

---

# PopCluster (Mac)

---

## 1. Introduction

This document briefly describes how to run PopCluster on a Mac. It should be read together with the PopCluster.pdf file, which gives much more details about the methods, inputs and outputs of PopCluster and is thus highly recommended to read before starting the use of PopCluster.

The zipped Mac package includes the following files in 3 subfolders:

- 1) PopClusterMac
- 2) Readme.pdf
- 3) PopCluster.pdf
- 4) PopClusterUpdateHistory.txt
- 5) ant377NoScale.dat
- 6) ant377NoScale.PcPjt

The first file is the binary compiled by Gfortran compiler (GNU fortran 6.3.0). It is for 64bit Mac, and may not work for 32bit OS. It is statically linked for serial run only. This binary file is placed in the subfolder "Bin".

The 2ed file is what you are reading. The 3rd is a detailed description of the use of PopCluster program. The 4th file is a brief of the updating history of PopCluster. These 3 documents are in the subfolder "Doc".

The 5th file is an example genotype data file, and the 6<sup>th</sup> is the corresponding parameter file. These 2 files are in the subfolder "Example".

Installation and run of PopCluster programs are all better conducted in Mac's x-terminal. It is suggested to open an x-terminal and then to obtain the administrator's privileges by typing

```
sudo -s
```

Otherwise, the shell operations (such as unzipping) and PopCluster run may have permission problems.

## 2. Installation

In x-terminal, unzip the downloaded file, and copy the 3 subfolders (with content files) into your PopCluster program folder on your Mac. Then go to the Bin subfolder, and use command "chmod u+x PopClusterMac" to make the binary executable on your Mac.

## 3. Running PopCluster

To make a serial run of PopCluster for the example dataset, simply go to the PopCluster program folder (using command cd) in the X terminal and type

```
./Bin/PopClusterMac INP:./Example/ant377NoScale.PcPjt
```

On finishing the run, analysis results will be in several output files located in the same folder as the parameter file “ant377NoScale.PcPjt”. That is subfolder “Example”, which acts as the project folder.

In the above command line, the key word “INP:” means INPut file. Other key words, as listed and explained in the main documentation PopCluster.pdf, are used to provide runtime parameters which override the parameter values set in the parameter file.

To run your own dataset, you need to make a project folder, to prepare a parameter input file and a genotype data file, and to save the 2 files in the folder. Suppose your project folder name with path is “MyPrj”, and in this folder you have a parameter file MyPar.PcPjt and a data file MyData.dat. Note the data file name “MyData.dat” must be specified in parameter file “MyPar.PcPjt”. You can run the analysis of your data from PopCluster program folder by the X-terminal command line

```
./Bin/PopClusterMac INP:MyPrj/MyPar.PcPjt
```

On finishing the run, analysis results will be in several output files stored in your project folder, MyPrj.

When executing the binary for the first time, you may see an error message “permission denied”. This means the binary needs to be unlocked for execution by issuing the X-terminal command line

```
chmod u+x ./Bin/PopClusterMac
```

in the PopCluster program folder.