

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032017\
 Data File : BG026346.D
 Acq On : 21 Mar 2017 3:37
 Operator : SJ/MA
 Sample : PB97680BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB97680BL

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-BG032017.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	344	350	360	rBV	268900	414495	5.24%	1.102%
2	5.620	424	431	446	rBV	1769964	2880036	36.39%	7.658%
3	7.206	694	701	715	rBV	1655395	2824698	35.69%	7.510%
4	7.577	756	764	776	rBV	1755215	3065886	38.73%	8.152%
5	8.047	837	844	855	rVB	334214	559256	7.07%	1.487%
6	8.370	890	899	910	rBV	1412380	2420937	30.59%	6.437%
7	9.204	1032	1041	1052	rBV	978696	1750108	22.11%	4.653%
8	10.843	1314	1320	1331	rBV	460797	835003	10.55%	2.220%
9	13.282	1727	1735	1746	rBV	3127226	4747864	59.98%	12.624%
10	14.651	1959	1968	1982	rBV2	696939	1153789	14.58%	3.068%
11	16.131	2213	2220	2236	rBV	2273132	3396616	42.91%	9.031%
12	17.383	2427	2433	2446	rBV2	981493	1491098	18.84%	3.965%
13	19.980	2869	2875	2886	rBV	5924157	7915223	100.00%	21.045%
14	21.631	3149	3156	3168	rVB	1300575	2099466	26.52%	5.582%
15	24.809	3687	3697	3709	rBV2	740551	2056061	25.98%	5.467%

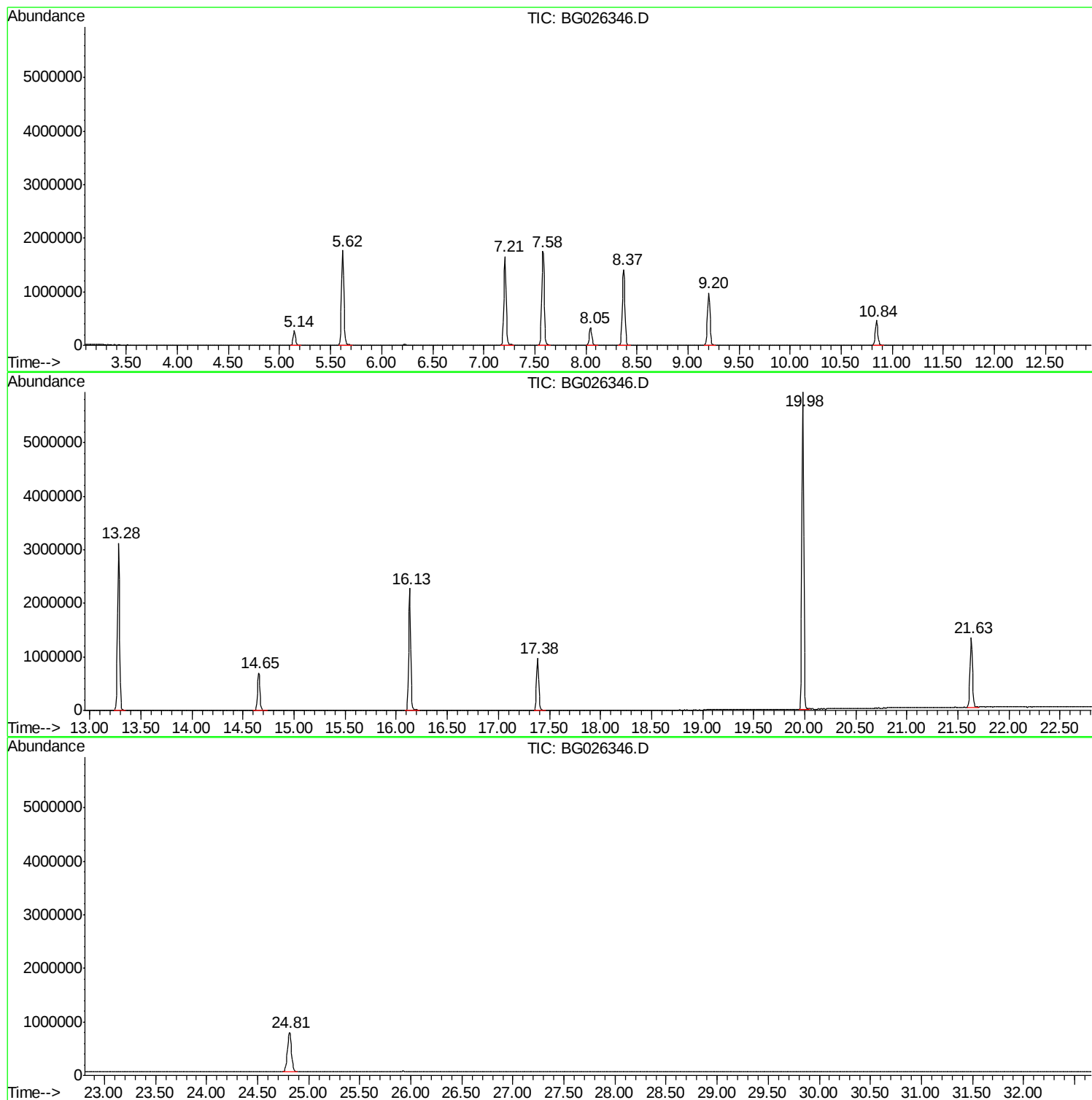
Sum of corrected areas: 37610536

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG032017.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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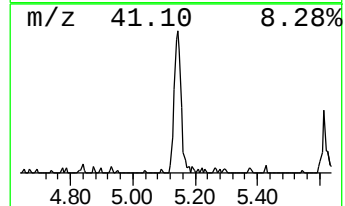
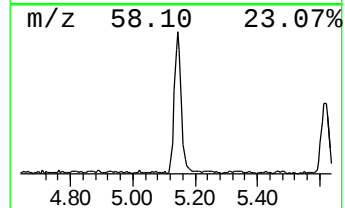
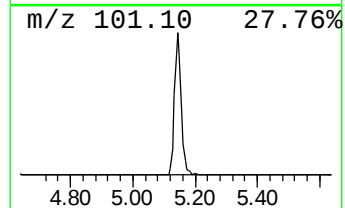
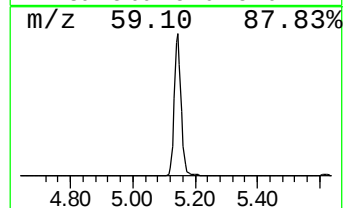
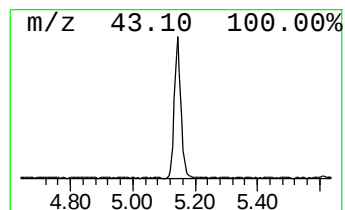
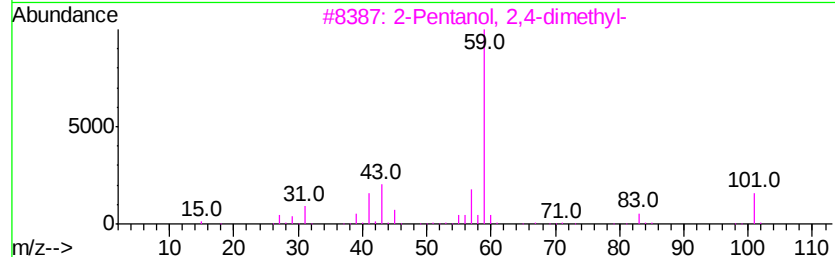
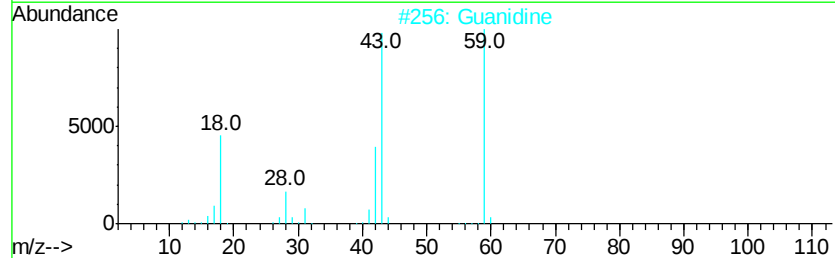
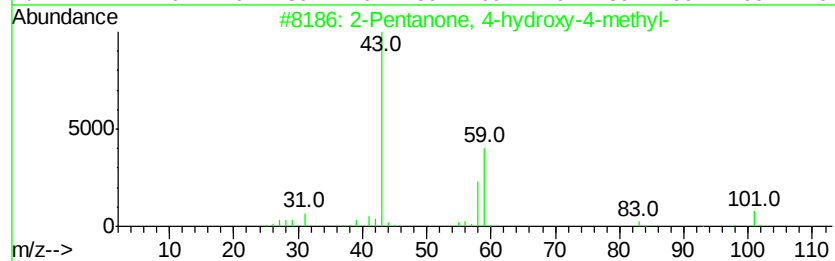
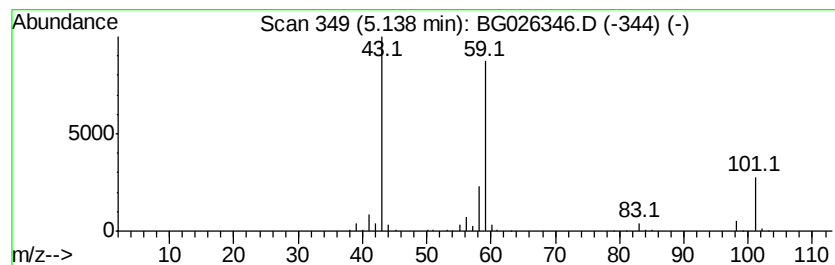
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	14.82 ng	414495	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
2			Guanidine	59	CH5N3	000113-00-8	47
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	36
4			Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	28
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25



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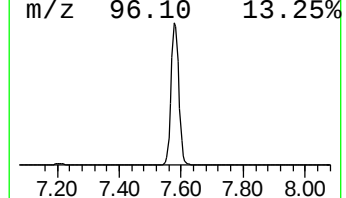
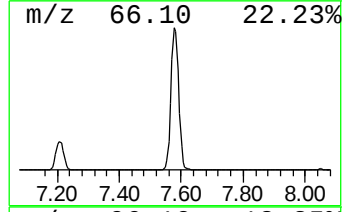
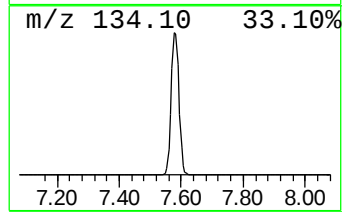
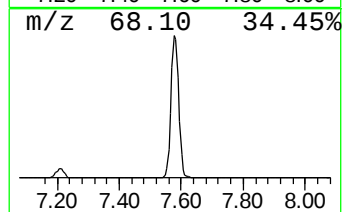
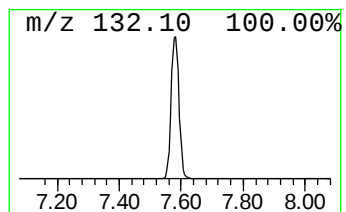
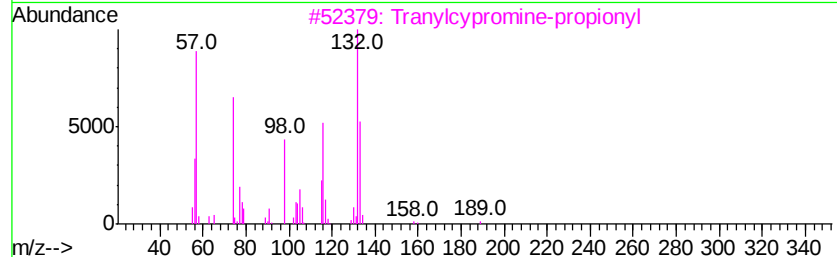
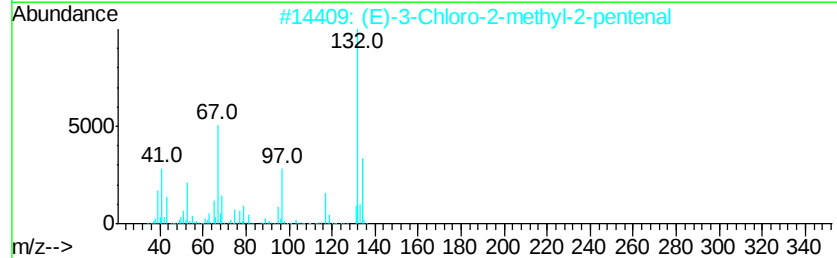
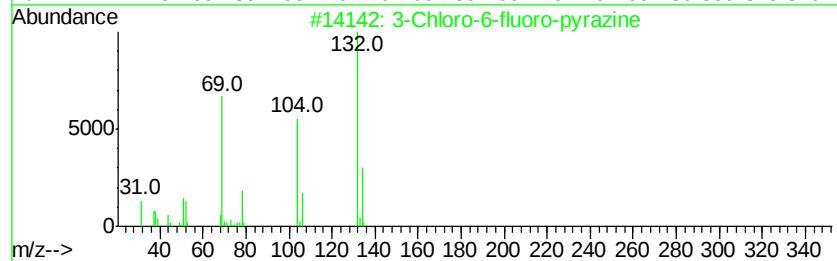
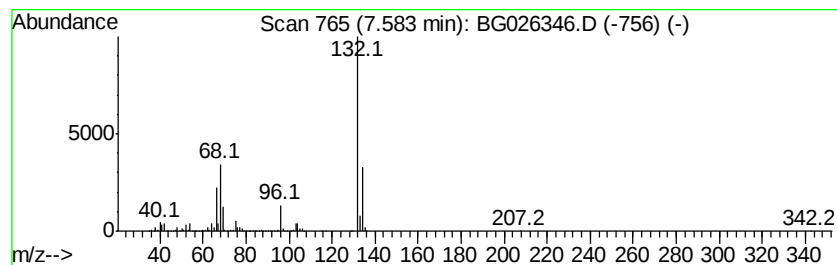
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	109.64 ng	3065890	1,4-Dichlorobenzene-d4	8.05

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	12
3		Tranvlcypramine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	14.8	ng	414495	1	8.05	559256	20.0
unknown7.58	7.58	109.6	ng	3065890	1	8.05	559256	20.0