

# EMSL: Molecular Science Computing

## Quick Reference Guide

WEB: <http://www.emsl.pnl.gov/capabilities/computing/>

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E-mail Address: [mscf-consulting@emsl.pnl.gov](mailto:mscf-consulting@emsl.pnl.gov)

### External users log on to Chinook

\$ ssh <UserID>@chinook.emsl.pnl.gov

First: enter PASSCODE from secureID

Second: enter Kerberos password

### Internal users log on to Chinook

\$ ssh <UserID>@chinook.emsl.pnl.gov

(enter Kerberos password)

#### Note:

Machines use Secure Shell Protocol 2.

You may have to use ssh2 or ssh -2

### System Information

Machine status

\$ sinfo

Nodes available

\$ showbf

### Job Commands

Submit job

\$ msub *jobscript.msub*

Show all jobs in queue

\$ showq

Status of all your jobs

\$ showq | grep <UserID>

Cancel job

\$ canceljob *jobid*

Estimate Job start

\$ showstart *jobid*

### File Systems

NFS - permanent home directory; /home/UserID

NWfs - archive file system; access using sftp nwfs.emsl.pnl.gov

Global File System - cleaned periodically; /dtemp/UserID

Scratch - cleaned after job runs; /scratch **Version 2b: Aug, 2008**

The EMSL is the Environmental Molecular Sciences Laboratory and is located at Pacific Northwest National Laboratory (PNNL), one of the U.S. Department of Energy multiprogram national laboratories.

### Allocation Information (Gold)

Account balance and account name

\$ gbalance -h -u <UserID>

### Using modules to set up your environment

List loaded modules

\$ module list

Available modules

\$ module avail

Removing modules

\$ module unload <module>

Adding new modules

\$ module load <module>

Swapping modules

{Currently not working on Phase-1}

Removing all modules

\$ module purge

### Available modules

Default environment: pnnl\_env

Note: Environment set default Integer\*4

Integer\*8 environment: pnnl\_env/i8

Use: module purge

Use: module load pnnl\_env/1.2.i8

{Additional options to be added later}

### Batch

Control flags for batch job script file, jobscript. See our website for sample scripts.

```
#!/bin/csh
#MSUB -A <your Gold account>
#MSUB -l "walltime=04:30:00" # (HH:MM:SS)
#MSUB -l "nodes=<number-of-nodes>;ppn=8" # (8 cores/node)
#MSUB -m ae # e-mail on "abort" or at the "end"
#MSUB -o <your output file>.%j
#MSUB -e <your error file>.%j
#MSUB -N <job name>
#MSUB -M <your email address>
```

To run a program, add the following lines to the jobscript file for a batch job.

```
crun -nodes <# of nodes> -cores <# of cores> <your program> <your args>
```

To copy a file or binary to /scratch on all the nodes (using one CPU per node):

```
pdsh -w "${SLURM_JOB_NODELIST}" /bin/cp <file> /scratch/<file>
```

### Execute Batch Job

Submit your job to the queue by typing: msub *jobscript.msub*

### Interactive Job

Start your interactive job/session (X-windows required) by typing:

```
isub -A <your Gold account> -N <# of nodes> -W <time limit HH:MM> -s <shell>
```

To run a program in the interactive X terminal, type:

```
crun -nodes <# of nodes> -cores <# of cores> <your program> <your args>
```

### Intel Compiler

Default compiler is Intel compiler:

ifort - fortran77 and fortran 90 compiler

icc - C and C++ compiler

### Useful Intel compiler flags

-i8 : integer\*8 ; -i4 : integer\*4

-O2 : default optimization level

-FR : free format fortran

-Vaxlib : link in portability library

### Intel MKL BLAS & LAPACK

Note: variables MLIB\_LIB and MLIB\_LDFLAGS

are set by the module environment

For Integer\*4 and Integer\*8 compiled code, include on the link line: \${MLIB\_LDFLAGS}

This environmental variable includes both the -L and -l arguments for MKL

### Linking MPI with Intel compiler

Automatic: use mpif90 or mpicc

For Details include -show: mpif90 -show or mpicc -show

Note: If these do not work, contact MSC Consultants [use e-mail address below]

For assistance, email [mscf-consulting@emsl.pnl.gov](mailto:mscf-consulting@emsl.pnl.gov)