

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041618\
 Data File : BG033826.D
 Acq On : 16 Apr 2018 21:15
 Operator : SJ/JU
 Sample : PB108262BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB108262BL

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	5.757	416	420	432	rBV2	39115	84478	2.16%	0.633%
2	6.286	504	510	530	rBV	167238	598681	15.31%	4.489%
3	7.989	790	800	819	rBV3	131297	622828	15.93%	4.670%
4	8.371	858	865	885	rBV2	198855	691785	17.69%	5.187%
5	8.859	940	948	956	rBV3	44320	133081	3.40%	0.998%
6	9.182	994	1003	1022	rBV	353687	905297	23.15%	6.787%
7	10.058	1144	1152	1168	rBV2	170797	616736	15.77%	4.624%
8	11.738	1430	1438	1457	rBV	80256	254580	6.51%	1.909%
9	14.094	1831	1839	1860	rBV	852072	1839006	47.03%	13.788%
10	15.469	2066	2073	2085	rBV2	190233	399084	10.21%	2.992%
11	16.944	2318	2324	2348	rBV	492935	1109399	28.37%	8.318%
12	18.201	2532	2538	2551	rBV2	280864	585188	14.97%	4.387%
13	20.722	2961	2967	2984	rBV	2688021	3909878	100.00%	29.314%
14	22.549	3270	3278	3297	rBV	392700	827087	21.15%	6.201%
15	26.315	3907	3919	3938	rBV2	206352	760685	19.46%	5.703%

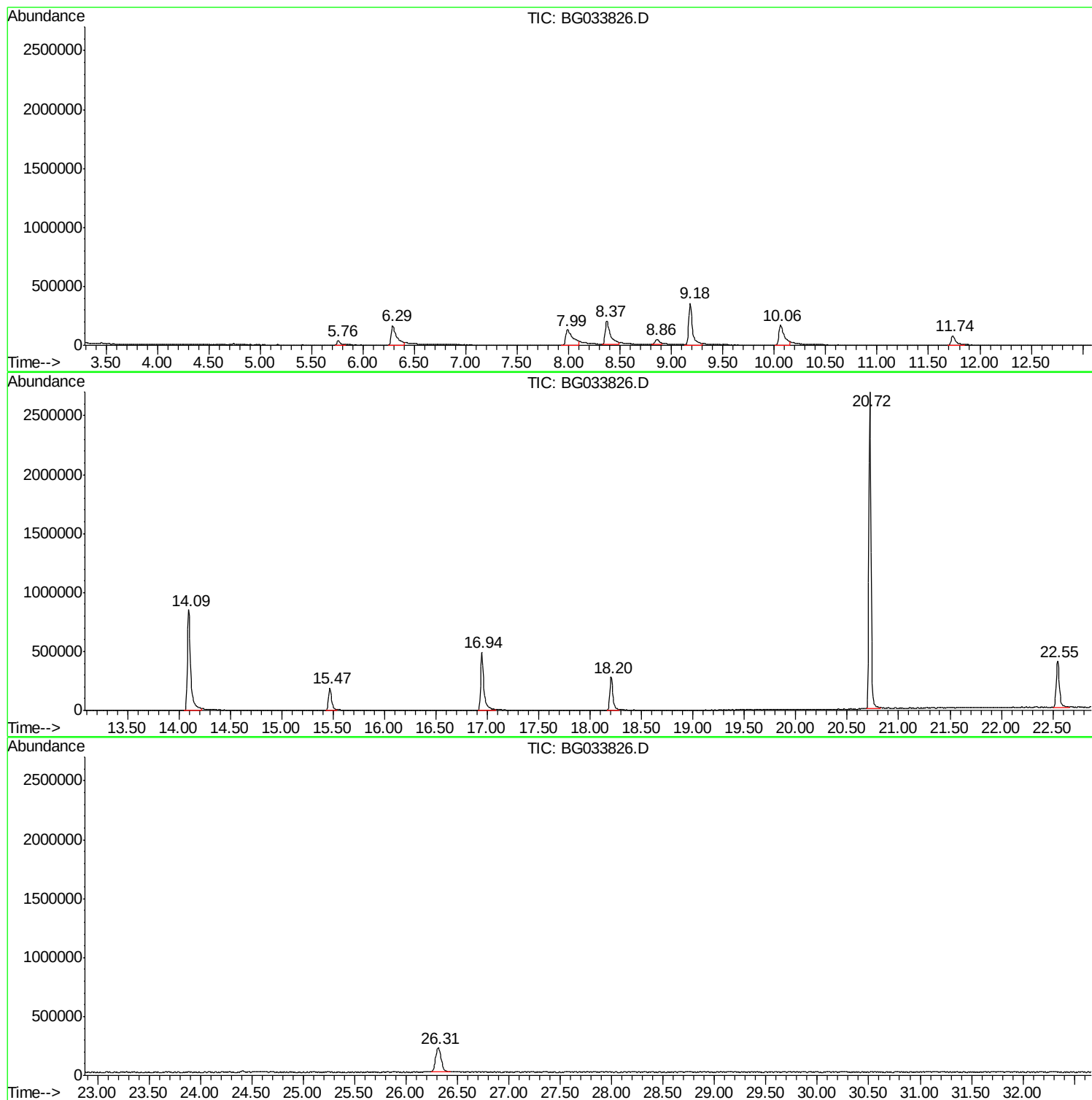
Sum of corrected areas: 13337793

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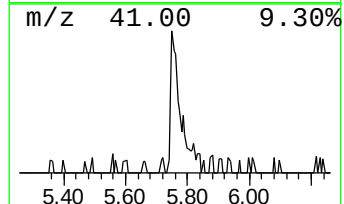
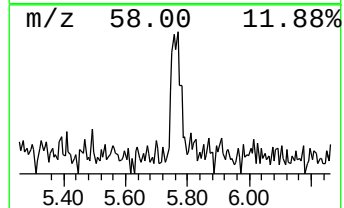
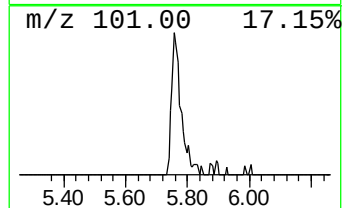
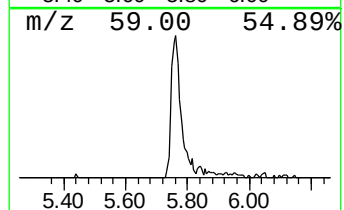
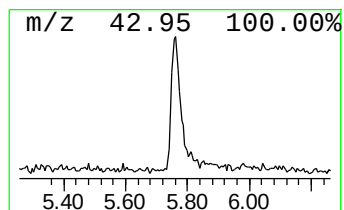
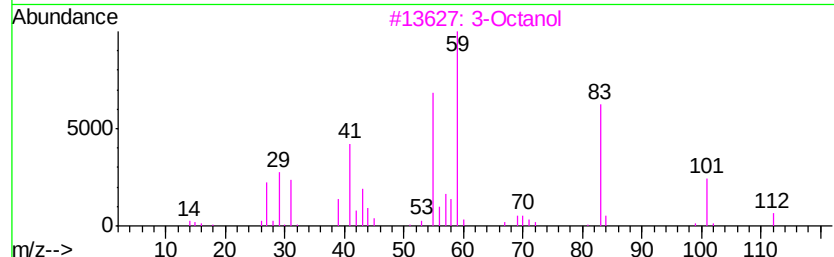
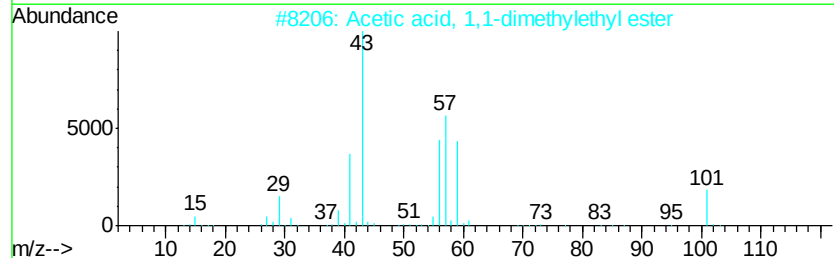
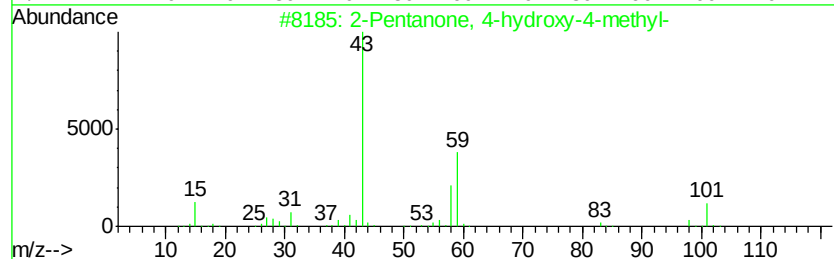
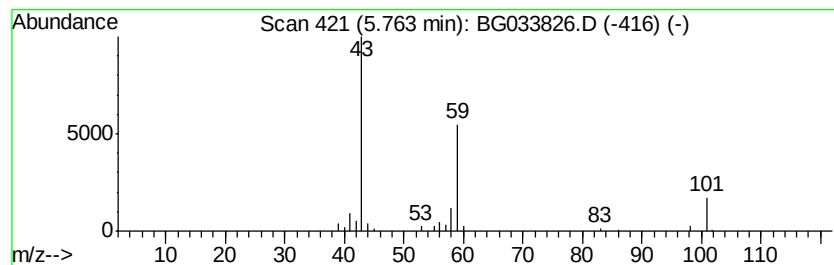
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.76	12.70 ng	84478	1,4-Dichlorobenzene-d4	8.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
3		3-Octanol	130	C8H18O	000589-98-0	33
4		Butane	58	C4H10	000106-97-8	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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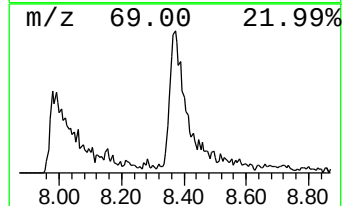
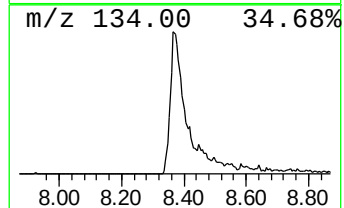
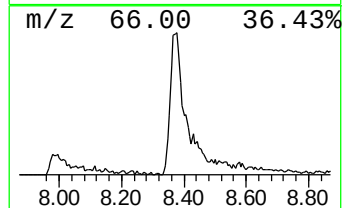
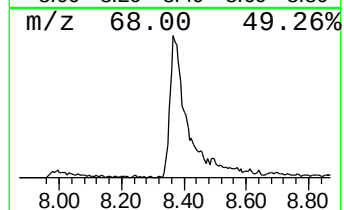
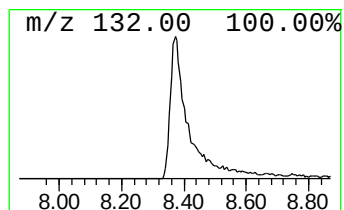
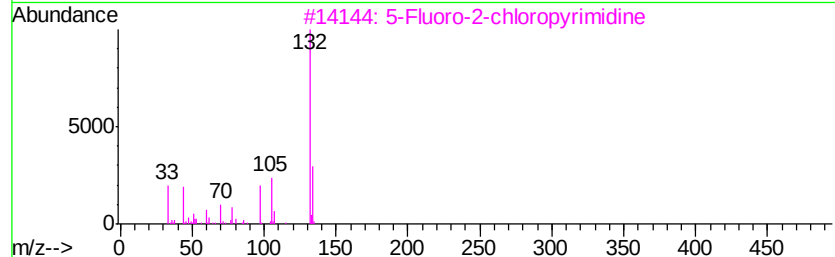
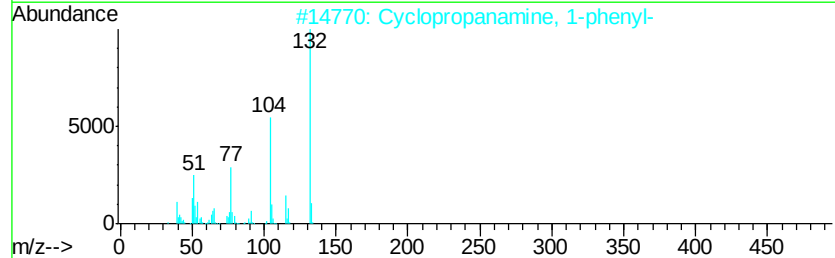
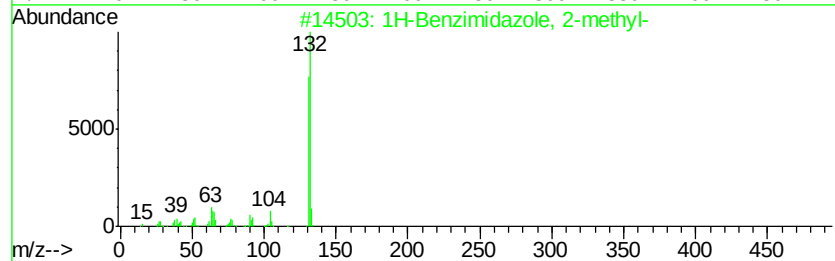
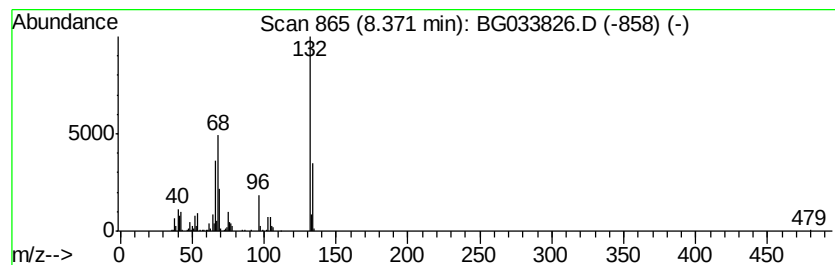
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Peak Number 2 unknown8.37 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.37	103.97 ng	691785	1,4-Dichlorobenzene-d4	8.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	12
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.76	12.7	ng	84478	1	8.86	133081	20.0
unknown8.37	8.37	104.0	ng	691785	1	8.86	133081	20.0