

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
 Data File : BG036427.D
 Acq On : 27 Aug 2018 12:10
 Operator : JU/SJ
 Sample : J4594-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1-3-RO

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	5.167	310	320	328	rBV	80957	131468	3.19%	0.700%
2	5.632	392	399	408	rBV	800327	1233346	29.92%	6.565%
3	7.218	661	669	679	rBV	726774	1204580	29.22%	6.412%
4	7.594	726	733	745	rVB	823550	1395305	33.85%	7.427%
5	8.064	807	813	821	rVB	188502	319220	7.74%	1.699%
6	8.387	861	868	876	rBV	605813	1047120	25.40%	5.573%
7	9.227	1002	1011	1018	rBV	499769	905559	21.97%	4.820%
8	10.872	1283	1291	1297	rBV	267446	479019	11.62%	2.550%
9	13.317	1699	1707	1716	rVB	1436126	2238242	54.30%	11.913%
10	14.692	1934	1941	1949	rVB2	432747	701616	17.02%	3.734%
11	16.172	2186	2193	2211	rBV	1147590	1741017	42.23%	9.267%
12	17.429	2400	2407	2416	rBV	649410	999582	24.25%	5.320%
13	20.038	2845	2851	2862	rBV	3159009	4122235	100.00%	21.941%
14	21.695	3126	3133	3141	rBV2	710373	1185961	28.77%	6.313%
15	24.915	3671	3681	3691	rBV3	371984	1083223	26.28%	5.766%

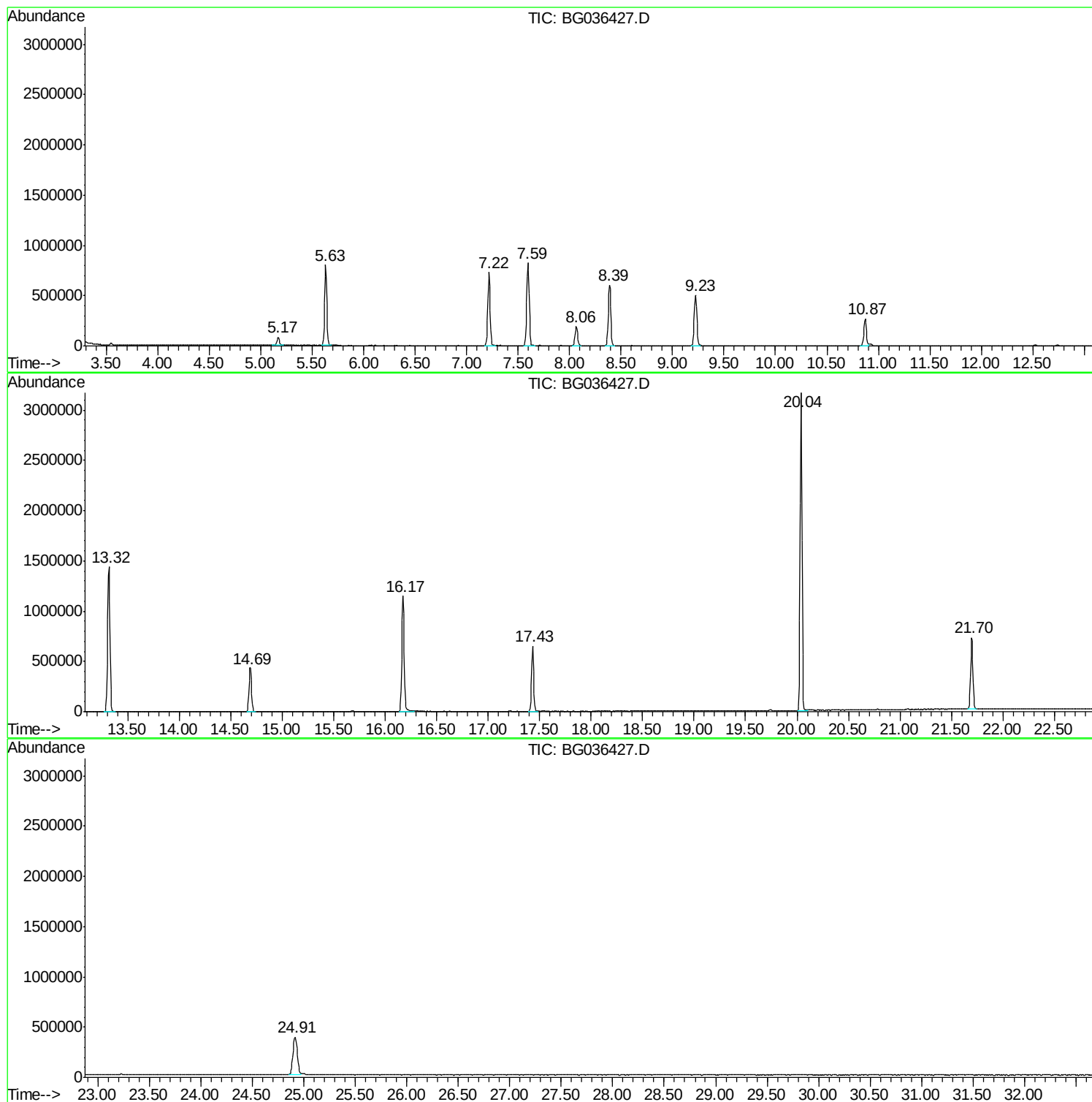
Sum of corrected areas: 18787493

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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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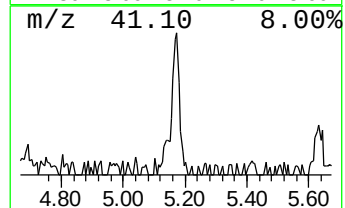
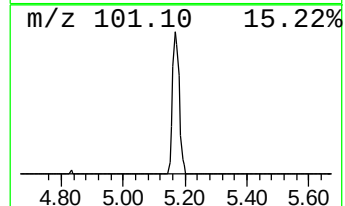
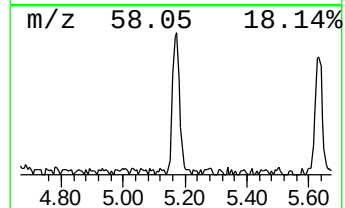
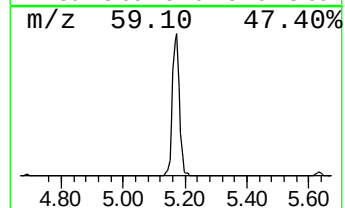
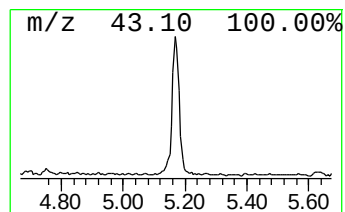
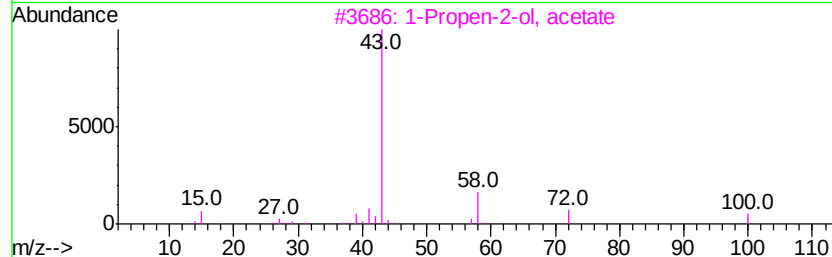
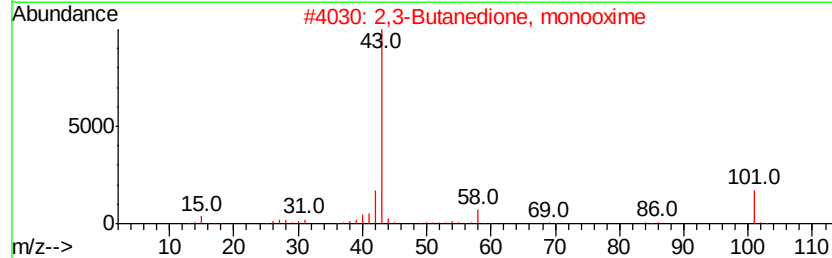
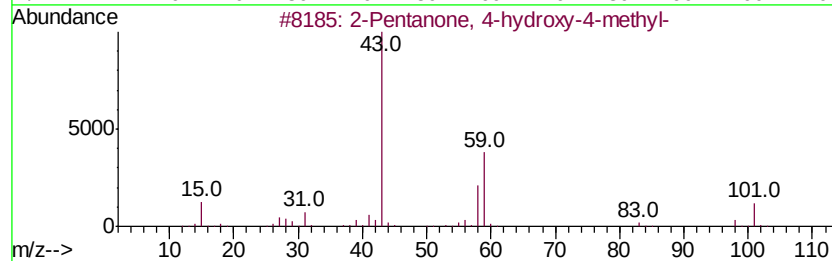
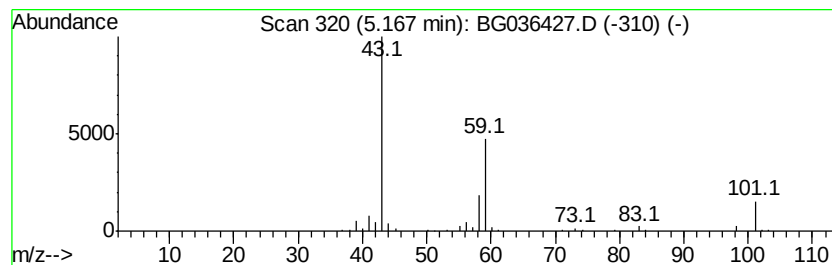
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	8.24 ng	131468	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	83
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		Butanoic acid, 4-amino-3-hydroxy...	119	C4H9NO3	000924-49-2	9



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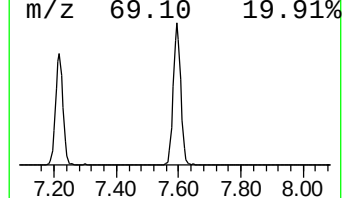
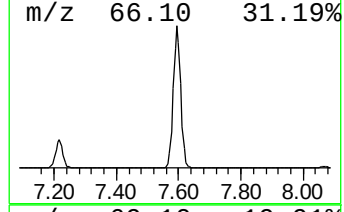
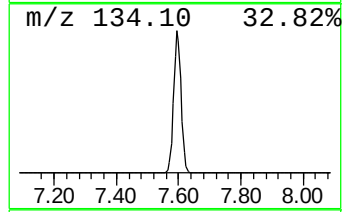
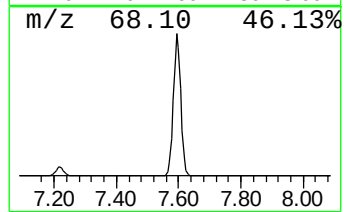
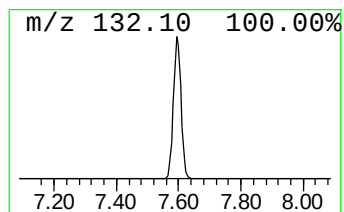
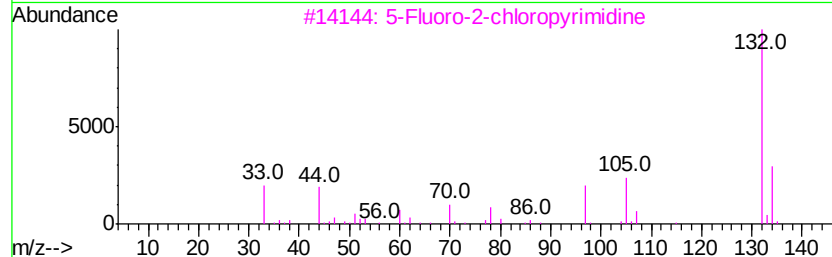
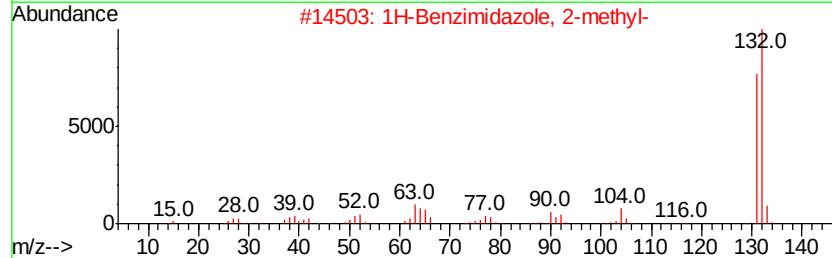
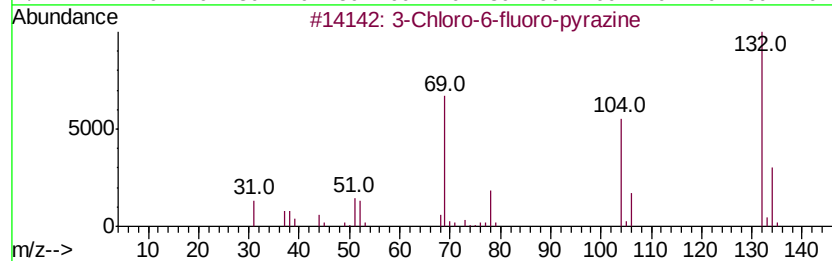
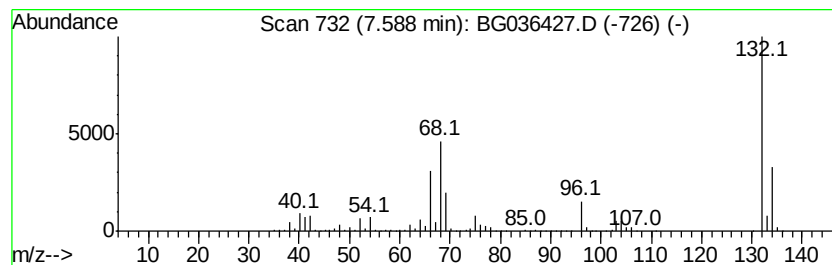
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	87.42 ng	1395310	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	25
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	8.2	ng	131468	1	8.06	319220	20.0
unknown7.59	7.59	87.4	ng	1395310	1	8.06	319220	20.0