Rules and Regulations of the 4<sup>th</sup> Annual IHPCSS Challenge



Trophy bears no relationship to reality.



# **Starting Point**

We give you working codes in MPI, OpenMP and OpenACC
 copy "challenge.tar" from /home/dsh/ihpcss18/.

- Set up to tackle a problem of size 672x672
  - challenge is to run a much larger problem

### Straightforward solutions

- no attempts at parallel (or serial) optimisation
- each only uses a single model

### No compilation instructions

you decide how best to compile



### **General Rules**

- Due Thursday midnight (!)
- 4 Nodes of Bridges
- Use any combination of MPI, OpenMP, OpenACC and Python
  - base versions (C and Fortran) for MPI, OpenMP and OpenACC on moodle
  - single tar file: challenge.tar
- How fast can you run a 10752 x 10752 Laplace code to convergence?
  - weird size chosen to decompose exactly on, e.g., 2, 4, 28 and 112 procs
  - can use smaller size of 672 x 672 for development



## **Some Specifics**

Can't change kernel (Must retain two core loops source)

- Can change number of MPI processes (Does not have to be 112 or 4)
- 1 Source File
- 1 Combined Environment/Compile/Submit/Execute script
  to make it easy for us to run your solutions!
- Mail to d.henty@epcc.ed.ac.uk by deadline



### **Rules For Lawyers**

#### No libraries

#### Don't mess with timer placement

• ?



# **Reality Checks**

Serial code converges at 3580 time steps. Yours should too.

- As we know, this is not enough to verify correctness. You should find point [8064][10702] in C and (10702,8064) in Fortran converges to 17.1 degrees.
- As discussed, the 10752 result differs from the 672 result.\*
  - smaller problem converges in 3264 time steps
  - check values: [504][622] in C, (622,504) in Fortran = 15.5 degrees
  - Plugging in Gauss-Seidel or Successive Over Relaxation (SOR) would be easy and interesting. But, not for our contest.

http://www.cs.berkeley.edu/~demmel/cs267/lecture24/lecture24.html is a brief analysis of these issues.



# Printing out the test point

- Straightforward in serial, OpenMP or OpenACC
  - single process, temperature stored in a single global array
- More complicated when you introduce MPI must locate owning process
  - if (8064/columns == mype+1) then
  - write(\*,\*) 'PE ', mype, ': T(10702,8064) = ', temperature(10702,columns)

end if

- if (8064/ROWS == my\_PE\_num+1)
- printf("PE %d: T(8064,10702) = %f\n", my\_PE\_num,Temperature[ROWS][10702]);
- This hacky piece of code requires at least 4 MPI processes!

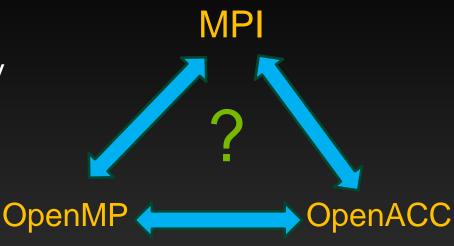
http://www.cs.berkeley.edu/~demmel/cs267/lecture24/lecture24.html is a brief analysis of these issues.



# **Suggested Things to Explore**

User Guide is your friend!

- Compiler flags -03
- Compiler
  - see Bridges documentation for how to use different modules
- MPI Environment Variables
  - 🔹 man mpi
- Thread placement google for KMP\_AFFINITY





# Decision

On Friday morning we will take the top self-reported speeds and run them in an interactive session

- Timings not within 10% of self-reported time will be disqualified
- Codes should print out "test point" at conclusion of run.
- Best of two runs for each finalist will determine winner



