

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011918\
 Data File : BG032153.D
 Acq On : 19 Jan 2018 15:58
 Operator : SJ/JU
 Sample : PB105755BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB105755BL

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-BG011218.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	5.658	397	403	411	rBV	26867	45235	1.07%	0.280%
2	6.169	482	490	505	rBV	550302	934247	22.05%	5.789%
3	7.849	767	776	790	rBV	550214	982460	23.19%	6.087%
4	8.255	837	845	854	rBV	558326	1033285	24.39%	6.402%
5	8.742	921	928	939	rBB	117413	208860	4.93%	1.294%
6	9.077	977	985	996	rBV	450451	839395	19.81%	5.201%
7	9.941	1124	1132	1141	rBV	308815	590778	13.94%	3.661%
8	11.621	1411	1418	1428	rBV	153811	295182	6.97%	1.829%
9	14.007	1816	1824	1835	rBV	1197231	1928876	45.52%	11.951%
10	15.376	2051	2057	2066	rVB2	285982	474864	11.21%	2.942%
11	16.851	2301	2308	2325	rBV	1069096	1682467	39.71%	10.425%
12	18.114	2516	2523	2532	rBV2	462957	732642	17.29%	4.540%
13	20.646	2948	2954	2965	rBV	2959419	4237075	100.00%	26.253%
14	22.444	3253	3260	3268	rBV2	578588	1069994	25.25%	6.630%
15	26.146	3873	3890	3904	rBV2	330150	1083876	25.58%	6.716%

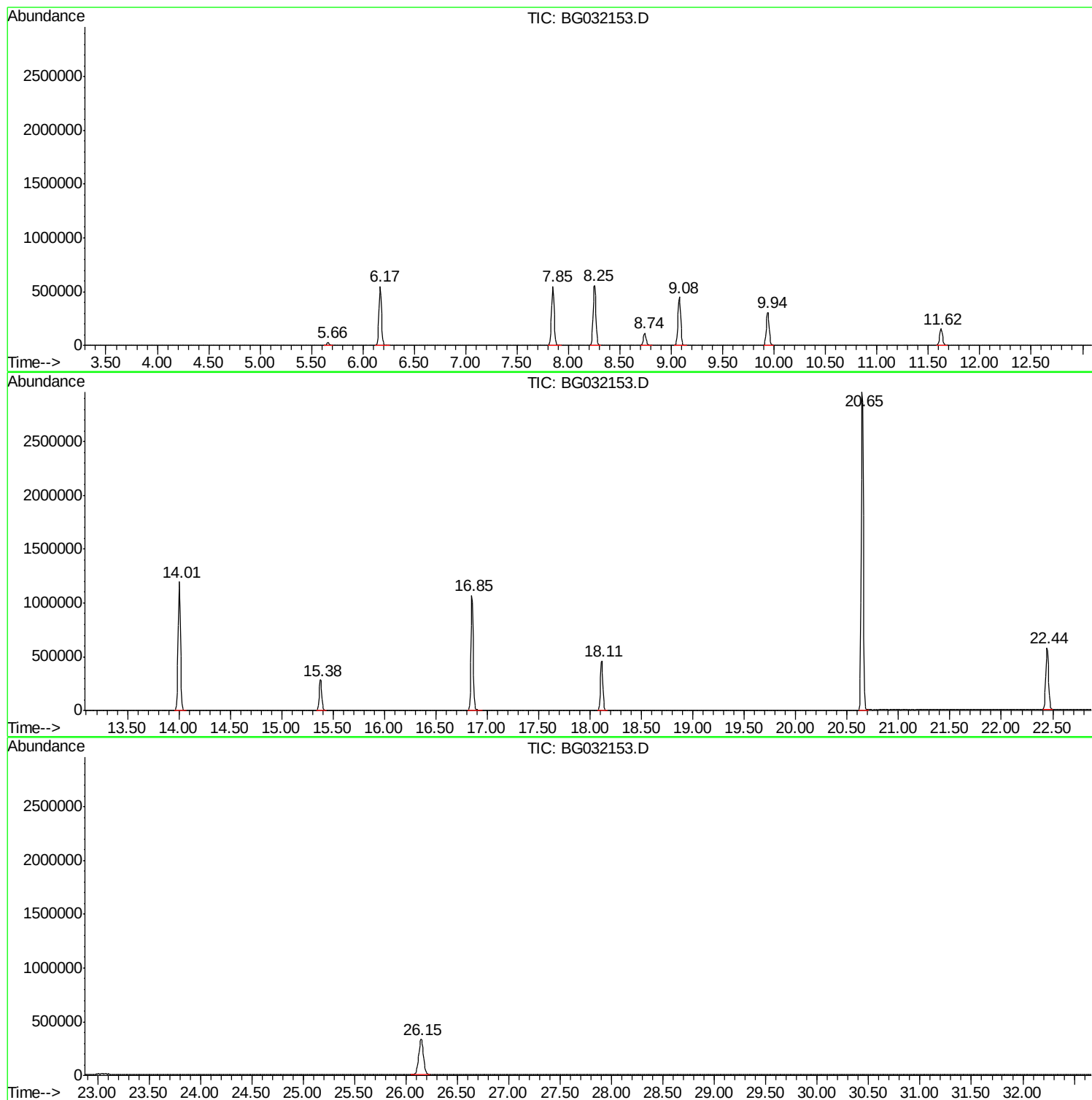
Sum of corrected areas: 16139236

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.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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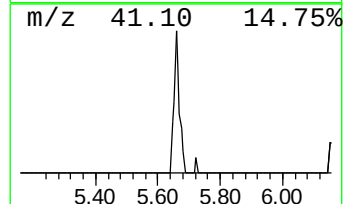
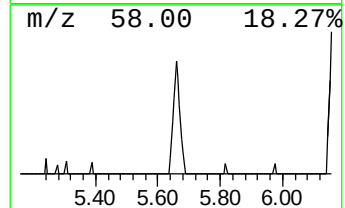
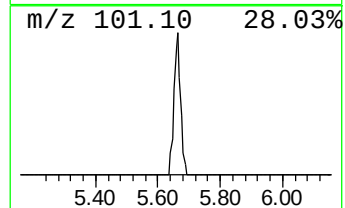
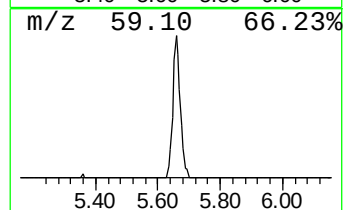
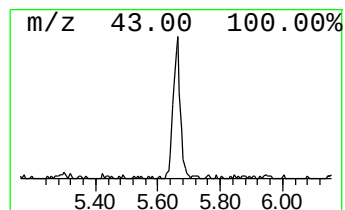
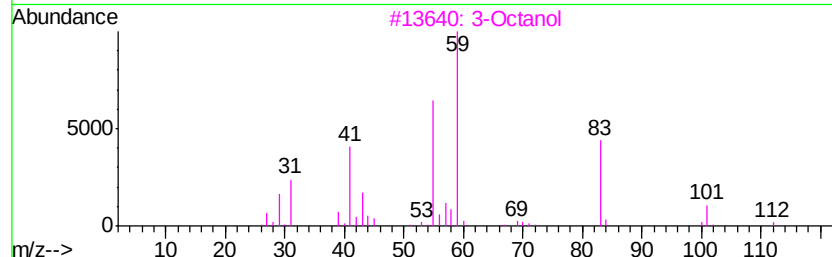
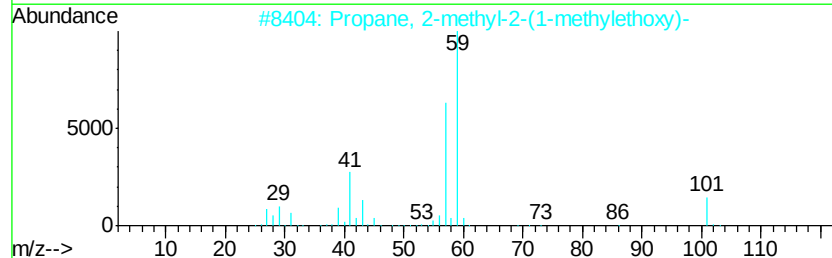
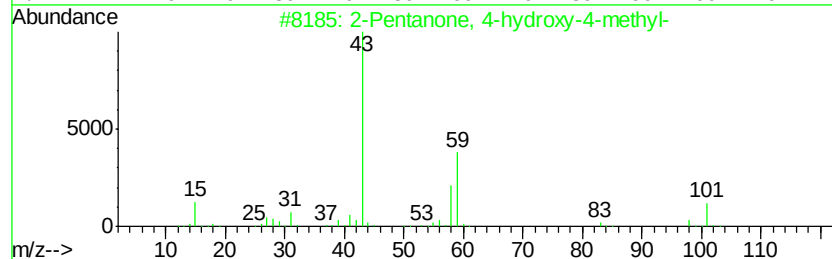
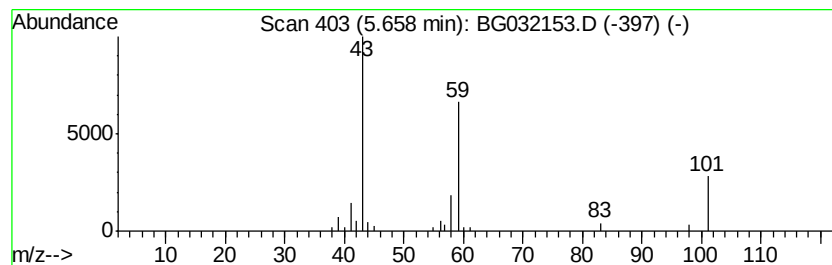
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.66	4.33 ng	45235	1,4-Dichlorobenzene-d4	8.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	39
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	38
3			3-Octanol	130	C8H18O	000589-98-0	25
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
5			3-Pentanol	88	C5H12O	000584-02-1	12



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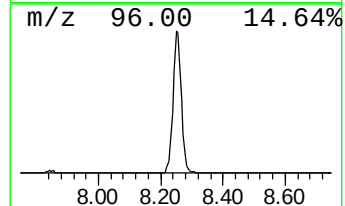
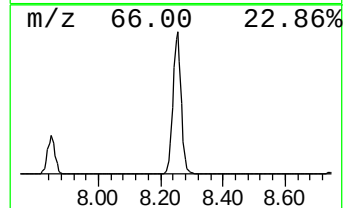
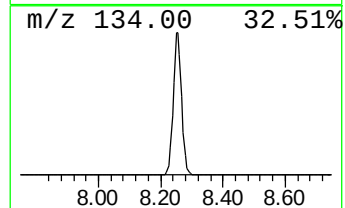
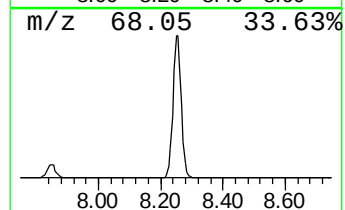
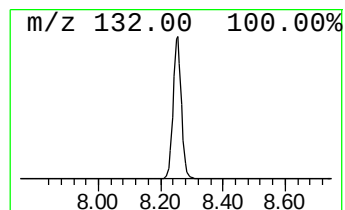
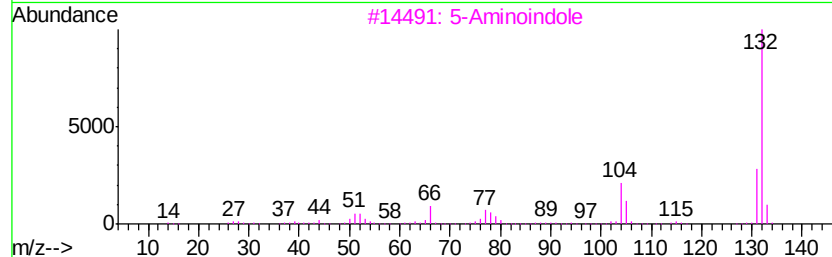
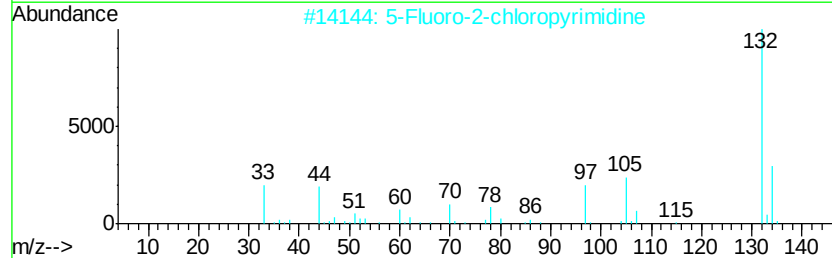
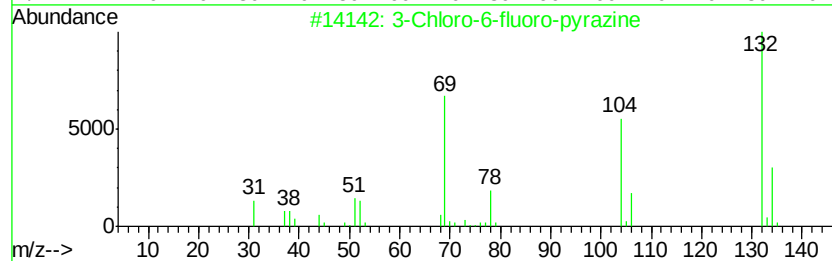
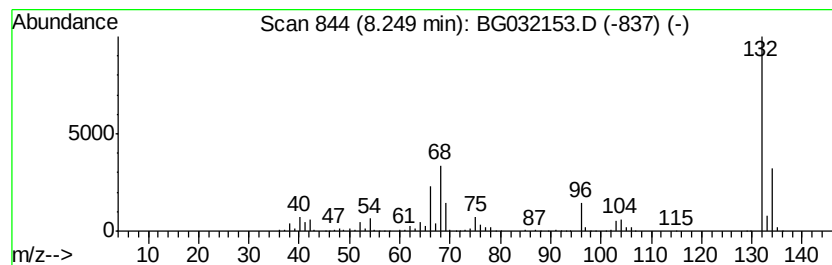
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Peak Number 2 unknown8.25 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.25	98.95 ng	1033290	1,4-Dichlorobenzene-d4	8.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	35
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.66	4.3	ng	45235	1	8.74	208860	20.0
unknown8.25	8.25	99.0	ng	1033290	1	8.74	208860	20.0