

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090418\
 Data File : BG036559.D
 Acq On : 5 Sep 2018 10:06
 Operator : JU/SJ
 Sample : J4761-06
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RBR-200031

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	5.165	309	319	328	rBV	124400	195360	3.29%	0.706%
2	5.630	390	398	409	rBV	1306009	2112742	35.63%	7.640%
3	7.216	659	668	682	rBV	1218643	2116366	35.69%	7.653%
4	7.592	723	732	741	rBV	1345235	2327089	39.25%	8.415%
5	8.056	803	811	819	rVB	220411	374829	6.32%	1.355%
6	8.379	858	866	875	rVB	1025567	1779442	30.01%	6.435%
7	9.220	1001	1009	1022	rVB	880705	1573140	26.53%	5.689%
8	10.865	1281	1289	1296	rBV	302762	546886	9.22%	1.978%
9	13.309	1697	1705	1717	rBV	2311672	3627757	61.18%	13.119%
10	14.678	1932	1938	1948	rBV2	470998	783039	13.21%	2.832%
11	16.164	2184	2191	2213	rBV	1753097	2774517	46.79%	10.034%
12	17.422	2399	2405	2417	rBV2	717059	1073876	18.11%	3.883%
13	20.030	2842	2849	2863	rBV	4466432	5929577	100.00%	21.443%
14	21.687	3124	3131	3143	rBV2	722590	1223237	20.63%	4.424%
15	24.895	3666	3677	3695	rBV2	333036	1150776	19.41%	4.162%
16	26.082	3871	3879	3890	rVB5	21376	63827	1.08%	0.231%

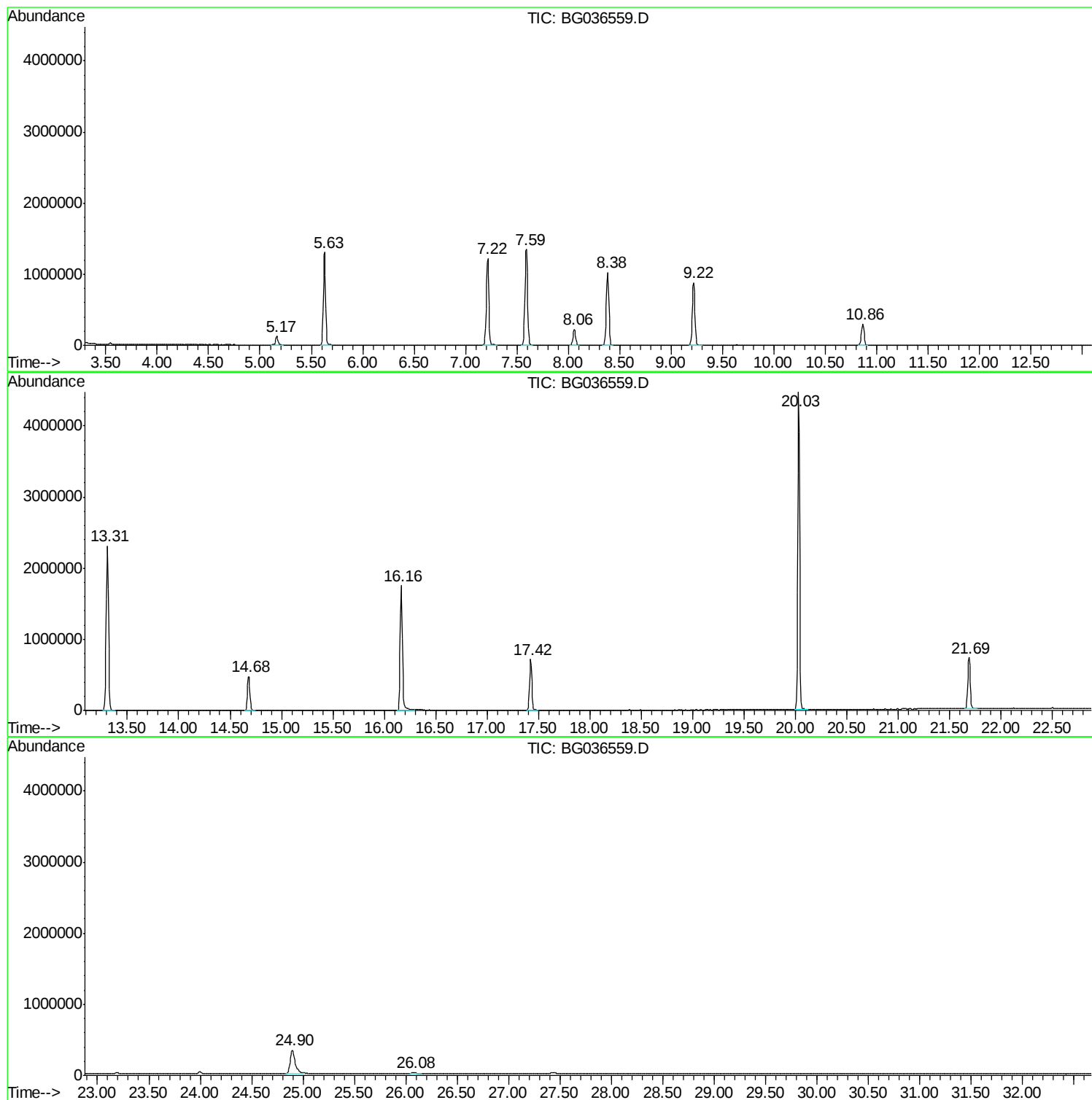
Sum of corrected areas: 27652460

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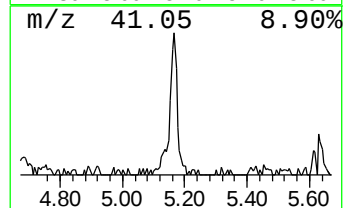
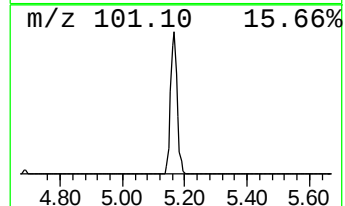
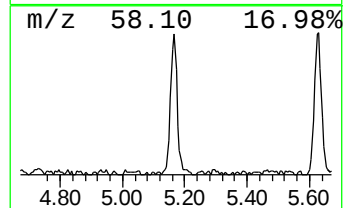
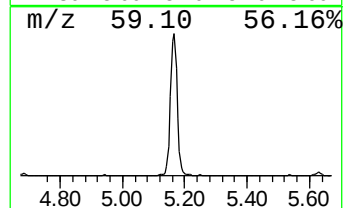
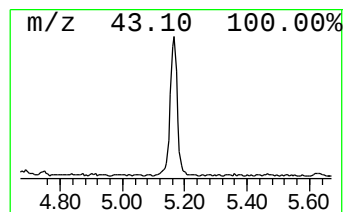
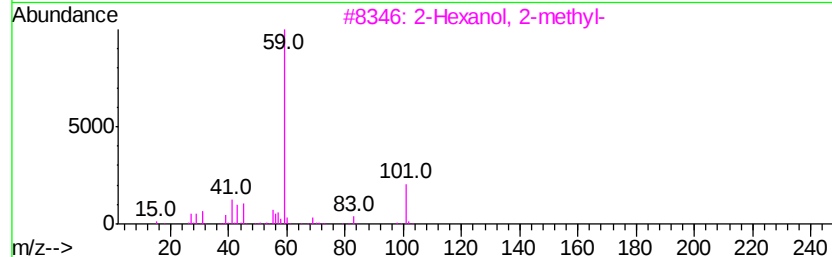
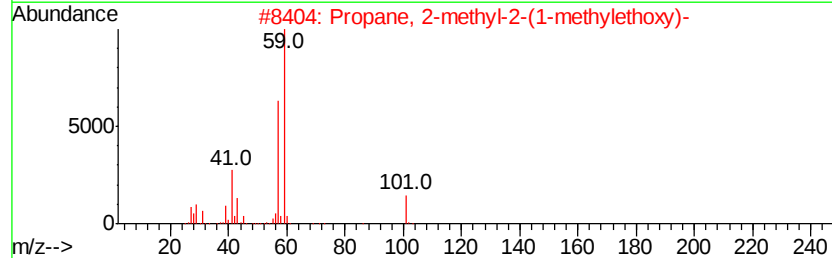
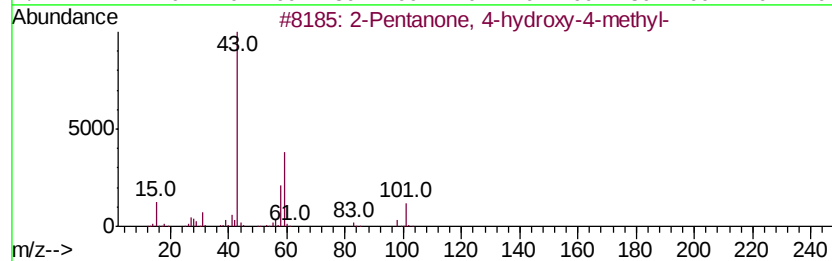
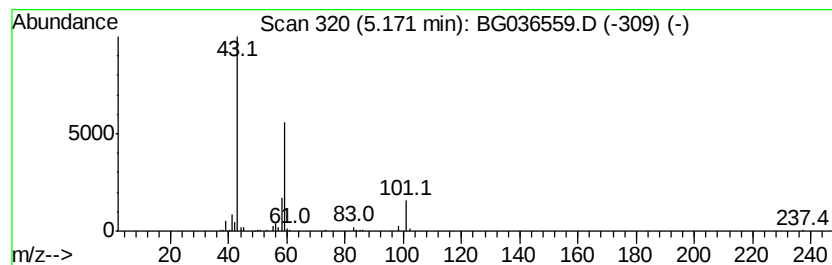
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	10.42 ng	195360	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	50
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Acetone	58	C3H6O	000067-64-1	9



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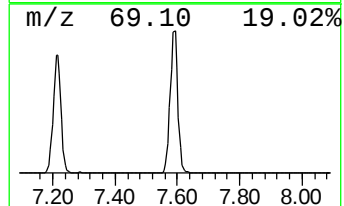
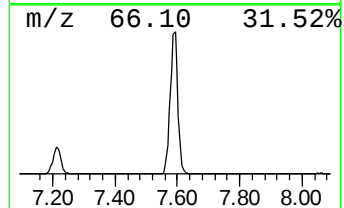
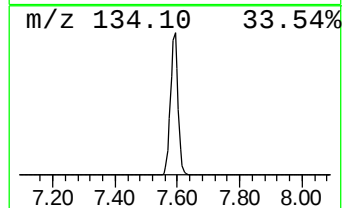
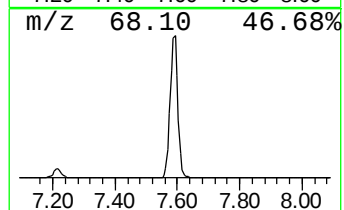
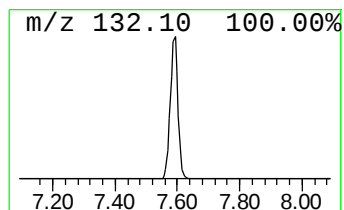
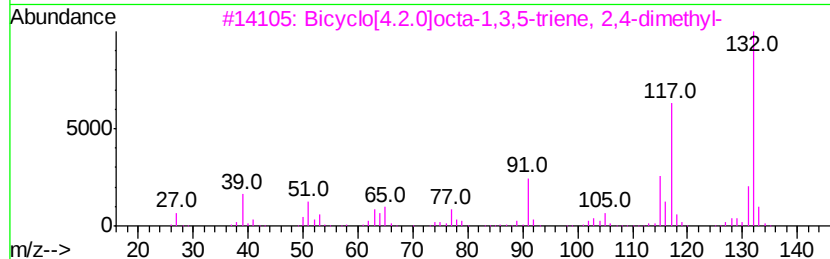
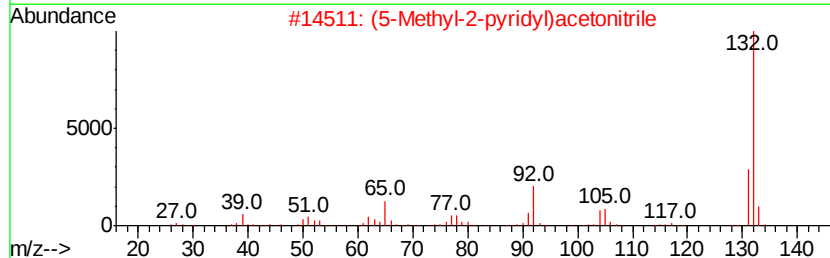
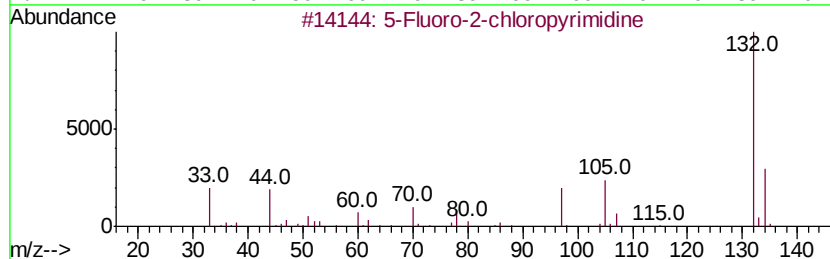
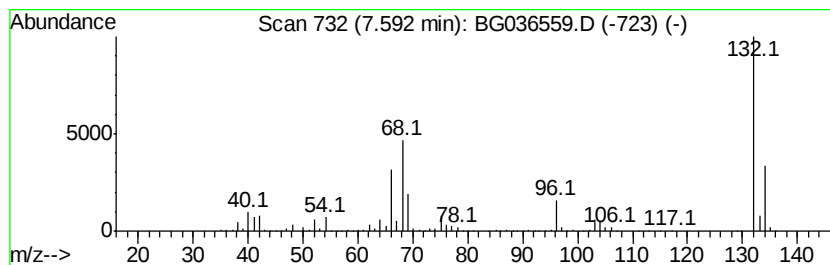
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	124.17 ng	2327090	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	10.4	ng	195360	1	8.06	374829	20.0
unknown7.59	7.59	124.2	ng	2327090	1	8.06	374829	20.0