

# Linux

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# Linux commands

Using ClustalW to align two sequences

# Step 1A: Accessing the IGB Biocluster

Open **Putty.exe**

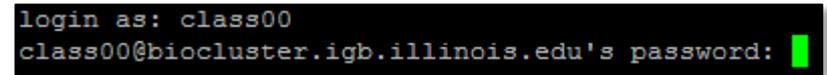
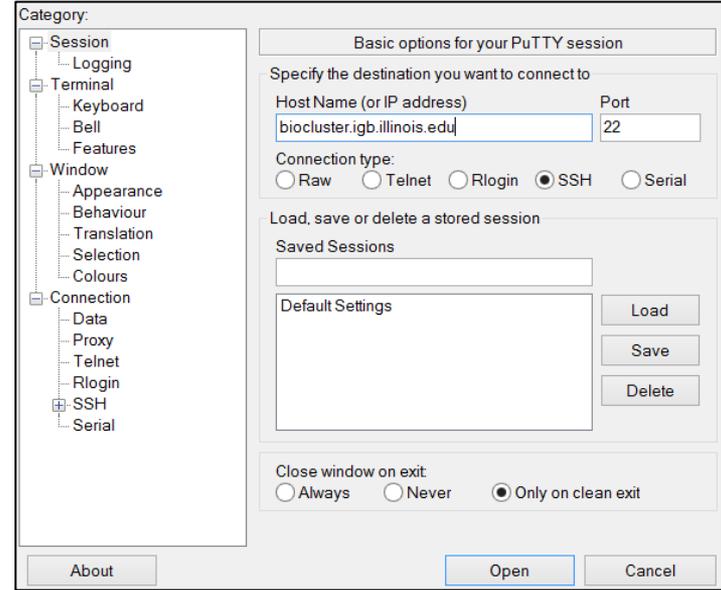
In the **hostname** textbox type:

`biologin.igb.illinois.edu`

Click **Open**

If popup appears, Click **Yes**

Enter login credentials assigned to you; example, user **class00**. You will not see any characters on screen when typing in password. Just type it.



**Now you are all set!**

# Step 1B: Listing files and directories (ls)

```
$ ls
```

```
# listing files in your current directory. When you first login, your  
directory is your home directory.
```

# Step 1C: Making Directories (mkdir)

```
$ mkdir ~/01_Linux
```

```
# create a subdirectory in your home directory. The tilde ~ character refers to your home directory.
```

```
$ ls
```

```
# to see the directory you just created.
```

## Step 1D: Changing directory (cd)

The lab is located in the following directory:

```
/home/classroom/mayo/2020/01_Linux
```

```
$ cd /home/classroom/mayo/2020/01_Linux
# tip: use "tab" for auto-completetion for path
$ ls
# to see the contents. You should see seqs.fa
```

## Step 1E: Print working directory (pwd)

```
$ pwd
# to see the full pathname. You should see
"/home/classroom/mayo/2020/01_Linux"
```

## Step 1F: Copying files (cp)

Copy `seqs.fa` from the data directory to your working directory.

```
$ cp /home/classroom/mayo/2020/01_Linux/seqs.fa ~/01_Linux/  
# tip: use "tab" for autocompletetion for path  
$ cd ~/01_Linux/
```

## Step 1G: Displaying the contents of a file on the screen (more)

```
$ more seqs.fa  
# you should see two sequences on your screen  
>seq1  
GATCGAGCGATCGTGCAGC  
GCAGAATGCGCGCTAG  
>seq2
```

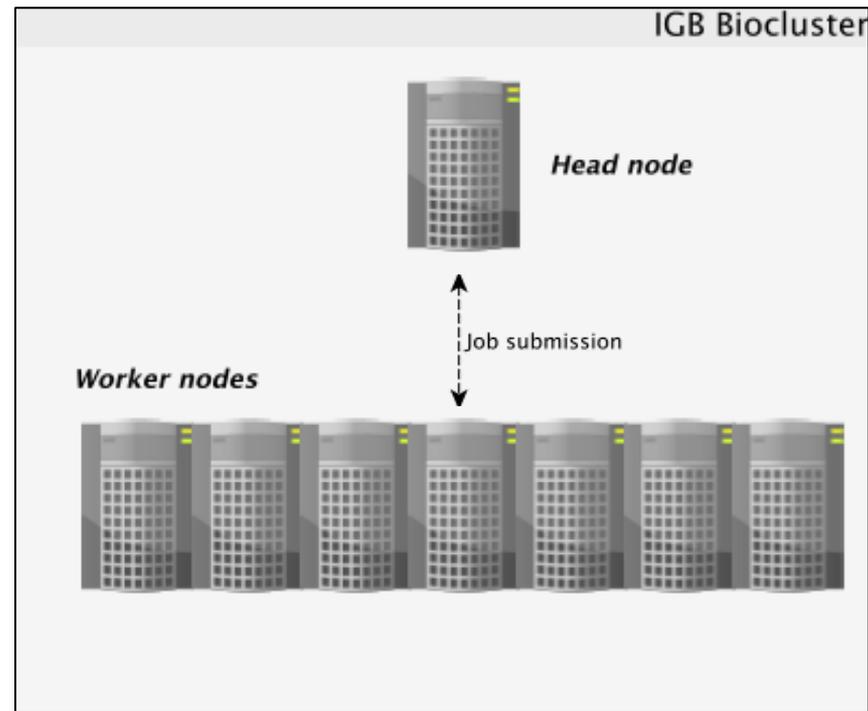
# Commands Summary

Command	Meaning
ls	list files and directories
mkdir directory	make a directory
cd directory	change to named directory
cd ~	change to home directory
cd ..	change to parent directory
pwd	display the path of the current directory
cp file1 file2	cp file1 and call it file2
more file	display the contents of a file

# Useful tips

Command	Meaning
tab	auto-complete path
↑	retrieve previous commands

# Step 1H: Run sequence alignment program



## Accessing the IGB Biocluster

# Step 1H: Run sequence alignment program

```
$ srun -p classroom -c 2 --mem 8000 --pty bash # SKIP IF DONE
# Open interactive session on biocluster with 2 cpus and 8G memory.

$ module load ClustalW2 # Load sequence aligner into the shell environment.

$ module list #See loaded tools

$ clustalw2 -INFILE=seqs.fa # Run the clustalW sequence aligner.
```

# Step 1H: Run sequence alignment program

You will see this on your screen, when the program is done.

```
CLUSTAL 2.1 Multiple Sequence Alignments
```

```
Sequence format is Pearson
```

```
Sequence 1: seq1          35 bp
```

```
Sequence 2: seq2          32 bp
```

```
Start of Pairwise alignments
```

```
Aligning...
```

```
Sequences (1:2) Aligned. Score: 21
```

```
Guide tree file created: [seqs.dnd]
```

```
There are 1 groups
```

```
Start of Multiple Alignment
```

```
Aligning...
```

```
Group 1:                      Delayed
```

```
Alignment Score 47
```

```
CLUSTAL-Alignment file created [seqs.aln]
```

# Step 1H: Run sequence alignment program

The alignment result is in `seqs.aln`. Use `more` command to see the result.

```
$ more seqs.aln
# You should see the following on your screen.
CLUSTAL 2.1 multiple sequence alignment
seq1          GATCGAGCGA-TCGTGCAGCGCAGAATGCGCGCTAG
seq2          GGTAGGGTAAATTGCCTACCGTCGATCGAGTA----
              * * * * * * * * * * * * * * * * * *
```

Exit putty by either closing the window or typing 'exit' in the command prompt.