

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036632.D
 Acq On : 10 Sep 2018 19:41
 Operator : JU/SJ
 Sample : J4769-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HR-05-090718-C

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	5.158	309	318	328	rBV	119555	190710	4.74%	0.989%
2	5.623	390	397	411	rBV	904249	1420997	35.32%	7.372%
3	7.209	660	667	684	rBV	859030	1493600	37.12%	7.749%
4	7.585	721	731	745	rBV	906514	1541165	38.30%	7.996%
5	8.055	804	811	819	rVB	186068	313109	7.78%	1.624%
6	8.378	858	866	874	rVB	676675	1154360	28.69%	5.989%
7	9.213	1001	1008	1019	rVB	578631	1031380	25.63%	5.351%
8	10.858	1280	1288	1294	rBV	254274	471650	11.72%	2.447%
9	10.911	1294	1297	1306	rVB2	43198	68659	1.71%	0.356%
10	13.302	1696	1704	1713	rBV	1643429	2492547	61.95%	12.932%
11	14.677	1931	1938	1946	rBV2	429123	692468	17.21%	3.593%
12	16.163	2184	2191	2215	rBV	1030595	1704727	42.37%	8.844%
13	17.421	2397	2405	2416	rBV2	595360	906938	22.54%	4.705%
14	20.029	2842	2849	2858	rBV	2989919	4023465	100.00%	20.874%
15	21.686	3122	3131	3150	rBV	561517	1031654	25.64%	5.352%
16	24.900	3668	3678	3694	rBV2	202954	737149	18.32%	3.824%

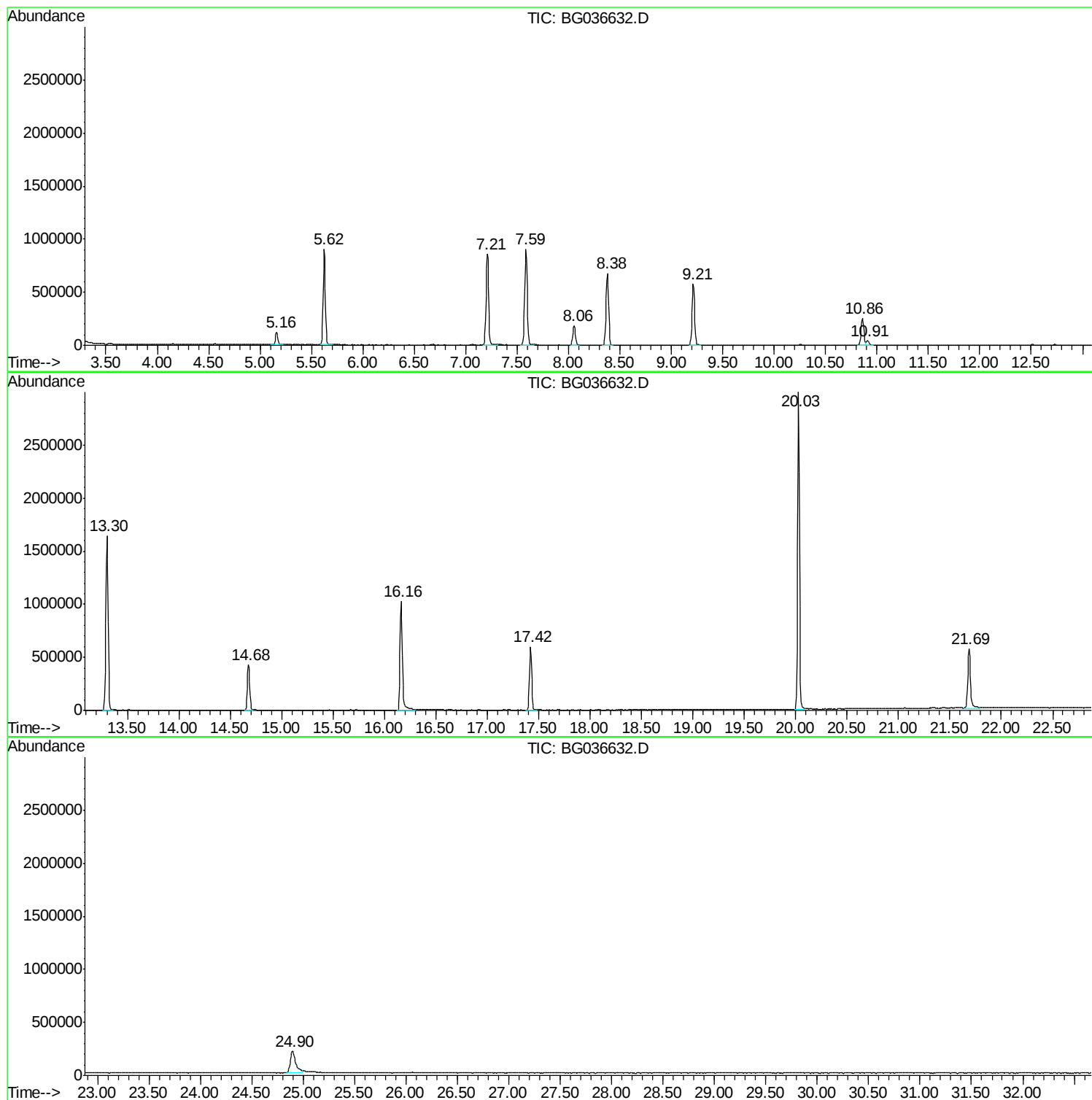
Sum of corrected areas: 19274578

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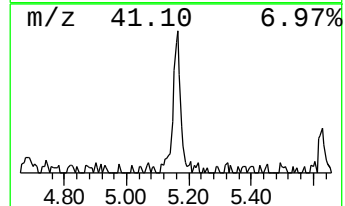
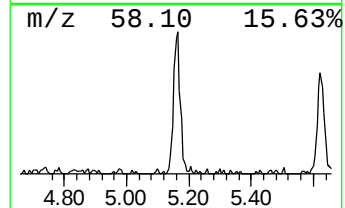
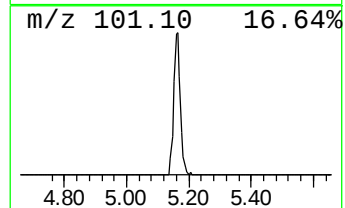
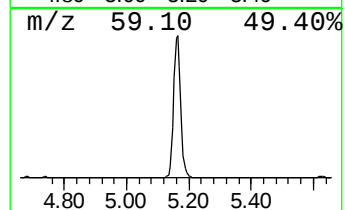
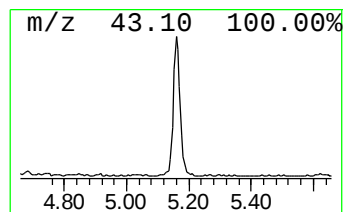
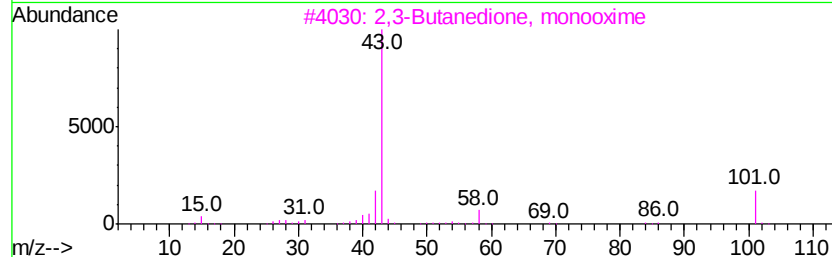
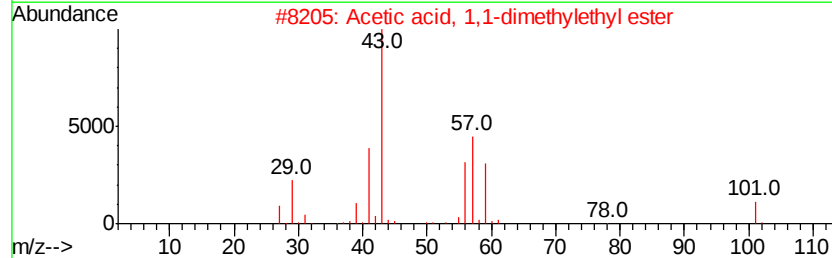
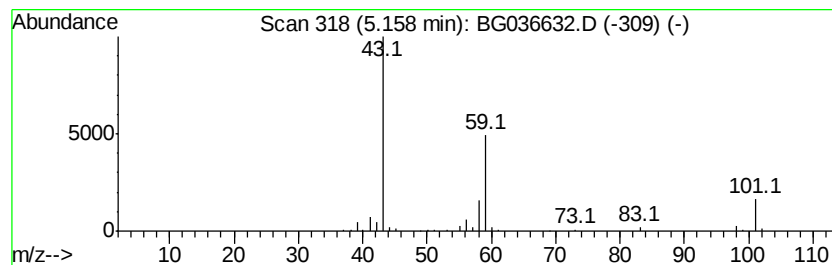
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	12.18 ng	190710	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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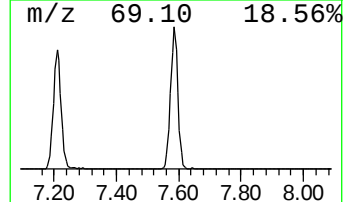
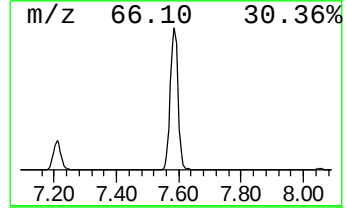
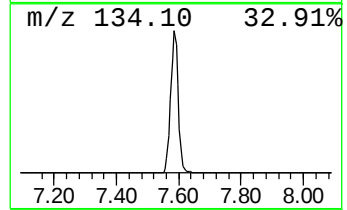
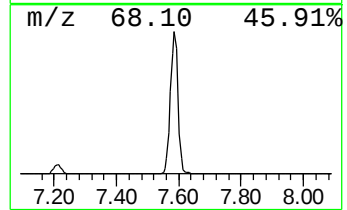
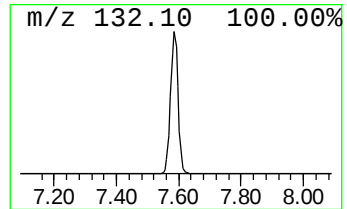
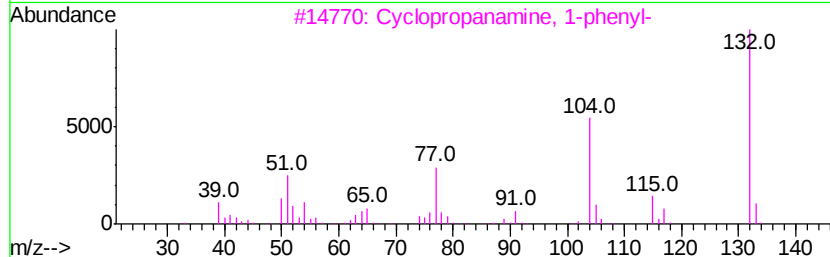
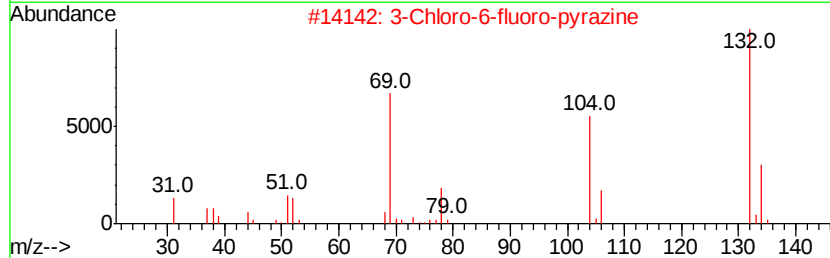
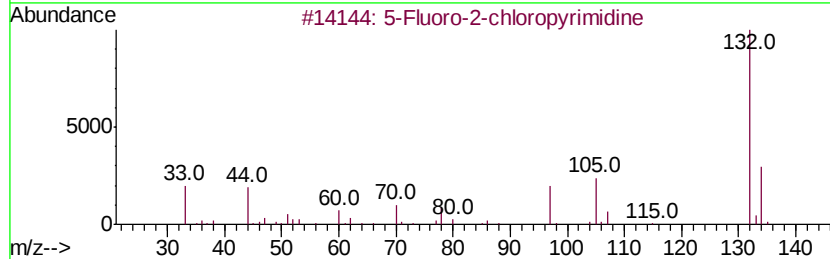
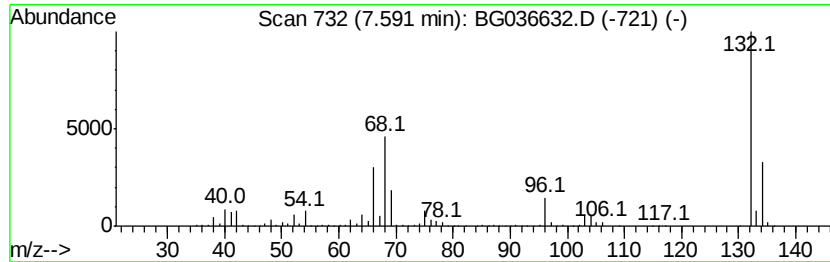
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	98.44 ng	1541170	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		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	12.2	ng	190710	1	8.06	313109	20.0
unknown7.59	7.59	98.4	ng	1541170	1	8.06	313109	20.0