OpenMP Lab

https://github.com/ResearchComputing/USGS_2016/

February 9, 2016 Timothy Brown



Source Code

All of the source code is online, in a github repository:

https://github.com/ResearchComputing/USGS_2016_02_09-10

Login to Yeti

1. Log in to Yeti.

laptop ~\$ ssh yeti.cr.usgs.gov



2. Clone the git repository

yeti-login01 ~\$ git clone \
https://github.com/ResearchComputing/USGS_2016_02_09-10

3. Start a compute job.



 Load the Intel parallel studio module compute80 ~\$ module load intel/psxe-2015

This contains the C, C++ and Fortran compilers as well as the Intel MPI library.

| Language | Compiler |
|----------|----------|
| С | icc |
| C++ | icpc |
| Fortran | ifort |

Number of threads

Figure out how many threads we have.
Using the omp_get_thread_num() library call.

| Example | Source |
|---------|-----------------------------|
| Fortran | thread_num/thread_num_f.f90 |
| С | thread_num/thread_num_c.c |

Compiling the program.

```
Intel icc -qopenmp -o thread_num_c thread_num_c.c
GCC gcc -fopenmp -o thread_num_c thread_num_c.c
PGI pgcc -mp -o thread_num_c thread_num_c.c
IBM xlc -qsmp=omp -o thread_num_c thread_num_c.c
```

- Execute the program, specifying different numbers of threads.
 - 1. ./thread_num_c
 - 2. env OMP_NUM_THREADS=1 ./thread_num_c
 - 3. env OMP_NUM_THREADS=64 ./thread_num_c
- What is the output?
 - Threads printed out their identification number.
 - Random order of numbers. Threads execute independently and in general order will be random.

Preprocessor

Using the pre-processor with C and Fortran code.

Using ifdef's within the source code file, means you can compile your code with OpenMP directives with out using OpenMP.

| Example | Source |
|--------------------------|--------------|
| Pre-processor in Fortran | ifdef/if.F90 |

Loops

| Example | Source |
|------------------------------------|--------------|
| Array addition using a for loop | loop/for.c |
| First private in a parallel region | loop/first.c |
| Last private in a parallel region | loop/last.c |
| Reduction operator | loop/total.c |

Sections

| Example | Source |
|-----------------------------|-----------------------|
| Distribute independent work | sections/sections.f90 |

Synchronization

| Example | Source |
|---------------------------------|----------------|
| Synchronization using a barrier | sync/barrier.c |
| An ordered reduction | sync/ordered.c |

Alignment

Data alignment is crucial when using SIMD directives for vectorization.

| Example | Source |
|----------------------|-------------------|
| Alignment in C | align/align_c.c |
| Alignment in Fortran | align/align_f.F90 |

Note icc, gcc and gfortran pre-processors have __BIGGEST_ALIGNMENT__ defined. However ifort does not (it is currently an open feature request). So the Makefile defines it for Fortran.

Target

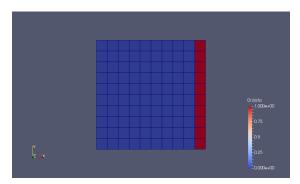
Using the OpenMP v4.5 target directives for off-loading to an accelerator.

| Example | Source |
|-------------|-------------------|
| Target in C | target/target_c.c |

Jacobi

The Jacobi method for stencil-based iterative solvers. The example is for a 3D diffusion equation.

| Example | Source |
|---------|--------|
| Fortran | jacobi |



- ► Load the HDF5 module for I/O. compute80 ~\$ module load tools/hdf5-1.8.15-intel
- Compile the serial jacobi program compute80 ~\$ make
- ► Execute the program with a $100 \times 100 \times 100$ domain and 1000 iterations. These are the deafults in the input.nl file.

```
compute80 ~$ ./jacobi -i input.nl
```

Questions?

Online Survey

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