

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
 Data File : BG036426.D
 Acq On : 27 Aug 2018 11:31
 Operator : JU/SJ
 Sample : PB112371BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112371BL

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	5.168	315	320	327	rBV	34739	53398	1.23%	0.260%
2	5.632	392	399	410	rBV	894571	1397733	32.08%	6.799%
3	7.219	661	669	679	rBV	948980	1588135	36.45%	7.725%
4	7.595	725	733	741	rBV	849719	1435576	32.95%	6.983%
5	8.065	806	813	824	rVB	197600	338582	7.77%	1.647%
6	8.388	860	868	876	rBV	636723	1071208	24.59%	5.211%
7	9.222	1002	1010	1020	rBV	510876	907278	20.83%	4.413%
8	10.873	1283	1291	1298	rBV	282124	504062	11.57%	2.452%
9	13.317	1696	1707	1715	rBV	1534710	2413368	55.40%	11.739%
10	14.686	1932	1940	1950	rBV2	481890	790691	18.15%	3.846%
11	16.173	2186	2193	2207	rBV	1235778	1839061	42.21%	8.946%
12	17.430	2401	2407	2418	rVB	734662	1113222	25.55%	5.415%
13	20.039	2845	2851	2858	rBV	3229790	4356642	100.00%	21.192%
14	21.696	3126	3133	3141	rBV	822318	1340620	30.77%	6.521%
15	24.916	3667	3681	3691	rBV3	447654	1291880	29.65%	6.284%
16	24.998	3691	3695	3702	rVB6	26227	55218	1.27%	0.269%
17	26.143	3884	3890	3900	rVB6	20234	61096	1.40%	0.297%

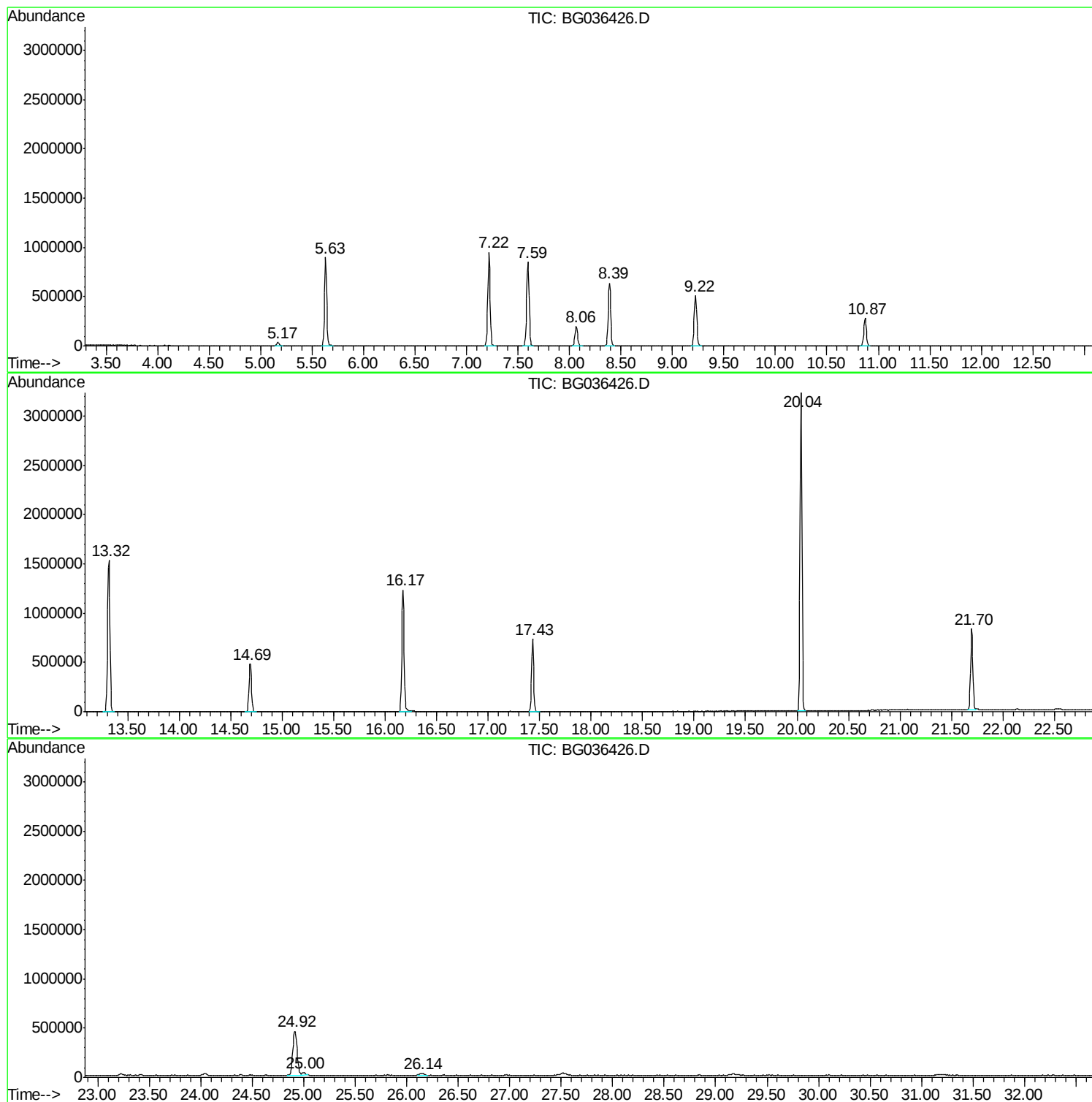
Sum of corrected areas: 20557770

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.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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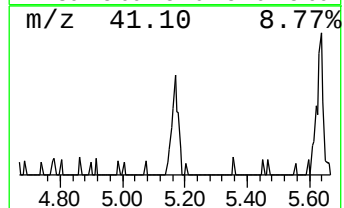
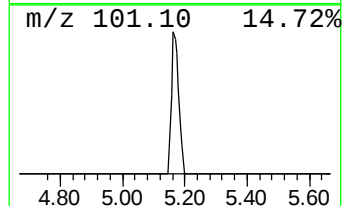
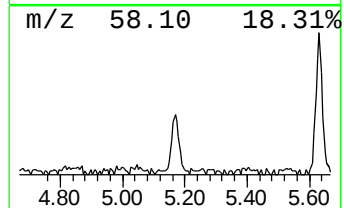
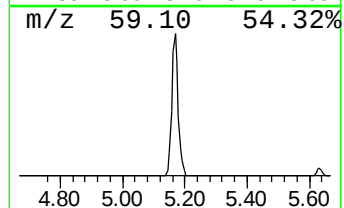
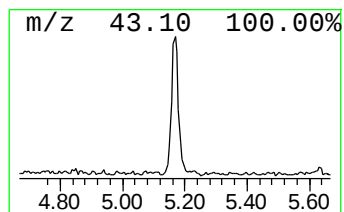
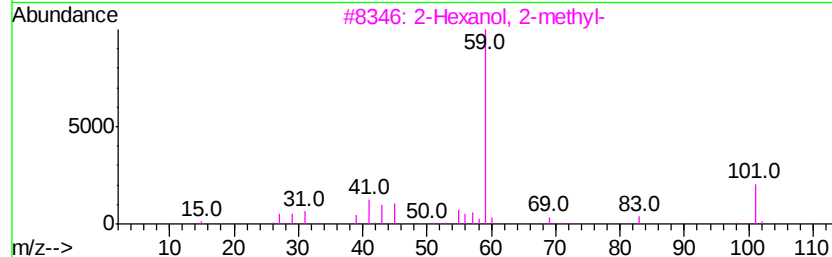
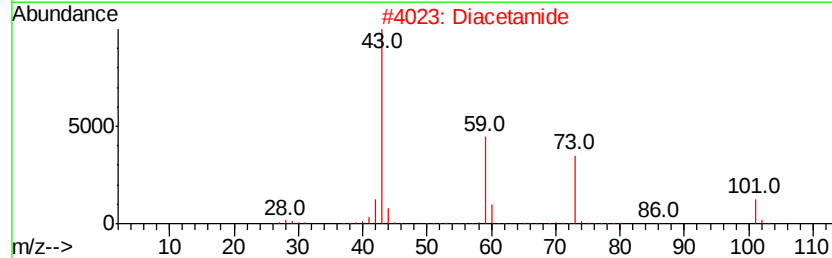
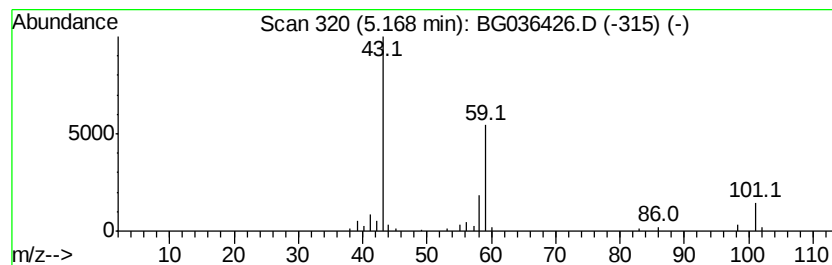
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.15 ng	53398	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Diacetamide	101	C4H7NO2	000625-77-4	38
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5			Acetamide	59	C2H5NO	000060-35-5	9



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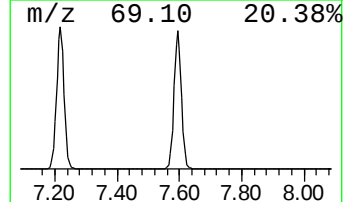
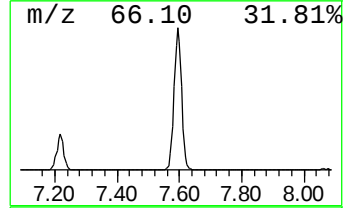
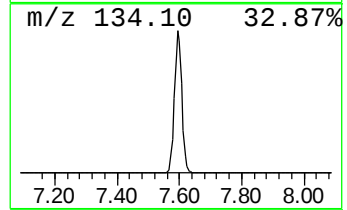
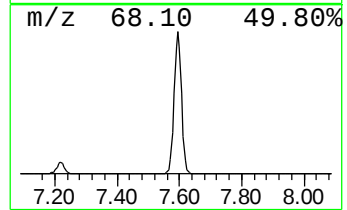
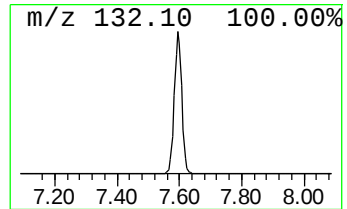
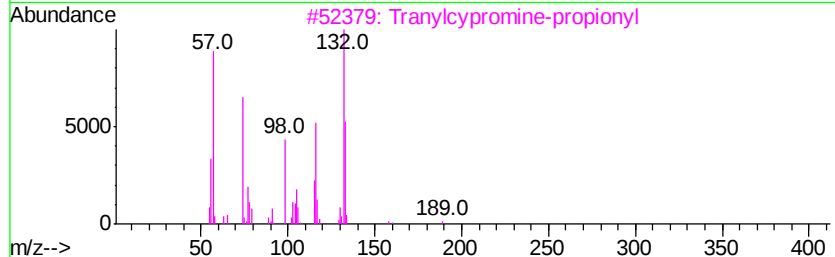
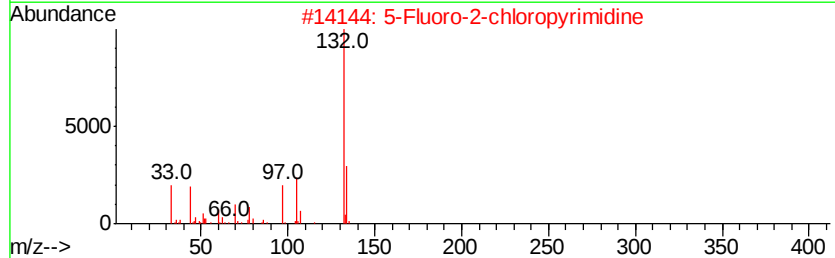
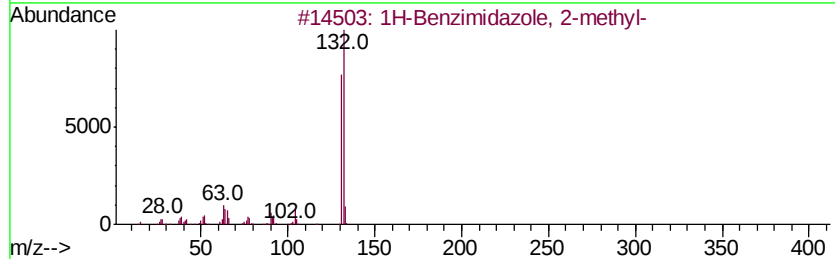
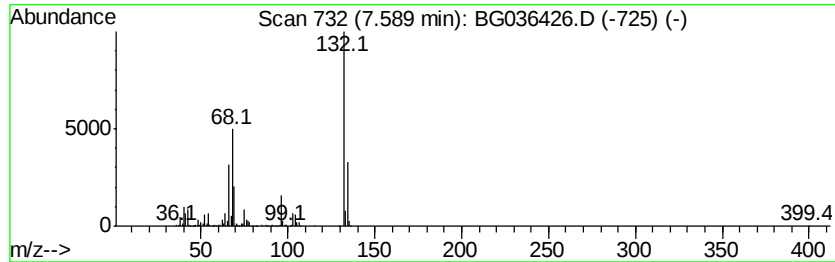
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	84.80 ng	1435580	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	3.1	ng	53398	1	8.06	338582	20.0
unknown7.59	7.59	84.8	ng	1435580	1	8.06	338582	20.0