

My completed code and data is contained in the folder **assembler/**.

1

The source code is contained in **src**. The program can be compiled by running **ant** and run with **bin/assembler <args>**. I set up a script **run_trials** which I used to easily playing with the parameters on my runs.

2

Results from the small first-week genome are contained in **reduced_genome/**.

3

Results from the second week reads are contained in the **full_genome** folders and labeled by coverage.

4

The approximate value of parameter a (coverage) required to get c contigs with the parameters $l = 50$ and $G = 10000$ is calculated as follows:

Given

$$c = Ne^{\frac{Nl}{G}}$$
$$a = \frac{Nl}{G}$$

we find

$$c = \frac{G}{l}ae^{-a}$$
$$ae^{-a} = \frac{cl}{G}$$

which cannot be easily simplified.

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