Introduction to Linux command line - Part 1

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2/10/2021

Preamble

For this first lesson, we are going to adopt the 'infant swimming' technique: it doesn't matter if you don't know the command line yet, you are going to write some code and hopefully get results;-) ... The deep understanding of the code will be revealed later on.

The \$ symbol at the beginning of the lines represents the *prompt* you have at the command line, i.e. the place where the cursor is located and ready to accept any characters you type. In the case of your terminal, some information are preceding the \$. We will see later what they are and how we can customize them. Note that depending on your system, the prompt could differ and the \$ replaced by another character such as >.

At the end of each line, you press enter, which sends the string to the *command line interpreter*, here the bash program.

In the following tutorial, some lines will start with a #. It means that the line is just a *comment*. You don't need to write it, but even if you did, it won't be interpreted.

Very very very important warnings

- 1. If something goes wrong, it is probably because you typed something wrong. Remember that every single character that you type will be interpreted and could have a special meaning depending on the context or position in the command line. This is especially the case for blank spaces or special characters such as but not limited to][)(}{|;&~.*\$#!. Follow very carefully what is written in the tutorial.
- 2. Some commands will use special codes starting with !, for example . . . cd !#:2. These special structures are related to the bash history expansion feature that will be explained in details later. This feature is really worth using because it improves greatly the ergonomy of the command line. That is the reason why I have tried to demonstrate its use as much as possible, along the exercise. However, since some of the codes really depend on the actual sequence the commands are entered, they may not work as expected if for some reason the flow of your commands differs slightly from what is proposed in the tutorial. So, you will always find the complete command line as a comment starting with # shortcut for . . .

Let's go...

Additional configuration

Open the shell configuration file ~/.bashrc with the text editor nano...

\$ nano ~/.bashrc

... and add 2 lines at the end.

```
\# add these lines at the end of the file (scroll down with the mouse or arrow key) PROMPT COMMAND='history -a'
```

```
shopt -s histverify
# Press Ctrl-x to quit, O or Y to accept changes and enter to validate the
# path to the file to be saved.

# Apply the changes
$ source ~/.bashrc

# If you did that correctly, the next command should display "history -a"
$ echo $PROMPT_COMMAND
```

Organize your workspace

```
$ cd ~
$ pwd
```

The last command should produce an output like /home/your_login_name

```
$ mkdir -p linux/1st_lesson && cd !#:2
# shortcut for mkdir -p linux/1st_lesson && cd linux/1st_lesson
$ pwd
```

When you press enter, the part !#:2 should be replaced by linux/1st_lesson or the exact string your entered just after mkdir and &&.

Download taxonomic data

```
$ wget "ftp://ftp.ncbi.nlm.nih.gov/pub/taxonomy/taxdump.tar.gz"

# In the next command, stop after typing 'tax' and press the tabulation key...
# That's a first example of auto-complete capabilities of the shell.
# In the following, this {TAB} notation will invite you to use this great feature
# as often as possible
$ tar -xzf tax{TAB}
$ ls -l
$ head na{TAB} no{TAB}
```

As you probably understood, the last command, head, displayed the first lines of each files, giving you a hint about their content. The next command will help you to better understand what they are.

```
$ less r{TAB}
# Should open readme.txt
# Scroll with mouse/tab gestures or arrows and quit with 'q'
```

One of the file contains data linking *taxids* with names (of species, genus, families, etc...) and the other one data linking *taxids* with taxonomy ranks. The objective is to retrieve the taxids and names for the 'family' and 'order' ranks. A new file containing 3 "columns" (taxids, ranks and names) will be created. But first, let's do some cleaning up to remove the files not needed.

```
# Start by doing some cleaning, you only need names.dmp and nodes.dmp. The other files
# can be deleted
$ ls -l
# Use the up arrow (and down arrow) to navigate in the history of the commands you typed...
```

```
# By doing this, you can modify the last command and add a pattern to filter the output
$ ls -l n[ao]*.dmp

# Use again the up arrow and edit the last command rather than re-typing everything...
$ ls -l !(n[ao]*.dmp)
$ rm !!:$
# shortcut for rm !(n[ao]*.dmp)
```

Data processing

In the next series of commands, you will process the data contained in *names.dmp* by using several commands acting sequentially. You will build what is called a *pipe* by adding one command at a time (separated by the pipe character |) and evaluating each intermediate result.

To make it easy, remember to use the up arrow key to displayed the last executed commands and simply adding the new command at the end.

```
$ head -n 40 na{TAB}
$ head -n 40 names.dmp | grep 'scientific name'
$ head -n 40 names.dmp | grep 'scientific name' | cat -T
$ head -n 40 names.dmp | grep 'scientific name' | tr -d '\t'
$ head -n 40 names.dmp | grep 'scientific name' | tr -d '\t' | cut -d '|' -f 1,2
# If you press 'enter' after a pipe symbol, the command is not executed because
# bash 'knows' it is not finished. The prompt changes to indicate that you are
# in a multiline command. You can test it if you like, but it is not necessary,
# I only use it here to avoid page overflow.
$ head -n 40 names.dmp | grep 'scientific name' | tr -d '\t' | cut -d '|' -f 1,2 |
> sort -t '|' -k 1b,1
# Save the result in a new file
$ head -n 40 names.dmp | grep 'scientific name' | tr -d '\t' | cut -d '|' -f 1,2 |
> sort -t '|' -k 1b,1 > file1
# Have a look at what you get
$ 1s -1h
$ wc -l f{TAB}
# Now, let's do it on the whole file
$ cat !head:3*
# shortcut for cat names.dmp | grep 'scientific name' | tr -d '\t' |
               cut -d '/' -f 1,2 / sort -t '/' -k 1b,1 > file1
$ ls -lh
$ wc -1 f{TAB}
```

From the nodes.dmp file, you will filter the lines related to *family* and *order* ranks (strictly!) and create a new file containing only taxid and rank columns.

```
$ head -n 10 no{TAB}
$ head -n 10 nodes.dmp | grep -E '\|family|\|order'
$ ^10^250^
# shortcut for $ head -n250 nodes.dmp | grep -E '\|family|\|order'
$ head -n 250 nodes.dmp | grep -E '\|family|\|order' | cut -d '|' -f 1,3
```

```
$ head -n 250 nodes.dmp | grep -E '\|family|\|order' | cut -d '|' -f 1,3 |
> sort -t '|' -k 1b,1

# If the result is ok, apply it on the whole nodes.dmp file and save it in a new file
$ cat !!:2* > file2
# shortcut for $ cat nodes.dmp | grep -E '\|family|\|order' | cut -d '|' -f 1,3 |
# > sort -t '|' -k 1b,1
$ ls -lh
$ wc -l f{TAB}2
```

Check quickly the result in both new files and make the join.

```
$ tail f*
$ join -t '|' -o '1.1 2.2 1.2' file* | head

# Some ordering
$ !!:- sort -t '|' -nk1
# shortcut for join -t '|' -o '1.1 2.2 1.2' file* | sort -t '|' -nk1

# OK ? So let's create the file
$ !! > taxonomy
# shortcut for join -t '|' -o '1.1 2.2 1.2' file* | sort -t '|' -nk1 > taxonomy

$ ls -lh
$ wc -l t{TAB} should produce taxonomy

# Add column names
$ sed -i '1i taxid|rank|name' t{TAB}
```

Now, consider you want to have a french version of this taxonomy file and to share those files (english or french) with colleagues who are using standard office software. You will do the translation of order to order and family to famille and change the field separator from | to , or ; which are the standard separators of .csv files recognized in common spreadsheet tools.

```
# The french version...
$ echo "taxid,rang,nom" > taxonomy_fr.csv
$ tail -n+2 t{TAB} >> !!:$
$ sed -i 's/|/,/g' !!:$
$ sed -i 's/order/ordre/' !!:$
$ ^order/ordre^family/famille^
# shortcut for sed -i 's/family/famille/' taxonomy_fr.csv

# The english version as .csv file
$ sed -i 's/|/,/g' t{TAB}
$ mv t{TAB} t{TAB}_en.csv
```