

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
 Data File : BG035787.D
 Acq On : 20 Jul 2018 16:51
 Operator : JU/SJ
 Sample : PB111042BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111042BL

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	5.144	311	316	323	rBV	40522	65682	1.56%	0.398%
2	5.614	389	396	412	rBV	551314	950480	22.51%	5.765%
3	7.194	658	665	677	rBV	560600	1014055	24.01%	6.151%
4	7.564	719	728	744	rBV	574871	996322	23.59%	6.043%
5	8.028	800	807	814	rBV	167946	282728	6.70%	1.715%
6	8.351	854	862	872	rVB	665920	1173458	27.79%	7.118%
7	9.191	997	1005	1018	rBV	492592	943059	22.33%	5.720%
8	10.830	1276	1284	1295	rBV	166809	344858	8.17%	2.092%
9	13.274	1693	1700	1715	rBV	1454559	2322111	54.99%	14.085%
10	14.654	1928	1935	1950	rBV	299386	528865	12.52%	3.208%
11	16.141	2181	2188	2200	rBV	665280	1148740	27.20%	6.968%
12	17.392	2395	2401	2416	rBV2	430583	740459	17.53%	4.491%
13	20.000	2839	2845	2862	rBV	3072769	4222960	100.00%	25.615%
14	21.657	3120	3127	3141	rBV2	503674	915784	21.69%	5.555%
15	24.858	3661	3672	3682	rBV2	271657	836883	19.82%	5.076%

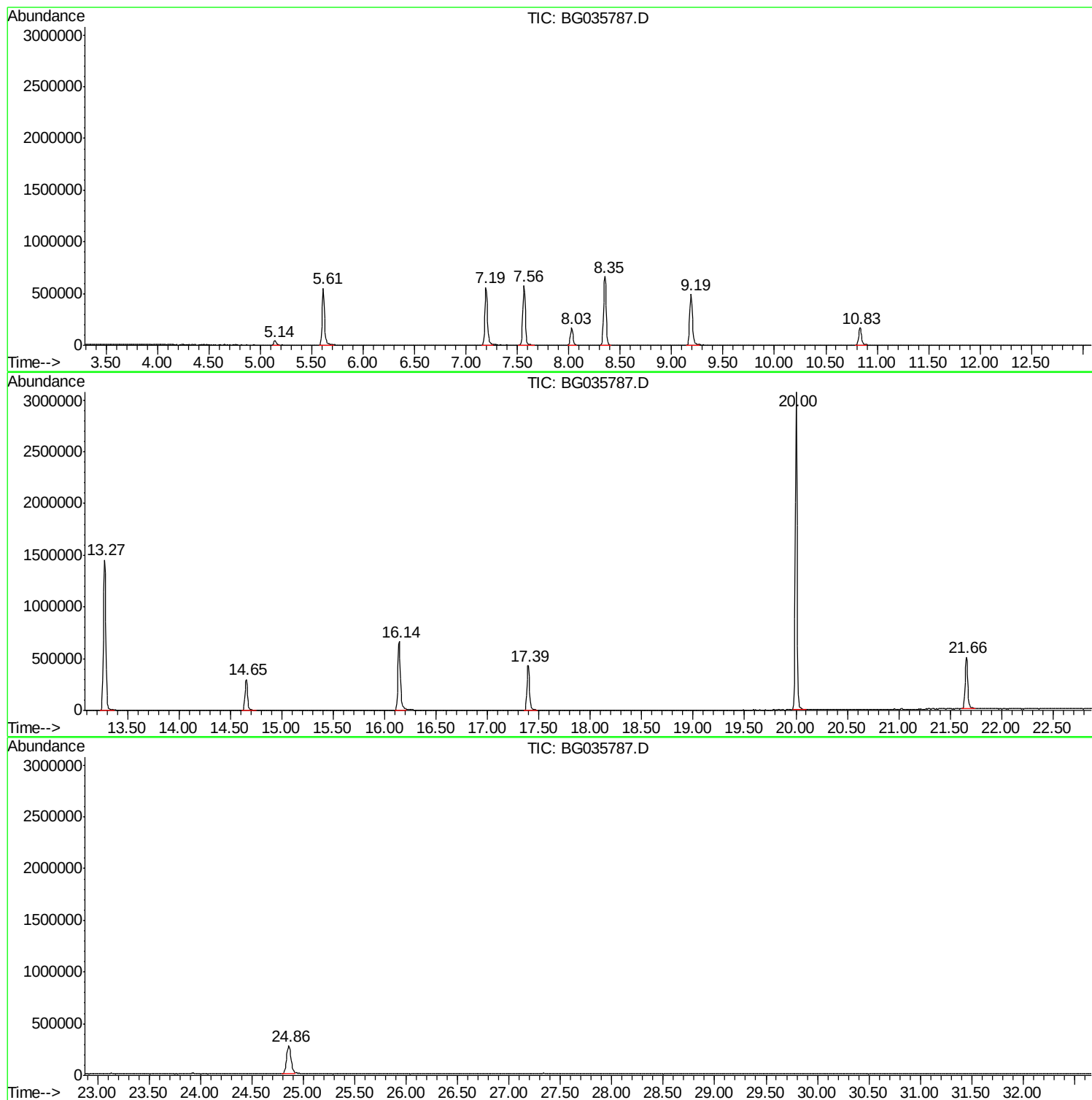
Sum of corrected areas: 16486444

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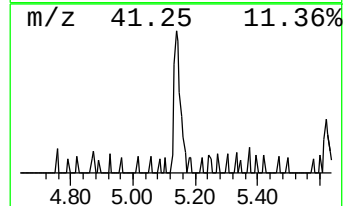
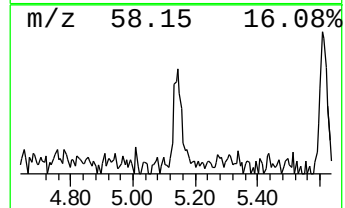
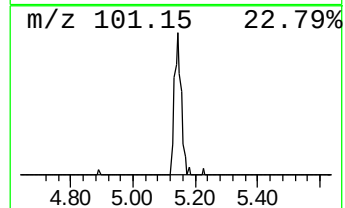
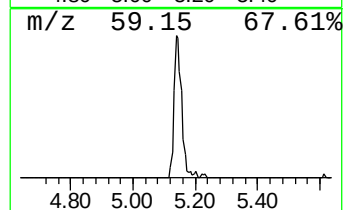
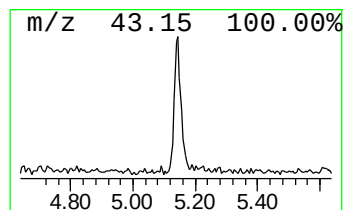
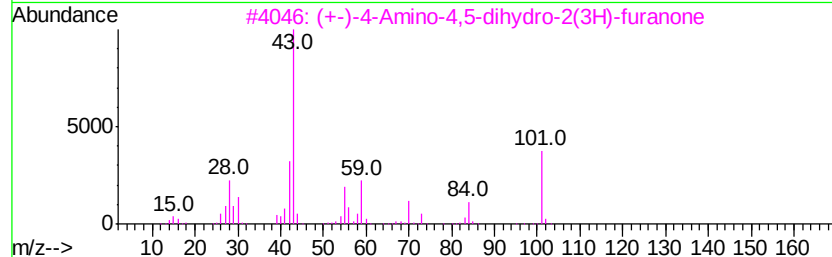
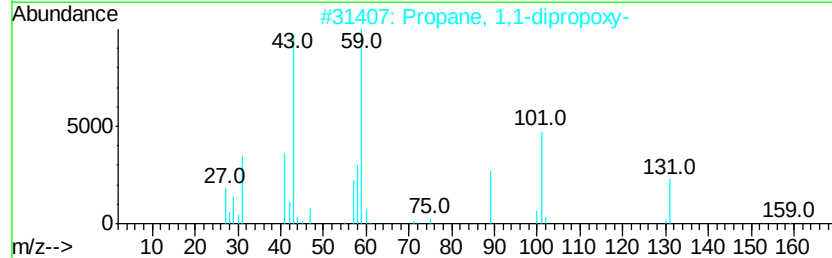
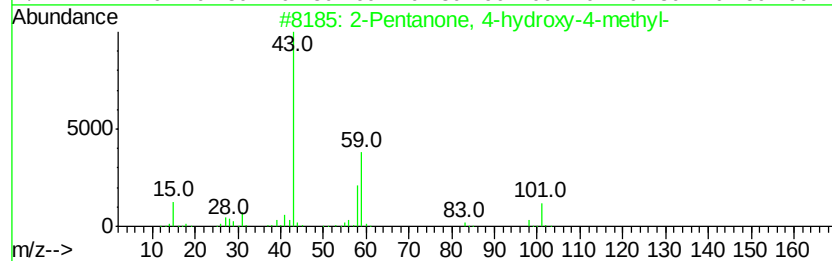
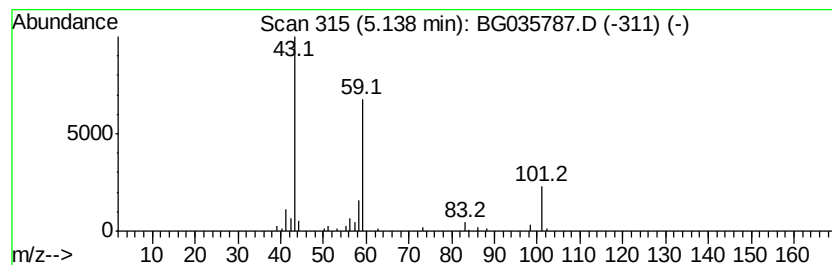
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.65 ng	65682	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38
3			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	38
4			2,3-Dimethyloxirane-2-carboxylic acid	130	C6H10O3	1000306-64-4	37
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	37



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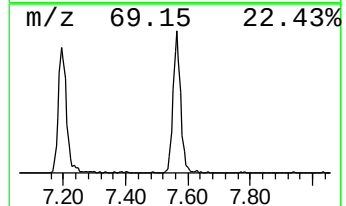
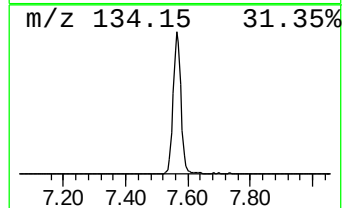
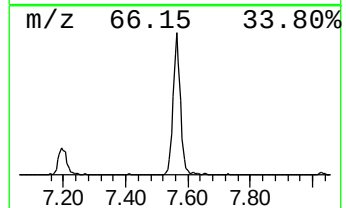
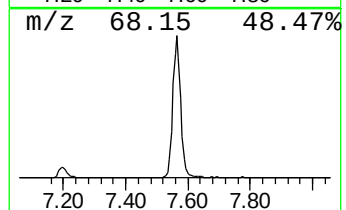
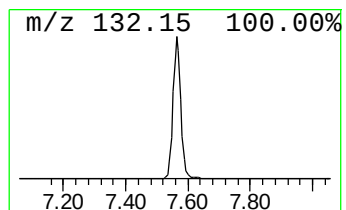
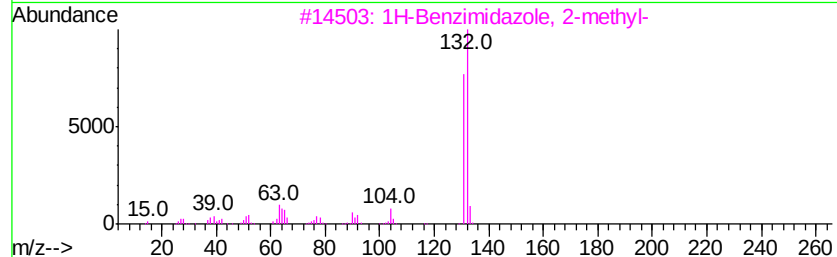
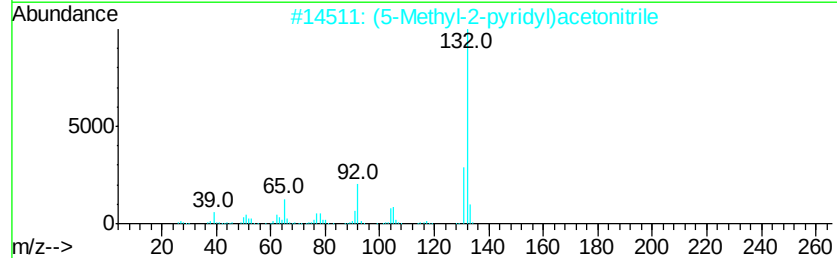
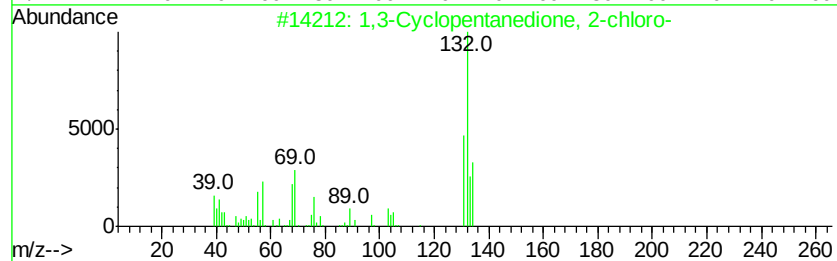
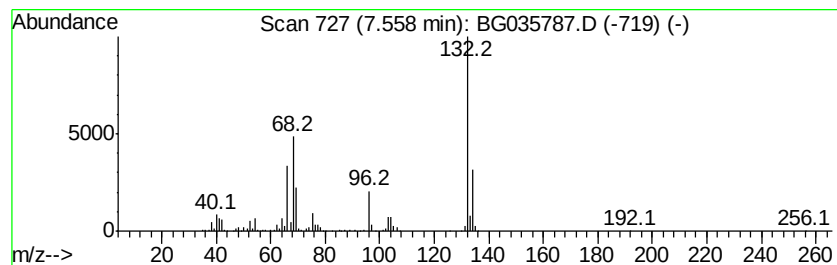
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Peak Number 2 unknown7.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	70.48 ng	996322	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
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...	5.14	4.7	ng	65682	1	8.03	282728	20.0
unknown7.56	7.56	70.5	ng	996322	1	8.03	282728	20.0