

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092017\
 Data File : BG028836.D
 Acq On : 20 Sep 2017 17:54
 Operator : SJ/JU
 Sample : PB102470BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB102470BL

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	5.392	298	307	316	rBV2	93445	151190	4.47%	0.978%
2	5.856	380	386	401	rBV	625995	1073107	31.75%	6.944%
3	7.442	649	656	672	rBV	618592	1113877	32.95%	7.207%
4	7.824	713	721	735	rVB	613061	1080356	31.96%	6.990%
5	8.288	793	800	808	rBV	130265	231684	6.85%	1.499%
6	8.611	848	855	866	rBV	442415	800817	23.69%	5.182%
7	9.458	991	999	1010	rBV	360271	672102	19.88%	4.349%
8	11.109	1272	1280	1292	rBV	177534	333850	9.88%	2.160%
9	13.529	1684	1692	1701	rBV	1102404	1780212	52.66%	11.519%
10	14.910	1919	1927	1939	rBV3	314386	532661	15.76%	3.447%
11	16.396	2173	2180	2194	rBV	971066	1629323	48.20%	10.543%
12	17.660	2389	2395	2405	rBV2	497871	782870	23.16%	5.066%
13	20.245	2829	2835	2844	rBV	2486254	3380289	100.00%	21.872%
14	21.984	3122	3131	3145	rBV	514044	952034	28.16%	6.160%
15	25.451	3702	3721	3741	rBV2	260450	940362	27.82%	6.085%

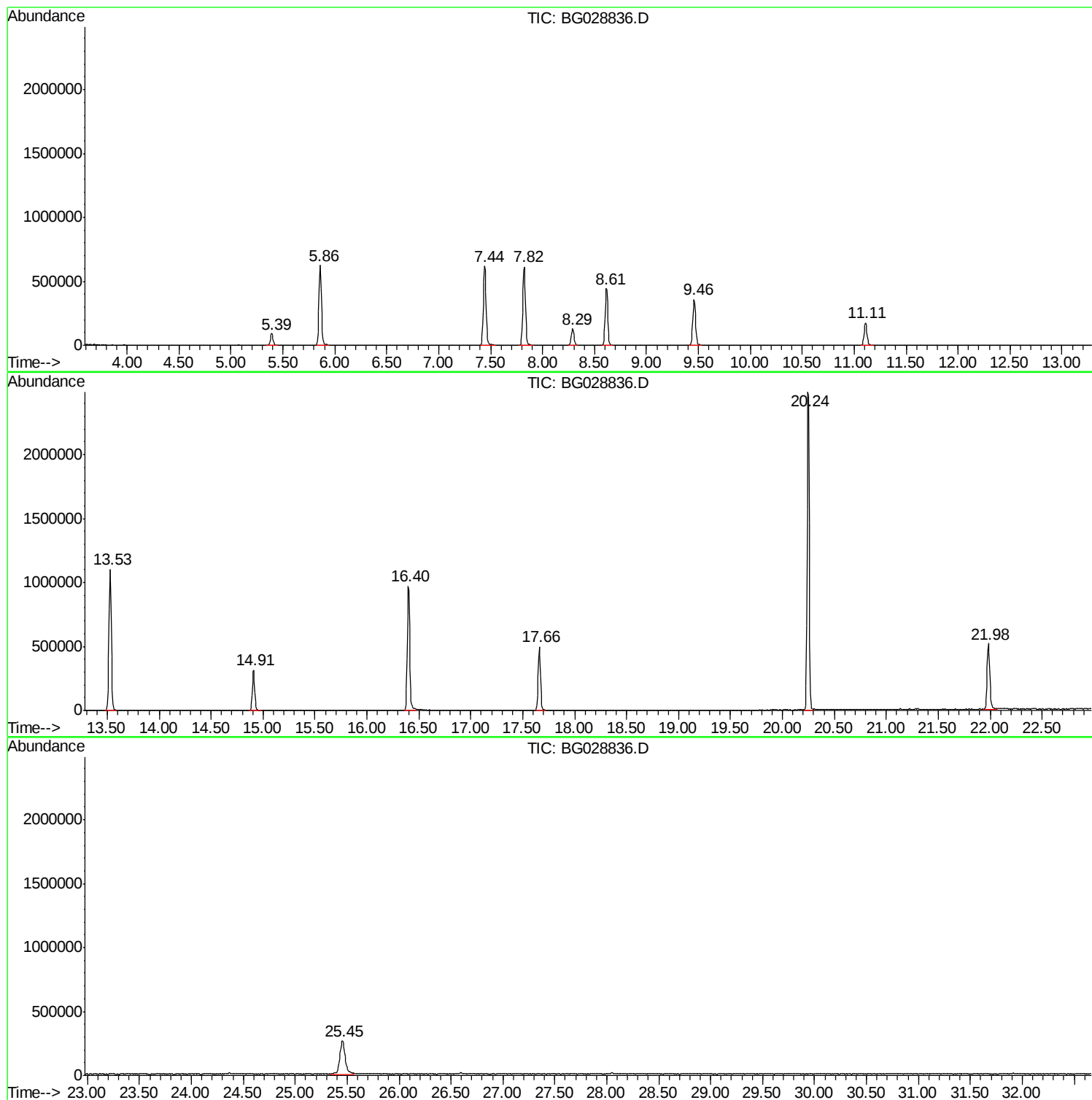
Sum of corrected areas: 15454734

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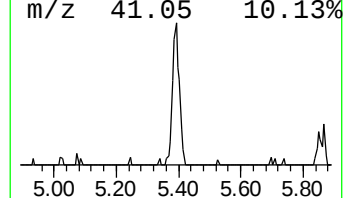
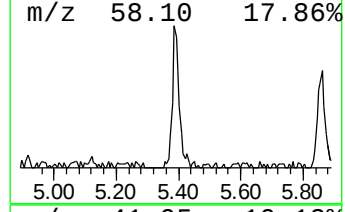
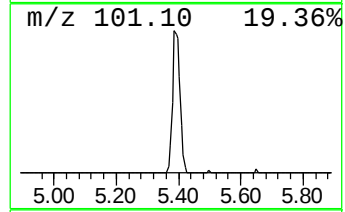
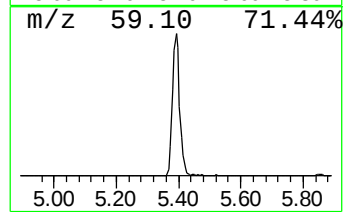
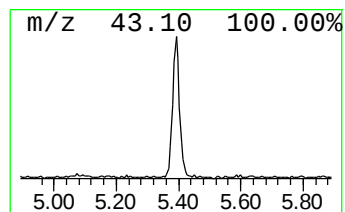
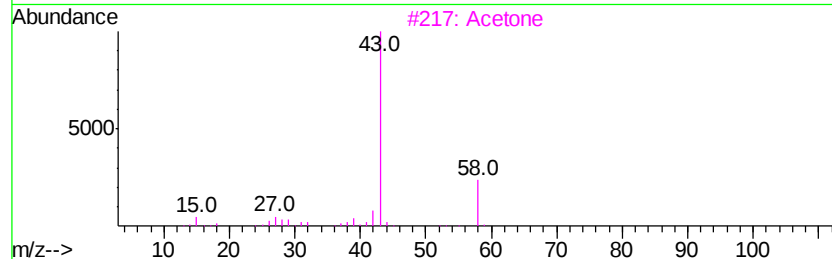
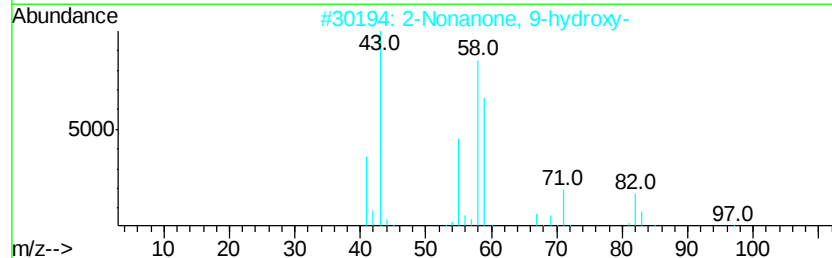
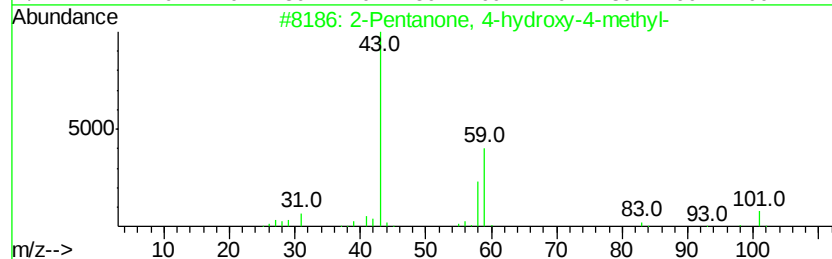
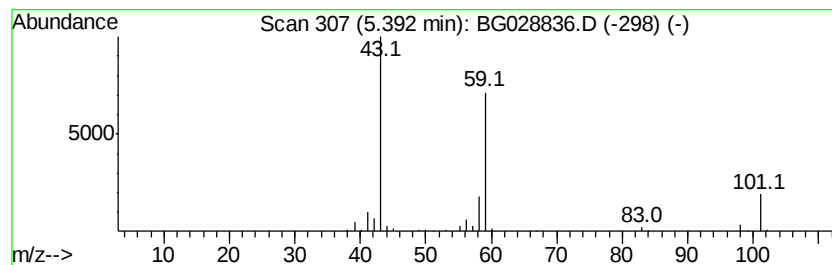
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.39	13.05 ng	151190	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	28
3		Acetone	58	C3H6O	000067-64-1	14
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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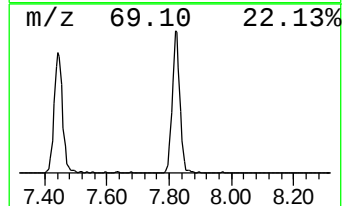
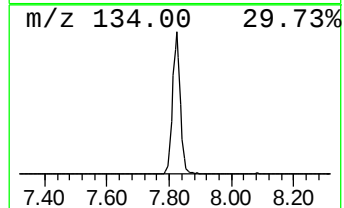
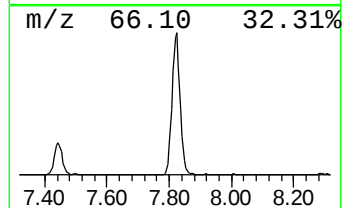
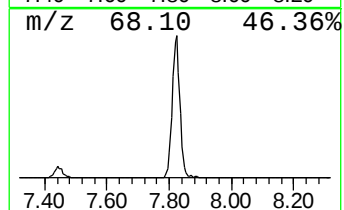
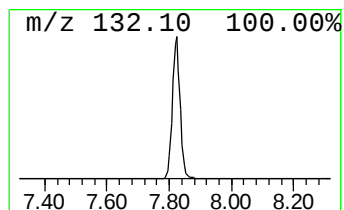
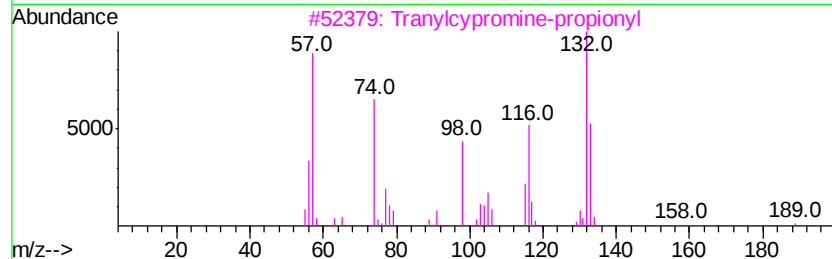
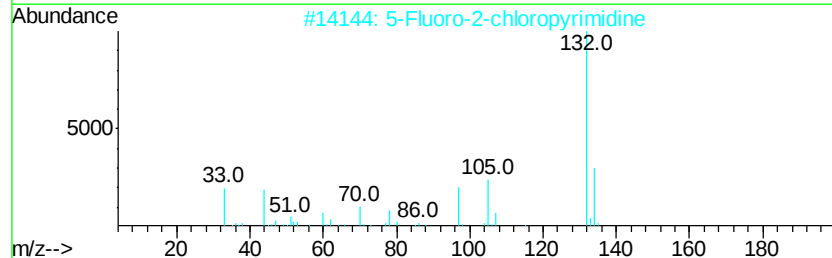
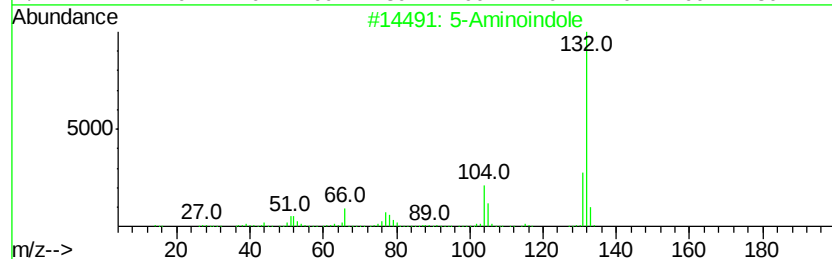
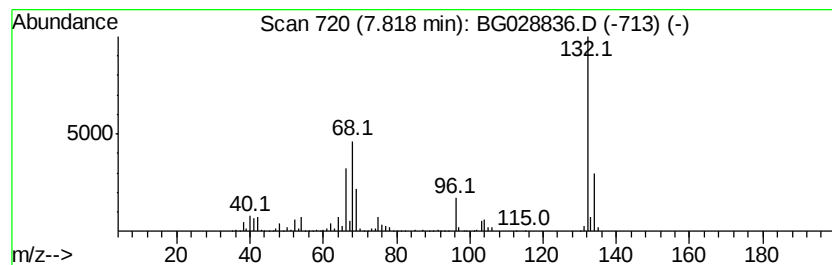
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	93.26 ng	1080360	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	12
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.39	13.1	ng	151190	1	8.29	231684	20.0
unknown7.82	7.82	93.3	ng	1080360	1	8.29	231684	20.0