# Basics of the Linux Command Line

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## **Getting Started**

Today, we'll be using JSLinux, a web-based terminal emulator, to practice commands in a UNIX environment. Access it here: https://bellard.org/jslinux/.

You'll be presented with a list of options; click "click here" under "Startup Link" in the first row (Alpine Linux 3.12.0).

#### **JSLinux**

Run Linux or other Operating Systems in your browser!

The following emulated systems are available:

CPU	os	User Interface	VFsync access	Startup Link	TEMU Config	Comment
x86	Alpine Linux 3.12.0	Console	Yes	<u>click here</u>	<u>url</u>	
x86	Alpine Linux 3.12.0	X Window	Yes	click here	<u>url</u>	Right mouse button for the menu.
x86	Windows 2000	Graphical	No	click here	url	Disclaimer.
x86	FreeDOS	VGA Text	No	click here	<u>url</u>	
riscv64	Buildroot (Linux)	Console	Yes	click here	url	
riscv64	Buildroot (Linux)	X Window	Yes	click here	url	Right mouse button for the menu.
riscv64	Fedora 33 (Linux)	Console	Yes	click here	url	Warning: longer boot time.
riscv64	Fedora 33 (Linux)	X Window	Yes	click here	<u>url</u>	Warning: longer boot time. Right mouse button for the menu.

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#### What is this?

A terminal - a window that allows you to interact with a remote computer via text-based commands. The localhost:~# is called the prompt - this will look different depending on the computer you're logged into, but generally, the first part will tell you the name of the server you're logged into, followed by the directory you're currently in, then ending with # or \$. After the prompt, you can type your commands. Hit Enter to run them.

```
Welcome to JS/Linux (i586)

Use 'vflogin username' to connect to your account.
You can create a new account at https://vfsync.org/signup .
Use 'export file filename' to export a file to your computer.
Imported files are written to the home directory.

localhost:-#
```

## Where am I?

By default, you'll be in your home directory. The pwd command will **print** the current **working directory**, telling you where you are. Usually, this will be something like /home/yourname

```
In [ ]: pwd
You can also establish who you are (your username).

In [ ]: whoami
```

To see the files and directories present in your current directory, use the ls command.

```
In [ ]: ls
```

You can get more detail on the files using ls -l.

```
In [ ]: ls -l
```

There are lots of other ways to customize the output of ls. You can use the man command to get a **man**ual for another command. *This won't work in our practice terminal, but it will on most systems.* 

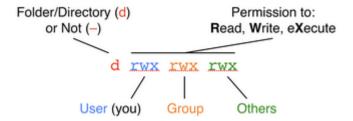
```
In [ ]: man ls
```

Now you can try something more complex.

```
In [ ]: ls -lah
```

## File permissions

So what's all this drwxrwxr-x? These are permissions, which tell you what you (and others) can do to a file or directory.



## Files and directories

In [ ]: | cd awesome data

You can create new directories, move to them, and see what's inside.

Note how your prompt changes as you move to a new directory.

```
In [ ]: ls
```

In [ ]: cd .. In [ ]: pwd You can chain together as many double dots as you like. This would take you up three directory levels: ../../... Another useful shortcut is ~, which represents your home directory. In [ ]: cd ~ In [ ]: pwd We can also create simple text files. In [ ]: echo "some fun text" > myfile.txt In [ ]: ls We can show the entire file with cat, scroll through it with less, or edit it with nano. In [ ]: cat myfile.txt In [ ]: less myfile.txt Type q to quit. In [ ]: nano myfile.txt Ctrl+x to exit, y / n to save or not save the file. Files and directories can be copied, moved, and removed. There is no recycle bin here; if you delete something, it's gone forever. Remove with care. In [ ]: cp myfile.txt newfile.txt In [ ]: mv newfile.txt awesome data In [ ]: cd awesome data In [ ]: mv newfile.txt delete this.txt In [ ]: rm delete this.txt In [ ]: cd ~

You can move to a directory above your current one using a double dot ( ... ).

To delete directories and their contents **recursively**, use the -r flag.

```
In [ ]: rm -r awesome_data
```

## Searching

We can use grep to search within text files. Move back to your home directory and try searching myfile.txt.

```
In [ ]: grep "fun" myfile.txt
```

To find a file, use the find command.  $\sim$  is the directory to search (recursively), -iname searches by name, case-insensitive, and "myfile.txt" is the name to search for. man find will get you more details.

```
In [ ]: find ~ -iname "myfile.txt"
```

## **Ending Things**

You can kill a running command with Ctrl+C.

Hint: you can paste this into the practice terminal by right-clicking on the flashing green cursor.

```
In [ ]: while :; do echo 'OH NO'; sleep 1; done
```

To terminate your connection to a server, use Ctrl+D.

## Other useful tips

You can use the mouse to select and copy text, but not to move your cursor. Use the left and right arrow keys to move around.

Copy and paste are a little different than you're used to - Ctrl+Shift+C and Ctrl+Shift+V.

You can use the up and down arrows to scroll through your command history - useful for editing previous commands.

You can use tab to autocomplete commands and filenames - type the beginning of the file name and then hit tab. Hit it twice to see a list of options if there's more than one possibility.

Asterisks (\*) are wildcard characters. For example, this would list all files in the current directory ending in .txt: ls \*.txt

If a text file is behaving strangely, the problem may be linebreaks. Windows, Mac, and Linux operating systems use different characters to tell a file the line has ended:  $\r\n$ ,  $\r$ ,  $\n$ ,  $\n$ . Use a decent text editor than can produce Unix linebreaks (not Notepad, never MS Word!). nano will work perfectly. Notepad++ is good for Windows users. You can fix incorrect linebreaks using the commands dos2unix and mac2unix.

You can see which directories will be searched when executing programs by examining your PATH variable: echo "\$PATH"

## Logging in using a real terminal

## Mac/Linux

You're using a UNIX-like system, so login using your local terminal using ssh (**s**ecure **sh**ell). The command will look something like this:

```
ssh your username@server address
```

Enter your password when prompted and hit Enter.

If you need to view graphics from the server, add -X or -Y. This can be finicky, so if one doesn't work, try the other.

```
ssh -X your_username@server_address or ssh -Y
your username@server address
```

For our current course server, you'll use a private key file (.pem) to prove your identity instead of a password:

```
ssh -i "your username md2023.pem" your username@18.222.38.194
```

For IU servers:

```
ssh your username@quartz.uits.iu.edu
```

OR

```
ssh your username@carbonate.uits.iu.edu
```

You'll need to complete 2-factor authentication with Duo to complete your login.

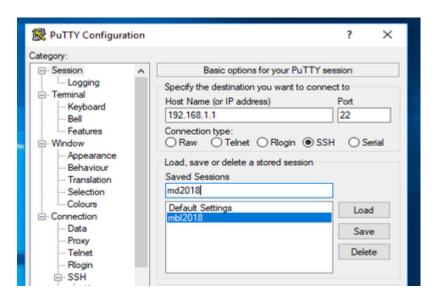
## Windows (using PuTTY)

#### For the course server:

- 1. Open PuTTY
- 2. Under Category, click on SSH > Auth
- 3. Click browse
- 4. Find your private key (your username md2023.pem) and select it
- 5. Under Category, click on SSH > Auth and check the "Enable X11 forwarding" box.
- 6. Under Category, click Session and input the server IP address (18.222.38.194) in the "host name" box
- 7. Type "md2023" in the box under saved sessions and click save.
- 8. Double-click on the "md2023" that appears under saved sessions.
- 9. Log in with your username. Your key should be used automatically.
- 10. For future logins, just double-click the "md2023" saved session.

#### For the IU servers:

- 1. Open PuTTY
- 2. If you need to view graphics from the server: Under Category, click on SSH > Auth and check the "Enable X11 forwarding" box.
- 3. Under Category, click Session and input the server address (quartz.uits.iu.edu or carbonate.uits.iu.edu) in the "host name" box
- 4. Type "md2023\_quartz" or "md2023\_carbonate" in the box under saved sessions and click save.
- 5. Double-click on the "md2023\_quartz" or "md2023\_carbonate" that appears under saved sessions.
- 6. Log in with your username and password.
- 7. You'll need to complete 2-factor authentication with Duo to complete your login.
- 8. For future logins, just double-click the "md2023\_quartz" or "md2023\_carbonate" saved session.





## Submitting jobs in a cluster environment with Slurm

## Batch scripts

Both quartz and carbonate are computing *clusters* made up of many individual computers, or *nodes*. Lots of people are constantly running lots of demanding programs, so we need a way to make sure everyone's jobs are managed fairly. Enter SLURM. SLURM is a scheduler that will add your job to a queue; it will run on an appropriate node when it's its turn.

To run anything other than basic commands, you'll need to create a script - a text file with some header lines that tells SLURM details about the job you want to run, followed by the commands themselves. Here's an example:

```
#SBATCH --job-name=your job name
                     ## (-N) number of nodes to use
                                     ## Name of the iob.
#SBATCH --nodes=1
#SBATCH --ntasks=1
launch. Typically 1 unless you're getting fancy.
#SBATCH --cpus-per-task=4
                           ## number of cores the iob needs
#SBATCH --error=%x-%j.err ## error log file. %x is replaced
by numeric job ID, %j is job name as above
#SBATCH --out=%x-%j.out ## out log file. %x and %j same as
with --error
#SBATCH --mail-type=fail,invalid depend,begin,end
you if the job starts, ends, fails, or can't start due to
invalid dependencies. optional.
#SBATCH --mail-user=your.email@wherever.edu ##replace with
your email. optional.
#SBATCH --time=1-02:03:04 ## Format is DD-HH-MM-SS - this job
would time out at 1 day, 2 hours, 3 minutes, 4 seconds.
optional, but choosing a shorter time will likely get your
job running sooner.
#SBATCH --mem=8G ##maximum memory usage. technically
optional but recommended.
#SBATCH -A slurm-account-name ##Name of SLURM account to
charge.
#SBATCH -p some partition ##Name of partition to run on
```

module load prokka

prokka --cpus 4 --outdir my\_prokka\_results --prefix my\_genome
--metagenome --kingdom archaea my seq.fa

Let's break that down.

#!/bin/bash denotes which *shell* you're using. Shells are what actually provide the command line interface. Most linux systems use bash. If you're on a mac, you may be using zsh. On quartz, carbonate, and most linux systems, we want bash.

Lines starting with #SBATCH are instructions to SLURM. They are used to set the name of your job, the amount of nodes, CPUs, and memory required for your job, where to send error and output logs, where to send an email if your job changes status, and more. The comments (##) above explain what each line does. It's to your advantage to request reasonable resources for your job - smaller requests are more likely to get run sooner in the queue, and you won't take up unnecessary resources that other users could have used.

The account (-A) and partition (-p) will be specific to the cluster you're using, if they are required at all.

Software is often installed as *modules* on clusters. You can load just the modules required for your job, which helps avoid conflicts between software packages. In this case, I was running prokka, so I activated that module with module load prokka.

To see what modules are available, run module avail.

To see what modules are currently loaded, module list.

To deactivate a module, module unload module name.

Finally, after all the setup, I can run my actual prokka command.

HOT TIP: Save this example and use it as a template when creating new job scripts.

## Running your jobs

To submit your script to the queue, use the command <code>qsub</code>:

```
qsub myscript.slurm
```

To check the status of your jobs, use the squeue command:

```
squeue -u your username
```

You should see something like this:

```
JOBID USER ACCOUNT NAME ST TIME_LEFT NODES CPUS TRES_PER_N MIN_MEM NODELIST (REASON)
6663845 my_username some_acct job_name_1 R 14:13:50
1 8 N/A 32G node1 (None)
6663849 my_username some_acct job_name_2 R 14:13:50
1 4 N/A 16G node2 (None)
```

JOBID: numerical job ID

USER: you

ACCOUNT: jobs are typically assigned to a specific account if you're on a cluster that tracks or bills for usage.

NAME: job name

ST: Job state. Some common states are:

- PD: pending (waiting its turn)
- R: running
- F: failed
- CG: completing (in the process of finishing; nearly done)
- CD: completed
- OOM: out of memory
- TO: timeout (job reached time limit)

TIME\_LEFT: remaining time until job reaches time limit

NODES: Number of nodes used

CPUs: Number of CPUs used

TRES\_PER\_N: Refers to **T**rackable **RES**ources - you won't need to worry about this.

MIN\_MEM: Memory you requested for your job

NODELIST: Names of node(s) being used by your job

REASON: Explains reasons for job state. If your job is running normally, this will be (None). If something has gone wrong, this may be helpful in figuring out why.

Once your job has completed, it will disappear from the queue. By default, jobs run in the directory you submitted the job from - look there for your output files.

If you need to cancel a job, use the scancel command. To cancel the first job above:

scancel 6663845

#### Interactive sessions

If you need to run a job interactively, rather than submitting a batch script, you can start an interactive session:

```
srun -p some-partition -A slurm-account-name --pty bash
```

This will open an interactive bash terminal. Make sure to end the session when you're done - Ctrl+D will disconnect you.

#### More info

IU's Slurm help page is here: https://kb.iu.edu/d/awrz