

Introduction to Bioinformatics

Git for Bioinformatics Analysis, Project Management and
Collaboration

Introduction

- When performing bioinformatic analysis and starting a bioinformatics project you're expected to write codes
- Not only that, sometimes you also have to track the results of your experimentation for the analysis
- Let's start with a story

First Scenario

Day 1

```
base_dir <- system.file("extdata", package = "minfiData")
base_dir <- "recfon_practice/data"

ann450k <- getAnnotation(IlluminaHumanMethylation450kanno.ilmn12.hg19)

targets <- read.metharray.sheet(base_dir, pattern="SampleSheet.csv")

# read in the raw data from the IDAT files
rgSet <- read.metharray.exp(targets=targets)
rgSet

# give the samples descriptive names
targets$ID <- paste(targets$Sample_Group, targets$Sample_Name, sep=".")
sampleNames(rgSet) <- targets$ID
```

Imagine you're writing codes in R for your analysis, the code works perfectly and returned the expected results

Day 2

```
base_dir <- system.file("extdata", package = "minfiData")
base_dir <- "recfon_practice/data"

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# give the samples descriptive names
targets$ID <- paste(targets$Sample_Group, targets$Sample_Name, sep=".")
sampleNames(rgSet) <- targets$ID
```

The next day you realize that you didn't need some lines of code, so you remove it

Day 14

```
base_dir <- system.file("extdata", package = "minfiData")
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```

Fast forward to day 14 of you're writing code it turned out that you have to go back to your initial code

Day 14

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sampleNames(rgSet) <- targets$ID
```

Well for a shorter code like something above its easier, we know that we removed ann450k and target variable

But the truth is

```
3703
3704   # Get all the CpG sites used in the analysis to form the background
3705   all <- DMPs$Name
3706   # Total number of CpG sites tested
3707   length(all)
3708
3709   par(mfrow=c(1,1))
3710   gst <- gometh(sig.cpg=sigCpGs, all.cpg=all, plot.bias=TRUE)
3711
3712   # Top 10 GO categories
3713   topGSA(gst, number=10)
3714   gst[,gst$ONTOLOGY == "BP"]
```

Turns out when you're doing bioinformatics project things can really get complicated

But the truth is

```
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```

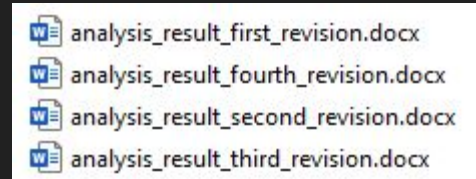
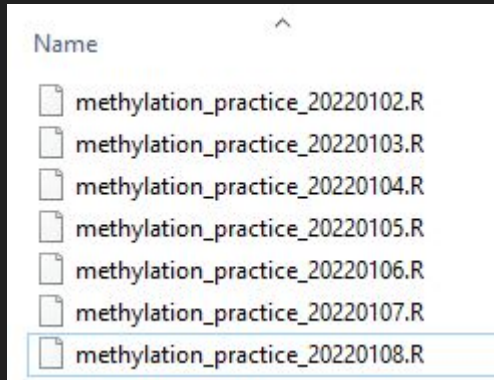
Number of files will increase significantly, so are the results and the processed data

But the truth is

```
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```

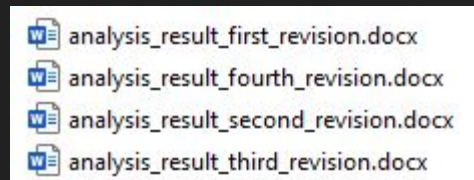
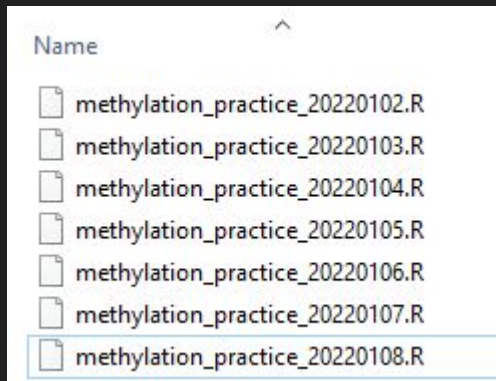
It will be hard to track the changes you made and to go back to that point if you're not using any version control system (VCS)

Well we have the conventional way of doing that



We can create multiple files and put a date suffix at the end of it's name right?

Well we have the conventional way of doing that



But this one is not practical! Why? Not only it increases the file size but it will also make the project directory less tidier

Second Scenario

Collaboration

```
base_dir <- system.file("extdata", package = "minfiData")
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targets <- read.metharray.sheet(base_dir, pattern="SampleSheet.csv")

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# give the samples descriptive names
targets$ID <- paste(targets$Sample_Group, targets$Sample_Name, sep=".")
sampleNames(rgSet) <- targets$ID
```

Mbak Dwi's task

```
par(mfrow=c(1,2))
barplot(colMeans(detP), col=pal[factor(targets$Sample_Group)], las=2,
        cex.names=0.8, ylab="Mean detection p-values")
abline(h=0.05, col="red")
legend("topleft", legend=levels(factor(targets$Sample_Group)), fill=pal,
        bg="white")

barplot(colMeans(detP), col=pal[factor(targets$Sample_Group)], las=2,
        cex.names=0.8, ylim=c(0,0.002), ylab="Mean detection p-values")
abline(h=0.05, col="red")
legend("topleft", legend=levels(factor(targets$Sample_Group)), fill=pal,
        bg="white")
```

Mbak Zahra's task

Sometimes the project is big enough that it requires collaboration

Collaboration

```
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Mbak Dwi's task

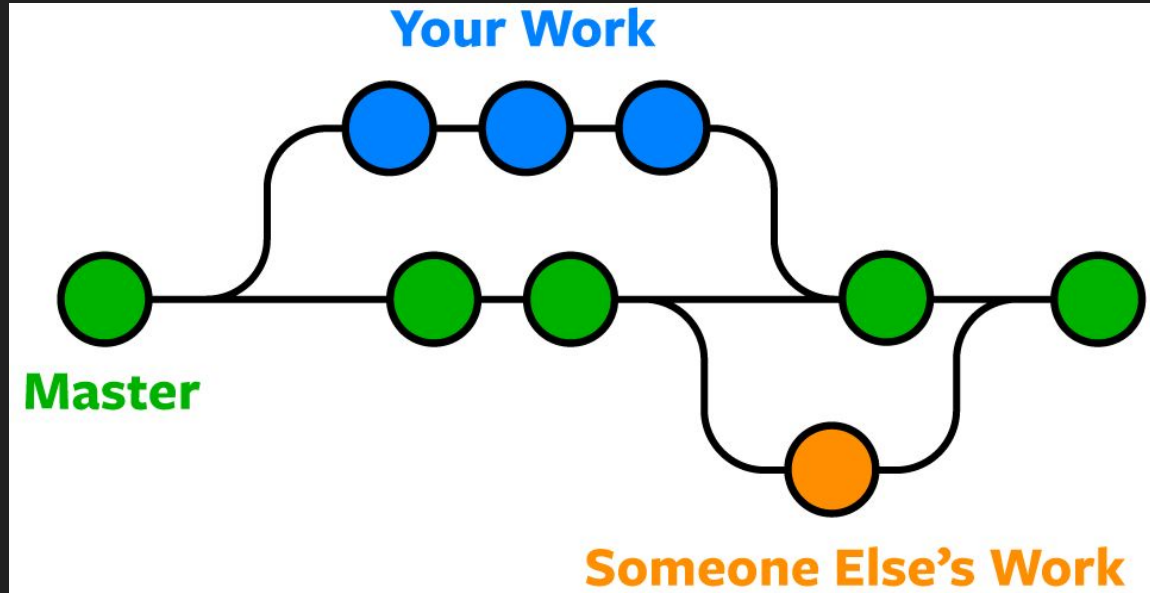
```
par(mfrow=c(1,2))
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barplot(colMeans(detP), col=pal[factor(targets$Sample_Group)], las=2,
        cex.names=0.8, ylim=c(0,0.002), ylab="Mean detection p-values")
abline(h=0.05, col="red")
legend("topleft", legend=levels(factor(targets$Sample_Group)), fill=pal,
        bg="white")
```

Mbak Zahra's task

Not only codes but all of the things that are related to the project, like report writing etc

Version Control System solves this problem



**So how do we overcome such
problems and doing the best
practice instead of the
traditional way?**

**Yes we have this thing called
Git**

What is Git

- Git is a Version Control System
 - Helps handling changes and maintaining history of our projects
 - Not only that it also has more features like **branch** and **merge**
- Git helps us revert back to the older version of our work in case something didn't go well
- Git provides easier access for collaboration
- Git has been widely used by bioinformaticians, software developers and etc.

Installation

- There are multiple ways on how to install git
- The easier way when we're working with Windows computer is to download it from <https://git-scm.com/downloads>
- If you're using UNIX based system you should follow this instruction <https://git-scm.com/book/en/v2/Getting-Started-Installing-Git>

GitHub account

- For this session we will be using GitHub to store our codes and analysis results
- To create an account go to github.com and register it with your email
- There are other providers for Git cloud but for now we will focus on using Github, they use the same command to so no worries with the commands

Repository

- A folder, usually used to organize a single project
- Repositories can contain folders and files, images, videos, spreadsheets, and data sets -- anything your project needs.
- Often, repositories include a README file, a file with information about your project. GitHub makes it easy to add one at the same time you create your new repository.

winter-internship-2017

Public

repo of my codes written during my internship at Gebze Technical University
and other internship-related stuff

☆ Star

internship

gebze-technical-university

rna-seq

python

r



Jupyter Notebook



1

Updated on Mar 18, 2017

**Let's start experimenting with
Git!**