**BIO334 Day 1: UNIX**

The UNIX operating system is made up of three parts; the kernel, the shell and the programs. The main control program in a UNIX OS is the kernel. The kernel does not allow the user to give it commands directly; instead when the user types commands on the keyboard they are read by another program in the OS called a shell which parses, checks and translates them in various ways and then passes them to the kernel for execution.

Everything in UNIX is either a file or a process. The UNIX filesystem is laid out as a hierarchical tree structure which is anchored at a special top-level directory known as the root (designated by a slash '/'). Because of the tree structure, a directory can have many child directories, but only one parent directory.

On Mac, you can open the terminal window by typing **Command-Space** and searching for "Terminal" (select the **Terminal.app**):

Text

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When you see "$" sign followed by the input cursor you are ready to go, so proceed to Exercise 1.

 means that it’s something you need to solve (solutions to these exercises can be read from the solutions file)

 useful tips or annotations

**Exercise 1: basic file and directory manipulation**

The dollar character ($) is a convention showing you are in the command line input, don’t type it

|  |  |
| --- | --- |
| $ **whoami** | prints out your user name (don’t worry if you do not recognize it, on your computer it will bear your name) |
| $ **pwd** | prints the current directory |
| $ **cd ~** | change directory command, the tide character is a shortcut to your home directory. In UNIX systems each user has a home directory. This will be your playground. |
| $ **pwd** | prints the current directory (notice the directory change). The last subdirectory is named after you, the user. |
| $ **ls –l** | ls means list, -l means long; this command lists all files, permissions and directories in the current directory |
| $ **mkdir** tmp | make directory; the command creates the directory “tmp” in your home directory |
| $ **cd** tmp | brings you to your newly created directory tmp |
| $ **pwd** | you can see that your current working directory changed |
| $ **cd ..** | move one level up |
| $ **ls -l** | you should see the newly created tmp directory |
| $ **press ↑** (arrow up key) | press arrow up and down to go through recent commands |
| $ **vi** tmp/poem.txt | start editing file tmp/poem.txt |
| **press i** | press i to switch to “insert” mode in the vi editor, while in this mode “insert” tag should be displayed at the bottom left corner of the page. |
| <http://www.poetryfoundation.org/poem/171621> | open web page and copy/paste the poem to the poem.txt file in your vi editor while in “insert”mode |
| press **ESC** | press ESC to switch to “command” mode in the vi editor. The “insert” tag on the bottom of the page should disappear. |
| type **:w** | Now you are back in the command mode. press **:** and then **w** and **ENTER** to write contents to file. The commands will appear in the lower left corner. |
| type **:q** | press **:** and then **q ENTER** to quit the vi editor and return to the shell |
| $ **cat** tmp/poem.txt | displays the poem |
| $ **cat** | what happens now? Try typing anything and press enter. |
| press **Ctrl** and **c** | (you are in the stdin mode where cat command listens to your keyboard input and outputs anything you type to the screen). Admittedly not very useful. Press the **Ctrl+c** combination to *kill* the current command.  Remember that combination. It is a useful trick to exit an unwanted command. |
| $ **head** tmp/poem.txt | displays first 10 lines of the poem |
| $ **tail** tmp/poem.txt | displays last 10 lines of the poem |
| $ **man** head | use **man** to get more information about the head command; find and write down the head command with the correct parameter to display not only the first 10 lines of the poem but the entire poem (20 lines) |
| $ **less** tmp/poem.txt | less is a convenient command for displaying file content. Large files that do not fit in the screen can be scrolled using “up” and “down” arrow. Mouse/trackpad won’t work. Return by pressing **q** |
| $ **grep** house tmp/poem.txt | if you are searching for something, you can use **grep** text to display only lines matching text How many lines you see? |
| $ **wc** -l tmp/poem.txt | displays the number of lines in the file |
| $ **wc** -w tmp/poem.txt | counts the words in the file |
| $ **cd** tmp | Enter the “tmp” directory. |
| $ **cp** poem.txt poem2.txt | **cp** command is used to make a copy the file. The second argument can be either a file name - this will make a copy of the file under the new name - or a path to a directory. The latter will create the copy of the file in the specified directory under the same name. For example  $ cp poem.txt ../ will create a copy of poem.txt in the directory above. |
| $ **ls** -l | you should see 2 files now |
| $ **cd** .. | go back one level |
| $ **cp** -r tmp poems | you just made a copy of the entire directory “tmp” as “poems” |
| $ **ls** -l | you should see now both directories “tmp” and “poems” |
| $ **ls** tmp/\*.txt | you can use the wildcard (**\***) to list files and directories that match your expression. In this case, all the files that ends with “.txt” extension. |
| $ **rm -r** tmp | **rm** removes files and folders. The -r flag stands for recursive. It will delete this folder and all files and folders within. |
| $ **clear** | clears the terminal display |

More often than not you won’t have to write down the whole command name or the whole file name, and it’s enough to write the first one or two letters and press [tab] and UNIX will fill the rest for you. If there is an ambiguity which command or file you have actually meant, press [tab] again and UNIX will display all the possibilities which start with your characters.

 Write **less p** in the command line and then press [tab] 4 times. What has happened each time you press [tab]?

Try to use [tab] as often as possible throughout the exercises. It should be a second nature to tap it all the time.

**Exercise 2: pipes and redirecting input / output**

The pipe operator takes the output from one command and uses it as the input for the next command:

$ command1 | command2 | command3 …

The standard output of command1 is redirected (piped) to the standard input of command2, etc.

It’s one of the most powerful tools at your disposal. This way you can combine an arbitrary number of programs in one command. In more technical terms, the pipe operator is used to create concurrently executing processes that pass data between them.

the output in UNIX is called *standard output*, as opposite to *standard error* output which is a way for the script to print any warnings or errors on the screen that won’t be passed to the next step.

For example, a common expression in bioinformatics that you may want to write is something like this:

read data.tsv file [then] select third column [then] select lines that say "sample2" [then] count how many lines there are

This would translate in command line to this expression:

$ cat data.tsv | cut -f 3 | grep "sample2"| wc -l

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| --- | --- |
|  | What is the standard **input** and standard **output** of the command **grep** in the example above? |
|  | Try it yourself, write a command that reads poem.txt file and prints only the lines that include the word “house” (use **cat** and **grep**) |

**Redirecting input and output.** The output from programs is usually written to the screen, while their input usually comes from the keyboard (if no file arguments are given or nothing is *piped in*). To redirect standard output to a file instead of the screen, we use the **>** operator:

|  |  |
| --- | --- |
| $ echo hello | displays “hello” on the screen |
| $ echo hello > hello.txt | writes “hello” to the file |
| $ cat hello.txt | displays contents of file; you should see “hello” |

In this case, the contents of the file hello.txt will be overwritten if the file already exists. If instead we want to append the output of the echo command to the file, we can use the >> operator:

|  |  |
| --- | --- |
| $ echo hello >> hello.txt | appends “hello” to the end of the file |
| $ cat hello.txt | displays contents of file |

Standard input can also be redirected using the < operator, so that input is read from a file instead of the keyboard:

$ wc -l < hello.txt

There are many ways to do the same thing in UNIX, you may have noticed that the above example is equivalent of the one that you have already learned:

$ cat hello.txt | wc -l

You can combine input redirection with output redirection, but be careful not to use the same filename in both places. For example:

$ wc -l < hello.txt > hello\_counts.txt

**Exercise 3: writing and executing a shell script**

You now know the basics of the **vi** editor from the first exercise. Try to copy/paste the below simple bash program into a file and execute it. The hello world bash script:

#!/bin/bash

echo "Hello World"

Copy the above 2 lines and save them to the file “hello.sh”. The first line tells the Unix how to run the program. In this case it says that the file is a shell command and it should run it using ‘bash’ shell (but it could be very well a Python or R script).

Make your file executable by typing:

$ chmod +x hello.sh

Every file and folder in UNIX system has a set of permissions which says who can read it, and who can write to it. One of the permissions states if the file can be executed. This command sets the permission so that you (or whoever have permissions to read the file) can also run it.

And now, to run it, you can simply type:

$ ./hello.sh

 What has been displayed?

Note the “**./”** at the start of the command. The “**.**” tells the system that the file is in the current directory (for comparison the “**..**” meant the directory above).

**Exercise 4: download FASTA file and count the number of proteins**

Now we will download a FASTA file. For this we will use a command called **curl**

$ curl <https://stringdb-static.org/download/protein.sequences.v11.0/9606.protein.sequences.v11.0.fa.gz> -o human\_proteome.fa.gz

The file will download in the current directory named human\_proteome.fa.gz. Unzip it using a command **gunzip**

$ gunzip human\_proteome.fa.gz

How the name of the file has changed when you unzipped it?

Count the number of proteins in the file using commands you have learned. In the fasta format file the first line of each protein record starts with “**>**” character.

**Exercise 5: sort file on column**

Download the protein abundance data (<https://pax-db.org/downloads/4.1/datasets/9606/9606-BRAIN-integrated.txt>) with curl, like you did in the previous exercise and name the file “abundance\_data.tsv”

 Explore the downloaded file using the **less** command. Note which column holds the abundance numbers.

Run the following command:

$ sort -k3nr abundance\_data.tsv | less

 Open manual of the **sort** command, look for option “**k**” and “**n**”, “**r**”. What do they mean in the context of the command above?

 Remove the “**n**” parameter and run the sort command again. Explain what has happened.

 **Exercise 6: write a simple bash script**

Often, you would like to run the same command with different parameters. As an exercise, write a simple bash script that will output numbers from 1 to 100. Use a for loop.

#!/bin/bash

for i in {1..100}

do

echo $i

done

Save the above code to a file (e.g. **script.sh**), make the file executable (**+x** flag) and run it.

What is the output?

 **Exercise 7: iterating over files**

Download the compressed folder containing a selection of proteomes with curl, and extract it using the following **tar** command. URL with the proteomes: <https://raw.githubusercontent.com/meringlab/Bio334/master/01_unix/exercises/data/proteomes.tar.gz>

$ tar xzvf proteomes.tar.gz

while **gzip** compresses one file, **tar** on the other hand is used to compress the whole directory structure and all the files in it. The arcane string **xzvf** translates to **X**tract **Z**ipped **V**isual **F**ile.

Enter the proteomes directory (**cd** command)

 Using the commands you have learned print out the number of files in the directory.

By utilizing the same concept of **for** loop from the previous exercise, can you try to create the shell script that iterates over all fasta files in the directory, print their names and the number of sequences in each file?

for filename in \*.fa

You can put more than one command between **do** and **done**, one line would print the file name, the other will print the sequence count.

**\* Exercise 8: exploring a FASTA format file**

*Dictyostelium discoideum* ([www.dictybase.org](http://www.dictybase.org)) is an interesting social amoeba and a well-studied model organism. The whole genome sequence is already available here:

<https://raw.githubusercontent.com/meringlab/Bio334/master/01_unix/exercises/data/dd.tar.gz>

Download the file with **curl** and extract it with **tar**. The FASTA format is widely used in sequence distribution, see the description at: <http://en.wikipedia.org/wiki/FASTA_format>.

Explore the file using **less**. Use **grep** to find out how many chromosomes are present in the file. Use **grep** (**-v**) to only print out the genomic sequence. How large is the genome?

Now look at the RNA-seq data sample stored in the rnaseq.fastq file. This file includes qualities for each nucleotide (see FASTQ description at <http://en.wikipedia.org/wiki/FASTQ_format>).

How many reads are in the file?

A picture containing text

Description automatically generated**\* Exercise 9: download and install bowtie2 software**

Bowtie2 is a short-sequence read aligner (e.g. 150nt long). The reads are aligned to a reference sequence (e.g. human genome). Here, we will use anaconda to install the package since curl does not work in this case due to many dependencies. Simply copy and paste the following commands into your terminal:

cd ~

. /opt/miniconda3/etc/profile.d/conda.sh

conda create --prefix ~/Documents/py38\_envs python=3.8

When asked to proceed, type y and press enter.

conda activate ~/Documents/py38\_envs

conda install -c bioconda bowtie2

Also here, press “y” and enter when asked if you want to proceed.

Now you can simply type “bowtie2” anywhere (in any directory) and the **shell** will find the **bowtie2** software.

**\* Exercise 10: searching for short sequences in the *Dictyostelium discoideum* genome**

Is the sequence “AAAAAGAGATACAT” present in the DD genome (dd/dd.fasta)?

 use **grep** to find out

\* You can also use **bowtie2,** which we previously installed. First build the index of the DD reference genome. The format is: “bowtie2-build <fasta\_file> <custom\_index\_name>”. In the dd folder, you could use:

$ bowtie2-build dd.fasta dd

Once the index is created (you do this only once for each reference genome, i.e. each FASTA file), you align one read (sequence) to the genome by typing:

$ bowtie2 -x dd -ac AAAAAGAGATACAT > dd.sam

Explore the SAM results file with **less -S**. The parameter “-S” prevents line wraps, so you can see one alignment per line.

Does **bowtie2** find more alignments compared to **grep**? Why could that be?

 **\* Exercise 11: alignment of RNA-seq sample reads to the *Dictyostelium discoideum* genome**

To align the RNA-seq reads in the rnaseq.fastq file, you first need to index the *Dictyostelium discoideum* genome. Use bowtie2-build to create an index with name **dd** (if you didn't already build the index in the previous exercise):

$ bowtie2-build fasta\_file custom\_index\_name

After you created the index, you can align the reads by typing:

$ bowtie2 -x dd -U rnaseq.fastq > dd.sam

The results are returned in SAM format and stored to the dd.sam file. How many reads align?

**System information, processes and other useful commands**

|  |  |
| --- | --- |
| **uname -a** | display system information |
| **man** *command* | display manual page of command |
| **df -h** | list mounted disks with available space |
| **du -h** *path* | show space usage |
| **top** | display running processes |
| **kill -9** *pid* | kill process |

**File and folder manipulation, compression**

|  |  |
| --- | --- |
| **pwd** | display current folder |
| **ls -l** *path* | list files and folders |
| **cd** *path* | change folder to path |
| **cd** ~ | change folder to home folder |
| **mkdir** *name* | make folder |
| **rmdir** *name* | remove folder |
| **cp** *source* *dest* | copy file/folder and all its contents |
| **less** *filename* | display file content |
| **wc** *filename* | count number of lines in file |
| **head** *filename* | shows first few lines of file |
| **tail** *filename* | shows last few lines of file |
| **gzip** *filename* | compress file with gzip (adds .gz extension) |
| **gunzip** *filename* | decompress filudecompress and remove .gz extension |
| **tar xfvz** *filename.tar.gz* | decompress files from tar.gz archive |
| **tar** **zcvf** *archive.tar.gz folder\_to\_compress* | creates archive.tar.gz |
| **unzip** filename.zip | unzip archive |

**Network and file transfer**

|  |  |
| --- | --- |
| **curl** URL -o filename | download URL to filename |
| **ssh** *username@host* | remote login to host with username |

**“vi” editor**

|  |  |
| --- | --- |
| **$ vi** *filename* | start editing file with vi |
| **i** | switch to “insert” mode |
| **ESC** | switch to “command” mode |
| **:w** | save |
| **:q** | quit |
| **:x** | save and quit |
| **/<pattern>** | search for pattern, <n> gives you the next match |
| **:q!** | quit without saving changes |