

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030218\
Data File : BG032956.D
Acq On : 2 Mar 2018 21:40
Operator : SJ/JU
Sample : PB107009BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB107009BL

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-BG022118.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	5.825	426	432	438	rBV2	23722	44124	1.25%	0.231%
2	6.330	509	518	531	rBV	975074	1659427	46.90%	8.678%
3	8.010	796	804	821	rBV	975023	1760479	49.75%	9.207%
4	8.428	867	875	884	rBV	934322	1717500	48.54%	8.982%
5	8.915	949	958	966	rBV	185421	332834	9.41%	1.741%
6	9.250	1007	1015	1025	rBV	689619	1293109	36.55%	6.762%
7	10.114	1153	1162	1176	rBV	550338	1042496	29.46%	5.452%
8	11.799	1441	1449	1458	rBV	264707	490123	13.85%	2.563%
9	14.161	1843	1851	1859	rBV	1471286	2373258	67.07%	12.411%
10	15.530	2077	2084	2093	rVB2	394521	653579	18.47%	3.418%
11	16.998	2327	2334	2349	rBV	1080534	1694990	47.90%	8.864%
12	18.261	2541	2549	2556	rBV2	512582	819336	23.16%	4.285%
13	20.776	2971	2977	2991	rBV	2603387	3538351	100.00%	18.504%
14	22.614	3282	3290	3303	rBV2	467081	893311	25.25%	4.672%
15	26.427	3927	3939	3959	rBV2	215632	809169	22.87%	4.232%

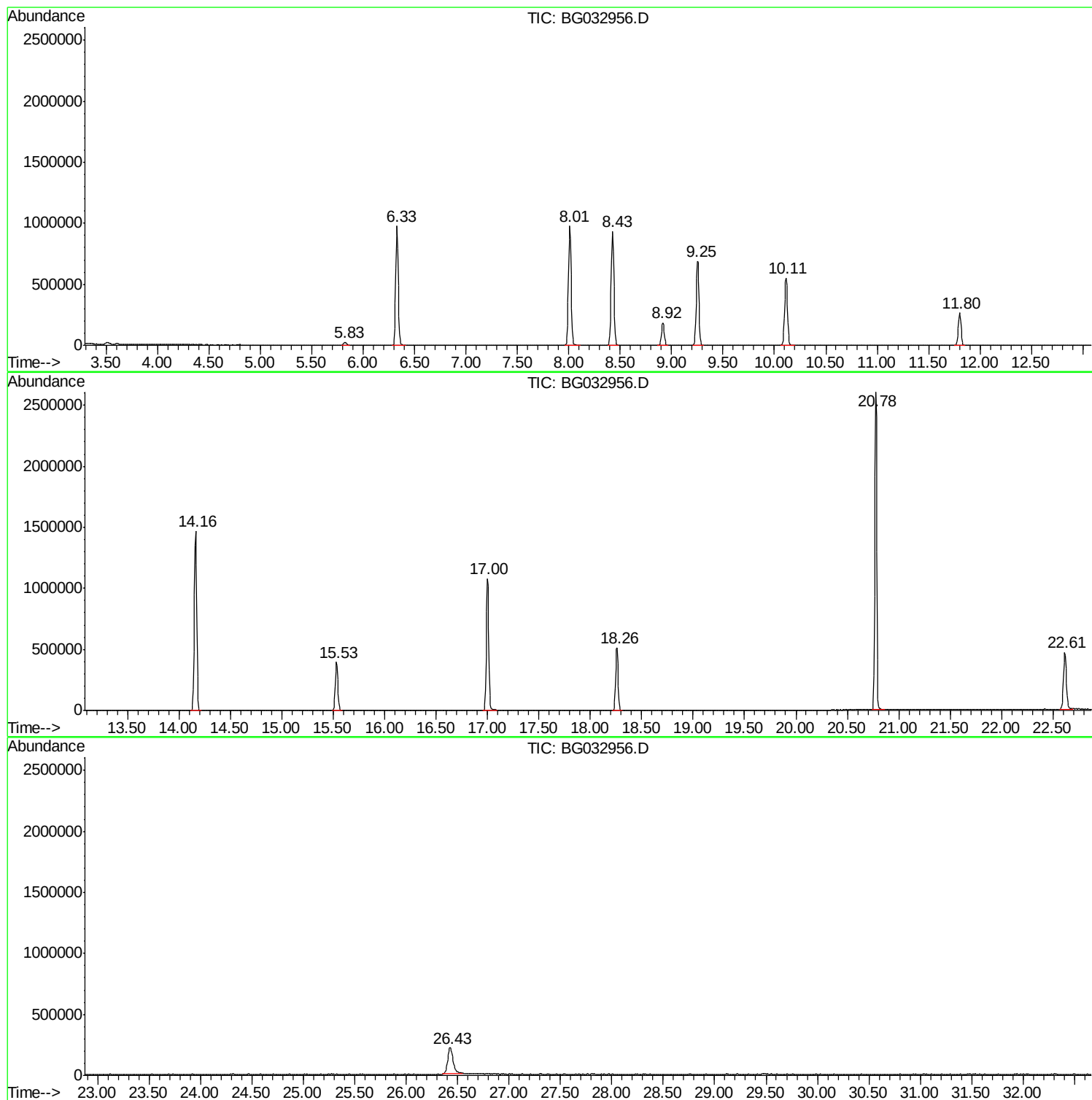
Sum of corrected areas: 19122086

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.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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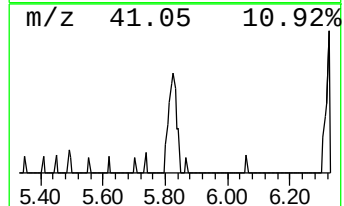
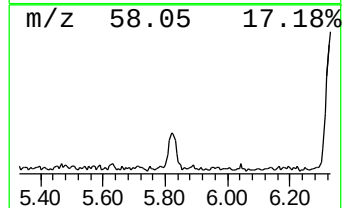
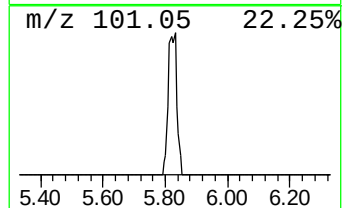
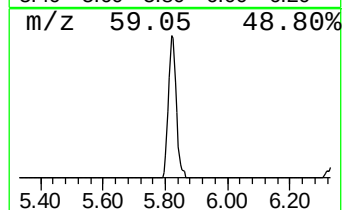
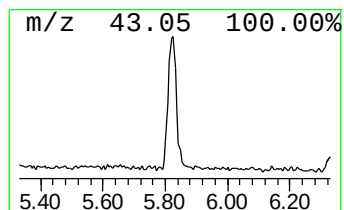
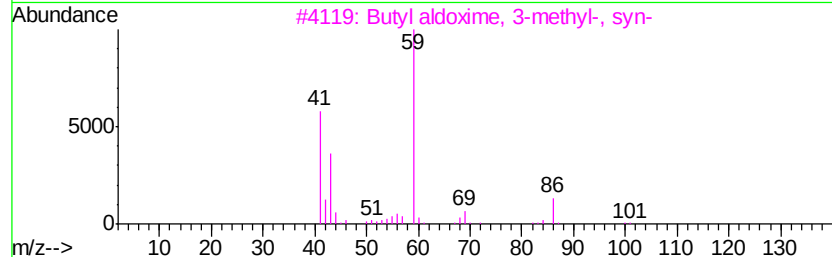
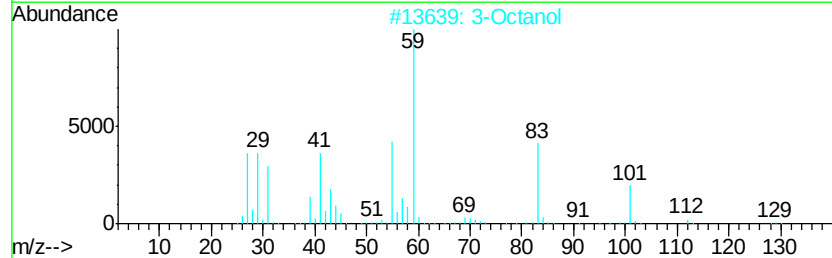
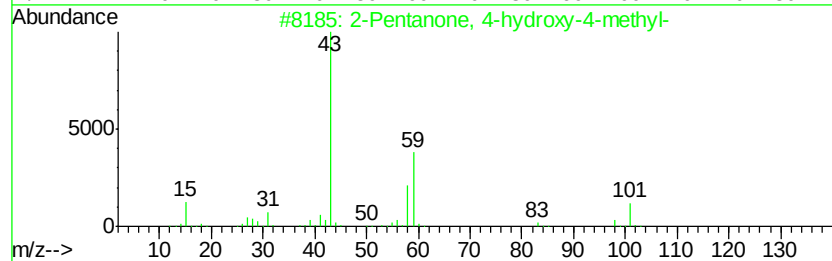
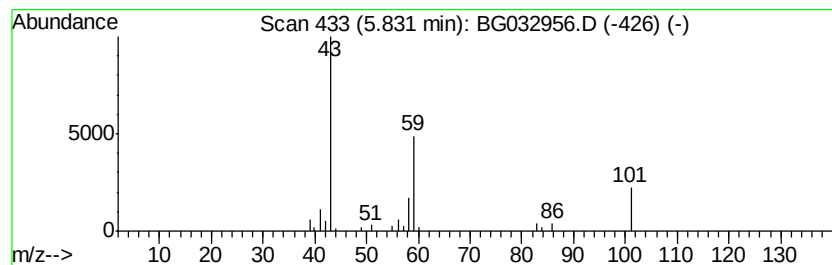
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.83	2.65 ng	44124	1,4-Dichlorobenzene-d4	8.92

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Octanol	130	C8H18O	000589-98-0	9
3		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7
5		Diacetamide	101	C4H7NO2	000625-77-4	7



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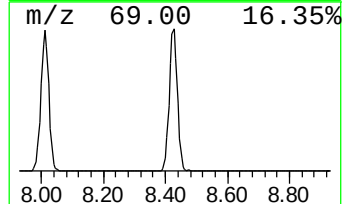
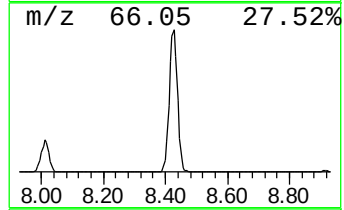
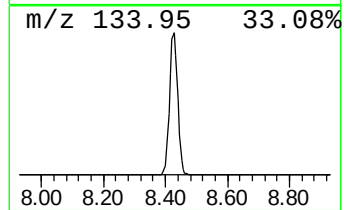
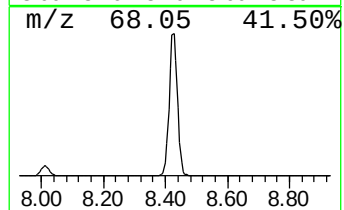
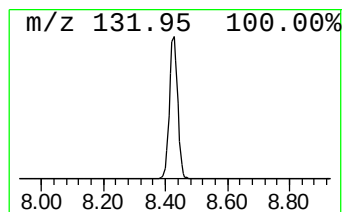
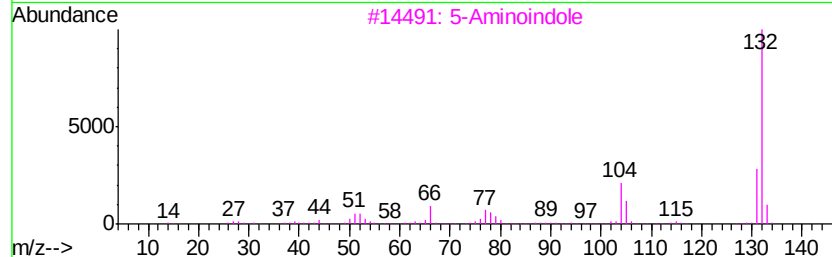
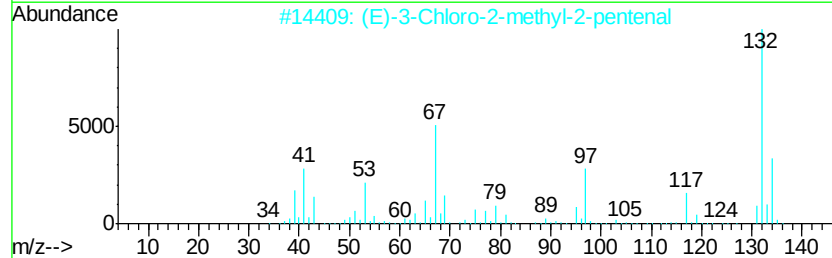
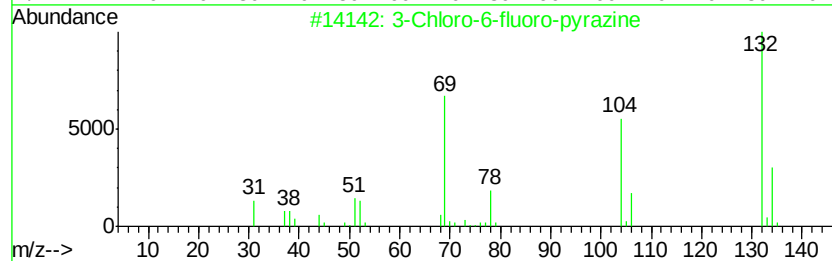
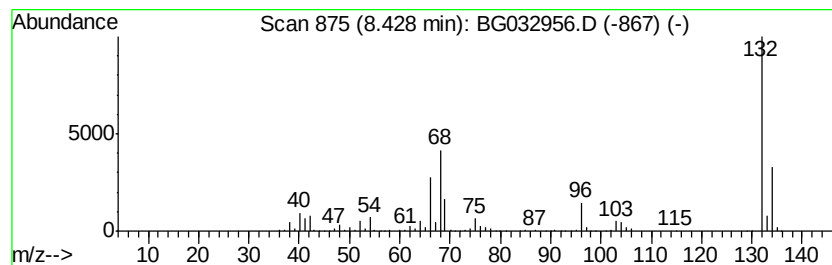
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Peak Number 2 unknown8.43 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.43	103.21 ng	1717500	1,4-Dichlorobenzene-d4	8.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.83	2.6	ng	44124	1	8.92	332834	20.0
unknown8.43	8.43	103.2	ng	1717500	1	8.92	332834	20.0