

### PARALLEL PROGRAMMING...

Copyright 2023 Patrick Lemoine. All rights reserved.

### **Parallel Programming: Overview**

#### **SESSION 2/3**



#### **Programming Interface for parallel computing**

MPI (Message Passing Interface)

OpenMP (Open Multi-Processing)

# Programming interface for parallel computing

병렬 컴퓨팅을 위한 프로그래밍 인터페이스

### **Programming interface...**

OpenMP



### Remember

CPU CPU Core Core Memory Memory Private Network Arrays Interconnect CPU CPU Core Core Memory Memory

CPU Core CPU Core CPU Core

Memory, Shared Arrays etc.

Typically less memory overhead/duplication.
Communication often implicit, through cache coherency and runtime

MPI (Message Passing Interface) is a multi-process model whose mode of communication between the processes is **explicit.** 

==> communication management is the responsibility of the user.

**OpenMP** (**Open Multi-Processing**) is a multitasking model whose mode of communication between tasks is **implicit** 

==> communications is the responsibility of the compiler.



## **Hybrid OpenMP-MPI**

### **Hybrid MPI and OpenMP**



Hybrid application programs using **MPI + OpenMP** are now common place on large HPC systems.

There are basically two main motivations:

- 1. Reduced memory footprint, both in the application and in the MPI library (eg communication buffers).
- 2. Improved performance, especially at high core counts where pure MPI scalability runs out.

### **Hybrid MPI and OpenMP**

#### **Parallel programming models**

Parallel execution is based on threads or processes (or both) which run at the same time on different CPU cores



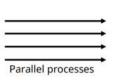
#### **Processes**

- Interaction is based on exchanging messages between processes
- MPI (Message passing interface)

#### **Threads**

- Interaction is based on shared memory, i.e. each thread can access directly other threads data
- OpenMP

### **Hybrid MPI and OpenMP**

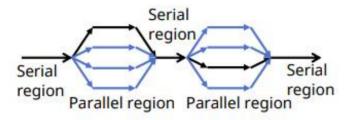


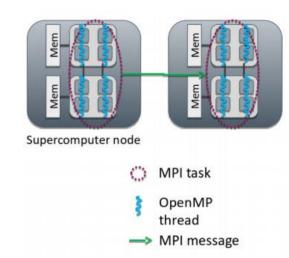
1: MPI: Processes

- Independent execution units
- Have their own memory space
- MPI launches N processes at application startup
- Works over multiple nodes

#### 2: OpenMP: Threads

- Threads share memory space
- Threads are created and destroyed (parallel regions)
- Limited to a single node





- **3: Hybrid programming:** Launch threads (OpenMP) within processes (MPI)
- Shared memory programming inside a node, message passing between nodes
- Optimum MPI task per node ratio depends on the application and should always be experimented.

### **Hybrid Programming**

In hybrid programming each process can have **multiple threads executing simultaneously** All threads within a process share all MPI objects Communicators, requests, etc.

#### MPI defines 4 levels of thread safety:

- ➤ MPI\_THREAD\_SINGLE : One thread exists in program
- ➤ MPI\_THREAD\_FUNNELED : Multithreaded but only the master thread can make MPI calls Master is one that calls MPI\_Init\_thread()
- > MPI\_THREAD\_SERIALIZED: Multithreaded, but only one thread can make MPI calls at a time
- ➤ MPI\_THREAD\_MULTIPLE: Multithreaded and any thread can make MPI calls at any time. Use MPI\_Init\_thread instead of MPI\_Init if more than single thread

### **Hybrid Programming**



#### Potential advantages of the hybrid approach

- Fewer MPI processes for a given amount of cores
  - Improved load balance
  - All-to-all communication bottlenecks alleviated
- Decreased memory consumption if an implementation uses replicated data
- Additional parallelization levels may be available
- Possibility for dedicating threads for different tasks
  - e.g. dedicated communication thread or parallel I/O
- Dynamic parallelization patterns often easier to implement with OpenMP



#### Disadvantages of hybridization

- Increased overhead from thread creation/destruction
- More complicated programming
  - Code readability and maintainability issues
- Thread support in MPI and other libraries needs to be considered

### **Hybrid Programming: Example**

```
#include <mpi.h>
#include <omp.h>
int main(int argc, char *argv[]) {
    int my_id, omp_rank;
    int provided, required=MPI_THREAD_FUNNELED;
    MPI_Init_thread(&argc, &argv, required,
                    &provided);
    MPI_Comm_rank(MPI_COMM_WORLD, &my_id);
#pragma omp parallel private(omp_rank)
    omp_rank = omp_get_thread_num();
    printf("I'm thread %d in process %d\n",
           omp_rank, my_id);
    MPI_Finalize();
```

```
$ mpicc -fopenmp hybrid-hello.c -o hybrid-hello
$ srun --ntasks=2 --cpus-per-task=4
    ./hybrid-hello

I'm thread 0 in process 0
I'm thread 0 in process 1
I'm thread 2 in process 1
I'm thread 3 in process 1
I'm thread 3 in process 1
I'm thread 1 in process 0
I'm thread 3 in process 0
I'm thread 1 in process 0
```

### **Hybrid Programming Example**

#### Numerical integration

$$\int_0^1 \frac{4}{1+x^2} \, dx = \pi$$

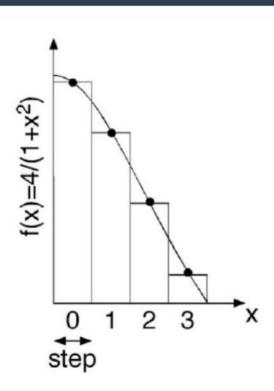
#### Discretization:

$$\Delta = 1/N: \text{ step} = 1/\text{NBIN}$$

$$x_i = (i+0.5)\Delta \ (i=0,...,N-1)$$

$$\sum_{i=0}^{N-1} \frac{4}{2}\Delta \cong \pi$$

```
#include <stdio.h>
#define NBIN 100000
void main() {
   int i; double step,x,sum=0.0,pi;
   step = 1.0/NBIN;
   for (i=0; i<NBIN; i++) {
      x = (i+0.5)*step;
      sum += 4.0/(1.0+x*x);}
   pi = sum*step;
   printf("PI = %f\n",pi);
}</pre>
```



```
#include <stdio.h>
#include <stdlib.h>
                        /* MPI header file */
#include <mpi.h>
#include <omp.h>
                            /* OpenMP header file */
#define NUM STEPS 100000000
#define MAX THREADS 4
int main(int argc, char *argv[]) {
  int nprocs, myid:
  int tid, nthreads, nbin;
  double start time, end time;
  double pi, Psum=0.0, sum[MAX_THREADS]={0.0};
  double step = 1.0/(double) NUM STEPS;
  /* initialize for MPI */
 MPI Init(&argc, &argv);
                           /* starts MPI */
  /* get number of processes */
  MPI Comm size(MPI COMM WORLD, &nprocs);
  /* get this process's number (ranges from 0 to nprocs - 1) */
  MPI Comm rank(MPI COMM WORLD, &myid);
 nbin= NUM STEPS/nprocs;
```

### **Hybrid Programming Example**

```
#pragma omp parallel private(tid)
     int i:
     double x:
     nthreads=omp get num threads();
     tid=omp get thread num();
     for (i=nbin*myid+tid; i < nbin*(myid+1); i+= nthreads) { /* changed*/
       x = (i+0.5)*step;
       sum[tid] += 4.0/(1.0+x*x);
for(tid=0; tid<nthreads; tid++) /*sum by each mpi process*/
     Psum += sum[tid]*step;
MPI Reduce(&Psum, &pi, 1, MPI DOUBLE, MPI SUM, 0, MPI COMM WORLD):/* added */
  if (myid == 0) {
    printf("parallel program results with %d processes:\n", nprocs);
    printf("pi = %g (%17.15f)\n",pi, pi);
  MPI Finalize():
  return 0;
```

#### Results

MPI

```
MPI uses 8 processes:
pi = 3.14159 (3.141592653589828)
```

OpenMP

```
OpenMP uses 8 threads:
pi = 3.14159 (3.141592653589882)
```

Hybrid

```
mpi process 0 uses 4 threads
mpi process 1 uses 4 threads
mpi process 1 sum is 1.287 (1.287002217586605)
mpi process 0 sum is 1.85459 (1.854590436003132)
Total MPI processes are 2
pi = 3.14159 (3.141592653589738)
```



Thank you for your attention!

