

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
 Data File : BG036429.D
 Acq On : 27 Aug 2018 13:28
 Operator : JU/SJ
 Sample : J4612-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RBR200027

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	310	320	329	rBV	89473	147794	3.65%	0.789%
2	5.632	392	399	408	rBV	765813	1200196	29.63%	6.410%
3	7.218	661	669	679	rBV	698279	1157095	28.57%	6.180%
4	7.594	724	733	745	rBV	834755	1395677	34.46%	7.454%
5	8.064	806	813	821	rVB	189975	319007	7.88%	1.704%
6	8.387	861	868	876	rBV	651093	1106048	27.31%	5.907%
7	9.222	1002	1010	1018	rBV	543769	963214	23.78%	5.145%
8	10.873	1283	1291	1296	rBV	259137	467874	11.55%	2.499%
9	10.920	1296	1299	1305	rVB2	27376	44322	1.09%	0.237%
10	13.317	1699	1707	1716	rBV	1546632	2347268	57.96%	12.537%
11	14.686	1933	1940	1948	rBV2	422187	691412	17.07%	3.693%
12	16.172	2186	2193	2207	rBV	1113557	1669093	41.21%	8.915%
13	17.430	2401	2407	2419	rVB2	630628	955013	23.58%	5.101%
14	20.038	2845	2851	2861	rBV	3057363	4049973	100.00%	21.631%
15	21.695	3126	3133	3142	rVB2	679946	1129700	27.89%	6.034%
16	24.915	3669	3681	3693	rBV	377630	1079109	26.64%	5.764%

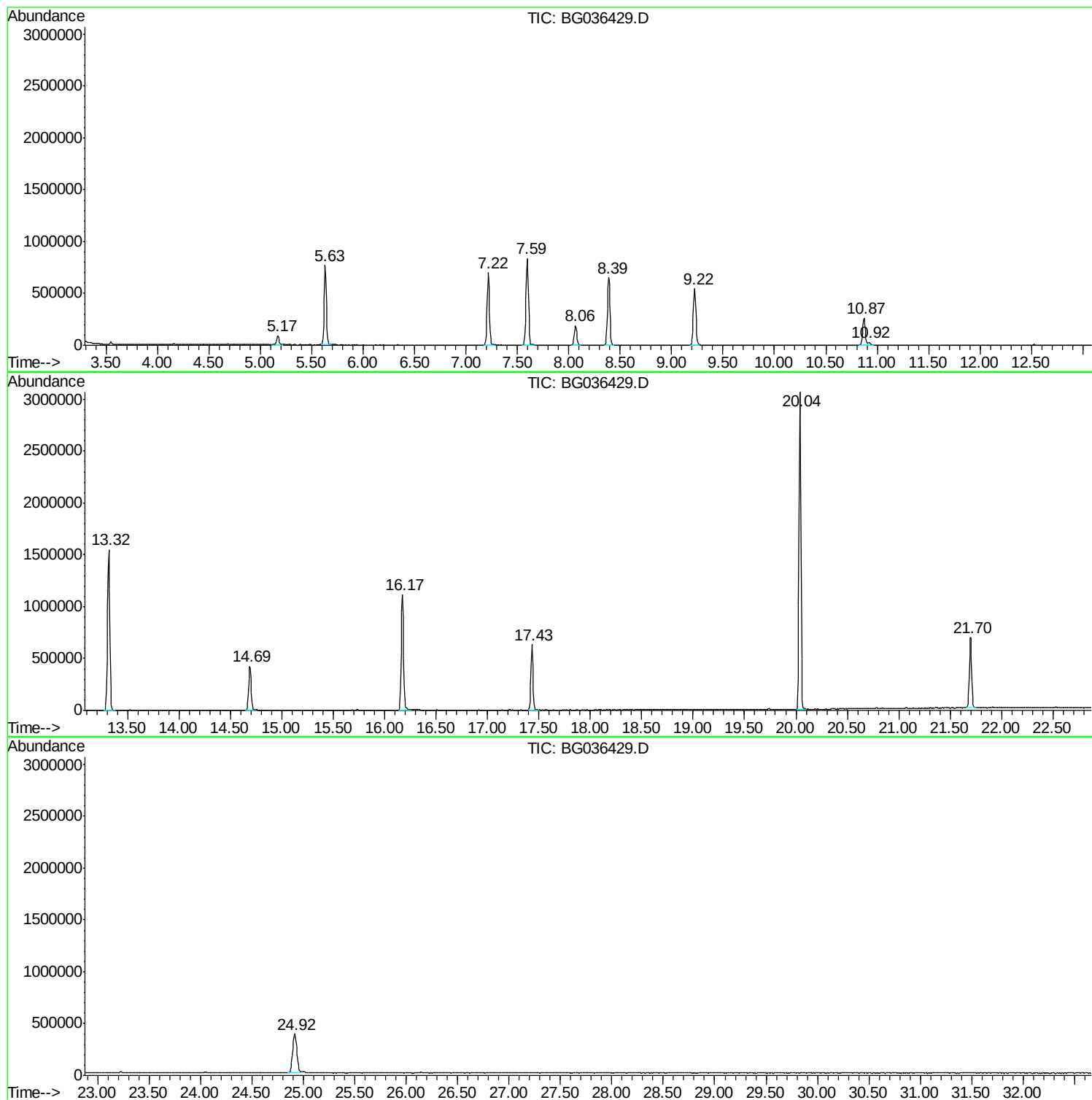
Sum of corrected areas: 18722795

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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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Instrument :
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ClientSampled :
RBR200027

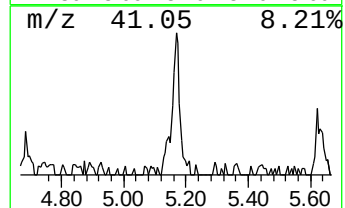
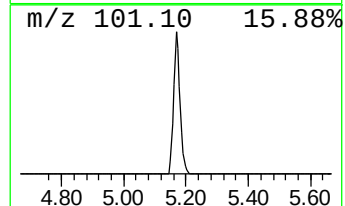
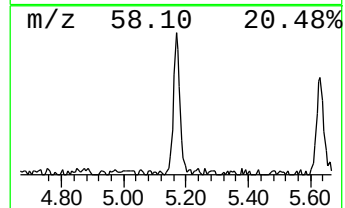
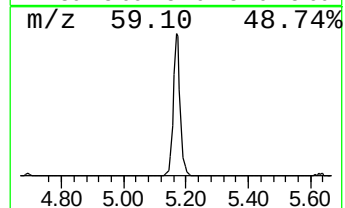
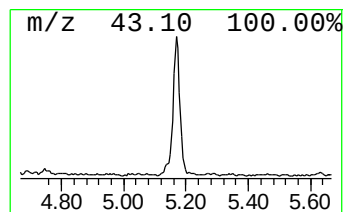
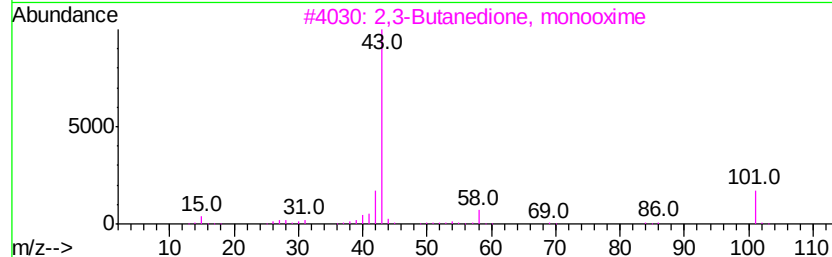
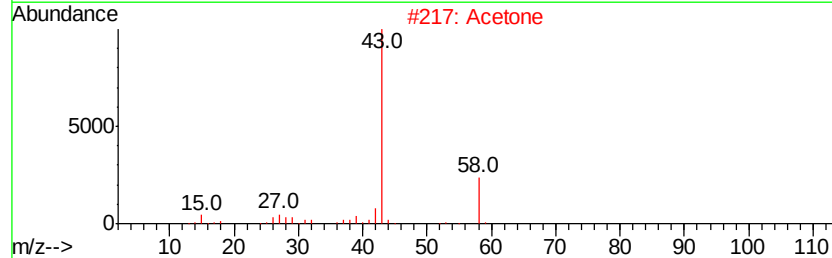
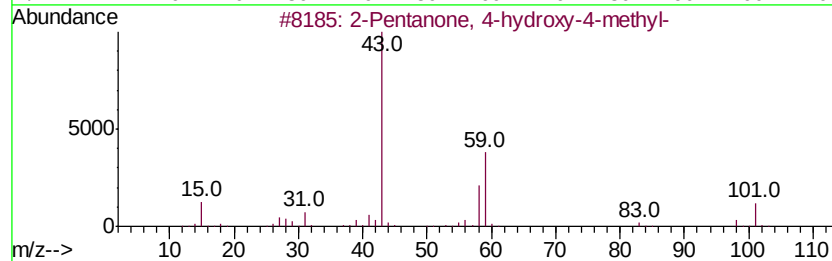
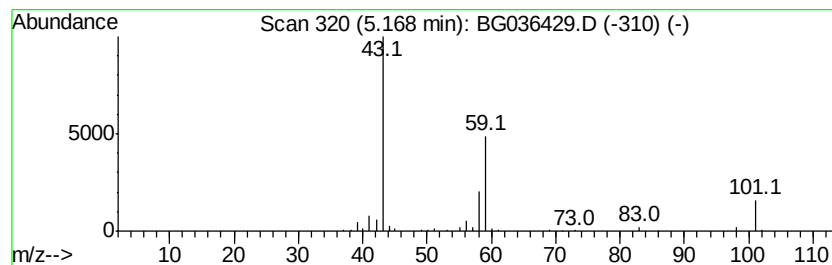
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	9.27 ng	147794	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2		Acetone	58	C3H6O	000067-64-1	27
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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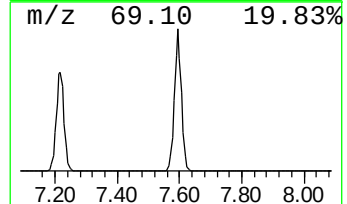
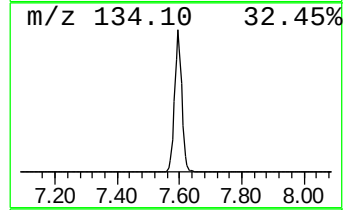
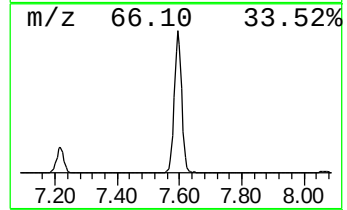
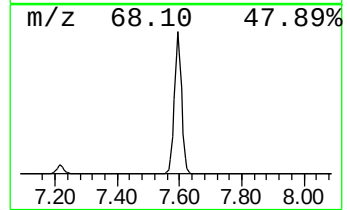
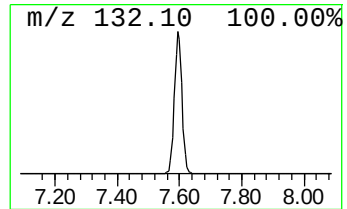
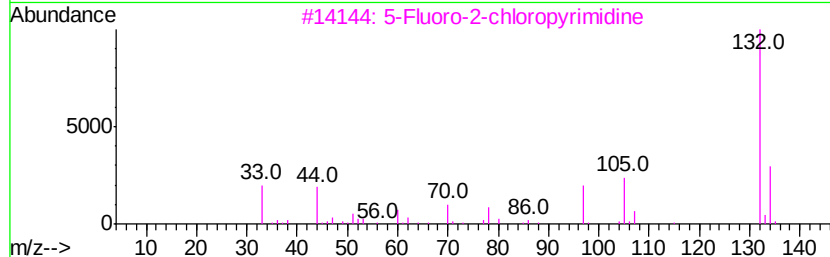
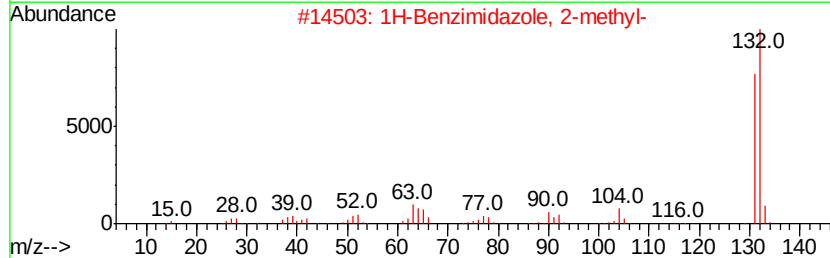
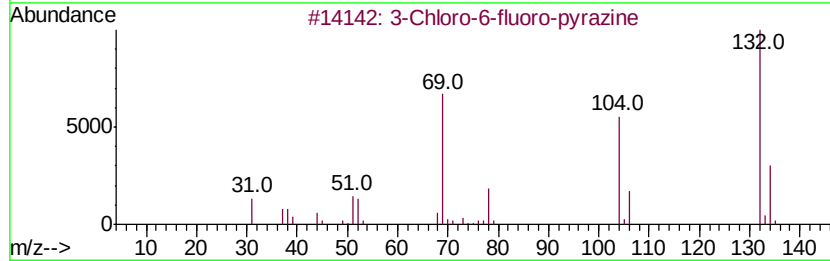
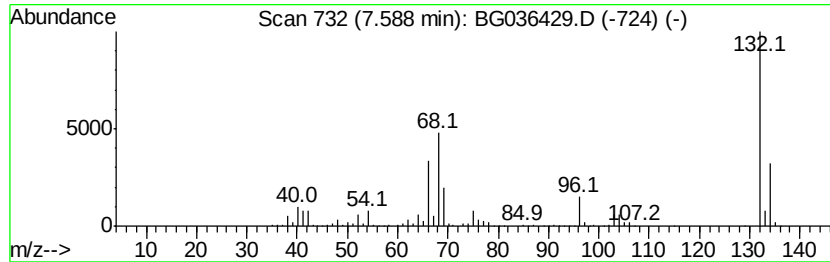
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	87.50 ng	1395680	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	1H-Benzimidazole, 2-methyl-		132	C8H8N2	000615-15-6	22
3	5-Fluoro-2-chloropyrimidine		132	C4H2ClFN2	062802-42-0	12
4	Cyclopropanamine, 1-phenyl-		133	C9H11N	1000351-26-2	10
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...	5.17	9.3	ng	147794	1	8.06	319007	20.0
unknown7.59	7.59	87.5	ng	1395680	1	8.06	319007	20.0