

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092517\
 Data File : BG028940.D
 Acq On : 26 Sep 2017 12:28
 Operator : SJ/JU
 Sample : I5385-06
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 092117-SIDI-F-U-S

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	5.380	299	305	310	rBV	27054	44488	1.14%	0.386%
2	5.844	377	384	399	rBV	160968	287486	7.36%	2.496%
3	7.430	648	654	666	rBV	95629	197858	5.07%	1.718%
4	7.812	712	719	729	rBV	277054	509723	13.05%	4.426%
5	8.277	791	798	808	rBV2	83448	148403	3.80%	1.289%
6	8.600	845	853	862	rBV	369518	685914	17.56%	5.956%
7	9.446	989	997	1012	rBV	305467	606046	15.51%	5.262%
8	11.097	1272	1278	1286	rBV	101710	200331	5.13%	1.740%
9	13.512	1681	1689	1702	rBV	973345	1587775	40.65%	13.787%
10	14.892	1916	1924	1932	rBV2	192435	322078	8.25%	2.797%
11	16.385	2170	2178	2194	rBV	670703	1134819	29.05%	9.854%
12	17.642	2386	2392	2400	rBV	336264	541361	13.86%	4.701%
13	18.500	2533	2538	2543	rBV5	22175	39366	1.01%	0.342%
14	20.233	2827	2833	2848	rBV	2913517	3906210	100.00%	33.918%
15	21.966	3121	3128	3136	rBV	371689	680500	17.42%	5.909%
16	25.415	3704	3715	3728	rBV	194177	624199	15.98%	5.420%

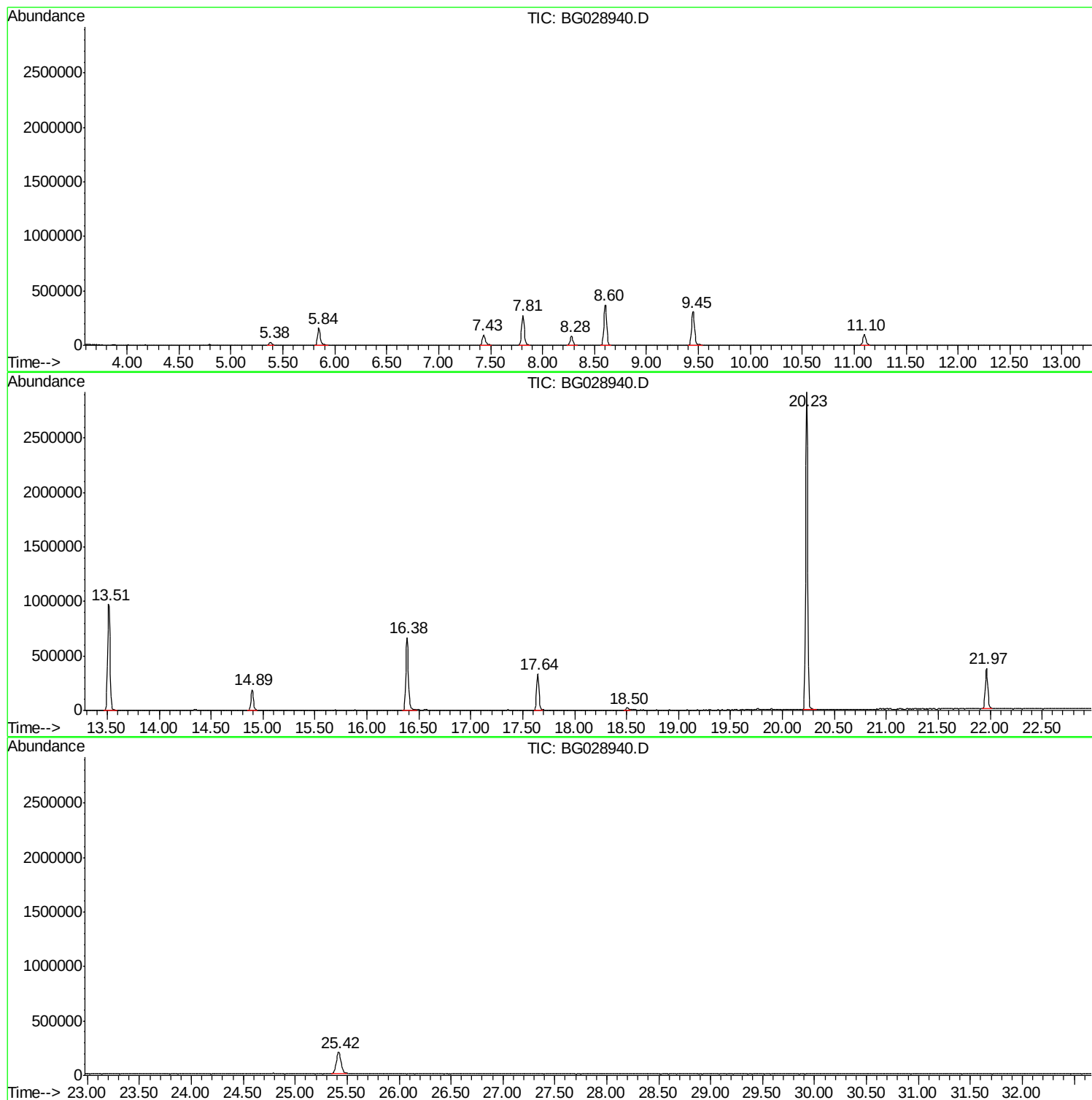
Sum of corrected areas: 11516557

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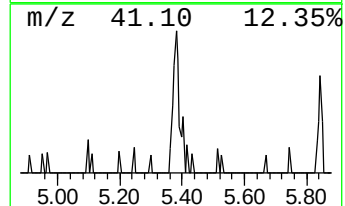
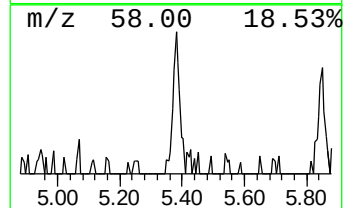
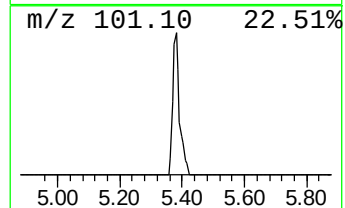
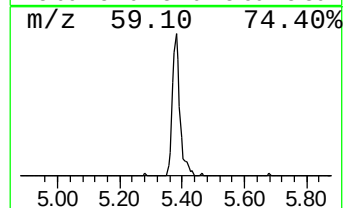
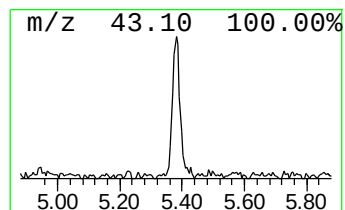
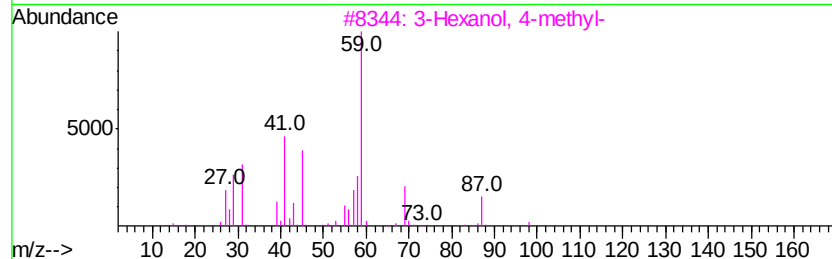
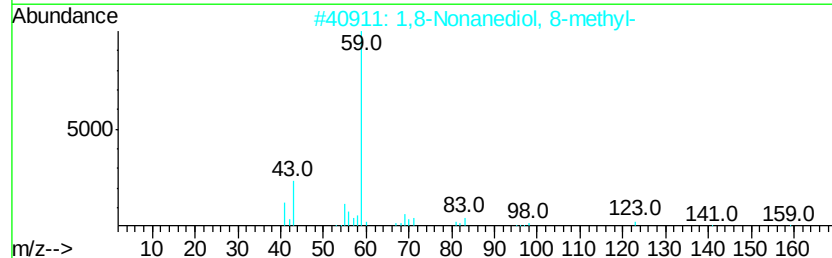
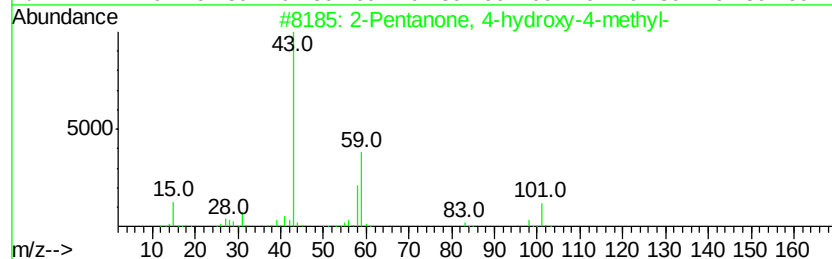
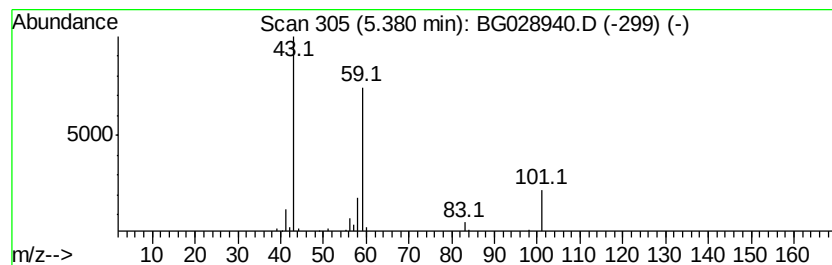
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.38	6.00 ng	44488	1,4-Dichlorobenzene-d4	8.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	40
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetone	58	C3H6O	000067-64-1	9
5			Guanidine	59	CH5N3	000113-00-8	9



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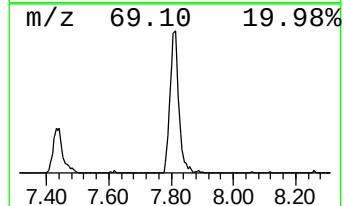
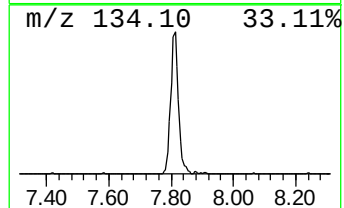
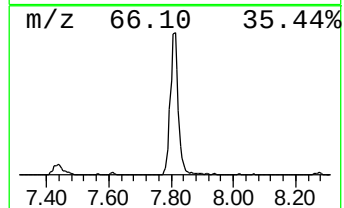
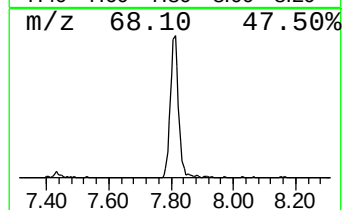
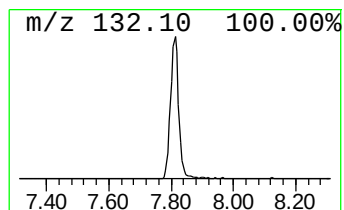
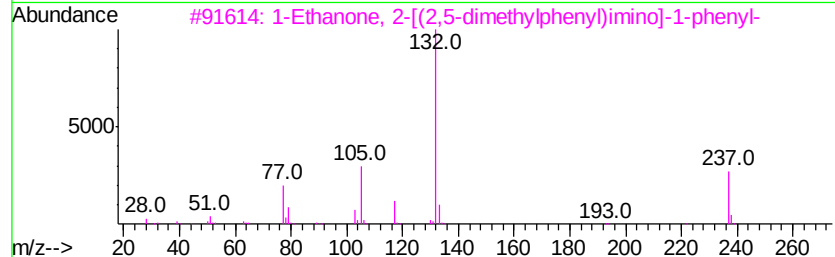
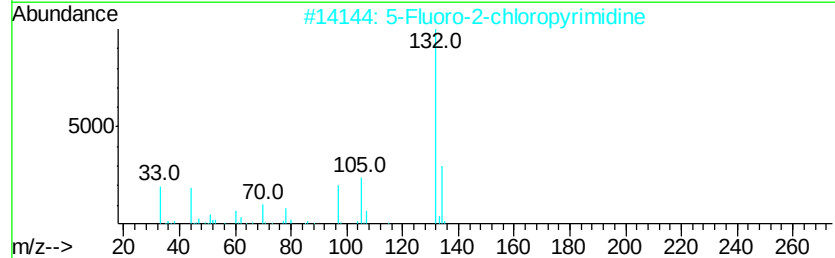
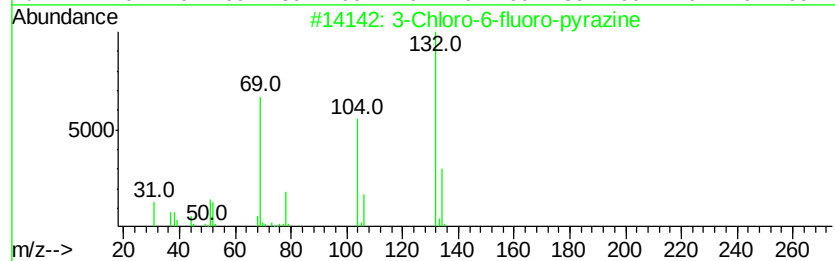
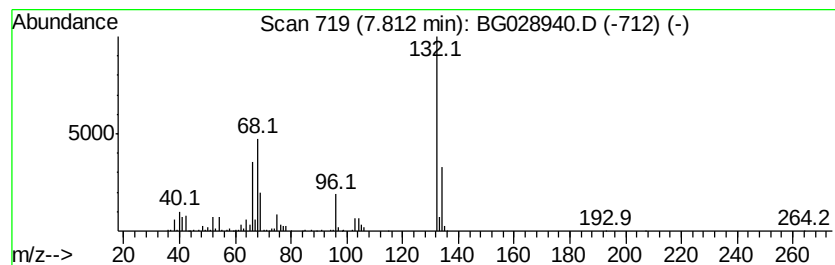
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Peak Number 2 unknown7.81 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	68.69 ng	509723	1,4-Dichlorobenzene-d4	8.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27		
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12		
3	1-Ethanone, 2-[(2,5-dimethylphen...	237	C16H15NO	1000319-31-2	12		
4	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12		
5	5-Aminoindole	132	C8H8N2	005192-03-0	10		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.38	6.0	ng	44488	1	8.28	148403	20.0
unknown7.81	7.81	68.7	ng	509723	1	8.28	148403	20.0