

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101717\
 Data File : BG029311.D
 Acq On : 17 Oct 2017 21:31
 Operator : SJ/JU
 Sample : PB103373BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103373BL

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-BG101617.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	5.245	325	333	341	rBV	130010	204870	6.74%	1.346%
2	5.721	406	414	422	rBV	410824	661893	21.78%	4.349%
3	7.319	677	686	696	rBV	447879	750865	24.71%	4.934%
4	7.701	742	751	761	rBV	414616	734761	24.18%	4.828%
5	8.171	824	831	838	rVB	194417	338587	11.14%	2.225%
6	8.494	878	886	896	rVB	477403	824424	27.13%	5.417%
7	9.334	1021	1029	1039	rBV	379981	686414	22.59%	4.510%
8	10.985	1299	1310	1318	rBV	311503	561137	18.47%	3.687%
9	13.412	1716	1723	1734	rBV	1269760	1922065	63.25%	12.629%
10	14.787	1950	1957	1970	rVB2	546331	876806	28.85%	5.761%
11	16.267	2201	2209	2220	rBV	630914	981286	32.29%	6.448%
12	17.525	2416	2423	2433	rBV2	740397	1107834	36.46%	7.279%
13	20.116	2857	2864	2874	rBV	2213716	3038741	100.00%	19.967%
14	21.796	3143	3150	3158	rBV	751089	1302299	42.86%	8.557%
15	25.104	3702	3713	3728	rBV2	373229	1227198	40.39%	8.063%

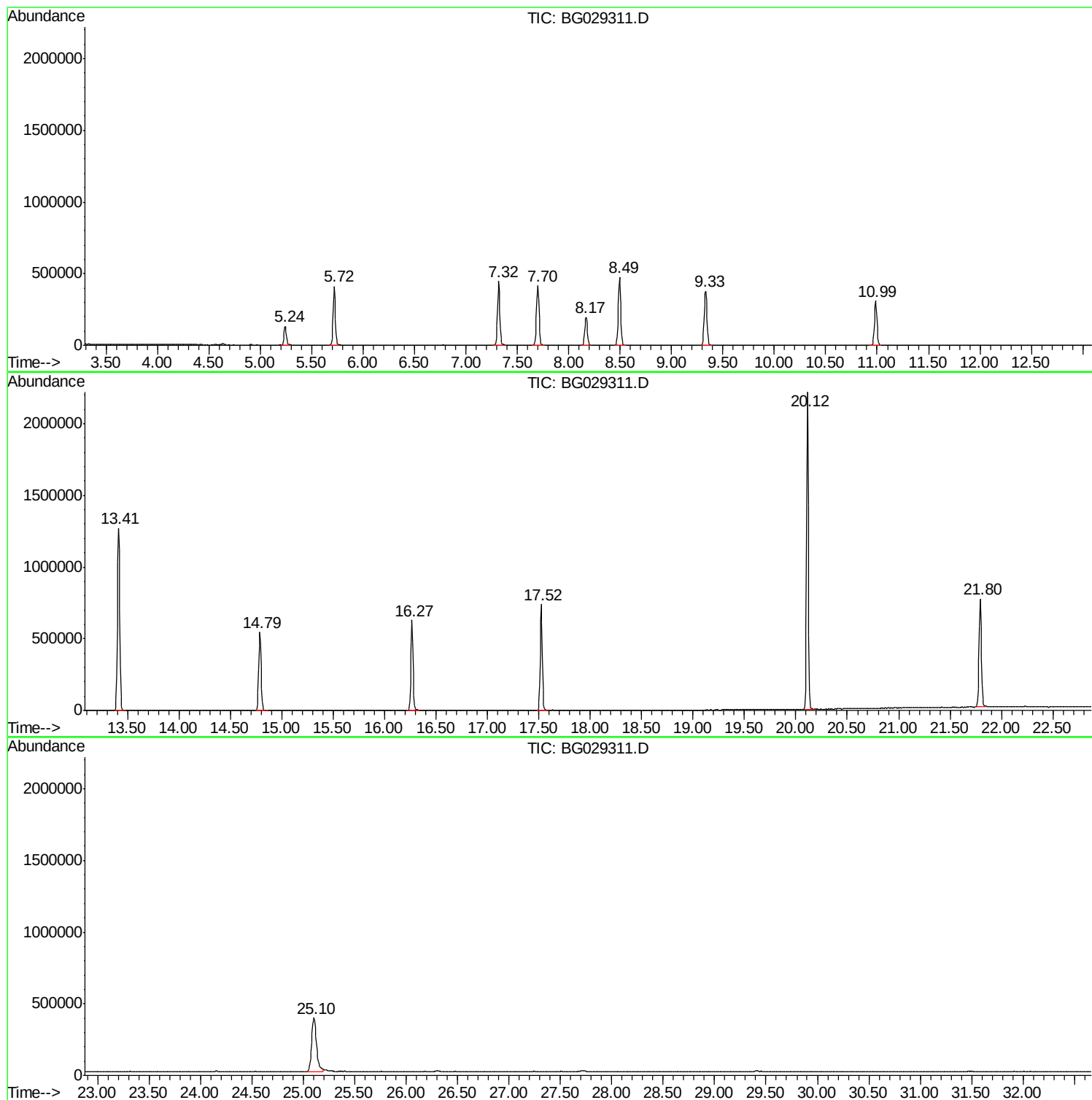
Sum of corrected areas: 15219180

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.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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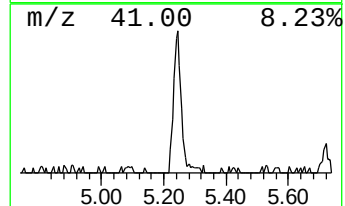
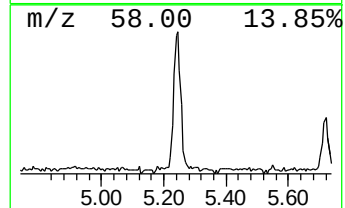
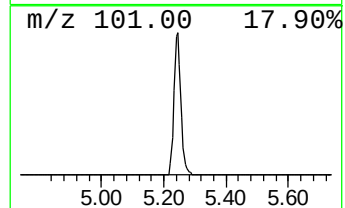
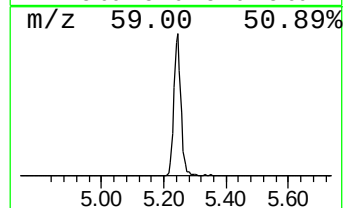
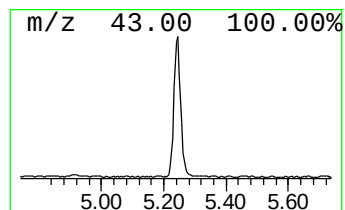
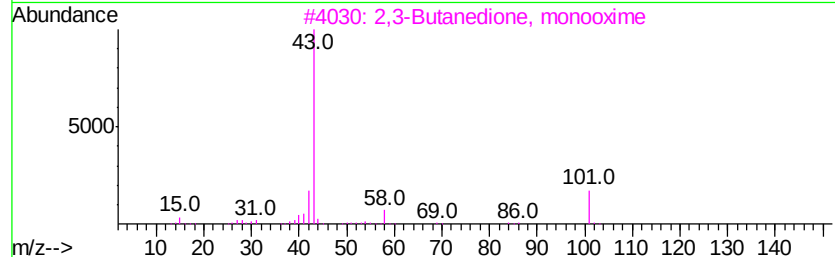
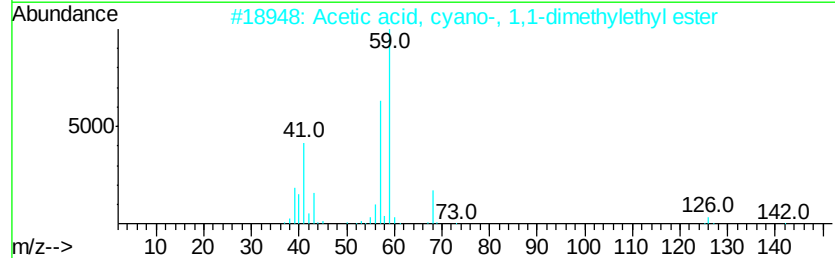
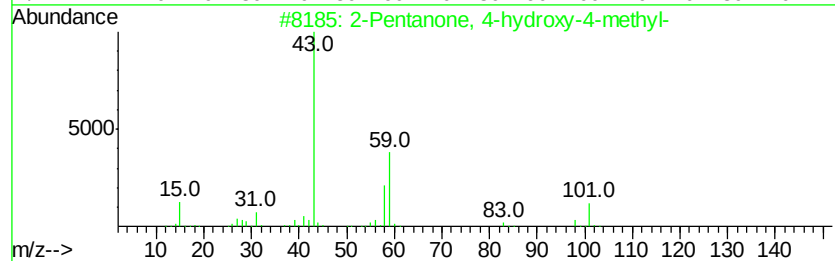
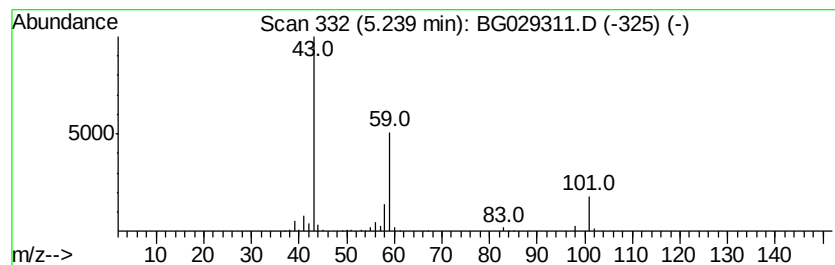
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	12.10 ng	204870	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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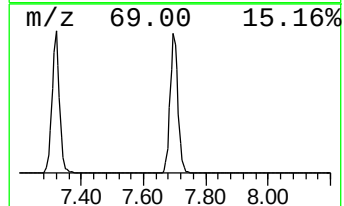
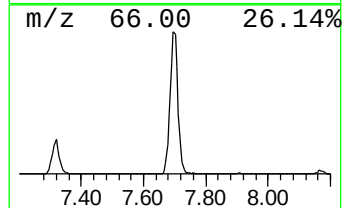
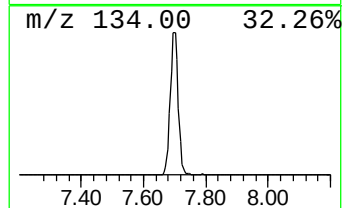
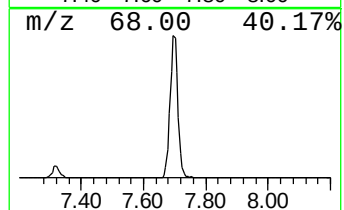
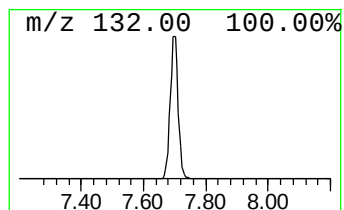
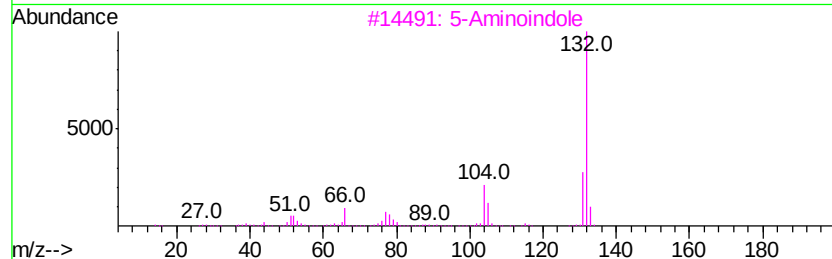
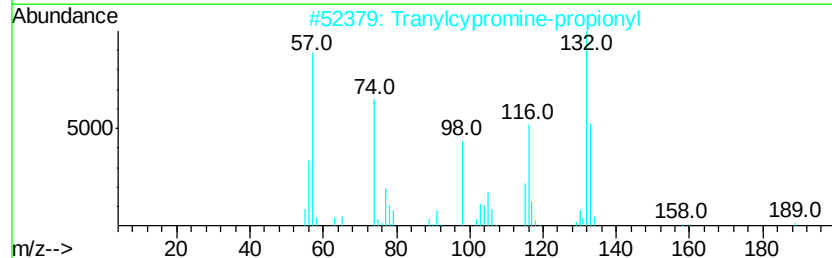
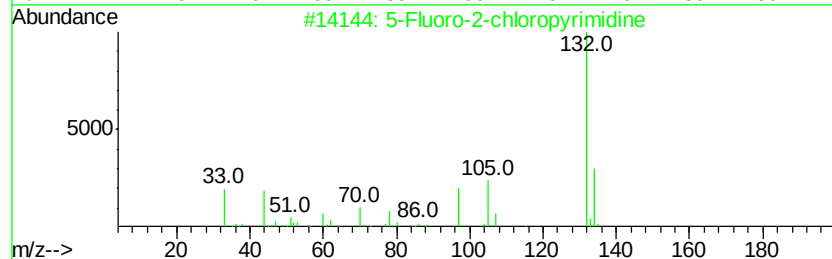
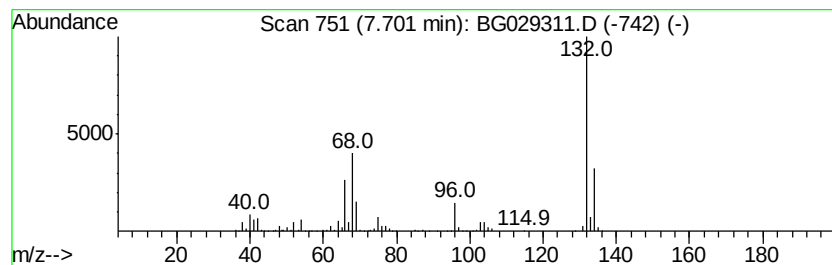
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Peak Number 2 unknown7.70 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	43.40 ng	734761	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
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...	5.24	12.1	ng	204870	1	8.17	338587	20.0
unknown7.70	7.70	43.4	ng	734761	1	8.17	338587	20.0