

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092118\
 Data File : BG036893.D
 Acq On : 22 Sep 2018 14:14
 Operator : JU/SJ
 Sample : J5019-03
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 091918-DBRI-E-D-S

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-BG092018.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	5.150	312	317	328	rBV	24959	41874	1.05%	0.286%
2	5.620	390	397	416	rBV	168601	343794	8.59%	2.352%
3	7.206	661	667	691	rBV	107567	267349	6.68%	1.829%
4	7.576	723	730	749	rBV	351939	699958	17.49%	4.788%
5	8.046	804	810	824	rBV	121501	240227	6.00%	1.643%
6	8.370	857	865	881	rBV	516237	987807	24.68%	6.757%
7	9.210	1000	1008	1026	rBV	437568	903886	22.59%	6.183%
8	10.849	1280	1287	1305	rBV	172561	356585	8.91%	2.439%
9	13.293	1696	1703	1721	rBV	1346847	2330019	58.23%	15.938%
10	14.122	1837	1844	1852	rBV	35571	58292	1.46%	0.399%
11	14.674	1931	1938	1950	rBV2	319113	561783	14.04%	3.843%
12	16.160	2184	2191	2210	rBV	742561	1267547	31.68%	8.670%
13	17.418	2398	2405	2420	rBV	468117	754853	18.86%	5.163%
14	20.026	2842	2849	2860	rBV	2894665	4001695	100.00%	27.372%
15	21.683	3124	3131	3141	rBV2	514957	874291	21.85%	5.980%
16	23.358	3410	3416	3423	rVB3	23287	47946	1.20%	0.328%
17	24.891	3666	3677	3691	rBV	298379	881707	22.03%	6.031%

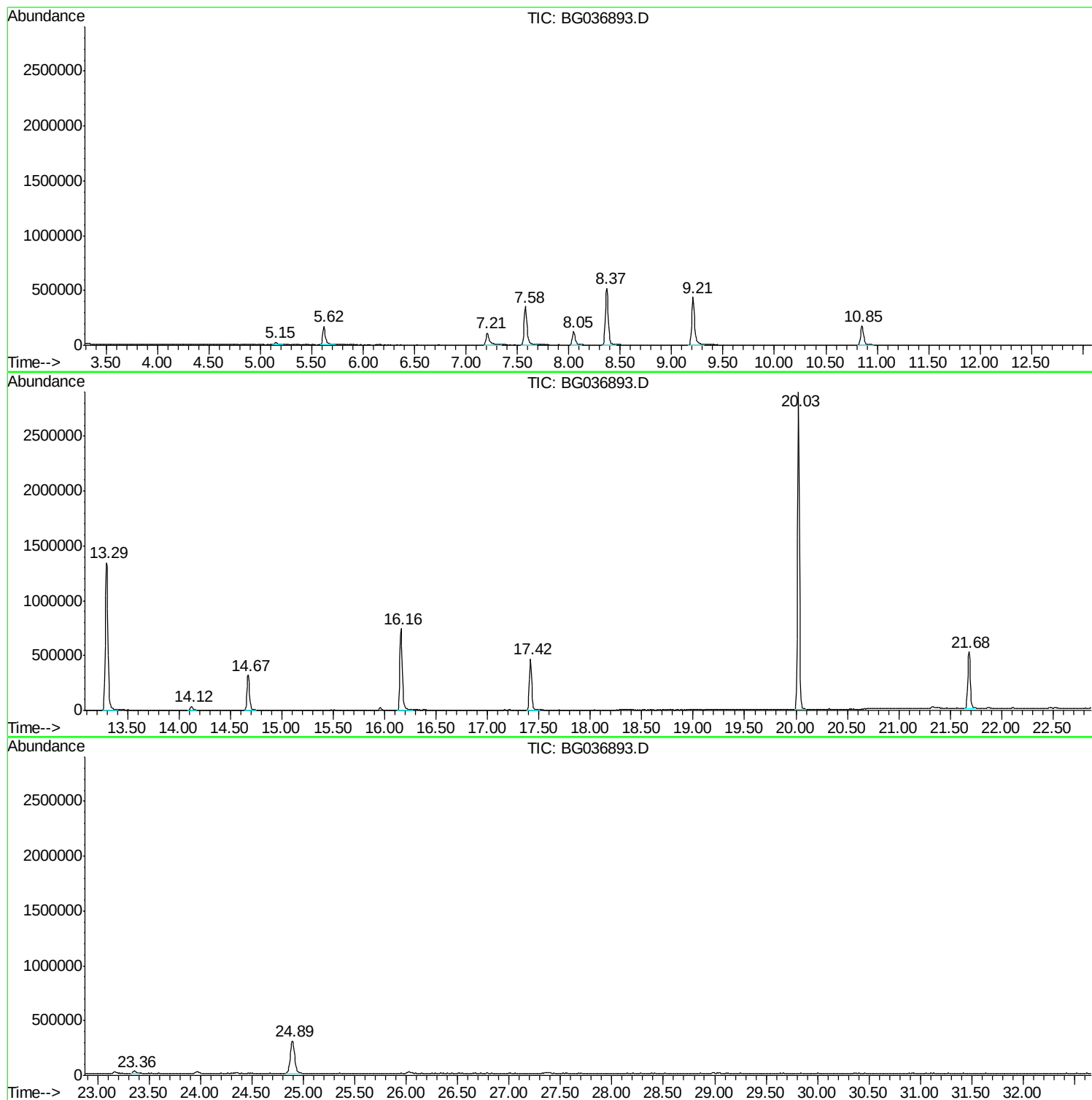
Sum of corrected areas: 14619613

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.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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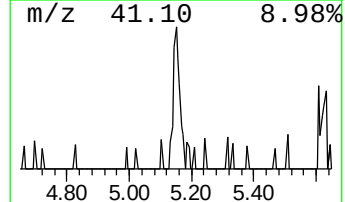
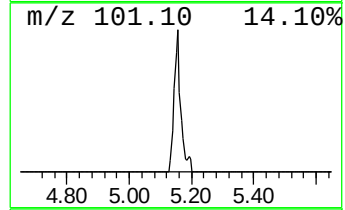
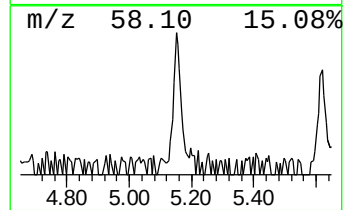
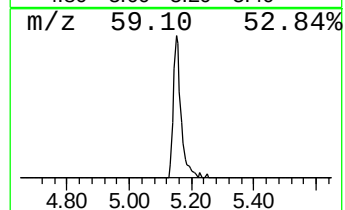
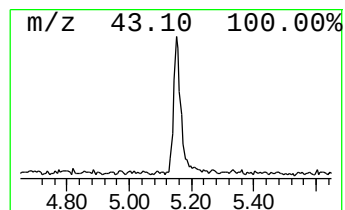
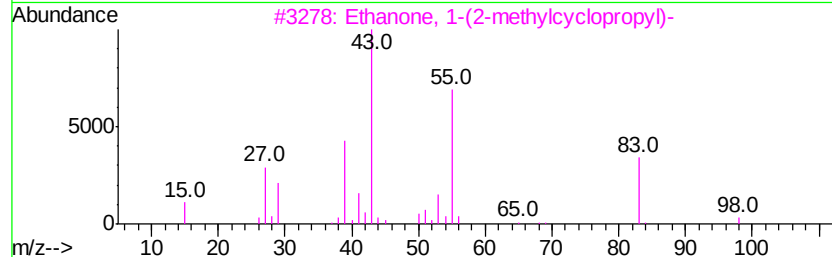
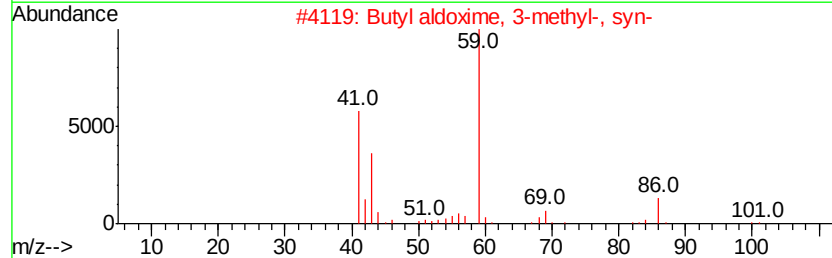
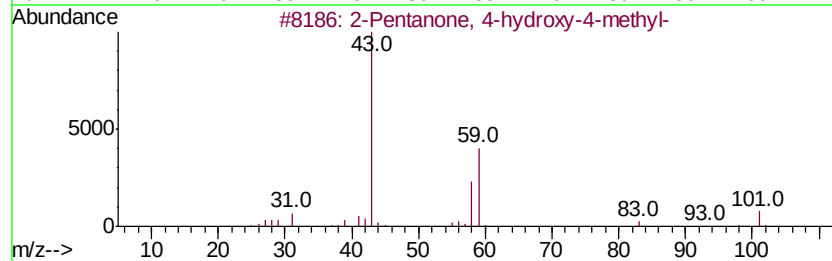
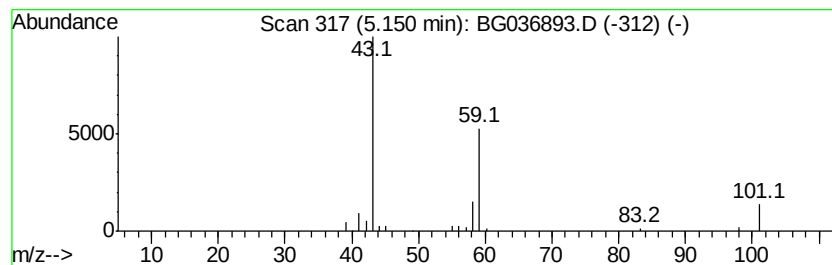
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.49 ng	41874	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3			Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			3-Hexanol	102	C6H14O	000623-37-0	9



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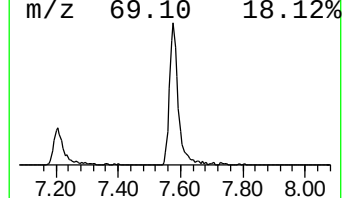
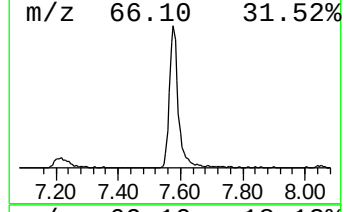
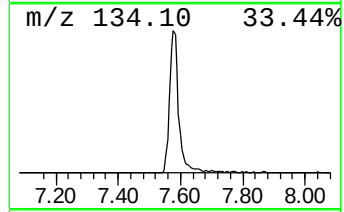
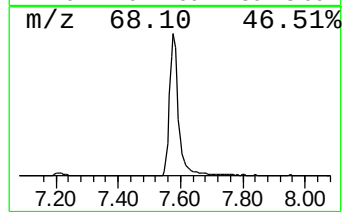
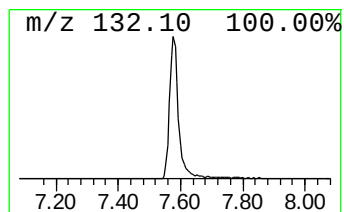
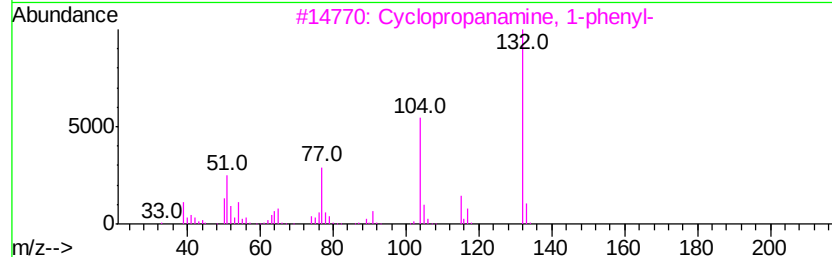
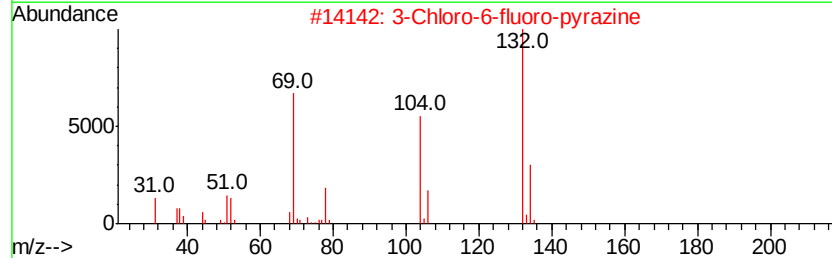
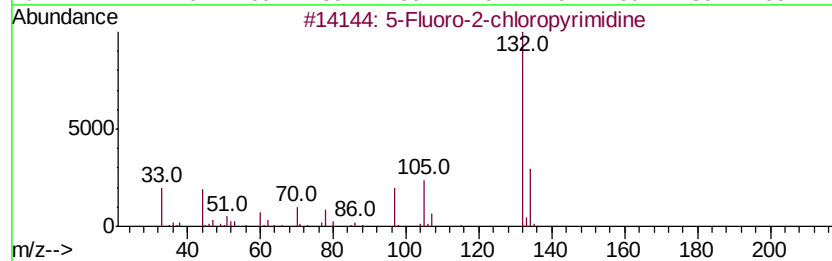
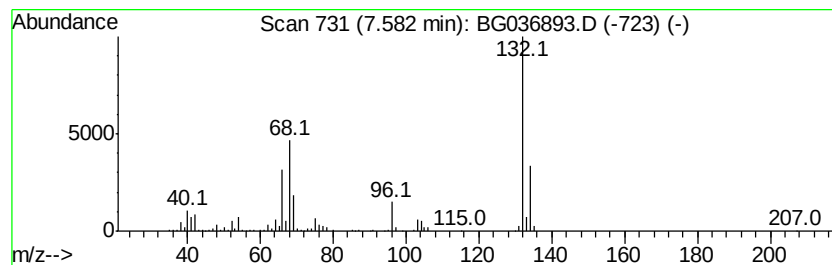
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	58.27 ng	699958	1,4-Dichlorobenzene-d4	8.05

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	16
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		Xenon	132	Xe	007440-63-3	9
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...	5.15	3.5	ng	41874	1	8.05	240227	20.0
unknown7.58	7.58	58.3	ng	699958	1	8.05	240227	20.0