

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081018\
 Data File : BG036279.D
 Acq On : 10 Aug 2018 22:40
 Operator : JU/SJ
 Sample : J4379-25
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-103D

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-BG071218.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	5.123	308	312	319	rBV	20403	31854	1.24%	0.332%
2	5.593	385	392	406	rBV	169214	333928	13.01%	3.481%
3	7.185	656	663	677	rBV2	97308	219598	8.56%	2.289%
4	7.543	717	724	740	rBV	382857	692546	26.99%	7.218%
5	8.008	796	803	811	rBV	92662	155069	6.04%	1.616%
6	8.331	848	858	871	rVB	385096	692925	27.01%	7.222%
7	9.171	992	1001	1013	rBV2	263187	550020	21.44%	5.733%
8	10.810	1272	1280	1293	rBV	82457	185944	7.25%	1.938%
9	13.253	1689	1696	1711	rBV	842776	1380451	53.80%	14.389%
10	14.634	1924	1931	1942	rBV2	168863	297853	11.61%	3.105%
11	16.120	2178	2184	2198	rBV	674780	1173622	45.74%	12.233%
12	17.377	2392	2398	2415	rBV2	237476	415552	16.20%	4.331%
13	19.985	2835	2842	2854	rBV	1881104	2565880	100.00%	26.744%
14	21.636	3117	3123	3134	rBV2	260298	476058	18.55%	4.962%
15	24.826	3656	3666	3684	rVB3	121427	422755	16.48%	4.406%

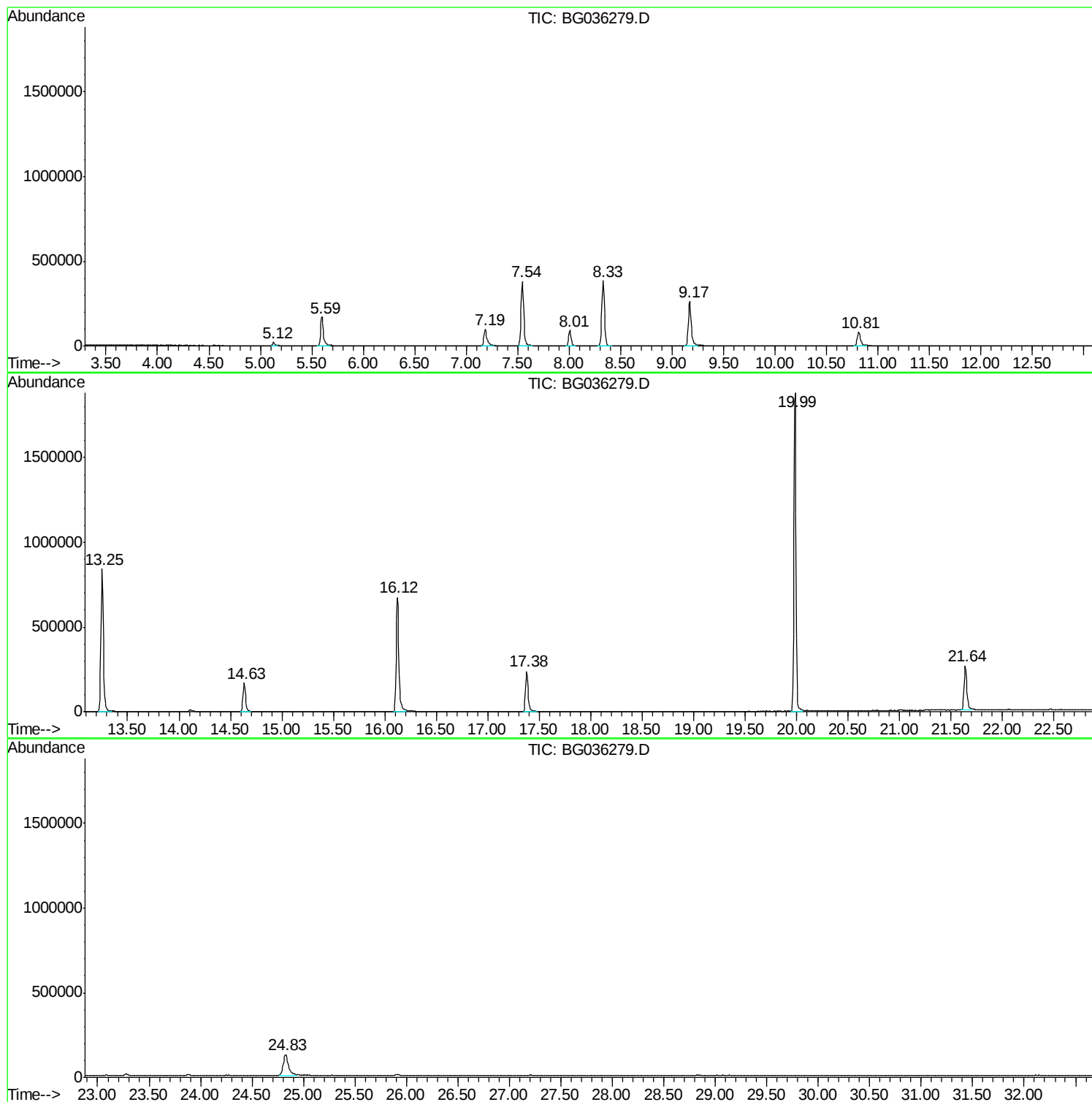
Sum of corrected areas: 9594055

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.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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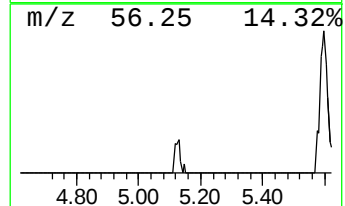
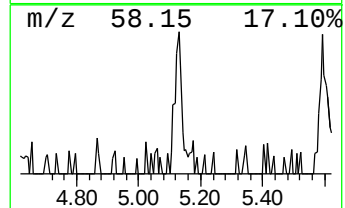
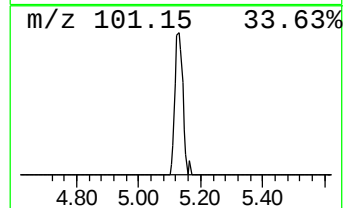
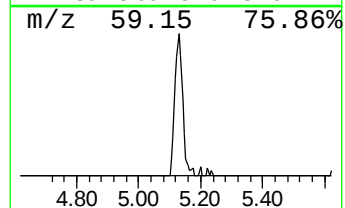
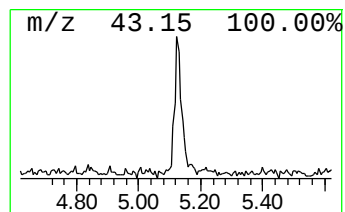
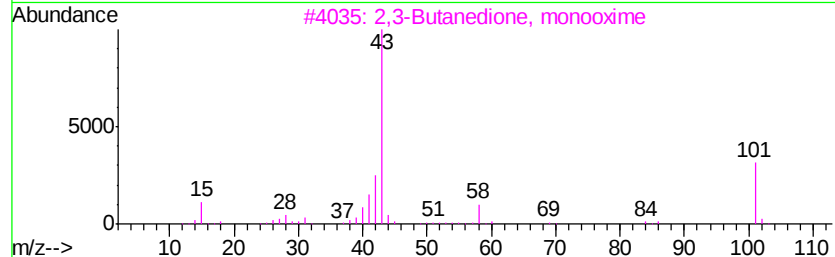
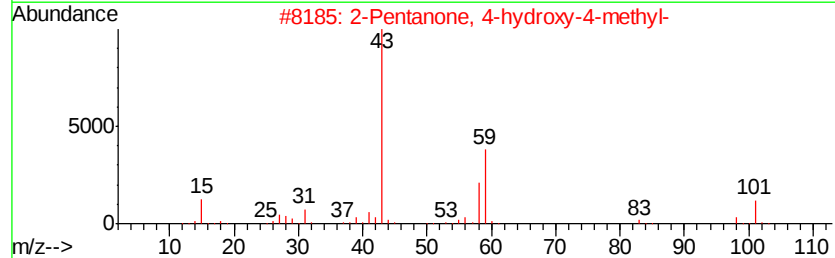
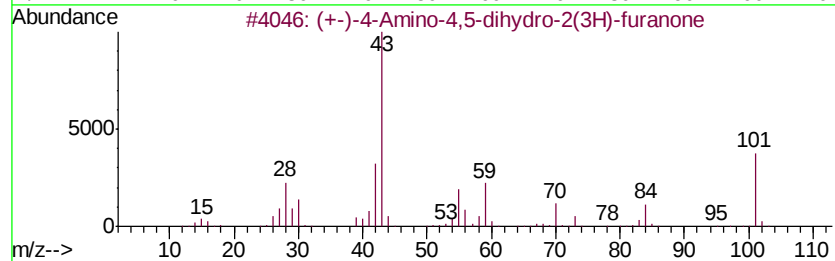
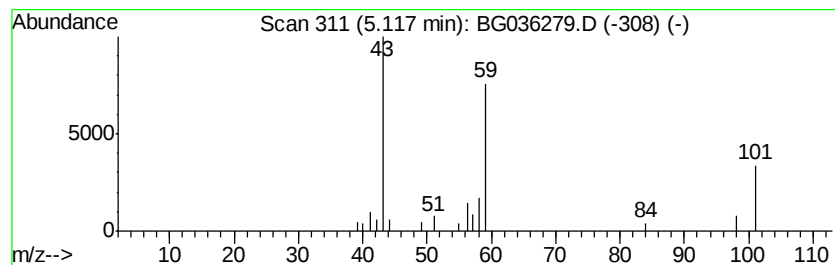
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Peak Number 1 (+)-4-Amino-4,5-dihydro-2(...) Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	4.11 ng	31854	1,4-Dichlorobenzene-d4	8.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7N02	016504-58-8	72
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
3	2,3-Butanedione, monooxime	101	C4H7N02	000057-71-6	32
4	Methyl 2-acetamidoacrylate	143	C6H9N03	035356-70-8	17
5	5-Hydroxypentanehydroxylamine, N...	245	C11H19N05	104556-13-0	17



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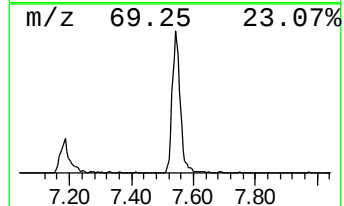
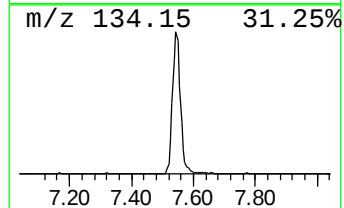
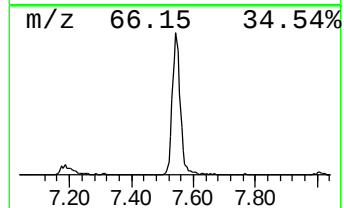
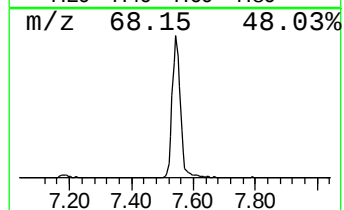
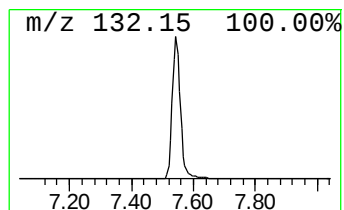
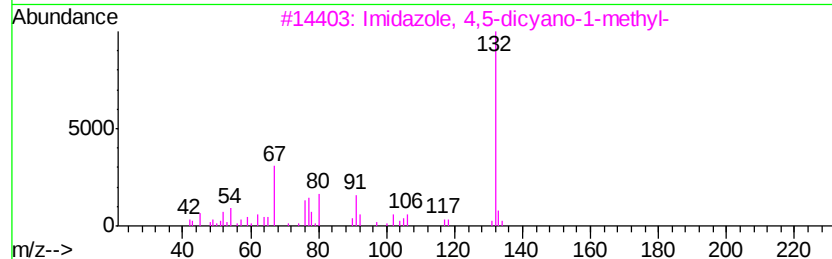
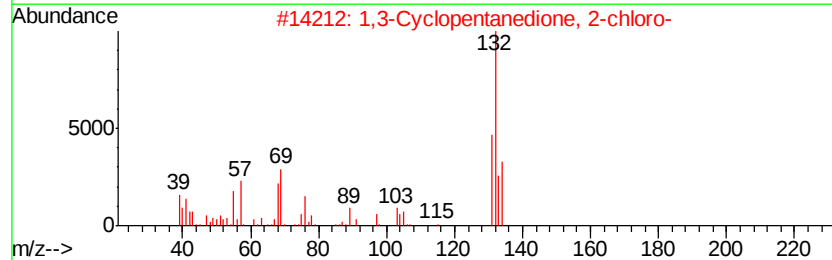
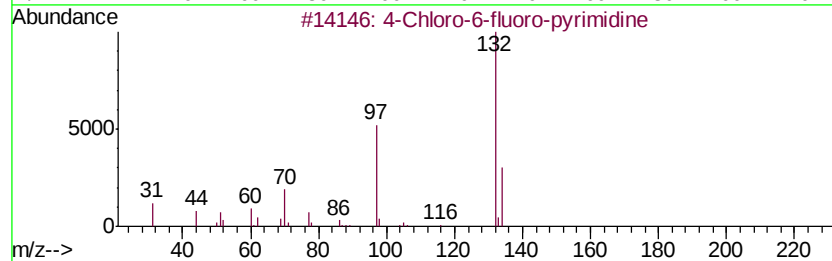
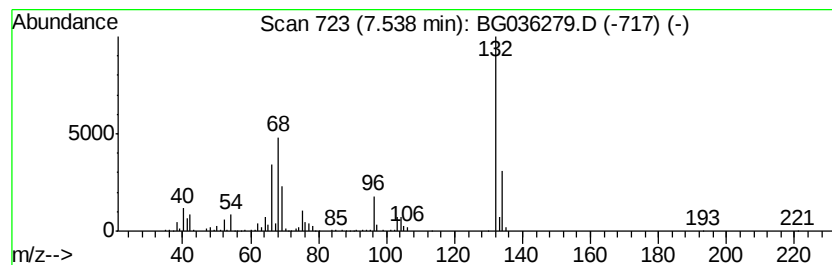
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Peak Number 2 unknown7.54 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	89.32 ng	692546	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(+)-4-Amino-4,5-...	5.12	4.1	ng	31854	1	8.01	155069	20.0
unknown7.54	7.54	89.3	ng	692546	1	8.01	155069	20.0