

Data Path : Z:\HPCHEM1\BNA G\DATA\BG040518\
 Data File : BG033608.D
 Acq On : 6 Apr 2018 00:49
 Operator : SJ/JU
 Sample : PB108007BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB108007BL

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-BG031918.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	5.768	416	422	431	rBV2	32458	56451	1.44%	0.332%
2	6.285	503	510	529	rBV	592062	1207466	30.80%	7.091%
3	7.971	790	797	826	rBV	641023	1423232	36.31%	8.358%
4	8.371	858	865	885	rBV	618532	1292436	32.97%	7.590%
5	8.858	941	948	963	rBV2	75613	185076	4.72%	1.087%
6	9.193	998	1005	1019	rBV	467250	940937	24.01%	5.526%
7	10.057	1146	1152	1172	rBV	352240	824531	21.04%	4.842%
8	11.743	1430	1439	1455	rBV	129290	294447	7.51%	1.729%
9	14.105	1833	1841	1861	rBV	1276045	2193615	55.96%	12.883%
10	15.474	2067	2074	2087	rBV2	285098	522420	13.33%	3.068%
11	16.949	2317	2325	2348	rBV	1153286	1944764	49.61%	11.421%
12	18.206	2530	2539	2553	rBV	411055	726029	18.52%	4.264%
13	20.727	2962	2968	2980	rBV	2794258	3919730	100.00%	23.020%
14	22.554	3272	3279	3291	rBV2	409310	806076	20.56%	4.734%
15	26.320	3909	3920	3937	rBB2	192860	690578	17.62%	4.056%

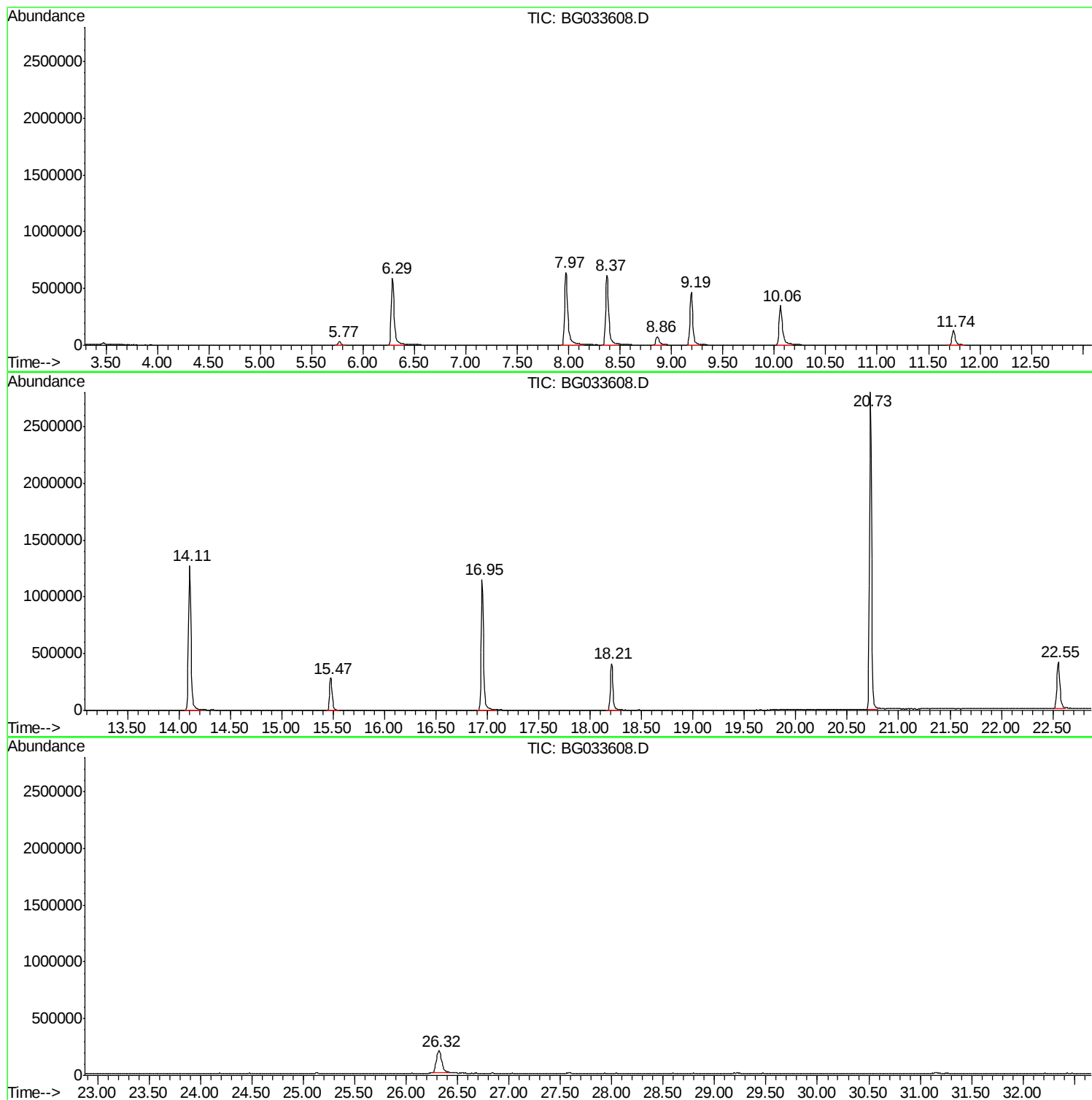
Sum of corrected areas: 17027788

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.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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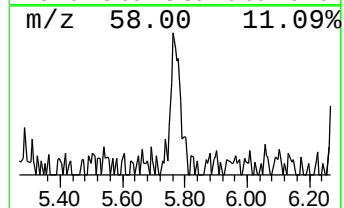
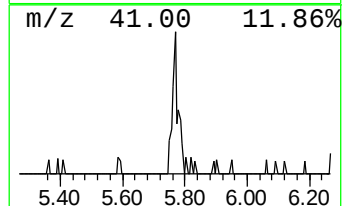
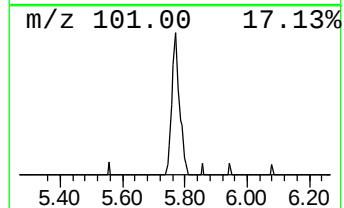
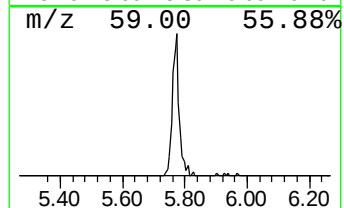
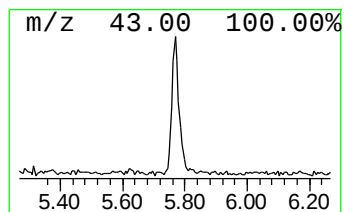
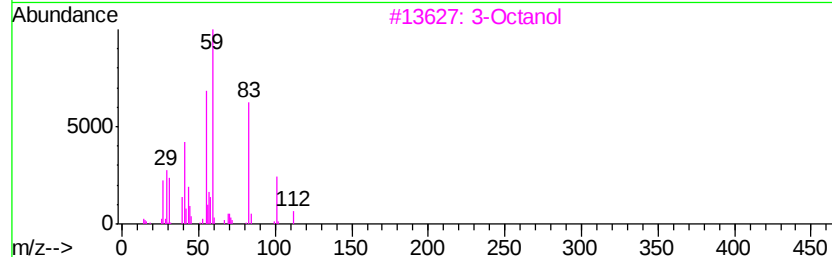
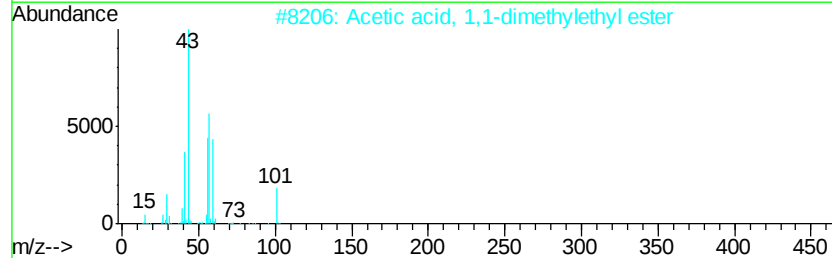
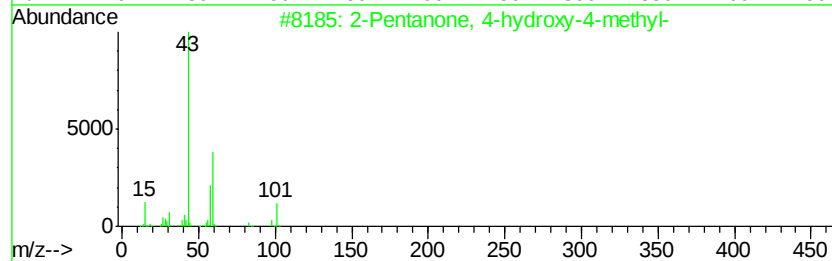
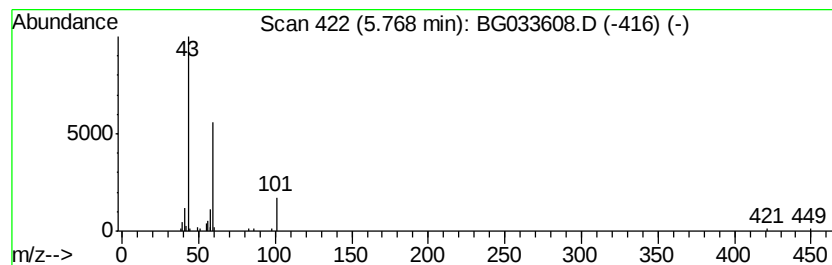
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.77	6.10 ng	56451	1,4-Dichlorobenzene-d4	8.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38	
3	3-Octanol	130	C8H18O	000589-98-0	33	
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28	
5	3-Pentanol	88	C5H12O	000584-02-1	28	



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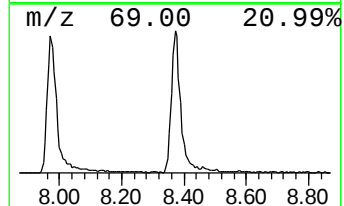
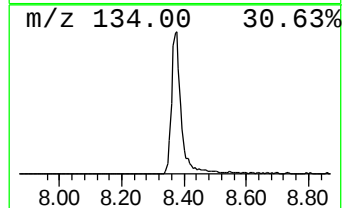
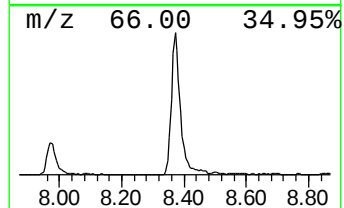
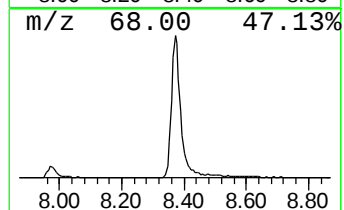
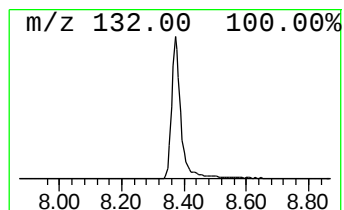
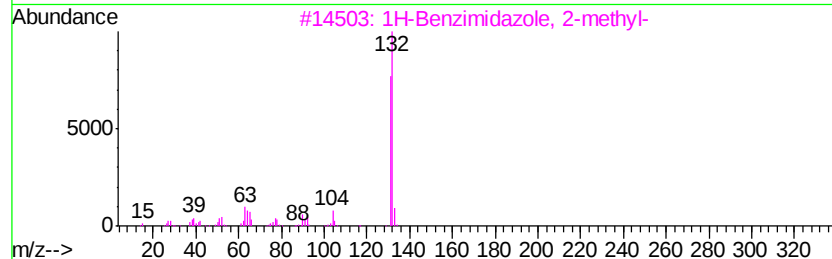
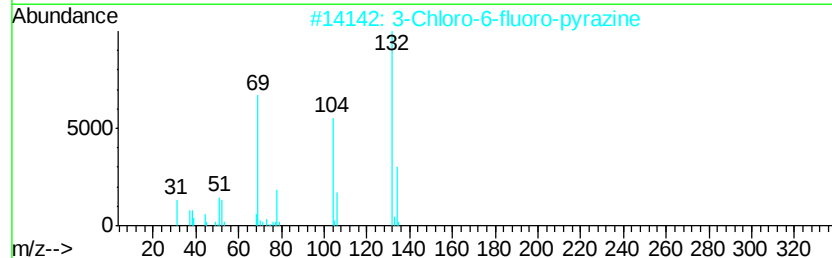
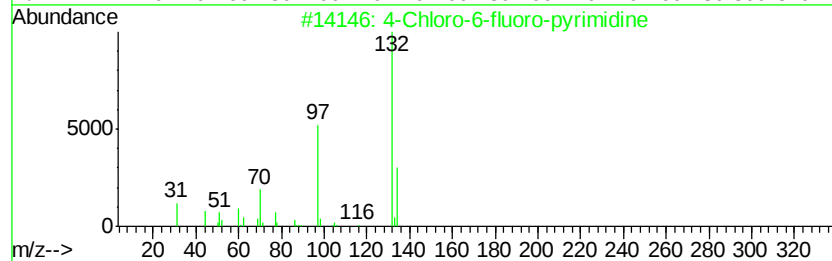
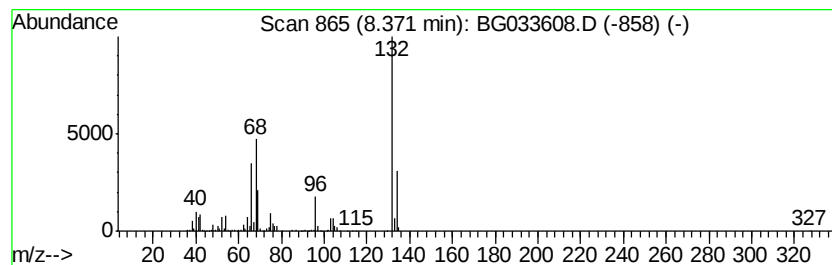
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Peak Number 2 unknown8.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.37	139.67 ng	1292440	1,4-Dichlorobenzene-d4	8.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.77	6.1	ng	56451	1	8.86	185076	20.0
unknown8.37	8.37	139.7	ng	1292440	1	8.86	185076	20.0