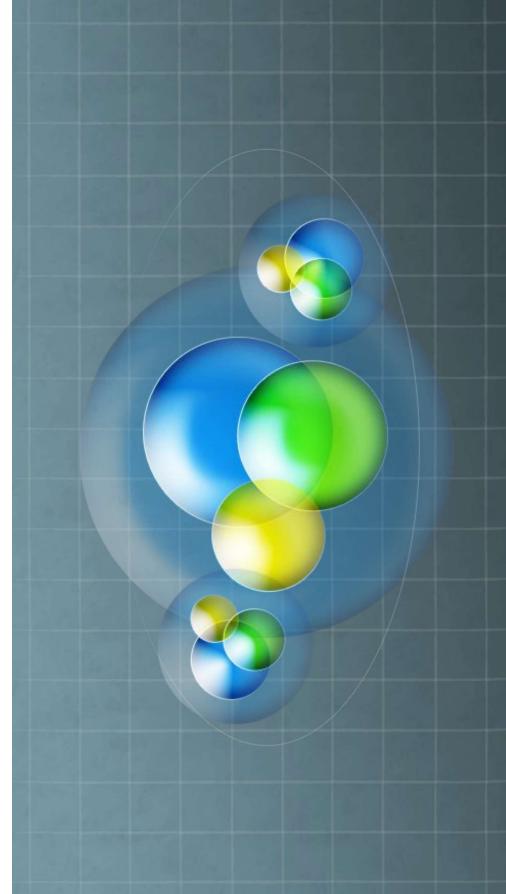
- Visit www.umich.edu/~cja/ and click "STATS 531 Introduction to Flux" to view these slides.
- Login to Flux using the instructions on the first few slides.



Dr Charles J Antonelli, LSA IT ARS Mark Montague, LSA IT ARS February, 2018



STATS 531 Introduction to Flux

Dr Charles J Antonelli, LSA IT ARS Mark Montague, LSA IT ARS March, 2018

Preliminaries

Connecting to Flux

(Windows)

- nstall the PuTTY/WinSCP installer for Windows: Select + under Get Going Get U-M PuTTY (UM_PuTTY_WinSCP.zip) Execute_the installer (UM_PuTTY_WinSCP.exe) If you don't have PuTTY installed Accept all defaults Compute at the U:
- /sourceforge.net/projects/xming/files/Xming-fonts/7.7.0.10/ Get the XMing fonts installer (Xming-fonts-7-7-0-10-setup.exe) Get the Xming installer (Xming-6-9-0-31-setup.exe) If you don't see this icon in your system tray Install the Xming X Server for Windows **Execute the installer Execute the installer** Accept all defaults Accept all defaults 7

Connecting to Flux

(Windows)

- Double-click "UM Internet Access Kit" icon on Desktop Double-click PuTTY application within
- Under Connection | SSH | X11, ensure Enable X11 Forwarding is checked In the Putty Application that appears: Enter "flux-login.arc-ts.umich.edu" in the Host Name box Click "Open" at bottom

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- Login with uniqname, Kerberos password, and Duo In the terminal window that appears: ψ
- 4. This creates an ssh session on Flux

Connecting to Flux

(Linux & macOS)

- . Start a terminal window
- ssh -X uniqname@flux-login.arc-ts.umich.edu In the terminal window that appears, type 7
- Login to Flux with Kerberos password and Duo ψ

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Roadmap

- **Preliminaries**
- Introduction to Flux
 The command line
 Flux mechanics
 Parallel R and pomp

ntroduction to Flux

XIII

Flux is a university-wide shared computational discovery , high-performance computing service.

- Provided by Advanced Research Computing at U-M
- Procurement, licensing, billing by U-M ITS
- Interdisciplinary since 2010



http://arc-ts.umich.edu/resources/compute-resources/

e Flux cluster

nodes

Compute nodes

Data transfer

node



Storage

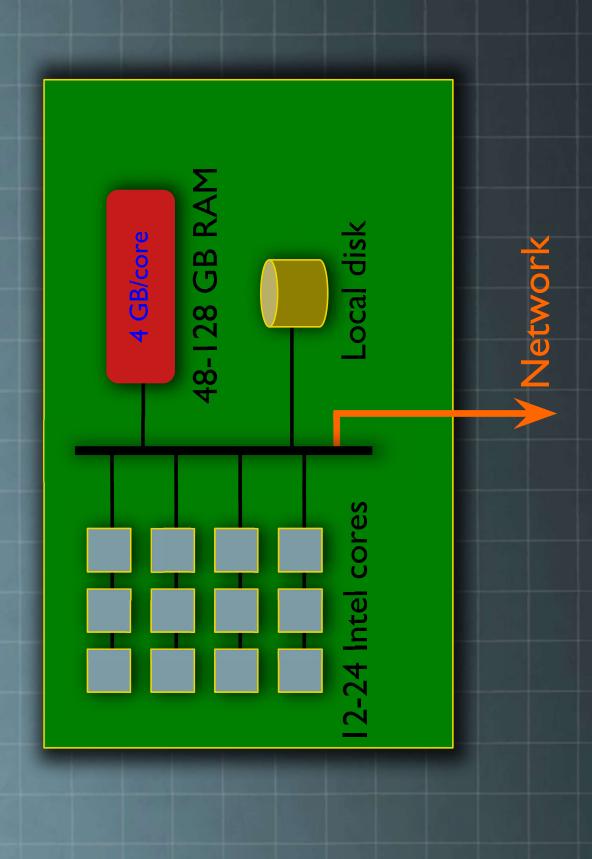




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A Standard Flux node



Using Flux

Three basic requirements:

A Flux login account https://arc-ts.umich.edu/fluxform

A Flux allocation

stats531w18_flux

http://its.umich.edu/two-factor-authentication A Duo app on your smartphone or tablet

Logging in to Flux

ssh -X *login@*flux-login.arc-ts.umich.edu

Campus wired or MWireless

Otherwise use VPN, or

ssh login.itd.umich.edu first

he command line

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~

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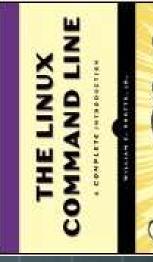
Command Line Reference

"The Linux Command Line: A Complete Introduction," No Starch Press, January 2012. William E Shotts, Jr.,

http://linuxcommand.org/tlcl.php.

Download Creative Commons Licensed version at

http://downloads.sourceforge.net/project/ inuxcommand/TLCL/13.07/TLCL-13.07.pdf





At the command prompt

- Basic input line editing commands
- **Backspace** erases previous character
- Left and right arrow move insertion point on the line
- Control-U erases the line to the insertion point, so you can start over
- **Enter** executes the line you typed
- Control-C interrupts whatever command you started and returns you to the shell prompt (usually)
- Up and down arrow will access your command history
- Type "exit" without the quotes to exit the shell

Flux mechanics

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Cluster batch workflow

- You create a batch script and submit it to PBS
- PBS schedules your job, and it enters the flux queue
- When its turn arrives, your job will execute the batch script
- Your script has access to all Flux applications and data
- When your script completes, anything it sent to standard output and error are saved in files stored in your submission directory
- You can ask that email be sent to you when your jobs starts, ends,
- You can check on the status of your job at any time, or delete it if it's not doing what you want
- A short time after your job completes, it disappears from PBS

```
#PBS -A youralloc_flux
#PBS -q flux
#PBS -I nodes=1:ppn=12,mem=47gb,walltime=00:05:00
#PBS -m abe
#PBS -N yourjobname
#PBS -V
```

cd \$PBS O WORKDIR R CMD BATCH --vanilla myscript.R myscript.out

#Your Code Goes Below:

Copying data

Using command line programs:

scp: copies files between hosts on a network over ssh scp *localfile uniqname@flux-xfer.arc-ts.umich.edu:remotefile* scp *localdir uniqname@flux-xfer.arc-ts.umich.edu:remotedir* scp *-r localdir uniqname@flux-xfer.arc-ts.umich.edu:remotefile localfile* scp *uniqname@flux-login.arc-ts.umich.edu:remotefile localfile*

Use "." as destination to copy to your Flux home directory: scp *localfile uniqname@*flux-xfer.arc-ts.umich.edu:.

... or to your Flux scratch directory:

scp localfile

*uniqname@*flux-xfer.arc-ts.umich.edu:/scratch/*allocname/uniqname*

sftp: an interactive file transfer program over ssh (a secure ftp) sftp uniquame@flux-xfer.arc-ts.umich.edu

Using graphical (GUI) applications: FileZilla (cross-platform):

http://filezilla-project.org/

Cyberduck (Mac): WinSCP (Windows):

nttps://cyberduck.io

nttp://www.itcs.umich.edu/bluedisc/

Basic batch commands

Once you have a script, submit it:
 qsub scriptfile

\$ qsub singlenode.pbs 6023521.nyx.engin.umich.edu You can check on the job status
 qstat jobid

qstat -u user
\$ qstat -u cja
nyx.engin.umich.edu:

SessID NDS Jobname hpc101i Username Queue flux cja 6023521.nyx.engi Job ID

Reg'd Elap TSK Memory Time S Time

00:05 Q

• To delete your job qdel jobid

Modules

The module command allows you to specify what Flux software you want to use

Search for string in module descrip Search for string in all module doc -- Add *path* to module search path Show versions of module *name* -- Load module name for use Show all available modules Show loaded modules -- Unload module name -- List all options string unload name key *string* spider *str*. load name use *path* avail avail module module module

Enter these commands at any time during your session

You can load multiple modules in the same login session

Software modules remain available throughout the session

Modules

- Once loaded, you can group a set of modules into a module save myset module set
- A module set can be restored at any time module restore myset
- List all module sets you've defined module savelist
- Create a module set to be loaded each time you log in module save

Lab

Task: Use the R multicore package

Copy sample code to your login directory

/scratch/data/workshops/stats/stats-samplecode.tar.gz .

-zxvf stats-sample-code.tar.gz

cd ./stats-sample-code

Examine labs. pbs and labs. R

•

Lab

Task: Use the R multicore package

- module load R
- Submit your job to Flux qsub lab3.pbs
- Watch the progress of your job qstat -u uniqname
- where uniqname is your own uniqname
- When complete, look at the job's output lab3.out

Parallel R and

Random numbers

Usually based on linear recurrences modulo m Generate stream of pseudorandom numbers Initialize the generator with some seed s Serial pseudorandom sequence

Performance, Reproducibility, Serializability Parallel pseudorandom sequence

Install pomp

Task: Install pomp package

- module load R/3.4.1
- install.packages ("pomp", repos="https: /cran.mtu.edu/")
- (answer y to "Would you like to use a personal library instead?" and "Would you like to create a personal library")

List of mirrors: https://cran.r-project.org/mirrors.htm

Multicore example

```
# Code shown below from: pomp-par.R
# To submit job: qsub pomp-par.pbs
```

```
cores <- as.numeric(Sys.getenv('PBS_NP', unset='8'))
                                                                  set.seed(2018,kind="L'Ecuyer")
                                                                                                                                       cl <- makeCluster(cores)
                                                                                                                                                                       registerDoParallel(cl)
                                library(doParallel)
rm(list=ls())
                                                                                                                                                                                                           trials <- 100
```

```
Multicore example
                                                                                                                                                                   .options.multicore=list(set.seed=TRUE)
                                                                                                                                                                                                                                                                     pompExample("gompertz")
                                                                                                                   r <- foreach(icount(trials),
                                                                 # pomp-par.R continued...
                                                                                                                                                                                                                                                                                                                                                                                                                       save(r,file="sims.Rda")
                                                                                                                                                                                                                                                                                          simulate(gompertz)
                                                                                                                                           .inorder=FALSE,
                                                                                                                                                                                                                                             for (i in 1:100) {
                                                                                                                                                                                                                     library(pomp)
                                                                                                                                                                                                                                                                                                                                                                                             stopCluster(cl)
                                                                                                                                                                                           ) %dopar% {
                                                                                             system.time(
```

MPI example

- Change the R module:module unload R/3.4.1module load Rmpi/3.4.1
- Install doMPI package:
- Rmpi
- install.packages("doMPI", repos="https://cran.mtu.edu/")

MPI example

```
# changes from multicore example shown in green
                                                    # Code shown below from: pomp-mpi.R
# To submit job: qsub pomp-mpi.pbs
```

```
cores <- as.numeric(Sys.getenv('PBS_NP', unset='8'))
                                                                set.seed(2018,kind="L'Ecuyer")
                                                                                                                                   cl <- startMPIcluster(cores)
                                                                                                                                                                     registerDoMPI(cl)
                                 library(doMPI)
rm(list=ls())
                                                                                                                                                                                                         trials <- 100
```

```
MPI example
                                                                                                                                                             .options.multicore=list(set.seed=TRUE)
                                                                                                                                                                                                                                                                pompExample("gompertz")
                                                                                                              r <- foreach(icount(trials),
                                                            # pomp-mpi.R continued...
                                                                                                                                                                                                                                                                                                                                                                                                                 save(r,file="sims.Rda")
                                                                                                                                                                                                                                                                                     simulate(gompertz)
                                                                                                                                      .inorder=FALSE,
                                                                                                                                                                                                                                       for (i in 1:100) {
                                                                                                                                                                                                               library(pomp)
                                                                                                                                                                                                                                                                                                                                                                                         closeCluster(cl)
                                                                                                                                                                                      ) %dopar% {
                                                                                       system.time(
```

Interactive jobs

You can submit jobs interactively:

```
qsub -I -X -V -l nodes=1:ppn=2 -l walltime=15:00
                                                 -A youralloc_flux -1 qos=flux -q flux
```

- This queues a job as usual
- Your terminal session will be blocked until the job runs
- When your job runs, you'll get an interactive shell
- When you exit the shell your job is deleted
- Interactive jobs allow you to
- Develop and test on cluster node(s)
- Execute GUI tools on a cluster node
- Utilize a parallel debugger interactively

Lab

Task: Use an interactive PBS session

- module load R
- qsub -I -V -1 nodes=1:ppn=2 -1 walltime=30:00 stats531w18_flux -q flux Start an interactive session

4-

Run R in the interactive shell
 cd \$PBS_O_WORKDIR
 R

- System-level
- freenodes
- [l-] sapousqd
- Account-level
- mdiag -a acct
- showq [-r][-i][-b][-w acct=acct]
- freealloc [--jobs] acct
- idlenodes acct [property]

shows available nodes for acct with property

with –jobs, shows resources in use

free resources in acct

with -r|i|b show more info for that job state

running/idle/blocked jobs for acct

cores & users for account acct

with -l, list only nodes marked down

nodes, states, properties

aggregate node/core busy/free

- User-level
- mdiag -u uniq
- showq [-r][-i][-b][-w user=uniq]

#running/idle/blocked jobs for uniq

allocations for user uniq

- Job-level
- qstat -f jobno
- gstat -n jobno
- checkjob [-v] jobno
- qpeek jobno

full info for job jobno # show nodes/cores where jobno running # show why jobno not running # peek at script output while jobno is running 35

Resources

- ARC User Guide http://arc-ts.umich.edu//flux-user-guide/
- http://arc-ts.umich.edu/flux/ ARC Flux pages
- Software Catalog http://arc-ts.umich.edu/software/

•

- Quick Start Guide
- http://arc-ts.umich.edu/flux/using-flux/flux-in-10-easy-steps/
- Flux FAQs

•

- http://arc-ts.umich.edu/flux/flux-faqs/
- For assistance, send email to:

•

Read by a team of people including unit support staff

hpc-support@umich.edu

- Can help with Flux operational and usage questions
- Programming support available

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