

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042015\
 Data File : BG016551.D
 Acq On : 20 Apr 2015 15:12
 Operator : TP/IZ
 Sample : PB82873BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB82873BL

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-BG040615.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	2.742	4	9	18	rVB2	32999	57036	1.70%	0.300%
2	4.764	348	353	364	rVB	133837	182875	5.46%	0.963%
3	5.245	428	435	454	rBV	1119680	1546458	46.20%	8.142%
4	6.844	700	707	721	rBV	1095478	1653475	49.40%	8.705%
5	7.190	759	766	783	rBV	1123995	1695245	50.65%	8.925%
6	7.654	838	845	852	rBV	225402	353721	10.57%	1.862%
7	7.972	892	899	909	rBV	811991	1266664	37.84%	6.669%
8	8.812	1034	1042	1050	rBV	588182	963840	28.80%	5.075%
9	10.439	1308	1319	1329	rBV	315012	530108	15.84%	2.791%
10	12.913	1733	1740	1753	rBV	1629628	2316844	69.22%	12.198%
11	14.294	1968	1975	1984	rBV2	485173	726388	21.70%	3.824%
12	15.786	2223	2229	2242	rBV	1007777	1481387	44.26%	7.799%
13	17.037	2435	2442	2454	rBV2	628136	907516	27.11%	4.778%
14	19.664	2883	2889	2899	rBV	2746849	3347063	100.00%	17.622%
15	21.215	3147	3153	3162	rBV	791670	989861	29.57%	5.212%
16	23.436	3523	3531	3540	rBV2	517502	975042	29.13%	5.134%

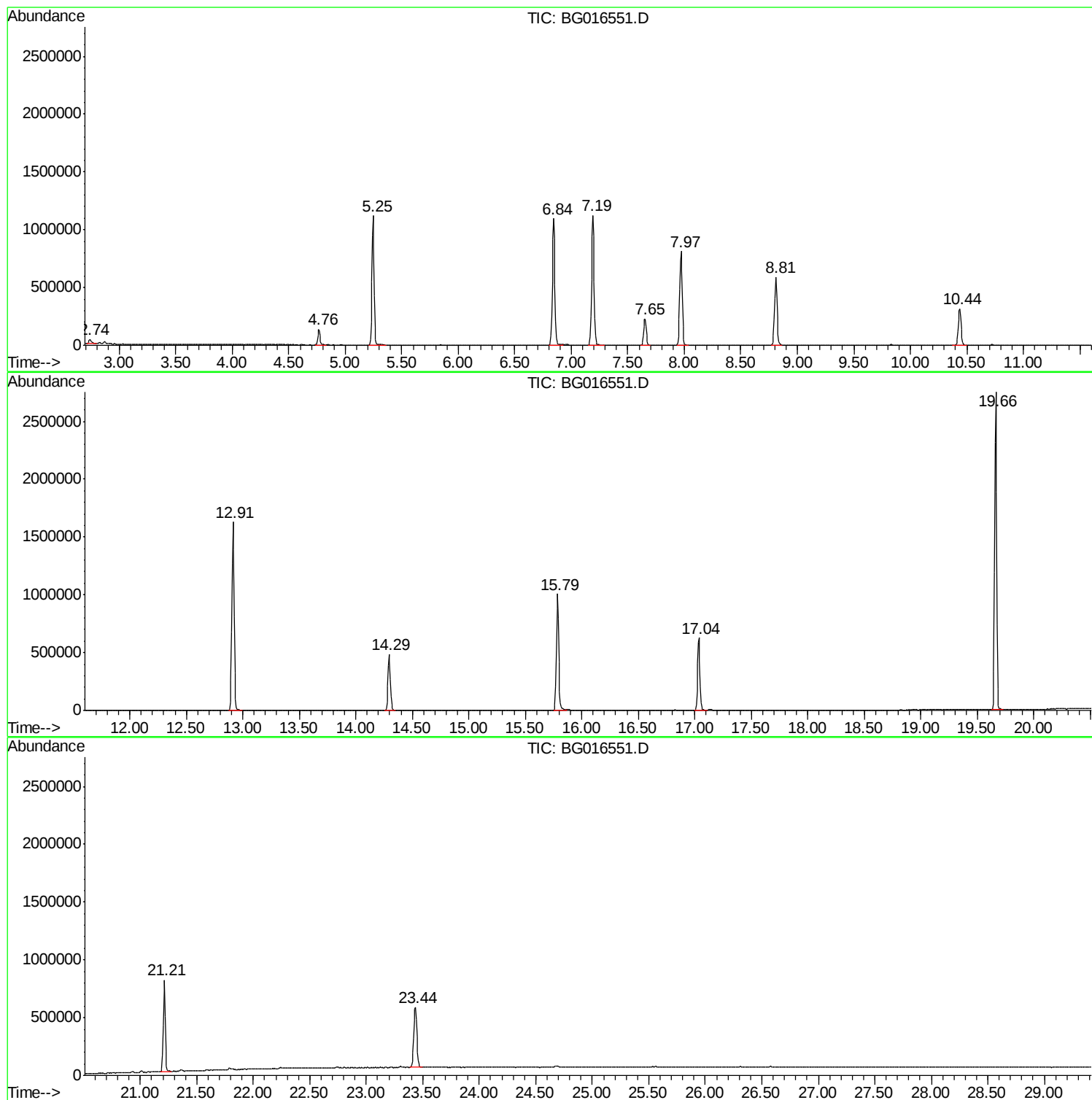
Sum of corrected areas: 18993523

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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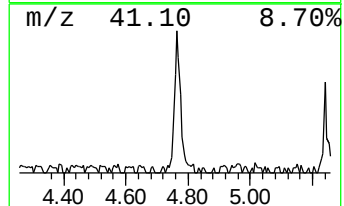
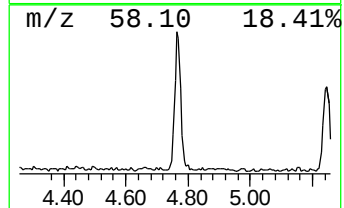
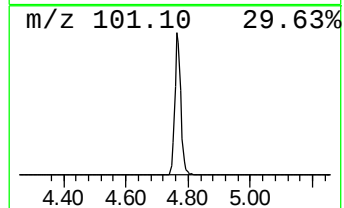
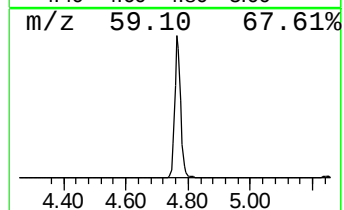
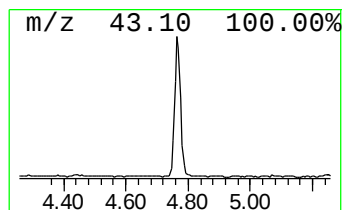
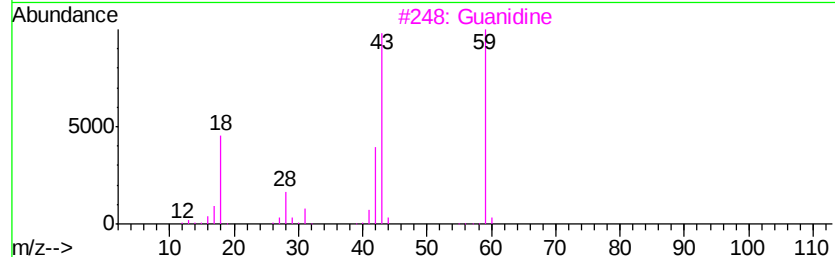
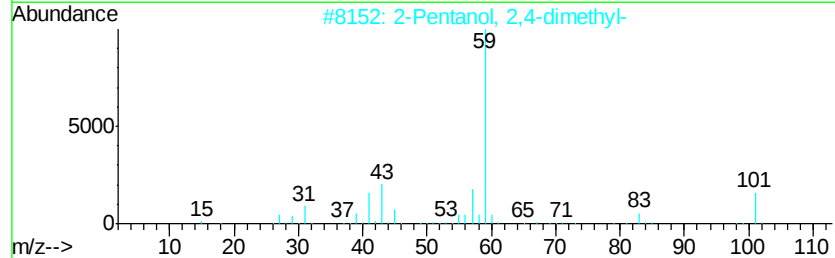
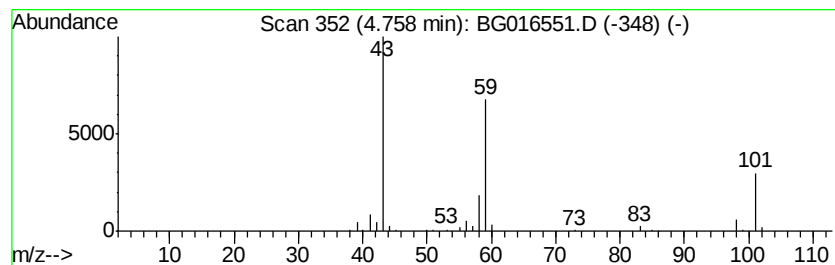
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.76	10.34 ng	182875	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40
3		Guanidine	59	CH5N3	000113-00-8	32
4		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17



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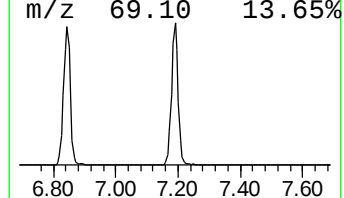
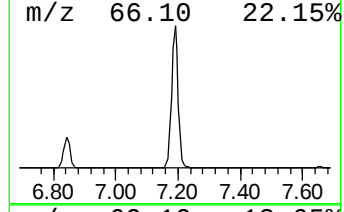
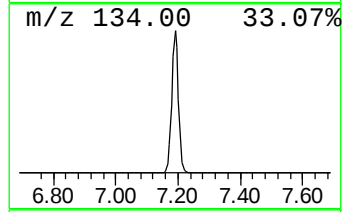
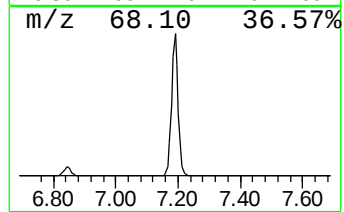
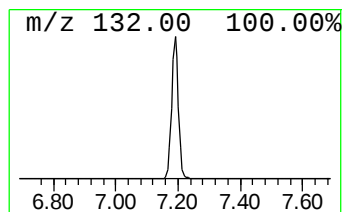
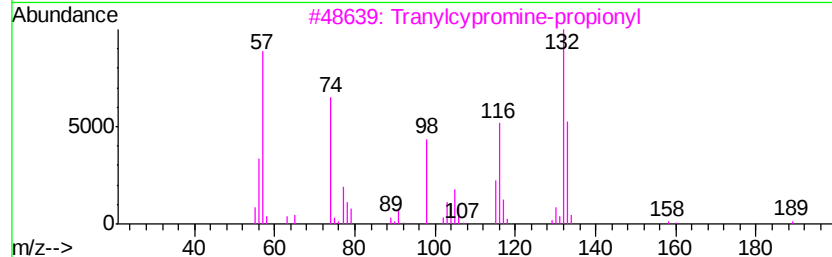
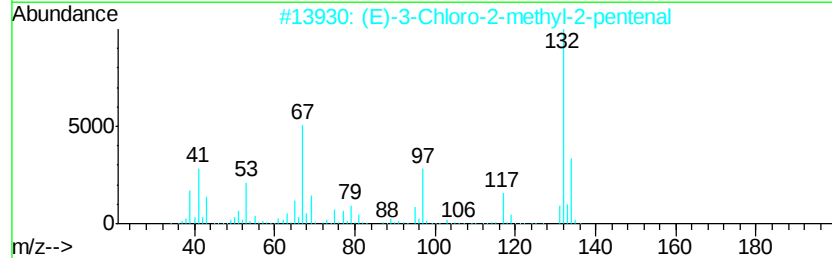
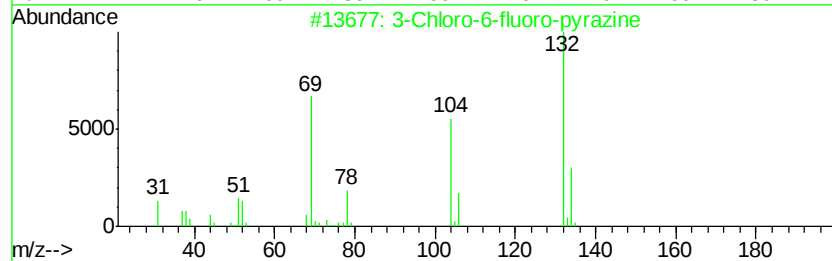
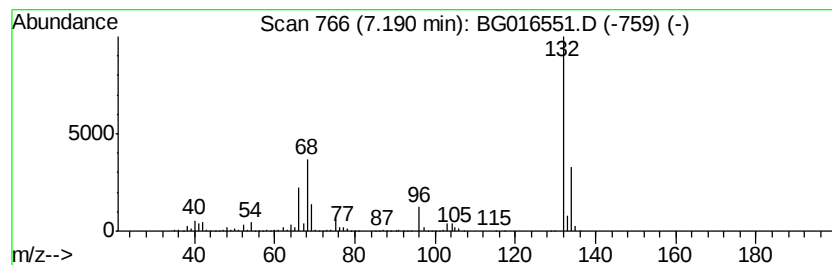
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Peak Number 3 unknown7.19 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	95.85 ng	1695250	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.76	10.3	ng	182875	1	7.65	353721	20.0
unknown7.19	7.19	95.8	ng	1695250	1	7.65	353721	20.0