

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
 Data File : BG024117.D
 Acq On : 24 Sep 2016 10:06
 Operator : UM/SJ
 Sample : H4902-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EB-01-092116

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	5.105	382	386	393	rVB	57722	84382	1.48%	0.414%
2	5.587	463	468	483	rBV	377186	691532	12.09%	3.390%
3	7.214	740	745	758	rVB	236521	452440	7.91%	2.218%
4	7.573	799	806	816	rBV	726369	1326106	23.19%	6.501%
5	8.037	880	885	892	rBV2	152965	264136	4.62%	1.295%
6	8.366	933	941	952	rBV	669608	1214218	21.24%	5.953%
7	9.223	1080	1087	1099	rBV	601553	1159977	20.29%	5.687%
8	10.874	1361	1368	1378	rBV	192309	391261	6.84%	1.918%
9	13.324	1778	1785	1795	rBV	1754767	2755169	48.18%	13.508%
10	14.716	2016	2022	2031	rBV	371741	625160	10.93%	3.065%
11	16.220	2271	2278	2297	rBV	1632696	2668681	46.67%	13.084%
12	17.483	2486	2493	2506	rBV	555482	897101	15.69%	4.398%
13	20.091	2931	2937	2947	rBV	4447944	5717992	100.00%	28.033%
14	21.789	3219	3226	3237	rVB	629132	1117318	19.54%	5.478%
15	25.125	3783	3794	3805	rBV2	329952	1031687	18.04%	5.058%

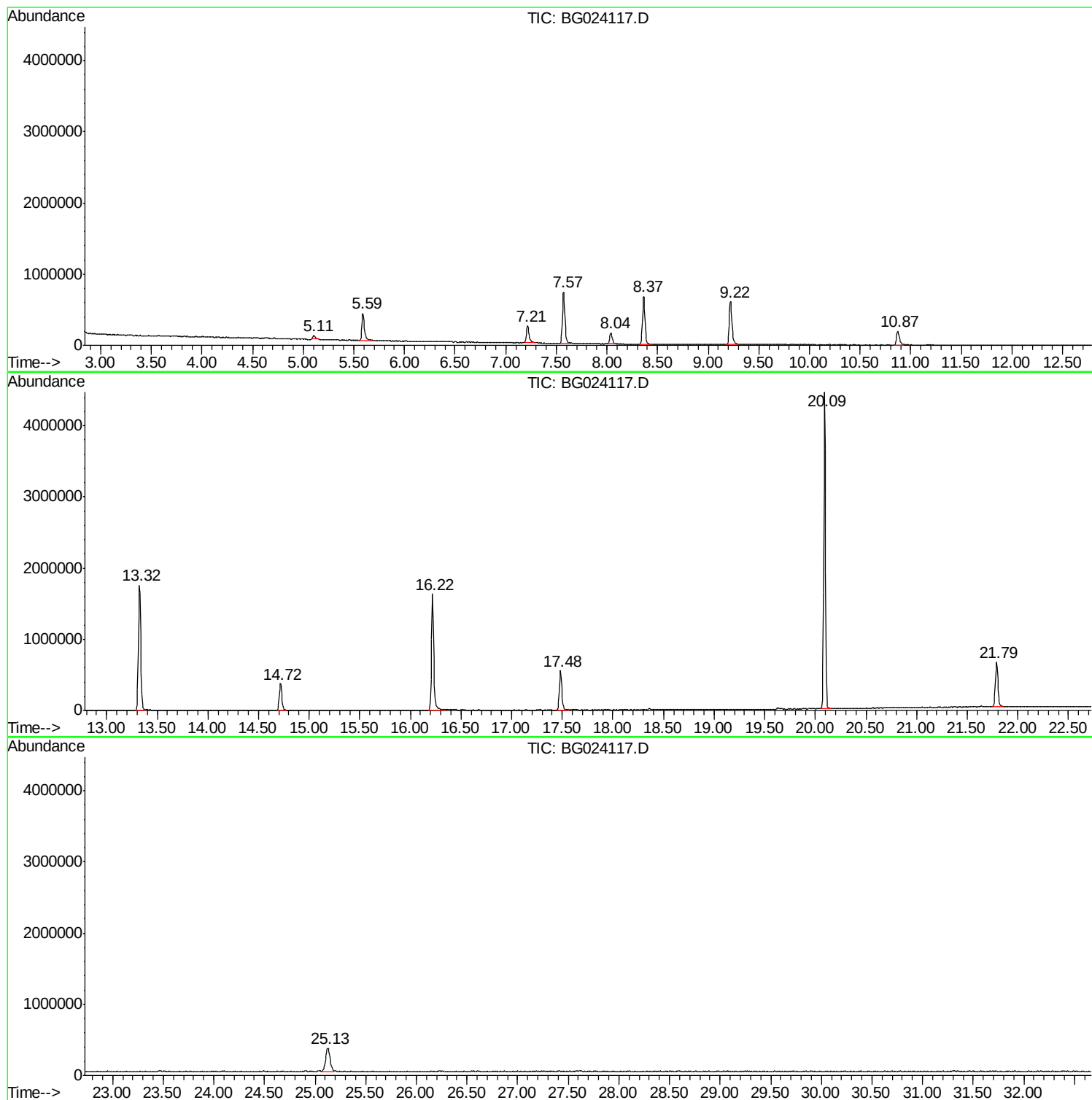
Sum of corrected areas: 20397160

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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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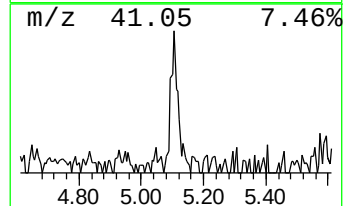
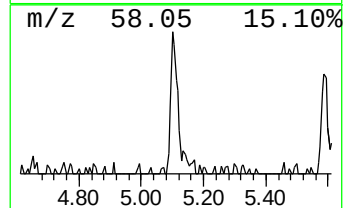
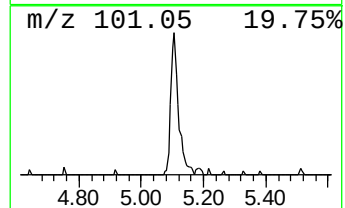
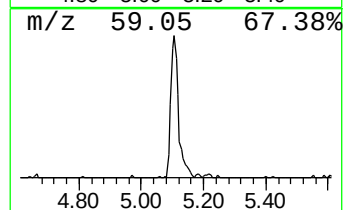
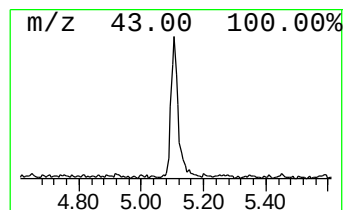
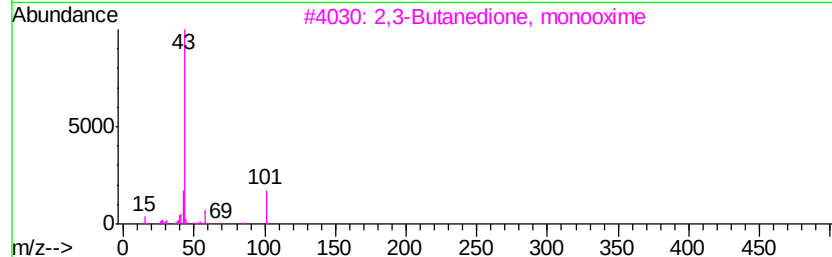
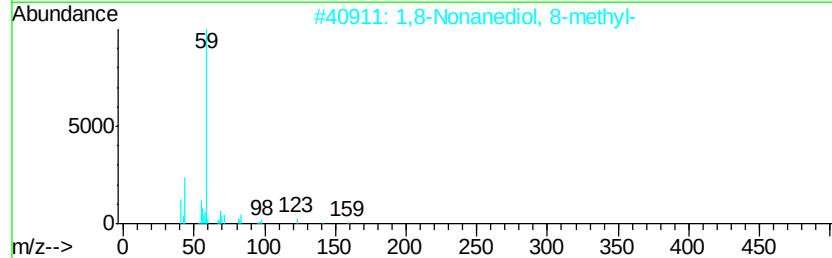
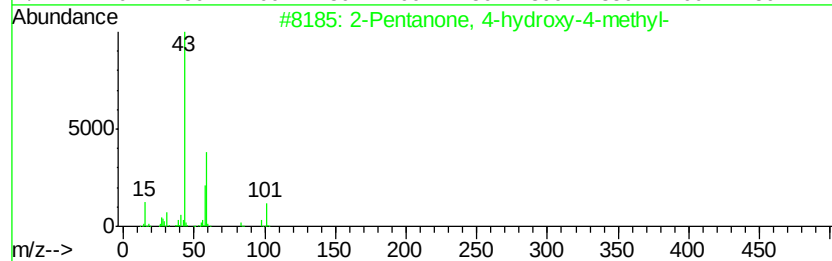
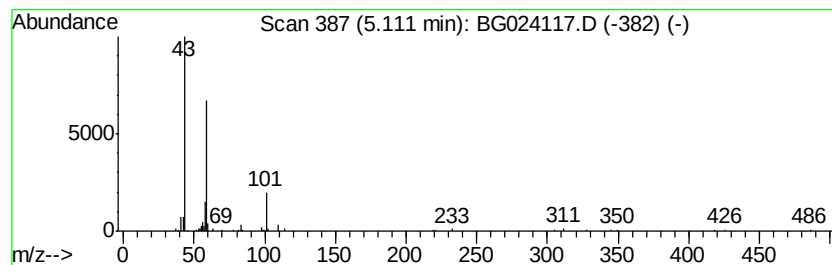
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	6.39 ng	84382	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			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	12
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			4-Ethyl-3-octanol	158	C10H22O	019781-26-1	9



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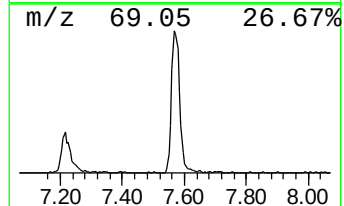
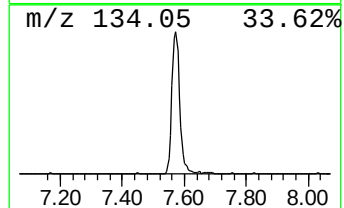
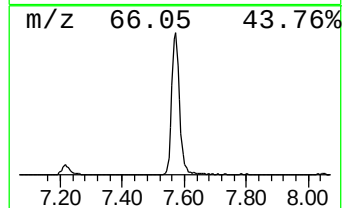
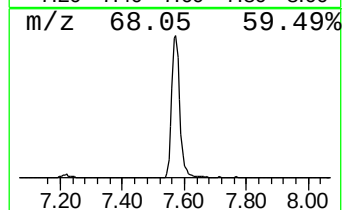
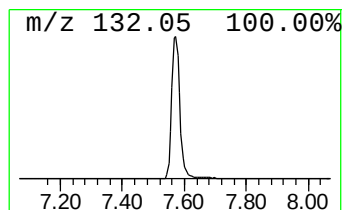
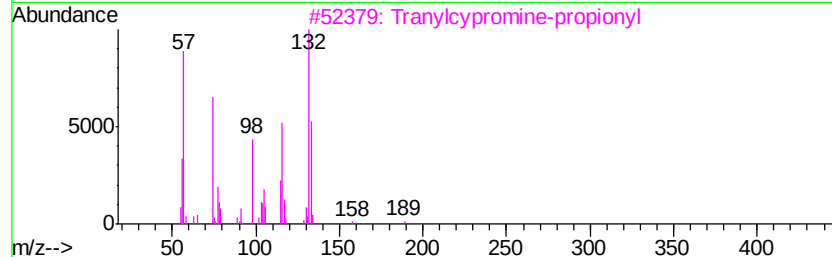
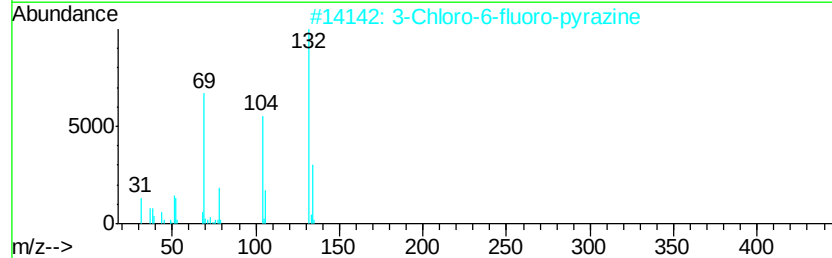
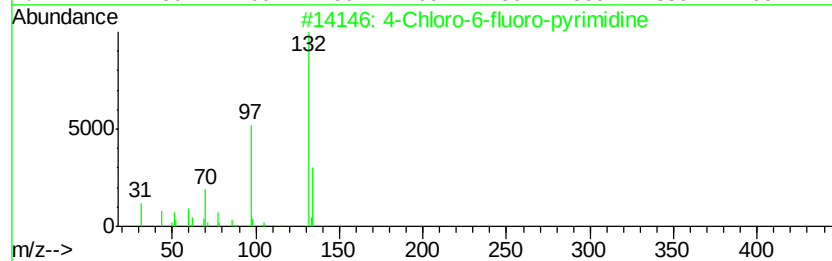
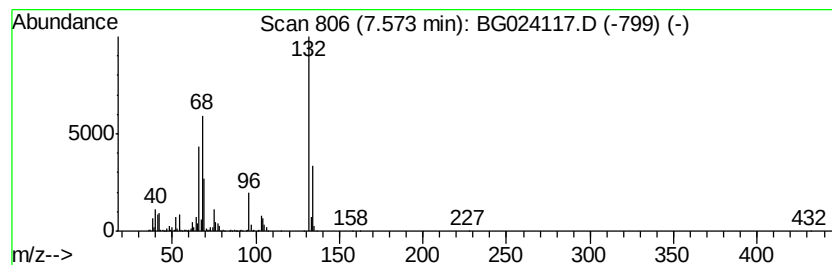
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	100.41 ng	1326110	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10
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...	5.11	6.4	ng	84382	1	8.04	264136	20.0
unknown7.57	7.57	100.4	ng	1326110	1	8.04	264136	20.0