

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG091417\
 Data File : BG028761.D
 Acq On : 15 Sep 2017 9:56
 Operator : SJ/JU
 Sample : I5223-03
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 090917-SIDI-E-U-B

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:\HPCHEM1\BNA_G\METHODS\8270-BG091217.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	5.435	307	314	321	rBV	55768	87588	2.78%	0.789%
2	5.900	387	393	407	rBV	242647	410998	13.03%	3.702%
3	7.480	655	662	676	rBV	191315	363137	11.51%	3.271%
4	7.862	718	727	735	rBV	292559	532069	16.87%	4.792%
5	8.326	799	806	820	rVB2	93106	170852	5.42%	1.539%
6	8.649	850	861	871	rBV	362164	658258	20.87%	5.929%
7	9.495	997	1005	1018	rBV	290960	554280	17.58%	4.992%
8	11.146	1279	1286	1294	rBV2	116998	227966	7.23%	2.053%
9	13.555	1687	1696	1706	rBV	934036	1495150	47.41%	13.466%
10	14.384	1831	1837	1845	rBV2	19679	37842	1.20%	0.341%
11	14.936	1924	1931	1941	rBV2	222454	377479	11.97%	3.400%
12	16.428	2178	2185	2197	rBV	579431	956493	30.33%	8.615%
13	17.692	2393	2400	2408	rBV2	354399	581396	18.44%	5.236%
14	19.936	2778	2782	2786	rVB2	41165	47686	1.51%	0.429%
15	20.277	2833	2840	2845	rBV	2389739	3153687	100.00%	28.403%
16	20.982	2956	2960	2964	rBV2	33120	36468	1.16%	0.328%
17	21.587	3058	3063	3072	rBV9	14675	38899	1.23%	0.350%
18	22.022	3129	3137	3147	rBV	381297	708618	22.47%	6.382%
19	25.512	3721	3731	3748	rBV3	205595	664320	21.06%	5.983%

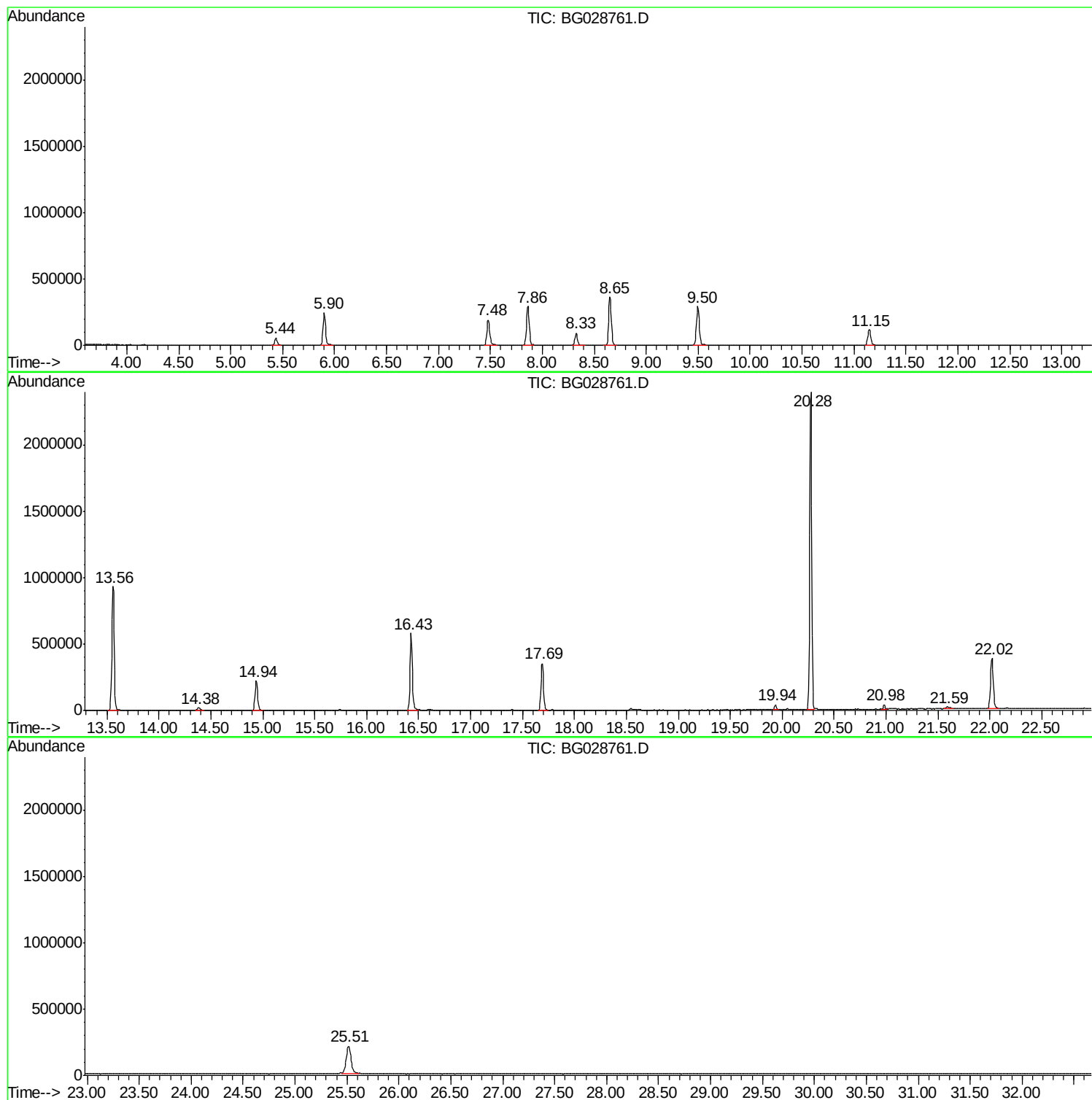
Sum of corrected areas: 11103186

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.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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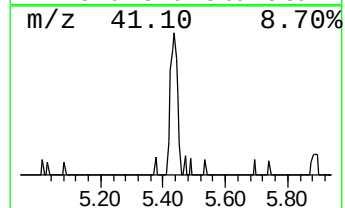
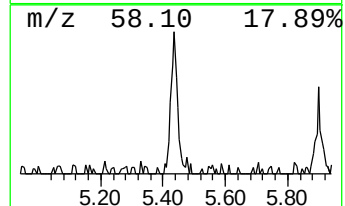
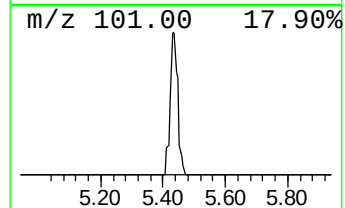
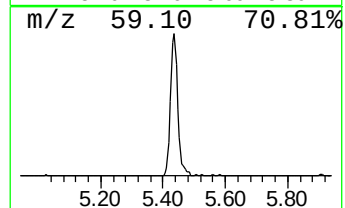
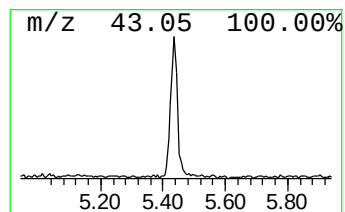
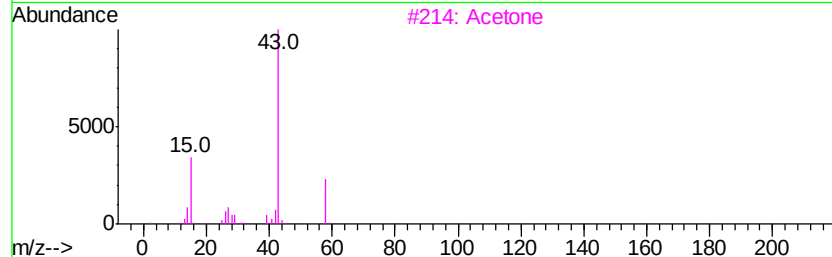
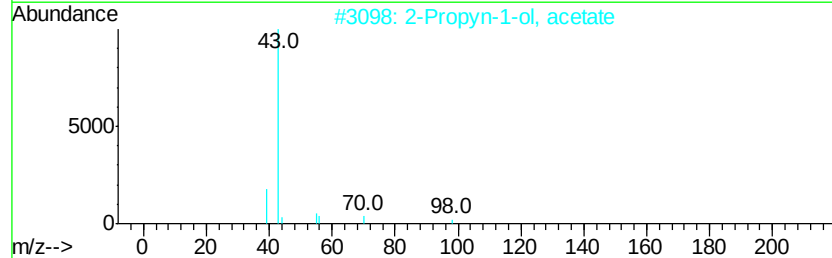
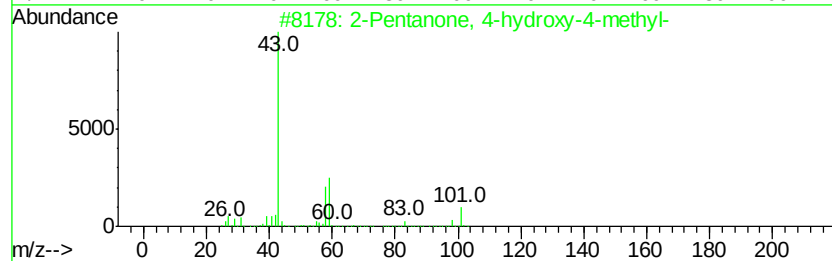
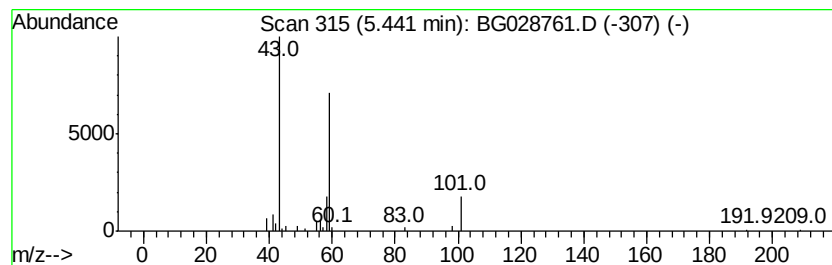
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.44	10.25 ng	87588	1,4-Dichlorobenzene-d4	8.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
3			Acetone	58	C3H6O	000067-64-1	9
4			Guanidine	59	CH5N3	000113-00-8	9
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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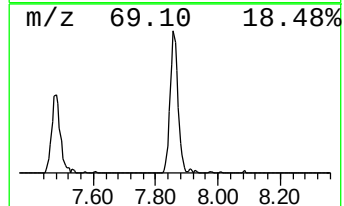
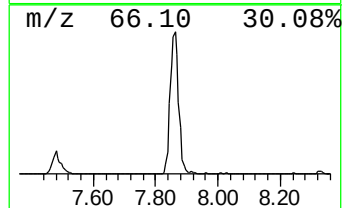
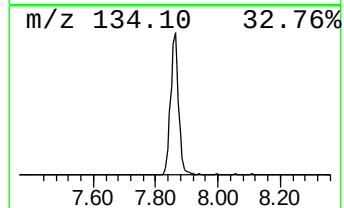
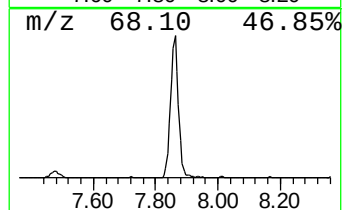
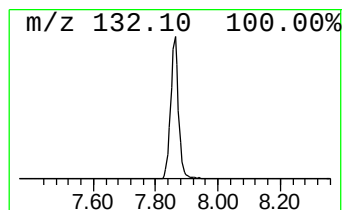
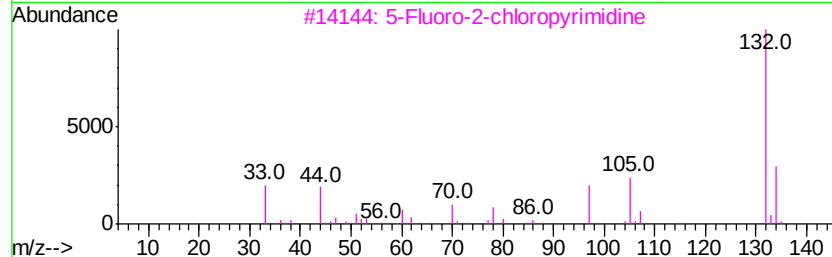
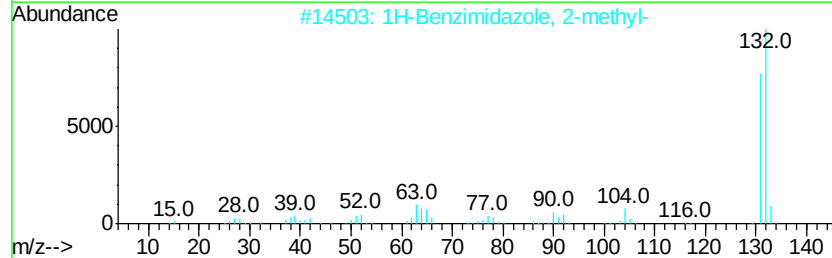
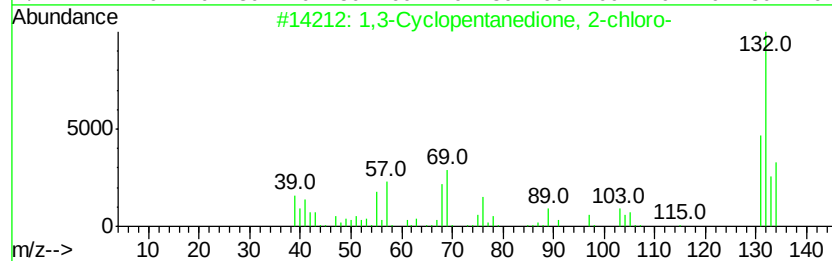
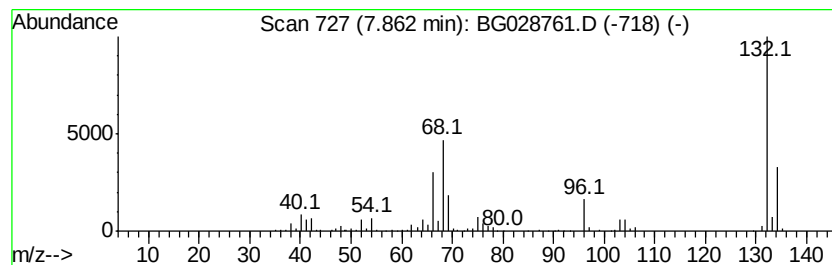
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Peak Number 2 unknown7.86 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	62.28 ng	532069	1,4-Dichlorobenzene-d4	8.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
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...	5.44	10.3	ng	87588	1	8.33	170852	20.0
unknown7.86	7.86	62.3	ng	532069	1	8.33	170852	20.0