

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092816\
Data File : BG024149.D
Acq On : 28 Sep 2016 16:39
Operator : UM/SJ
Sample : PB93653BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB93653BL

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:\HPCHEM1\BNA_G\METHODS\8270-BG092316.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	5.106	382	386	394	rVB	163113	266950	5.30%	1.309%
2	5.594	463	469	478	rBV	556332	1009094	20.05%	4.949%
3	7.215	739	745	768	rBV	580822	1158303	23.01%	5.681%
4	7.574	800	806	824	rVB	633817	1155084	22.95%	5.665%
5	8.044	881	886	892	rBV2	183872	321805	6.39%	1.578%
6	8.367	935	941	950	rVB	696757	1225197	24.34%	6.009%
7	9.224	1081	1087	1099	rBV	545585	1103779	21.93%	5.414%
8	10.875	1362	1368	1383	rBV	212423	470926	9.35%	2.310%
9	13.331	1779	1786	1807	rBV	1614367	2752225	54.67%	13.499%
10	14.723	2014	2023	2036	rBV	439357	806847	16.03%	3.957%
11	16.227	2272	2279	2301	rBV	753850	1523588	30.27%	7.473%
12	17.490	2485	2494	2510	rBV	600935	1132389	22.49%	5.554%
13	20.092	2931	2937	2946	rBV	3760585	5034116	100.00%	24.691%
14	21.790	3220	3226	3239	rBV	671460	1297959	25.78%	6.366%
15	25.126	3784	3794	3809	rBV2	340408	1129887	22.44%	5.542%

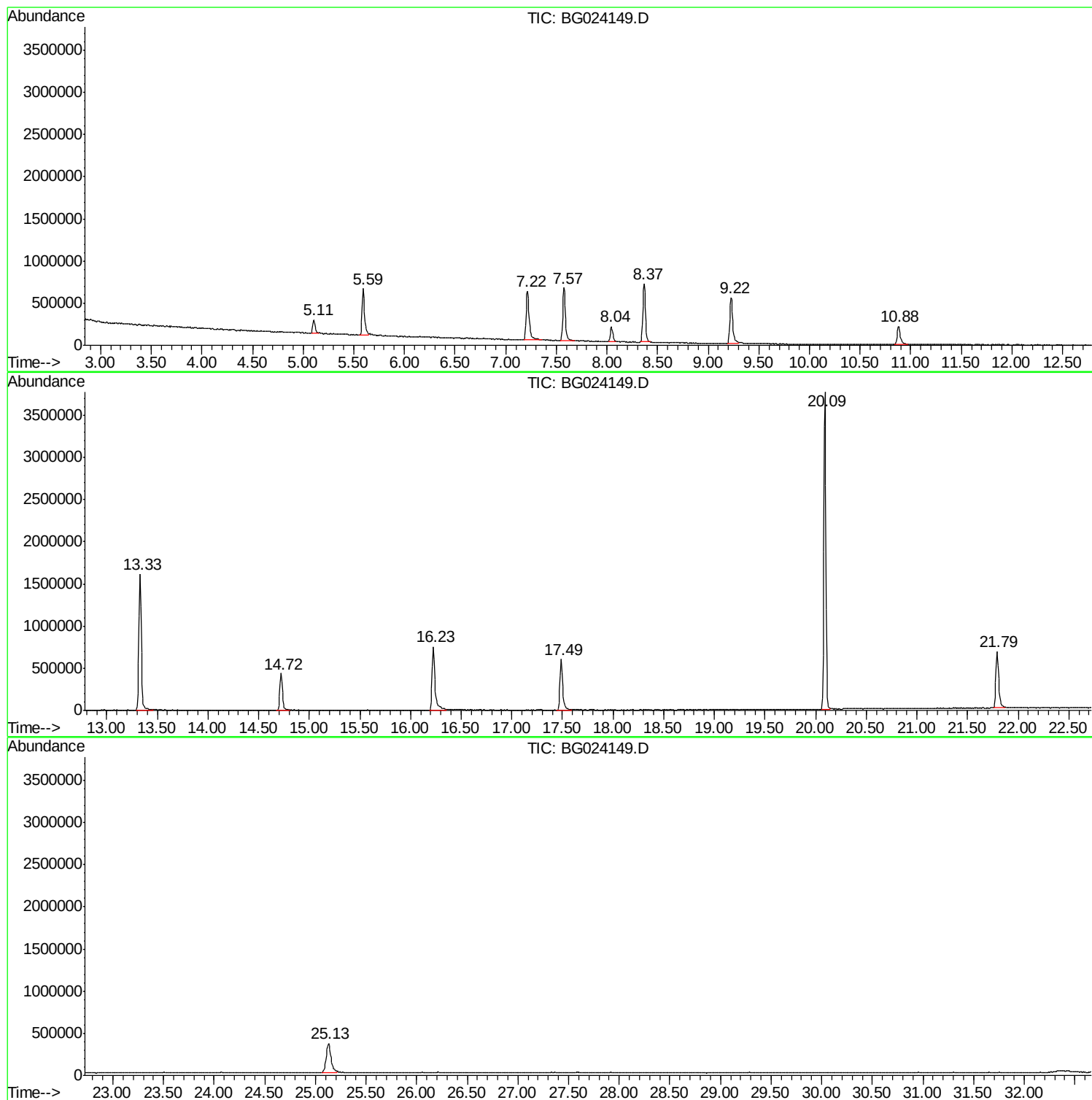
Sum of corrected areas: 20388149

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG092316.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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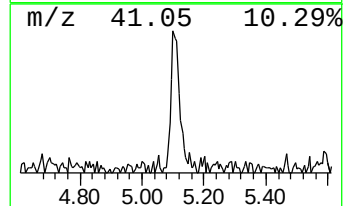
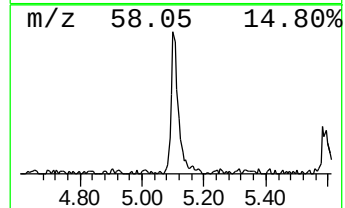
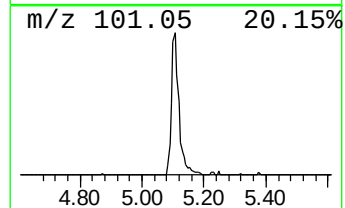
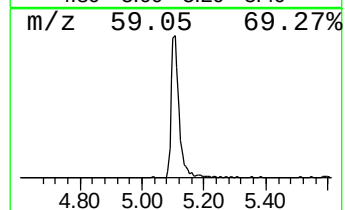
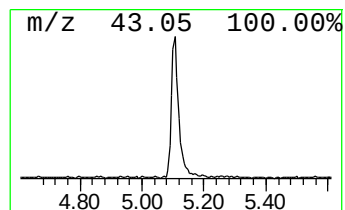
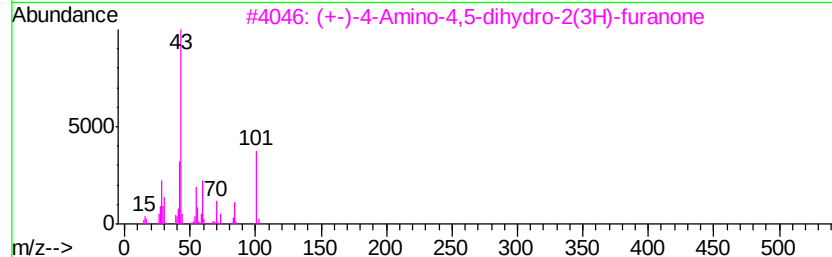
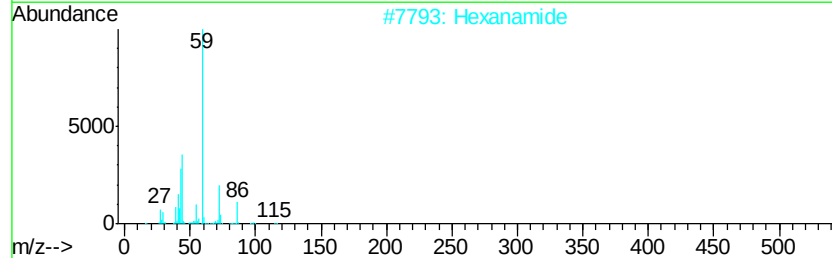
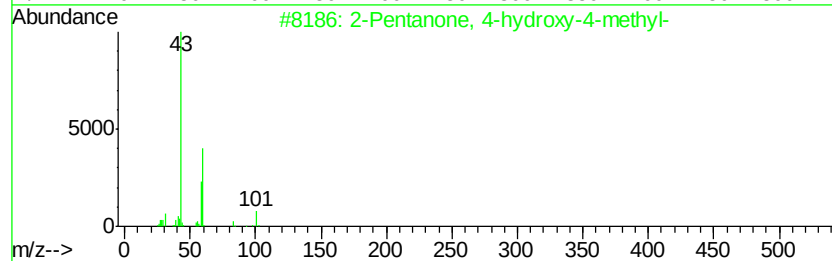
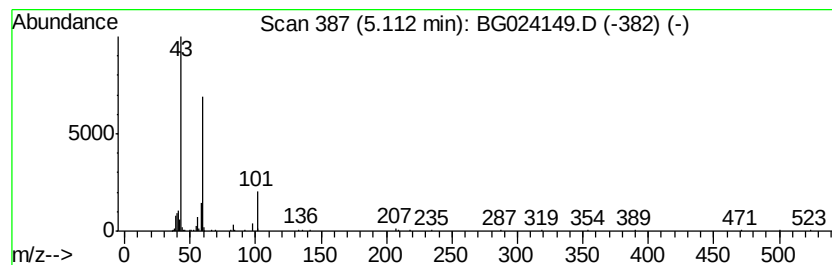
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	16.59 ng	266950	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Hexanamide	115	C6H13NO	000628-02-4	9
3			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	9



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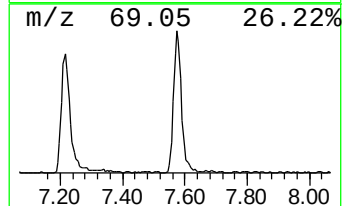
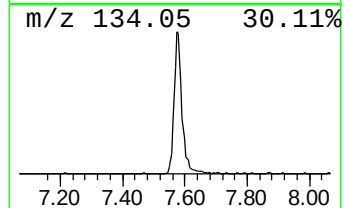
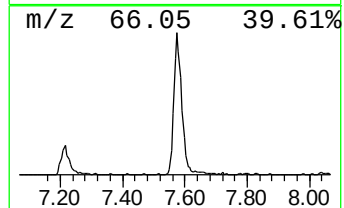
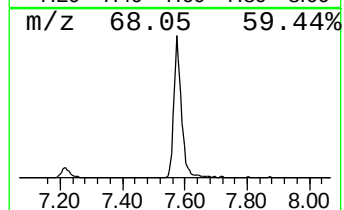
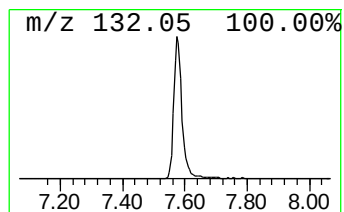
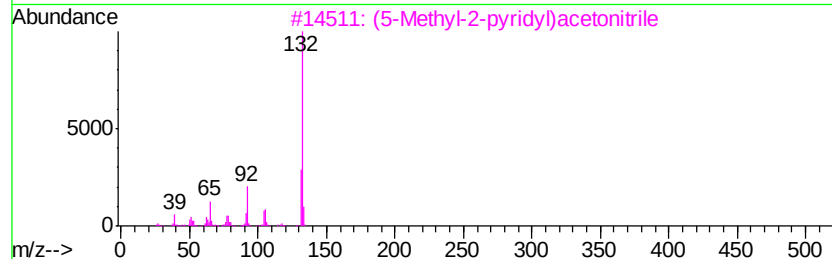
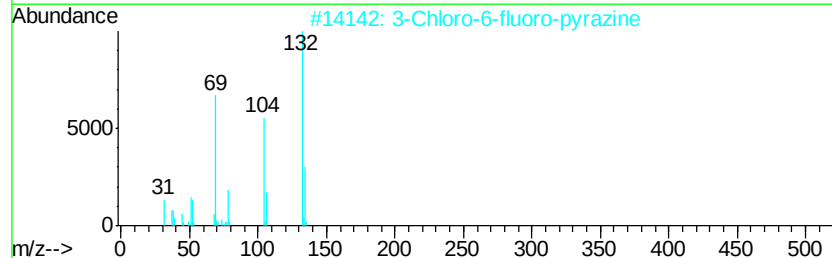
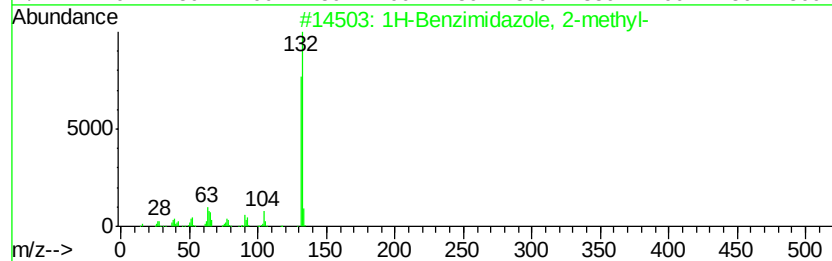
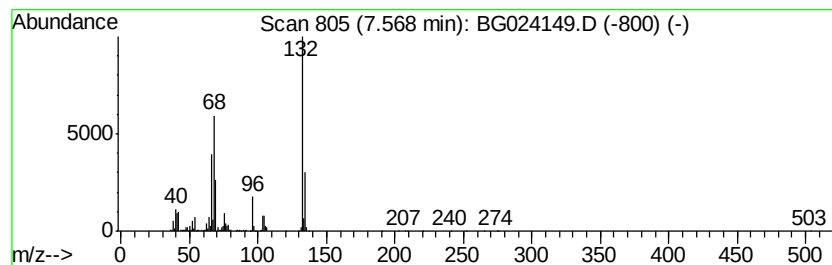
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Peak Number 2 unknown7.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	71.79 ng	1155080	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.11	16.6	ng	266950	1	8.04	321805	20.0
unknown7.57	7.57	71.8	ng	1155080	1	8.04	321805	20.0