SLAC Home | Computing Home | Unix at SLAC | High Performance Computing at SLAC | Parallel Computing at SLAC

Mixing MPI and OpenMP

Please make sure you read the SLAC specific instructions about MPI and OpenMP before you continue here.

Mixed "Hello World"

```
Name this little "Hello World" program hello.c:
#include <stdio.h>
#include "mpi.h"
#include <omp.h>
int main(int argc, char *argv[]) {
  int numprocs, rank, namelen;
  char processor_name[MPI_MAX_PROCESSOR_NAME];
  int iam = 0, np = 1;
  MPI Init(&argc, &argv);
  MPI Comm size(MPI COMM WORLD, &numprocs);
  MPI Comm rank(MPI COMM WORLD, &rank);
  MPI Get processor name(processor name, &namelen);
  #pragma omp parallel default(shared) private(iam, np)
    np = omp_get_num_threads();
    iam = omp_get_thread_num();
    printf("Hello from thread %d out of %d from process %d out of %d on %s\n",
           iam, np, rank, numprocs, processor name);
  }
  MPI Finalize();
}
```

Compiling and Linking Mixed MPI and OpenMP Programs

Once you have your example program, you can compile and link it with

- Linux: /afs/slac.stanford.edu/package/OpenMPI/bin/mpicc -openmp hello.c -o hello
- **Running Mixed Programs**

Solaris: N/A

• Interactively

```
alfw@morab> export OMP_NUM_THREADS=4
alfw@morab> /afs/slac.stanford.edu/package/OpenMPI/bin/mpirun --mca pls_rsh_agent ssh \
-np 2 -machinefile machinefile.morab -x OMP_NUM_THREADS ./hello
Hello from thread 0 out of 4 from process 0 out of 2 on morab006
Hello from thread 1 out of 4 from process 0 out of 2 on morab006
Hello from thread 2 out of 4 from process 0 out of 2 on morab006
Hello from thread 3 out of 4 from process 0 out of 2 on morab006
Hello from thread 0 out of 4 from process 1 out of 2 on morab001
Hello from thread 3 out of 4 from process 1 out of 2 on morab001
Hello from thread 1 out of 4 from process 1 out of 2 on morab001
Hello from thread 2 out of 4 from process 1 out of 2 on morab001
```

Note that you have to tell OpenMPI to set the OMP_NUM_THREADS environment variable for OpenMP for each process it starts using the -x OMP_NUM_THREADS command line argument.

• Executing via LSF Batch System

LSF's bsub command will pick up the OMP_NUM_THREADS and other environment variables and set them before starting your job on the different hosts.

alfw@morab> export OMP_NUM_THREADS=4 alfw@morab> bsub -a openmpi -q mpiq -n 4 ./hello

References

- SLAC specific MPI tutorial
 SLAC specific OpenMP tutorial
 Parallel Computing at SLAC

Alf Wachsmann. Last Modified: Mar 22, 2006.