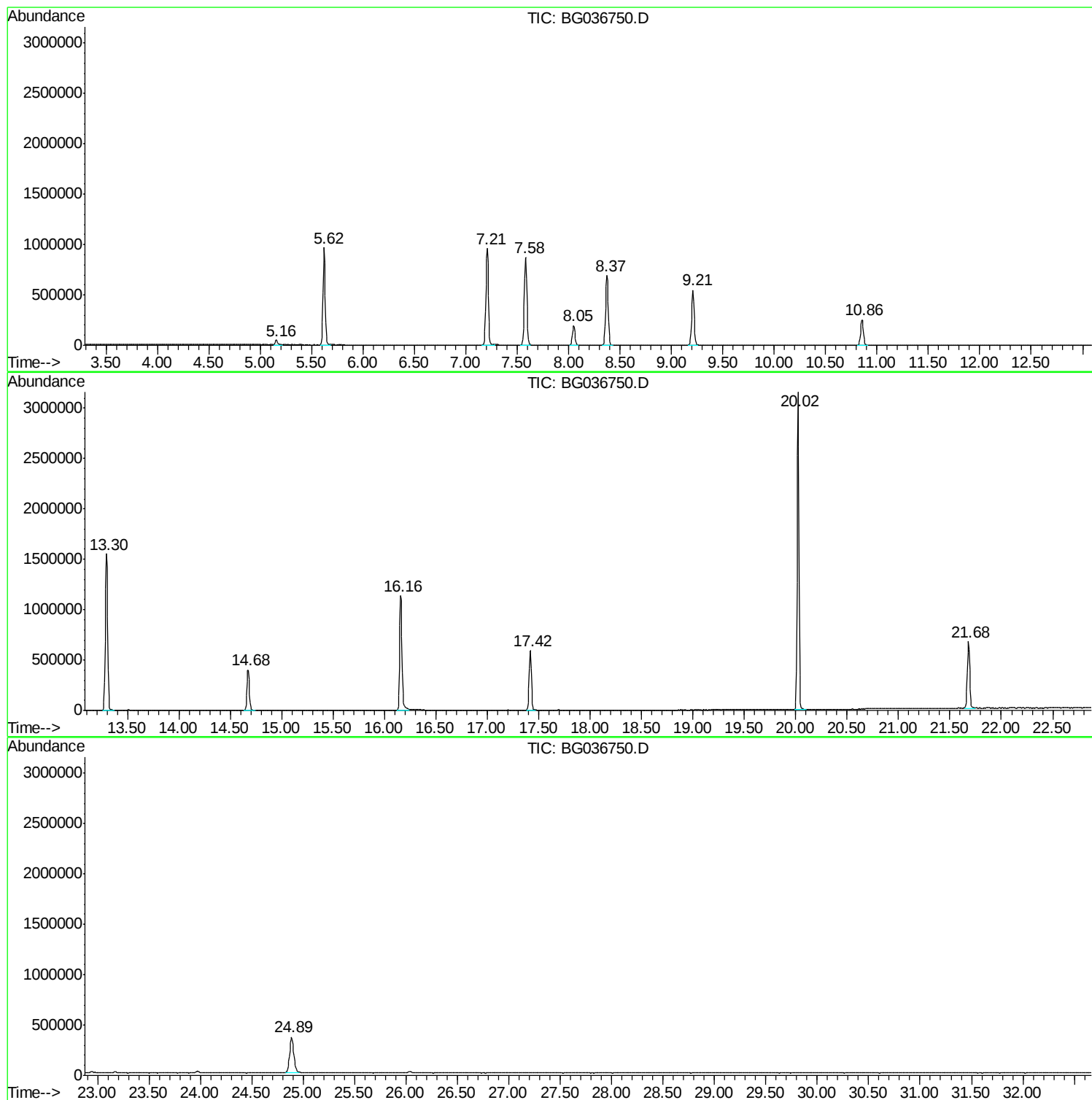
Sum of corrected areas: 19683639

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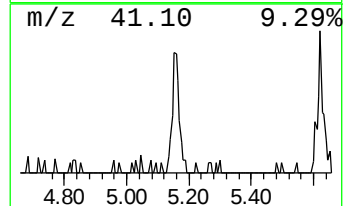
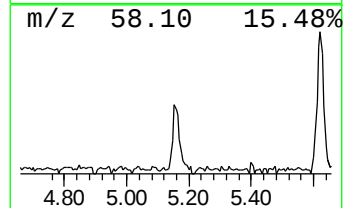
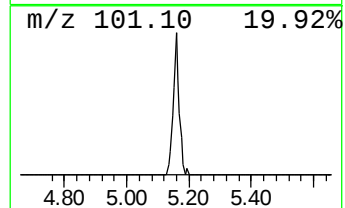
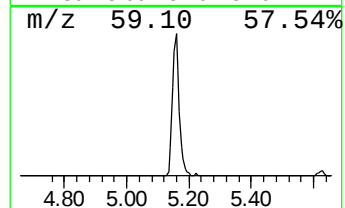
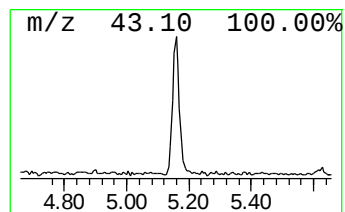
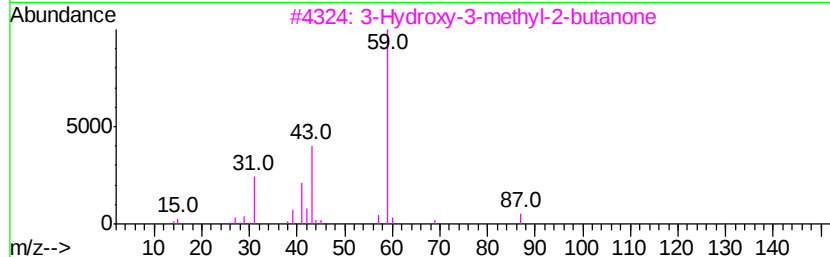
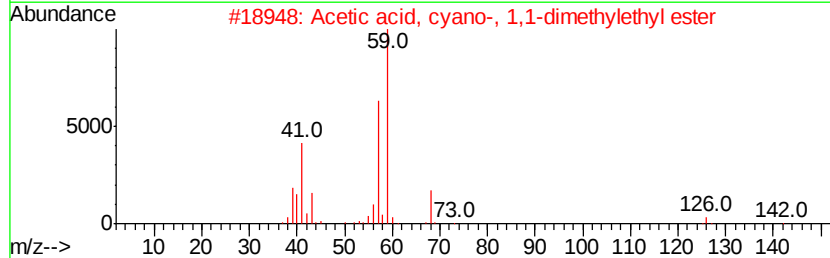
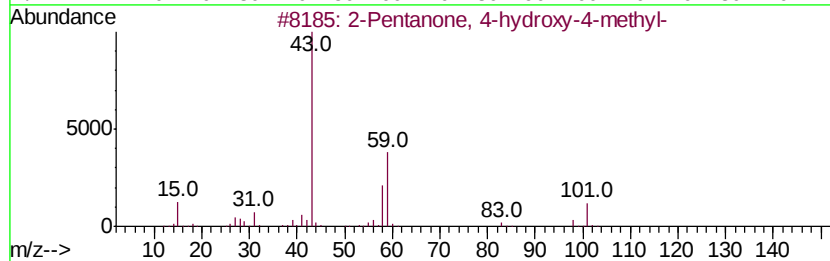
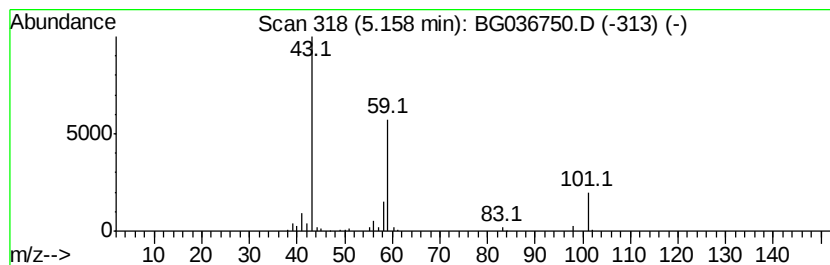
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.52 ng	74000	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	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
4			Propanoic acid, 2-hydroxy-2-meth...	132	C6H12O3	000080-55-7	9
5			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9



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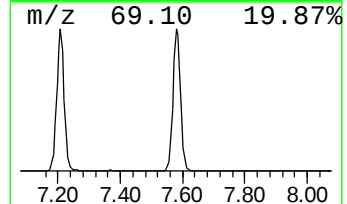
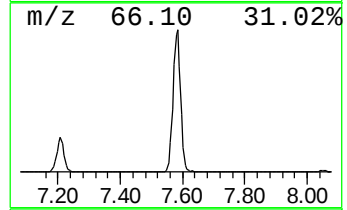
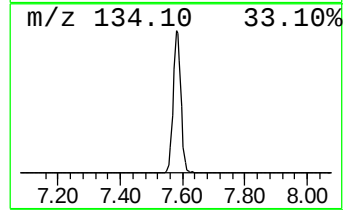
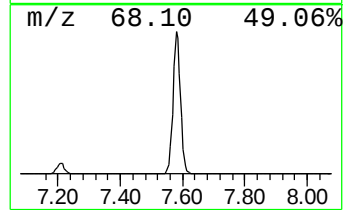
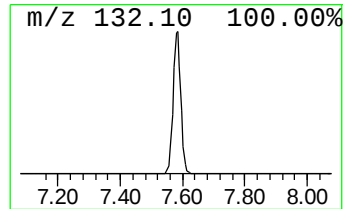
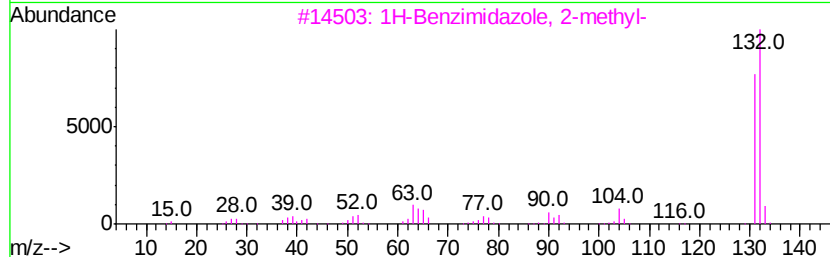
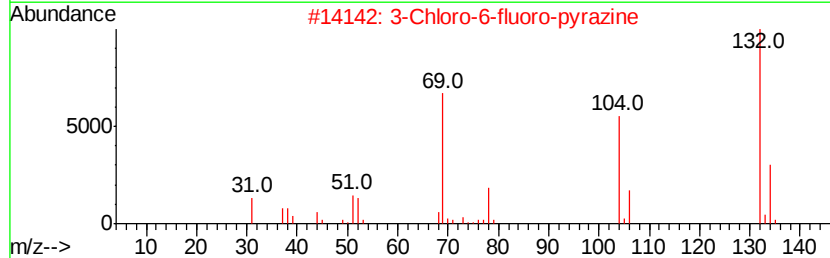
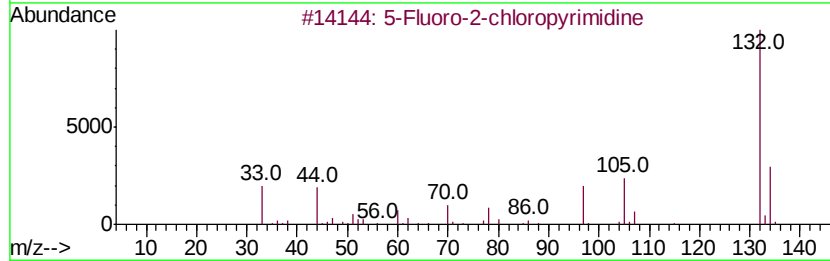
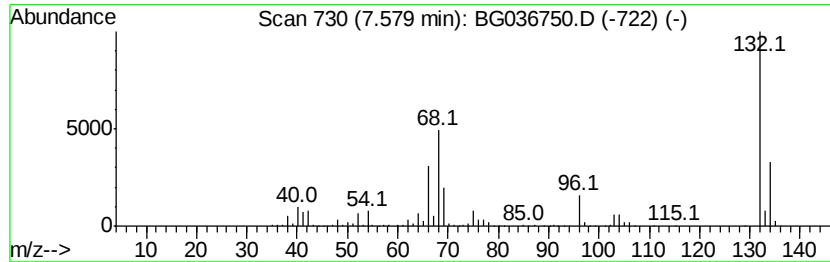
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	92.51 ng	1513160	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
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	4.5	ng	74000	1	8.05	327143	20.0
unknown7.58	7.58	92.5	ng	1513160	1	8.05	327143	20.0