

# ClusterJob

# An automated system for massive computational experiments

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# **Getting Started**

ClusterJob (CJ) is an automation system for massive computational experiments. Using CJ, one can conduct large-scale computational studies on remote clusters in a painless and reproducible manner. CJ builds 'reproducible' computational packages that are easy to reuse and share with others. CJ is written mainly in perl and has simple commands that are easy to learn. The project started at Stanford University by Hatef Monajemi and his PhD advisor David L. Donoho with the goal of making large-scale scientific computing simpler and more efficient. Current implementation allows submission of Matlab and Python jobs.

# CJ on Github

You may access the CJ Github repository at https://github.com/monajemi/clusterjob
This book is a work in progress. You may find the AsciiDoc source code for this book at https://github.com/monajemi/CJ-book.

A PDF copy of this documentation can be found at http://clusterjob.org/documentation/book.pdf.

### Setting up CJ

Installing a CJ agent is very simple. Follow the two steps below to set one up!

#### Step 1: Setting up authentication key

CJ assumes that the secure shell (ssh) to cluster is handled in a password-free manner. There are various ways to achieve this. Some clusters use Kerberos and some others might use ssh-keygen. We will explain how you could setup a key using ssh-keygen in this document. ssh-keygen is a Unix utility that is used to generate, manage, and convert authentication keys for ssh authentication. If you already have a password-free connection to your cluster, you may skip this step.

#### 1. Check SSH setup

You can check to see if you already set up the public ssh key. Open a terminal and enter:

```
$ ls -al ~/.ssh
```

if you see any of the following, you probably have already setup the ssh-keygen

- id\_rsa.pub
- id\_dsa.pub
- id\_ecdsa.pub
- · id\_ed25519.pub

#### 2. Generate a SSH key on your machine

If you have not setup ssh key, do not worry. It is very simple to set up. Open terminal, and enter,

```
$ ssh-keygen -t rsa -C "your_email@example.com"
# This will generate a SSH key. Just use the default setting if
# you are asked questions in the process of key generation
```

This will generate a key in '~/.ssh/id\_rsa.pub'

3. Copy the key to remote server

The last step is to copy the content of your \$HOME/.ssh/id\_rsa.pub to \$HOME/.ssh/authorized\_keys located on the remote server

4. Check connection

You should check your authentication keys by trying to connect to server. In your terminal enter,

```
$ ssh username@cluster.stanford.edu
```

Some clusters use the so-called Multi-Factor Authentication (e.g., Duo Two-Factor). In this case, each time you log on, you need to provide an additional code. In order to avoid it, you need to add the following lines for each cluster (here corn.stanford.edu) to your local \$HOME/.ssh/config file. This will create a tunnel on your first login, and uses the same tunnel for your subsequent connections.



```
Host corn corn?? corn.stanford.edu corn??.stanford.edu
ControlMaster auto
ControlPath ~/.ssh/%r@%h:%p
ControlPersist yes
```

If you are using Stanford clusters, you may check https://web.stanford.edu/group/farmshare/cgi-bin/wiki/index.php/Advanced\_Connection\_Options for more info.

### Step 2: Installing CJ

1. Clone ClusterJob from GitHub

Clone ClusterJob to a directory where you would like to install it, say ~/CJ\_install

```
cd ~/
rm -rf monajemi-clusterjob-* CJ_install
curl -sL https://github.com/monajemi/clusterjob/tarball/master | tar -zx - && mv
monajemi-clusterjob-* CJ_install
```

You may also use the following alternative command to clone CJ with the entire history

```
git clone https://github.com/monajemi/clusterjob.git ~/CJ_install
```

#### 2. Install perl dependencies

ClusterJob is written in perl and so depends on other perl modules. You can install CJ dependencies via cpan or cpanm. Copy and paste the following lines into your terminal:

```
sudo cpan -i Data::Dumper Data::UUID FindBin File::chdir File::Basename File::Spec IO::Socket::INET IO::Socket::SSL Getopt::Declare Term::ReadLine JSON::PP JSON::XS Digest::SHA Time::Local Time::Piece Moo HTTP::Thin HTTP::Request::Common JSON URI
```

#### 3. Provide configuration info

You will need to update the contents of ~/CJ\_install/cj\_config file to reflect your own information.

This file contains the following information:

```
CJID <YOUR CJ ID>
CJKEY <YOUR CJ KEY>
SYNC_TYPE auto
SYNC_INTERVAL 300
```

1. To use CJ remotely, you need to provide your unique CJID and CJKEY obtained from http://clusterjob.org.



- 2. If you plan to use CJ locally without syncing to CJ remote database, you must provide an arbitrary CJID.
- 3. If you do not include a CJKEY, your meta data will not sync to CJ remote database even if you have registered an acount on http://clusterjob.org.



#### 4. Provide cluster info

You will need to update the contents of /CJ\_install/ssh\_config file to reflect your own server setup.

The CJ convention for each remote machine is

```
[MACHINE_ALIAS]
           your host
Host
User
           your username
           batch queue system
Bqs
           CJ remote repository
Repo
           Matlab version
MAT
MATlib
           Matlab libraries
Python
           Python version
Pythonlib Python packages/channels
[MACHINE_ALIAS]
```

#### Example:

```
[sherlock]
Host
            login.sherlock.stanford.edu
            monajemi
User
            SLURM
Bqs
Repo
            /scratch/users/monajemi/CJRepo Remote
MAT
            matlab/R2017a
MATlib
            ~/BPDN/CVX/cvx:~/mosek/7/toolbox/r2013a
Python
            python/3.6
Pythonlib
            scipy:cuda80:torchvision:pytorch:-c soumith
[sherlock]
```



You can add as many machines as you want to ssh\_config.

#### 5. Build an alias for CJ

For easy use, you may want to add an alias for calling src/CJ.pl to your ~/.profile or ~/.bashrc:

```
$ echo "alias cj='perl ~/CJ_install/src/CJ.pl'" >> ~/.bashrc
$ echo "alias cj='perl ~/CJ_install/src/CJ.pl'" >> ~/.profile
```

#### 6. Initialize your CJ agent

You may now initialize your CJ agent by

```
$ cj init
```

You may check if the agent is installed by



\$ cj who

This should print out something like

agent: E9078FA4-8423-11E6-B9A8-DFE0D454C74A



You may install as many CJ agents as needed on a single machine or different machines.

# **CJ Basics**

After you have successfully started a CJ agent, it is time to have some fun! In what follows, we will demonstrate how you can use CJ to make a reproducible computational package from your MATLAB code and run it on a remote cluster in a hassle-free way.

## A simple Matlab example

We will now go over a simple example of running a Matlab code on a cluster to demonstrate the usage of ClusterJob.

#### **Running Jobs**

Suppose we would like to run the following simple matlab code. The code includes two for loops indexed by i and j, and for each such combination it write a line to the file results.txt.

simpleExample.m

This code runs on our personal machine without any error. We now wish to run this code on a cluster named **sherlock** using ClusterJob. To run this code serially, once we are in the folder containing **simpleExample.m**, we simply type the following command in a terminal.

```
$ cj run simpleExample.m sherlock -m "This is a simple test for run command"
```

This command should result in the following output:

```
CJmessage::runing [simpleExample.m] on [sherlock]
CJmessage::Sending from: /Users/hatef/github_projects/clusterjob/example/MATLAB
CJmessage::Creating reproducible script(s) reproduce_simpleExample.m
CJmessage::Sending package 07264a5d
CJmessage::Submitting job
CJmessage::1 job(s) submitted (10097640)
```

The run command uses only one processing core. We could however run our code in parallel by simply changing the command to parrun:

\$ cj parrun simpleExample.m sherlock -m "This is a simple test for parrun command"

This time, you should see the following output informing you of submitting 15 jobs instead.

CJmessage::parruning [simpleExample.m] on [sherlock]

CJmessage::Sending from: /Users/hatef/github\_projects/clusterjob/example/MATLAB

CJmessage::Creating reproducible script(s) reproduce simpleExample.m

CJmessage::Sending package 0ed00c68

CJmessage::Submitting job(s)

CJmessage::15 job(s) submitted (10097772-10097786)



The parallelization of your code happens automatically with no further effort from you.

As you can see, each instance of the run command produces a reproducible package with a distinct PID (for Package IDentifier). A PID is a SHA1 code, which is essentially a 40 digits long hexadecimal number. Though PIDs are 40 digits long, when using CJ commands you can provide only a short version of it, which contains the initial 8 characters only.



A PID is valid as long as it is a hexadecimal number of length 8 or more



The full PID can be retrieved using \$ cj info short\_pid command.

#### **Checking Status**

You can check the status of your jobs using state command.

\$ cj state 07264a5d

pid 07264a5d33bab71c1463f651b1ff920f6d32bb1c
remote\_account: monajemi@sherlock.stanford.edu

job\_id: 10109624
state: COMPLETED

Or for a parallel case,

```
$ cj state 0ed00c68
pid 0ed00c6851c504af7d8064a954aba44cf1da40f2
remote_account: monajemi@sherlock.stanford.edu
                      COMPLETED
1
       10097772
2
       10097773
                      COMPLETED
3
       10097774
                      COMPLETED
4
       10097775
                      COMPLETED
5
       10097776
                      COMPLETED
6
       10097777
                      COMPLETED
7
       10097778
                      COMPLETED
8
       10097779
                      COMPLETED
9
       10097780
                      COMPLETED
10
       10097781
                      COMPLETED
11
       10097782
                      COMPLETED
12
       10097783
                      COMPLETED
13
       10097784
                      COMPLETED
14
       10097785
                      COMPLETED
15
       10097786
                      COMPLETED
```

#### **Getting Results**

You can pull your computational package entirely or partially using get command. this command will pull the package associated with a particular PID into a temporary directory on your local machine:

```
$ cj get 07264a5d

CJmessage::Getting results from 'sherlock'
CJmessage::Please see your last results in
/Users/hatef/CJ_get_tmp/07264a5d33bab71c1463f651b1ff920f6d32bb1c
```

If we now look at the contents of this package, we would see the following. You may want to take a moment to see what is inside each file.

#### CJ CONFIRMATION.TXT

contains confirmation that this package was run by ClusterJob.

#### CJrandState.mat

contains the information about the random seed that generated the results so that other people can reproduce your results.

#### results.txt

is the output of your code

#### reproduce\_simpleExample.m

reproduces these results upon execution.

For a parallel package, if we use get, we will see the following output:

```
$ cj get 0ed00c6851c504
   CJmessage::Getting results from 'sherlock'
   CJmessage::Run REDUCE before GET for reducing parrun packages
   CJmessage::Please see your last results in
   /Users/hatef/CJ_get_tmp/0ed00c6851c504af7d8064a954aba44cf1da40f2
```



For parallel packages, one typically needs to use reduce before get

Clearly, we are asked to use reduce before get for a parallel package. We will discuss reduce command in the next section. If we now look at the folder containing this parallel package, we see:

```
$ ls -lt /Users/hatef/CJ_get_tmp/0ed00c6851c504af7d8064a954aba44cf1da40f2
total 80
-rwxr-xr-x 1 hatef staff
                           2429 Oct 7 16:06 CJ_CONFIRMATION.TXT
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:28 13
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:26 15
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:26 14
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 11
                            306 Oct 6 22:25 10
drwxr-xr-x 9 hatef staff
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 6
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 7
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 8
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 3
                            306 Oct 6 22:25 5
drwxr-xr-x 9 hatef staff
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 9
                            306 Oct 6 22:25 4
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 12
drwxr-xr-x 9 hatef staff
drwxr-xr-x 9 hatef staff
                            306 Oct 6 22:25 2
                            306 Oct 6 22:24 1
drwxr-xr-x 9 hatef staff
-rw-r--r-- 1 hatef staff
                            435 Oct 6 22:23 qsub.info
-rw-r--r-- 1 hatef staff 11591 Oct 6 22:23 master.sh
-rw-r--r-- 1 hatef staff
                            308 Oct 6 22:23 simpleExample.m
```

We now see 15 directories for 15 jobs submitted. If we further look into one of these directories:

```
$ ls -lt /Users/hatef/CJ_get_tmp/0ed00c6851c504af7d8064a954aba44cf1da40f2/1 total 40
-rw-r--r-- 1 hatef staff 2865 Oct 6 22:24 CJrandState.mat
-rw-r--r-- 1 hatef staff 6 Oct 6 22:24 results.txt
drwxr-xr-x 5 hatef staff 170 Oct 6 22:23 logs
drwxr-xr-x 3 hatef staff 102 Oct 6 22:23 scripts
-rw-r--r-- 1 hatef staff 1953 Oct 6 22:23 bashMain.sh
-rw-r--r-- 1 hatef staff 515 Oct 6 22:23 reproduce_simpleExample.m
-rw-r--r-- 1 hatef staff 196 Oct 6 22:23 simpleExample.m
```

### **Reducing Results**

The reduce command is designed for reducing the results of a parallel run into a single file. This is because the parrun command executes the for loops in the main script in parallel by generating independent sub-directories and submitting a separate job for each index combination.

If we look at the results.txt file in sub-directories 1 and 2 for instance, we see:

```
$ cat /Users/hatef/CJ_get_tmp/0ed00c6851c504af7d8064a954aba44cf1da40f2/1/results.txt
1,1,2
$ cat /Users/hatef/CJ_get_tmp/0ed00c6851c504af7d8064a954aba44cf1da40f2/2/results.txt
1,2,3
```

We certainly want these results to be reduced to one single results.txt file. This is done via the reduce command.

```
$ cj reduce results.txt 0ed00c68
CJmessage::Checking progress of runs...
CJmessage::Reducing results...
CJmessage::Do you want to submit the reduce script to the queue via srun?(recommneded
for big jobs) Y/N?
n
SubPackage 1 Collected (6.67%)
SubPackage 2 Collected (13.33%)
SubPackage 3 Collected (20.00%)
SubPackage 4 Collected (26.67%)
SubPackage 5 Collected (33.33%)
SubPackage 6 Collected (40.00%)
SubPackage 7 Collected (46.67%)
SubPackage 8 Collected (53.33%)
SubPackage 9 Collected (60.00%)
SubPackage 10 Collected (66.67%)
SubPackage 11 Collected (73.33%)
SubPackage 12 Collected (80.00%)
SubPackage 13 Collected (86.67%)
SubPackage 14 Collected (93.33%)
SubPackage 15 Collected (100.00%)
CJmessage::Reducing results done! Please use "CJ get 0ed00c68 " to get your results.
```

This will produce the new file results.txt in the main directory, which contains:

```
$ cat /Users/hatef/CJ_get_tmp/0ed00c6851c504af7d8064a954aba44cf1da40f2/results.txt
1,1,2
1,2,3
1,3,4
1,4,5
1,5,6
2,1,3
2,2,4
2,3,5
2,4,6
2,5,7
3,1,4
3,2,5
3,3,6
3,4,7
3,5,8
```



The file to reduce can be any acceptable MATLAB output, for instance output.mat.

## CJ commands

To see a full list of options

```
cj -help
```

In what follows, we review important commands.

#### RUN

```
cj run <script> <machine> -dep <DEPENDENCY_FOLDER> -m <MESSAGE> -mem <MEM_REQUESTED>
```

#### **DEPLOY**

```
cj deploy <script> <machine> -dep <DEPENDENCY_FOLDER> -m <MESSAGE> -mem
<MEM_REQUESTED>
```

#### **PARRUN**

parellel run for embarssingly parallel problems.

```
cj parrun <script> <machine> -dep <DEPENDENCY_FOLDER> -m <MESSAGE> -mem
<MEM_REQUESTED>
```

#### RERUN

To rerun a previously failed package.

```
cj rerun <PACKAGE> <FOLDER_NUMBER>
```

#### STATE

To see the state of the last job submitted through CJ,

```
cj state
```

To see the state of a particular package,

```
cj state <PACKAGE>
```

To see the state of a particular folder in a parrun package,

```
cj state <PACKAGE> <FOLDER_NUMBER>
```

#### LOG

To see the last N instances of CJ run,

cj log <N>

To see detailed log of a particular package,

cj log <PACKAGE>

To see all the packages

cj log all

#### INFO

To see the information of the last call to CJ,

cj info

To see the information of a particular package,

cj info <PACKAGE>

#### **REDUCE**

To reduce the results of the last parrun call,

cj reduce <RESULTS\_FILENAME>

To reduce the results of a particular parrun package,

cj reduce <RESULTS\_FILENAME> <PACKAGE>

#### GET

To get the results of the last call back,

cj get

To get the results of a particular package call back,

cj get <PACKAGE>

#### **CLEAN**

To remove the last package and its associated jobs,

cj clean

To remove a particular package and its associated jobs,

cj clean <PACKAGE>

#### **SHOW**

To show the program run by CJ at the last call,

cj show

To show the program run by CJ for a particular package,

cj show <PACKAGE>