
Usersguide of TUHEP cluster

Here is the usersguide of TUHEP cluster. Please read this document carefully before using the cluster and network in the office. If you have problems, please contact Zhenwei Yang (9498, zhenwei.yang@gmail.com).

Some important reminders:

- 1、 Please close the tunet client or web-based logins in advance. If you forget it, please close them directly WITHOUT clicking “Disconnect” button. Otherwise, the whole network would be lost for several minutes.
- 2、 The main purpose of TUHEP cluster is data analysis of HEP and other scientific calculations. The gate node is “hepfarm” (hostname), with ip 166.111.32.48 and domain name <http://server.hep.tsinghua.edu.cn>. The gate node is merely for login from outside, and no other services are provided. One cannot run any resource time/resource-consuming program either.
If one need to submit a job or run other analysis computing, please login the computing nodes, hepfarm21/22/23 (details followed in next page).
- 3、 Disk quota is applied to hepfarm21/22/23. One cannot exceed 2GB in home directory. If you have large-size data, please put them in subdirectories of /projects.
- 4、 Please change your password frequently for safety.
- 5、 p2p applications are strictly banned both in the cluster and office. Net service in the office is only for research and teaching, and one cannot provide your account and password to others. Otherwise, your account would be banned.
- 6、 Attention please, for important private data, especially codes, please backup by yourself. There’s always potential risk of data lose.

Well, before going to next page, make sure again that you’ve read these reminders carefully.

5. Comments for the job script example

First of all, the first 4 lines are essential, and the “#”s here do NOT mean comments!!!

Line 1: to specify shell of the script.

Line 2: to specify the job queue. If the job is expected to be finished in 20 minutes, you may change the S queue: #PBS -q S

Line 3: to specify the job name.

Line 4: **usually just keep it as this default format**. PBS server would run your job in the most unoccupied nodes, and each job uses one CPU core. If you do want your job to be run in some specified node, like hepfarm22, change this line like this:

```
#PBS -l nodes=hepfarm22:ppn=1
```

Do NOT change the number “1” in “ppn=1”, unless you are sure that your program support parallel computing!

Line 5 and later: to go to the working directory, set up environments and run your programs.

Reminders: the submitted job will be run in batch mode, please do not add “&” in the end of qsub command.

Frequently used job commands:

qsub <jobscript>: to submit a job

qstat -n : to scan submitted jobs

qdel <jobID> : to delete a job