

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073018\
 Data File : BG035983.D
 Acq On : 30 Jul 2018 19:34
 Operator : JU/SJ
 Sample : PB111568BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111568BL

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	5.132	309	314	320	rBV	21901	38064	1.04%	0.272%
2	5.602	387	394	409	rBV	426944	739690	20.21%	5.278%
3	7.188	657	664	677	rBV	458391	819449	22.38%	5.847%
4	7.558	719	727	737	rBV	446522	780294	21.32%	5.568%
5	8.022	800	806	816	rVB3	128925	217157	5.93%	1.549%
6	8.345	854	861	873	rVB	516928	901263	24.62%	6.431%
7	9.185	994	1004	1019	rBV	375844	740591	20.23%	5.284%
8	10.824	1276	1283	1302	rBV	134702	285219	7.79%	2.035%
9	13.268	1690	1699	1712	rBV	1223704	1976686	54.00%	14.104%
10	14.648	1926	1934	1943	rBV2	268592	465637	12.72%	3.322%
11	16.135	2181	2187	2205	rBV	538089	995743	27.20%	7.105%
12	17.392	2394	2401	2416	rBV2	429432	740001	20.21%	5.280%
13	20.000	2838	2845	2856	rBV	2633646	3660721	100.00%	26.120%
14	21.651	3119	3126	3138	rBV	473703	879121	24.01%	6.273%
15	24.846	3660	3670	3684	rBV2	232181	775357	21.18%	5.532%

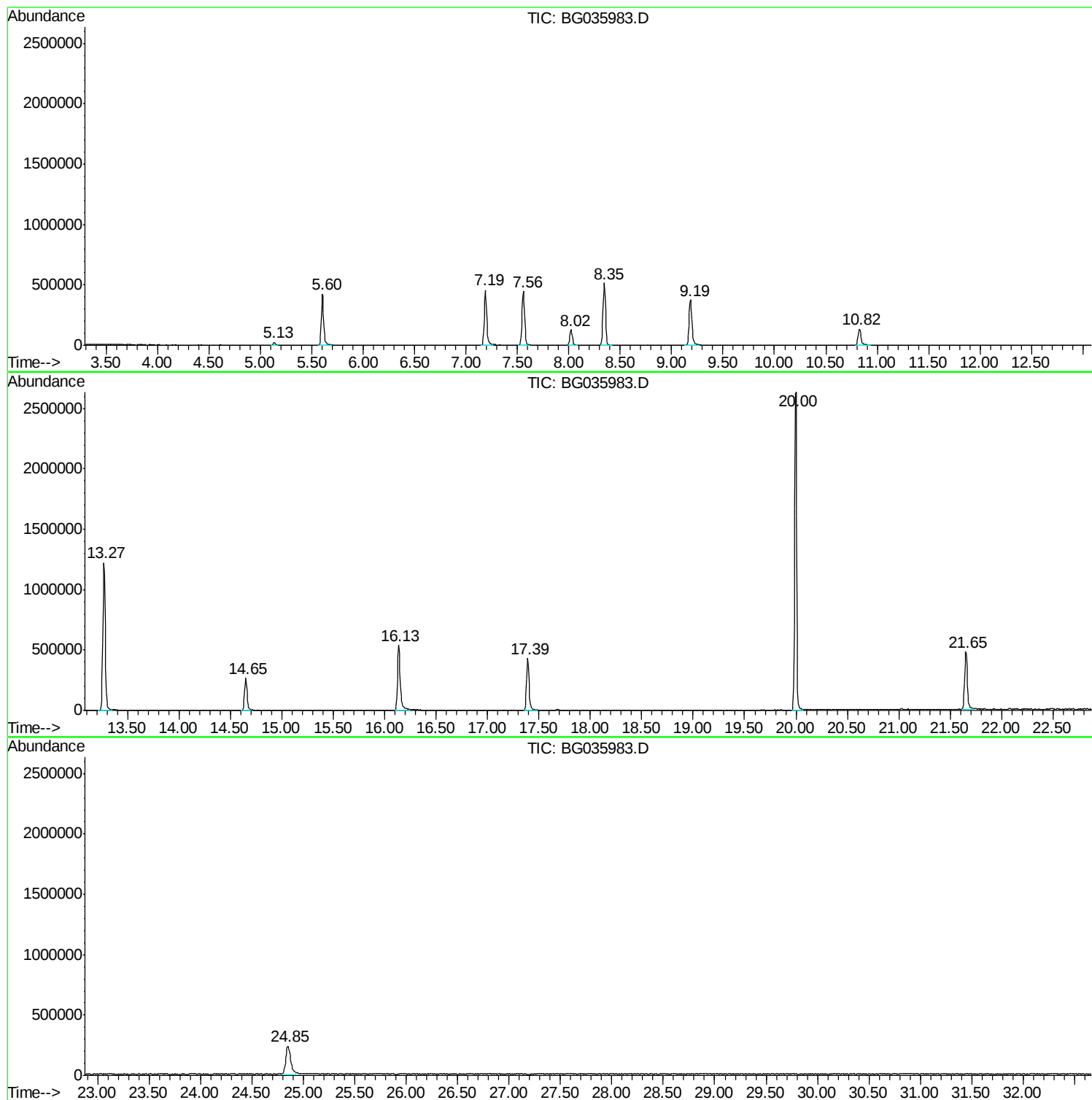
Sum of corrected areas: 14014993

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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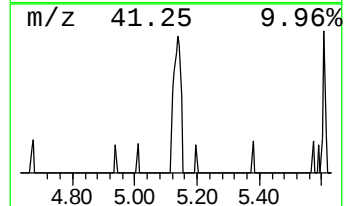
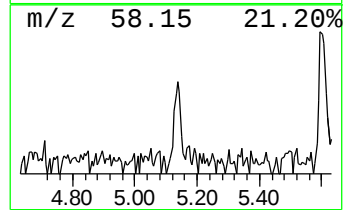
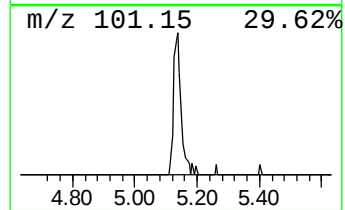
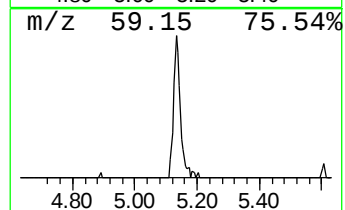
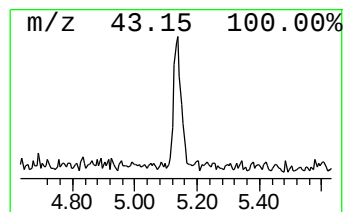
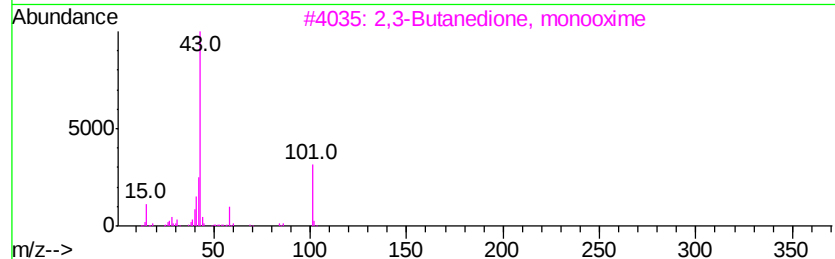
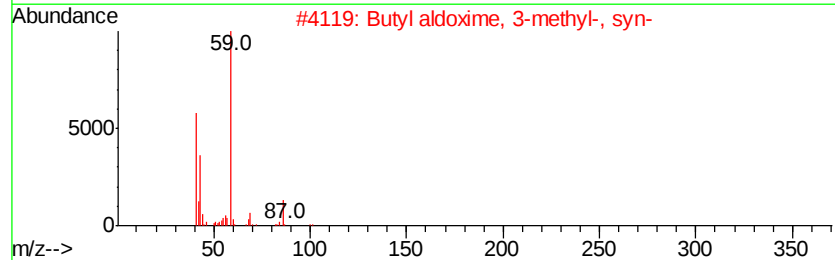
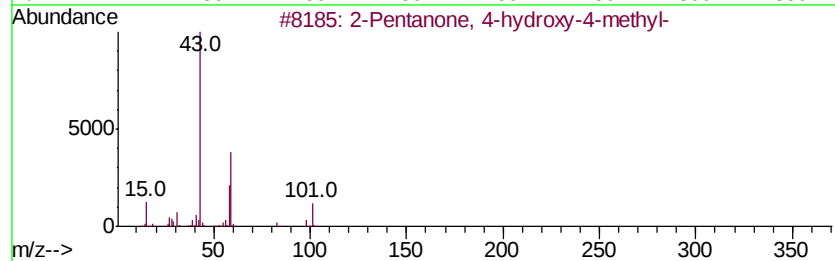
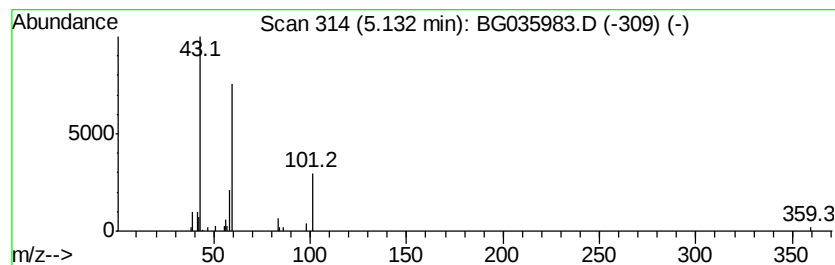
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	3.51 ng	38064	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	9
5			.alpha.-D-Galactopyranoside, met...	249	C10H19NO6	017296-07-0	9



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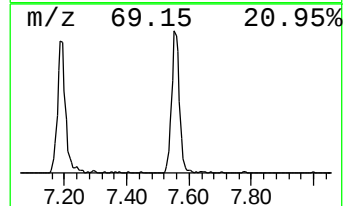
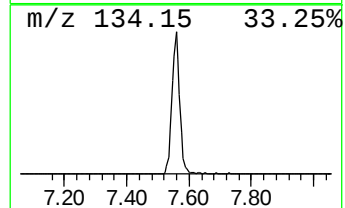
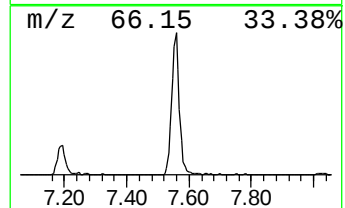
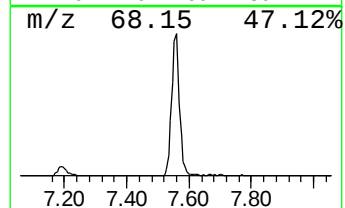
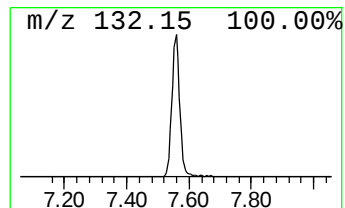
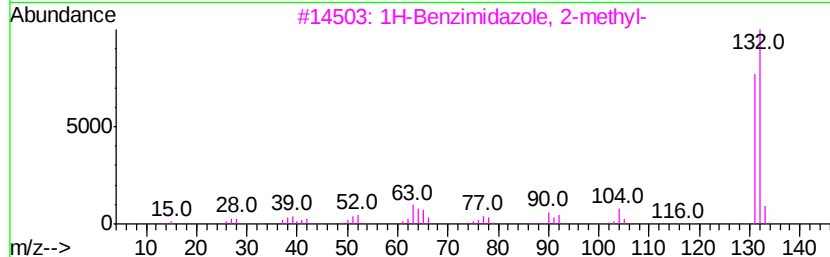
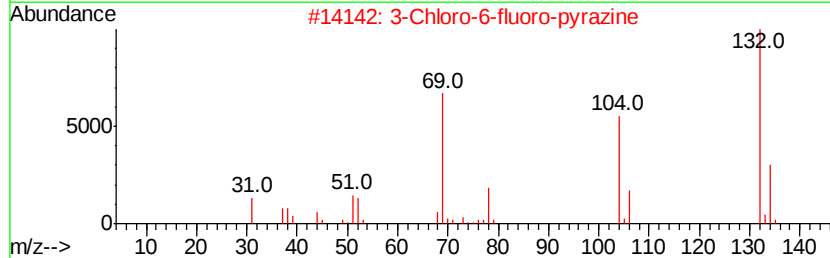
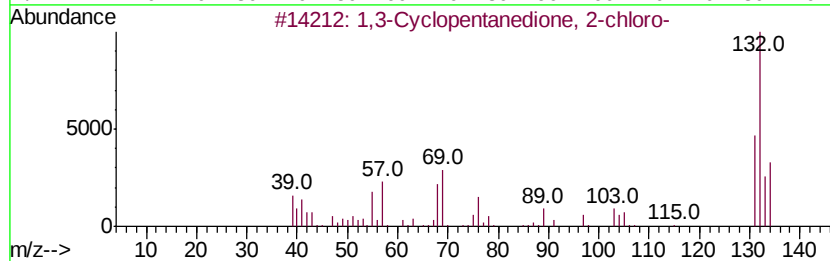
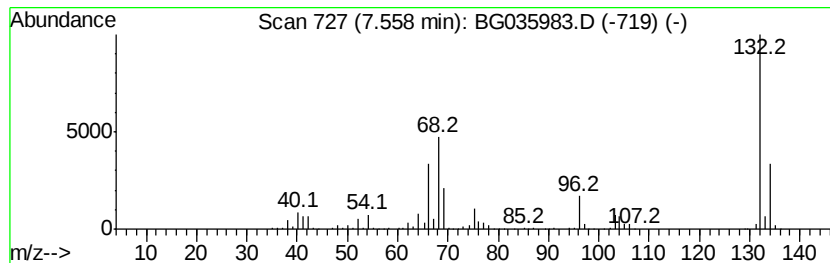
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Peak Number 2 unknown7.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	71.86 ng	780294	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
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...	5.13	3.5	ng	38064	1	8.02	217157	20.0
unknown7.56	7.56	71.9	ng	780294	1	8.02	217157	20.0