

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060718\
 Data File : BG034970.D
 Acq On : 8 Jun 2018 2:30
 Operator : SJ/JU
 Sample : PB110034BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110034BL

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-BG060718.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	5.187	317	323	330	rVB	48966	77082	2.20%	0.485%
2	5.646	394	401	413	rVB	713722	1126565	32.17%	7.087%
3	7.226	662	670	678	rBV	734317	1268757	36.23%	7.981%
4	7.608	726	735	743	rBV	683667	1162078	33.19%	7.310%
5	8.078	808	815	822	rBV	159019	259286	7.40%	1.631%
6	8.401	862	870	879	rBV	508094	885599	25.29%	5.571%
7	9.235	1003	1012	1022	rBV	413397	735059	20.99%	4.624%
8	10.886	1285	1293	1301	rVB	197180	343703	9.82%	2.162%
9	13.329	1702	1709	1717	rBV	1183501	1880057	53.69%	11.827%
10	14.698	1933	1942	1952	rVB2	337101	538646	15.38%	3.388%
11	16.184	2187	2195	2208	rBV	1029235	1597064	45.61%	10.047%
12	17.441	2402	2409	2416	rBV	500535	747091	21.34%	4.700%
13	20.056	2847	2854	2859	rBV	2728258	3501580	100.00%	22.028%
14	21.712	3129	3136	3145	rBV	562457	922030	26.33%	5.800%
15	24.955	3676	3688	3702	rBV2	305269	851714	24.32%	5.358%

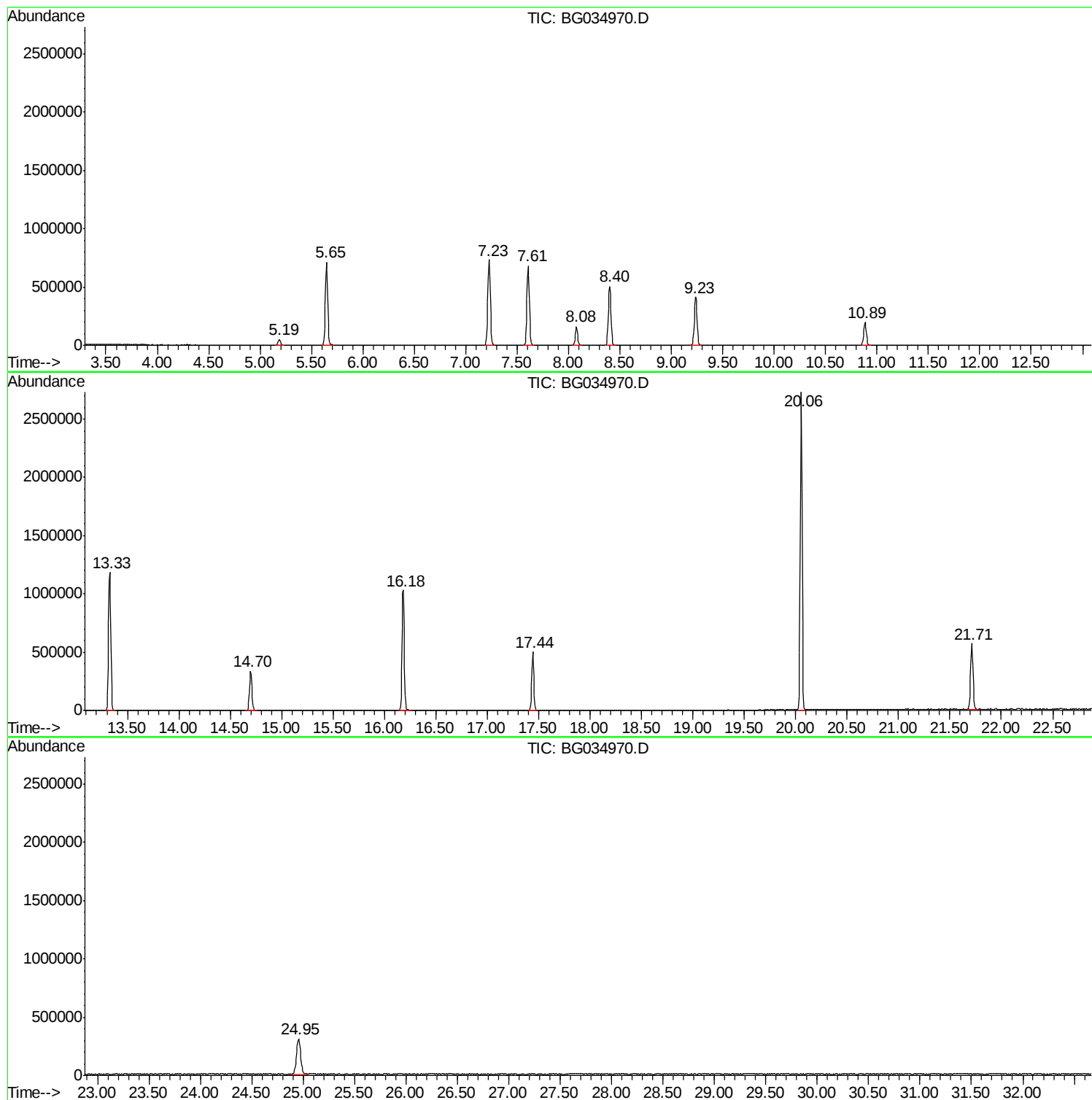
Sum of corrected areas: 15896311

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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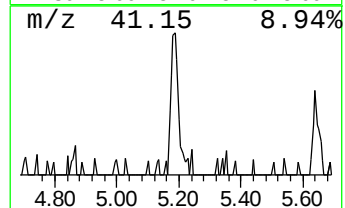
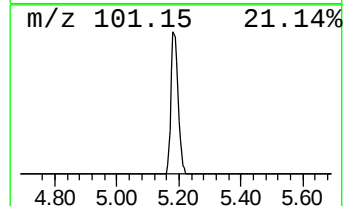
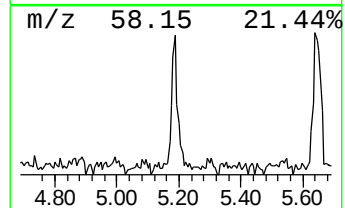
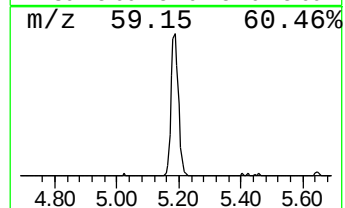
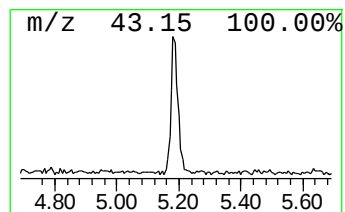
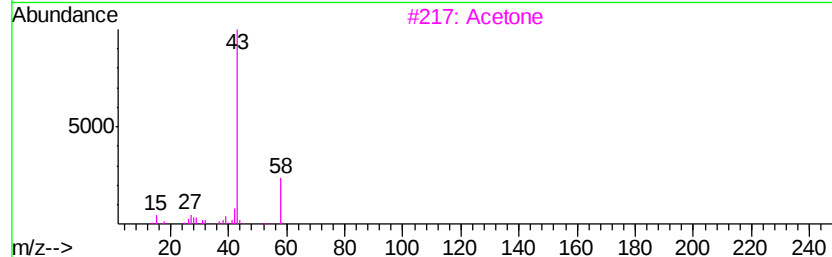
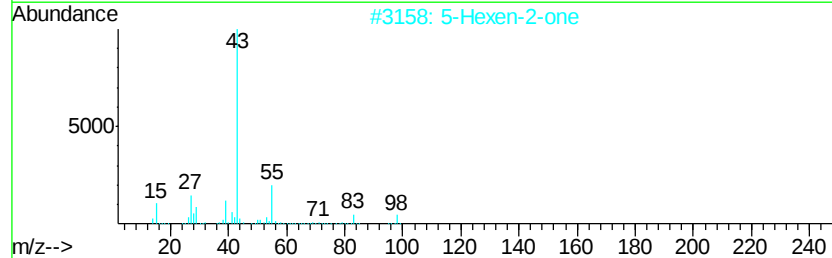
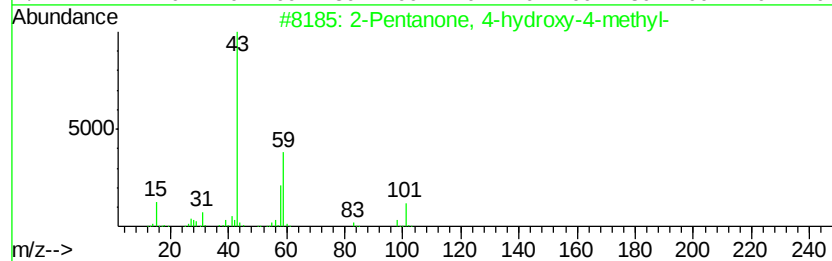
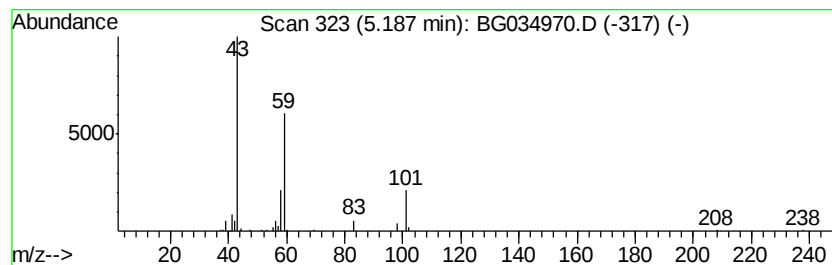
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	5.95 ng	77082	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
3	Acetone	58	C3H6O	000067-64-1	9		
4	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9		
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		



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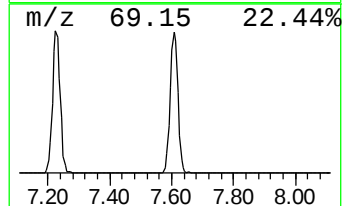
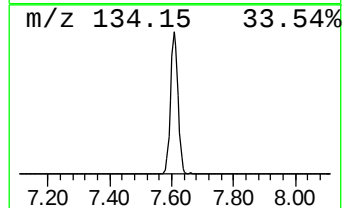
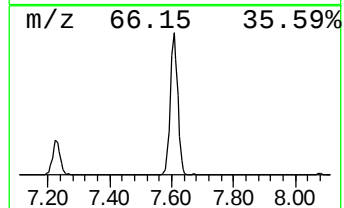
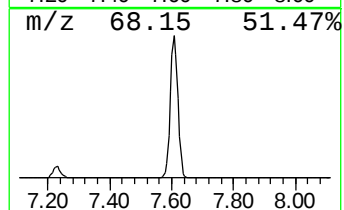
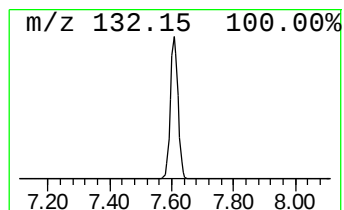
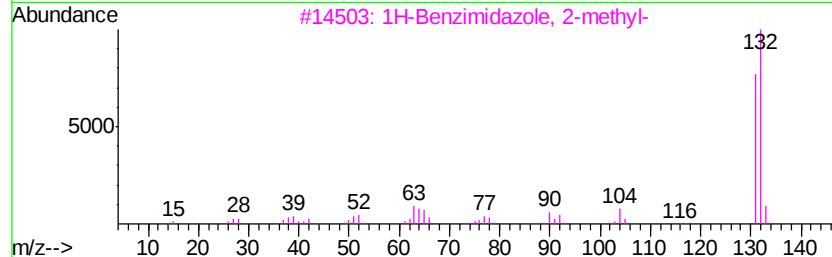
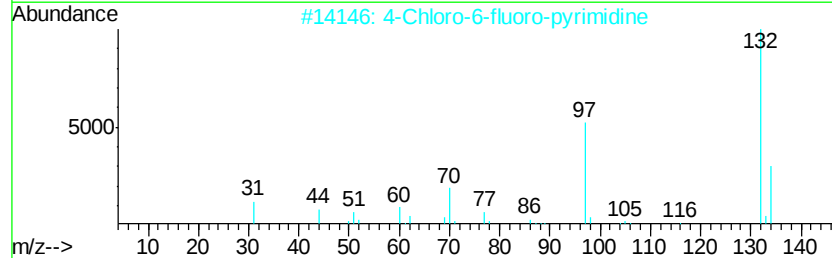
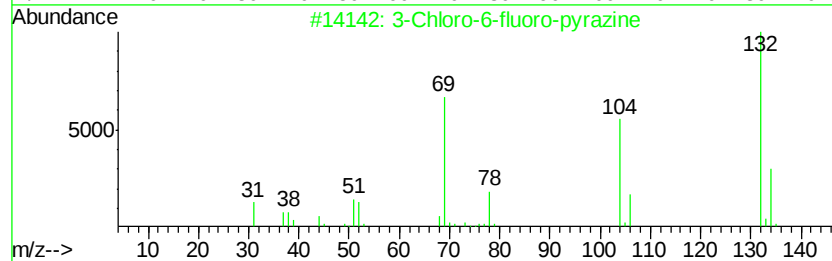
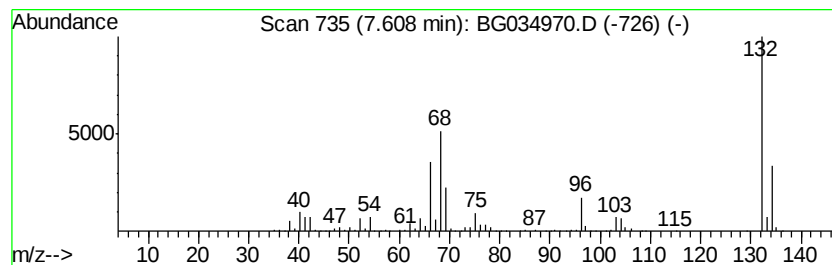
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	89.64 ng	1162080	1,4-Dichlorobenzene-d4	8.08

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		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	6.0	ng	77082	1	8.08	259286	20.0
unknown7.61	7.61	89.6	ng	1162080	1	8.08	259286	20.0