

## Lab 1: Getting Started

For this and all future labs, you will be working on the CS linux systems. You can either use a lab machine, or remotely log on to one (connect to `linuxlab.cs.pdx.edu`).

Download and unzip the file `lab1.zip` from D2L. You'll see a `lab1` directory with a `mpihosts` file and multiple versions of the sum program:

```
sum.c sum-pthd.c sum-omp.c sum-mpi.c sum1.chpl sum2.chpl
```

### Checking Your Environment

- Check to make sure you have the latest version of `gcc`:

```
linux> gcc --version
gcc (Ubuntu 4.8.4-2ubuntu1~14.04.1) 4.8.4
```

- Check to make sure you have access to the MPI compiler, `mpicc`:

```
linux> which mpicc
/usr/bin/mpicc
```

- Use `addpkg` to add the Chapel compiler to your environment:

```
linux> addpkg
...
```

Select `chapel-1.12`; exit, logout, then re-login.

### Compiling the Programs

- To compile Pthreads programs, use `gcc` with `-pthread` flag:

```
linux> gcc -pthread -g -o sum-pthd sum-pthd.c
```

- To compile OpenMP programs, use `gcc` with `"-fopenmp"` flag:

```
linux> gcc -fopenmp -g -o sum-omp sum-omp.c
```

- To compile MPI programs, use `mpicc`:

```
linux> mpicc -g -o sum-mpi sum-mpi.c
```

- To compile Chapel programs, use `chpl`:

```
linux> chpl -g -o sum1 sum1.chpl
```

### Running Pthreads and OpenMP Programs

These programs are run just like regular C programs:

```
linux> ./sum-pthd
linux> ./sum-omp
```

Note that for these two examples, the number of threads is hardwired in the program. In the future, we'll see how to make that adjustable.

**Exercise** Add a `printf` statement in the `worker()` routine in `sum-pthd.c` to show the id and the work range of each individual thread.

You may want to place this statement under a control flag:

```
#ifdef DEBUG
    printf(...);
#endif
```

This way, the same program can be compiled to two different versions:

```
linux> gcc -pthread -g -o sum-pthd sum-pthd.c           # regular version
linux> gcc -DDEBUG -pthread -g -o sum-pthd-debug sum-pthd.c # debug version
```

It is not as easy to do the same in the OpenMP program. We'll defer it to the future.

## Running MPI Programs

Before running MPI programs, you need to setup a host file. Copy `mpihosts` to your home directory, and set the following env variable:

```
setenv OMPI_MCA_orte_default_hostfile ~/mpihosts
```

You should include this line in your shell startup file (*e.g.* `.cshrc`) to avoid typing it in every time.

An MPI program is run with `mpirun`, with a flag `-n <#copies>` indicating the number of copies you'd like to execute:

```
linux> mpirun -n 4 ./sum-mpi    // running 4 copies of the program
```

Note that for this program, the number of program copies is specified externally at the time of execution.

**Exercise** Add a `printf` statement in `sum-mpi.c` to show the rank and size of the current copy of the program.

## Running Chapel Programs

For running Chapel programs, you need to set the following env variables in your shell startup file:

```
setenv CHPL_HOME /pkgs/chapel/chapel-1.12.0
setenv CHPL_HOST_PLATFORM linux64
setenv CHPL_COMM gasnet
setenv GASNET_SPAWNFN S
setenv GASNET_SSH_SERVERS "bevatron boson ..." # list of host names
setenv SSH_CMD ssh
```

A Chapel program is run with a flag `-nl <#locales>` indicating the number of locales you'd like to use:

```
linux> ./sum1 -nl 1           // running the program over 1 locale
```

**Exercise** In both programs, the problem domain size `N` is a configurable constant. Try to change it at the time of execution:

```
linux> ./sum1 --N=10 -nl 1
```