# Introduction to Parallel Programming

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## Outline

• Part1:Introduction to parallel programming

- Part2:Parallel Programming Tutorials
  - MPI
  - Pthreads
  - OpenMP

# Outline

- Models for Parallel Systems
- Parallelization of Programs
- Levels of Parallelism
  - Instruction level
  - Data Parallelism
  - Loop Parallelism
  - Functional/task Parallelism
- Parallel Programming Patterns
- Performance Metrics

# **Models for Parallel System**

- Machine Model
  - Describe machines at lowest level of abstraction of hardware, e.g., registers,
- Architecture Model
  - Describe architecture at the level of how processing units, memory organization, interconnect organized and execution model of instructions
- Computation Model
  - Formal model of AM for designing and analyzing algorithm, e.g. RAM, PRAM
- Programming Model
  - Describe the machine from the programmer point of view: how the programmer to code

# Parallel Programming Model

- Parallel Programming model Influenced by
  - Architecture Design
  - Programming Language
  - Compiler
  - Runtime
- Several criteria make them different
  - level of parallelism (instruction level, data level, loops level, procedural level)
  - implicit or use-defined explicit specified parallelism
  - how parallel program parts are specified
  - the execution model of parallel units (SIMD, SPMD, Sync, Async)
  - how to communicate (explicit comm or shared variables)

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- Typical Parallelization Steps
  - Partition/Decomposition: Algorithm is split into tasks with dependencies
  - Scheduling: Tasks are assigned to processes
  - Mapping: Processes are mapped to physical processors



- Partition/Task Decomposition
  - Task: a sequence of computation unit of parallelism, can be at different levels: instruction level, loop level, functional level
  - Task Granularity: Coarse grained and fine grained
  - Compromise between number of tasks and granularity: enough tasks to keep all processors busy and enough granularity to amortize the scheduling/mapping overhead

- Assign tasks to processes/threads
  - A process normally executes multiple different tasks
  - The goal is load balance: each process should have about the same number of computations to perform
  - Static (at the initialization phrase at program start) or Dynamic (during program execution)

- Mapping processes to physical processor/cores
  - Each process or thread is mapped to a separate processor or core
  - Goal: get an equal utilization of the physical processors or cores while keeping communication between the processors as small as possible.

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## **Parallelism at Instruction Level**

- Task unit: one instruction
- Dependency: True, Anti, Out between instructions
- Scheduling: schedule instructions to execute in different function unit

```
For i (1...n)
C[i]=A[i]+B[i]-A[i]*B[i]
```

Loop:

LD R1, @A LD R2, @B **ADD R3, R1, R2 MUL R4, R1, R2** SUB R5, R3, R4 ST R5, @C JNZ Loop



Loop: LD R1, @A LD R2, @B **ADD R3, R1, R2 MUL R4, R1, R2** SUB R5, R3, R4 ST R5, @C JNZ Loop

#### **Parallelism at Instruction Level**

- How to program Instruction Level Parallelism?
  - Write assembly code by hand—find the dependencies and schedule by hand
  - Hardware helps you automatically—Superscalar processor
  - Compiler helps you automatically—scheduling techniques for VLIW processor

#### **Data Parallelism**

- Task unit: one operation on single element
- Dependency: None
- Schedule: Pack different data element into SIMD instruction

for(i=0;i<n;i++) C[i]=A[i]+B[i]-A[i]\*B[i]



for i=0;i<n;i+=4){ C[i...i+3]=A[i...i+3]+B[i...i+3]-A[i...i+3]\*B[i...i+3]

#### **Data Parallelism**

- How to program Data Parallelism?
  - Write assembly code by hand—using SIMD instructions, e.g. MMX, SSE
  - Let compiler help you automatically—Using autovectorization, e.g. gcc "-ftree-vectorize -msse2", but not as high efficient as hand coded
  - Using data-parallel programming language

# **Loop Parallelism**

- Task unit: one iteration of the loop
- Dependency: dependencies between loop iterations
- Schedule: schedule different loop iterations to execute on different processors/cores

```
for(i=0;i<n;i++)

C[i]=A[i]+B[i]-A[i]*B[i] for(i=0;i<n/2;i++)

C[i]=A[i]+B[i]-A[i]*B[i] Core0

without any dependencies

between iterations for(i=n/2+1;i<n;i++)

C[i]=A[i]+B[i]-A[i]*B[i] Core1
```

## Loop Parallelism

- How to program Loop Parallelism?
  - Write multithread code by hand—Decompose the loop into different threads
  - Using high level programming language—e.g., OpenMP

#pragma omp parallel
for(i=0;i<n;i++)
 C[i]=A[i]+B[i]-A[i]\*B[i]</pre>

Compiler do it—Under research, some experimental compilers, e.g. PLUTO

## **Functional Parallelism**

- Task unit: code segments (statements, basic block, loops, functions)
- Dependency: dependencies between tasks
- Schedule: schedule different tasks to execute on different processors/cores

Fib(n)=Fib(n-1)+Fib(n-2)



f1=Fib(n-1)
f2=get\_result\_from\_core1 Core0
return f1+f2;

f=Fib(n-2) return f

# **Functional Parallelism**

- How to program Functional Parallelism?
  - Write multithread code by hand—Decompose the computation into different threads
  - Using high level programming language—e.g., OpenMP, Cilk, Codelet

```
int fib(int n)
{
    if (n < 2)
        return n;
    int x = cilk_spawn fib(n-1);
    int y = fib(n-2);
    cilk_sync;
    return x + y;
}</pre>
```

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# **Parallel Programming Patterns**

- Parallel programs consists of a collection of tasks that are executed by processes/threads
- Patterns: provide specific coordination structures for processes/threads
  - Fork-Join
  - SPMD and SIMD
  - Master–Slave
  - Client–Server
  - Pipelining
  - Task Pools
  - Producer–Consumer

# **Fork-Join**



- Initial time, there is only one main thread do sequential work
- Fork all work threads to do the work in parallel
- Join all the threads and continue to do sequential work

# **Single Program and Multiple Data**



- Each processor executes the same copy of the program
- Each processor has a logical copy of data
- Each processor uses p\_id to find their own part data

#### **Master and Slave**



- Master control the main function of program execution
- Slave does the actual computation which is assigned by Master thread

# Task Pool



- Tata structure in which tasks to be performed are stored and from which they can be retrieved for execution
- A fixed number of threads is used for the processing of the tasks
- a thread can generate new tasks and insert them into

#### **Producer-Consumer**



Producer threads produce data which are used as input by consumer threads

 Common data buffer is used, which can be accessed by both of threads

 Synchronization has to be used to ensure a correct coordination between producer and consumer threads

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### Performance Metrics for Parallel Programs

- Execution Time
  - Sequential execution time:  $T_s$
  - Parallel execution time: *T*<sub>p</sub>
  - Overhead:  $To = pT_p T_s$
- Speedup
  - Speedup= $T_s/T_p$

# Adding N numbers using N processing elements



#### Amdahl's law

- A program execution time  $T_s$  is composed of a fraction of sequential execution time  $T_s * f$  and a fraction of parallel execution time  $T_s * (1-f)$
- Speedup= $\frac{T_s}{T_s * f + T_s * \frac{1-f}{p}} = \frac{1}{f + \frac{1-f}{p}} \pounds \frac{1}{f}$

#### Amdahl's law

• Do you really need parallel computing for your program?

 Speedup is limited by the sequential part of your program

• What is the bottleneck, how many benefits can you get if you try to parallelize it?

#### References





Introduction to Parallel Computing http://wwwusers.cs.umn.edu/~karypis/parbook/

#### Parallel Programming: for Multicore and Cluster Systems

http://www.amazon.com/Parallel-Programming-Multicore-Cluster-Systems/dp/364204817X

# **MPI Programming: A Tutorial**

#### Haitao Wei

#### Thanks to slides from Robert Pavel

#### and Daniel Orozco

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#### MPI

- Stands for Message Passing application programmer Interface.
- It is a specification. There is not one MPI.
- The specification describes primitives that can be used to communicate and program.
- Inspired by the Communicating Sequential Processes paper.



Belfast, Northern Ireland

# Why Learn MPI?

- MPI is the *de facto* standard to program MIMD systems.
- It can be used in SMP systems as well.
- Very versatile, can run on:
  - Symmetric or asymmetric systems
  - Local networks or over the internet
  - On serial processors

# Why Learn MPI?

- Comparatively easy to use.
- All communication is explicit.
- Easy to learn
- Reasonably good performance
- Most importantly: Everyone already uses it
# What is MPI?: An example

- MPI is like exchanging emails with your advisor
- Your advisor gets hundreds of emails per day
  - If he doesn't know an email is coming, he can't respond
- But if he is expecting an email, he'll read it

# **MPI: Execution Model**

- There are P processes that are created at the beginning.
- All processes execute the same program.
- Processes communicate and synchronize using send and receive operations.
- Operations can be **blocking** or **nonblocking**.

# The MPI Programming Model

- Write ONE program that everybody runs.
- Initialize the MPI library:
   MPI\_Init
- Clean the MPI library at the end:

– MPI\_Finalize

## **The Basics**

### MPI\_Init

Initialize the MPI execution environment

int MPI\_Init( int \*argc, char \*\*\*argv )

### **Input Parameters**

#### argc

Pointer to the number of arguments

#### argv

Pointer to the argument vector

### **MPI\_Finalize**

Terminates MPI execution environment

### **Synopsis**

int MPI\_Finalize( void )

## **The Basics**

### MPI\_Comm\_rank

Determines the rank of the calling process in the communicator

**Synopsis** 

int MPI\_Comm\_rank( MPI\_Comm comm, int \*rank )

### **Input Argument**

comm

communicator (handle)

### **Output Argument**

rank

rank of the calling process in the group of comm (integer)



#### MPI\_Comm\_size

Determines the size of the group associated with a communicator

### Synopsis

int MPI\_Comm\_size( MPI\_Comm comm, int \*size )

### **Input Parameter**

comm

communicator (handle)

### **Output Parameter**

size

number of processes in the group of comm (integer)

# **Code Example**

### Hello World!

```
1 #include <mpi.h>
```

```
2
```

```
3 main(int argc, char *argv[])
```

```
4
```

```
5 int npes, myrank;
```

```
6
```

- 7 MPI\_Init(&argc, &argv);
- 8 MPI\_Comm\_size(MPI\_COMM\_WORLD, &npes);
- 9 MPI\_Comm\_rank(MPI\_COMM\_WORLD, &myrank);
- 10 printf("From process %d out of %d, Hello World!\n",
- 11 myrank, npes);
- 12 MPI\_Finalize();

13 }

./mpicc hello.c -o hello ./mpirun -np 3 hello From process 2 out of 3, Hello World! From process 1 out of 3, Hello World! From process 0 out of 3, Hello World!

# Sending and Receiving Data

- Now let's actually do something useful
- MPI, at it simplest, is a series of matched sends and receives
- Host A sends a message to Host B. Host B receives the message
- These sends and receives are blocking by default
- What is blocking?













Should I use red paint?









Working



Yay, I got an ack. Back to work...



## Yes. Use red Ack









## Now with code?

First, let's learn us some syntax!

IPI_Send erforms a blocking send
<pre>ynopsis nt MPI_Send(void *buf, int count, MPI_Datatype datatype, int dest,</pre>
put Parameters
uf
initial address of send buffer (choice)
ount
number of elements in send buffer (nonnegative integer)
atatype
datatype of each send buffer element (handle)
est
rank of destination (integer)
ng
message tag (integer)
omm
communicator (handle)

### **MPI\_Recv** Blocking receive for a message **Synopsis** int MPI Recv(void \*buf, int count, MPI Datatype datatype, int source, int tag, MPI Comm comm, MPI Status \*status) **Output Parameters** buf initial address of receive buffer (choice) status status object (Status) **Input Parameters** count maximum number of elements in receive buffer (integer) datatype datatype of each receive buffer element (handle) source

rank of source (integer)

#### tag

message tag (integer)

#### comm

communicator (handle)

# Now for Examples

### The right way

```
MPI_Comm_rank (comm, &myrank);
if (my rank == 0) {
```

```
MPI_Send (sendbuf, count, MPI_INT, 1,
tag, comm);
```

```
MPI_Recv (recvbuf, count, MPI_INT, 1, tag, comm, &status);
```

```
}
```

```
else if (my rank == 1) {
```

```
MPI_Recv (recvbuf, count, MPI_INT, 0, tag, comm, &status);
```

```
MPI_Send (sendbuf, count, MPI_INT, 0, tag, comm);
```

### .

### the wrong way

```
MPI_Comm_rank (comm, &my rank);
if (my rank == 0) {
    MPI_Recv (recvbuf, count, MPI_INT, 1, tag,
    comm, &status);
    MPI_Send (sendbuf, count, MPI_INT, 1, tag,
    comm);
    }
    else if (my rank == 1) {
        MPI_Recv (recvbuf, count, MPI_INT, 0, tag,
        comm, &status);
        MPI_Send (sendbuf, count, MPI_INT, 0, tag,
        comm);
    }
```

# Huh?

- Why didn't the wrong way work?
  - ...
- Deadlock
- Both processes are waiting for a message to arrive.

## Can we work around that?

Yes!

Non-blocking communication is one way

# MPI\_Isend and MPI\_Irecv

- You can look online for the syntax
  - And it is really long...
- The simple logic is
  - Start a send/recv
  - Do some work
  - Only check when you need the data





Should I use red paint?





Working with the green





Working with the green



Yes, use the red paint. Ack



Working with the green







Oh, cool. I can use red





Working with the red



# A Code Example

### The Wrong Way done Right!

```
MPI_Comm_rank (comm, &my rank);
if (my rank == 0) {
MPI_IRecv (recvbuf, count, MPI_INT, 1, tag, comm, &recv_request);
MPI_ISend (sendbuf, count, MPI_INT, 1, tag, comm, &send_request);
MPI_Wait(&recv_request, &status);
MPI_Wait(&send_request, &status);
}
else if (my rank == 1) {
MPI_IRecv (recvbuf, count, MPI_INT, 0, tag, comm, &recv_request);
MPI_ISend (sendbuf, count, MPI_INT, 0, tag, comm, &send_request);
MPI_Wait(&recv_request, &status);
MPI_Wait(&recv_request, &status);
```

### }

## **The Last Must-Have Tool**

- What if we need to ensure one phase is complete before starting the next
  - Finish washing your hands before you leave the restroom
- How to guarantee that in MPI?
  - A series of blocking sends and receives?
  - Something else?

# Or just use Barrier

- A Barrier halts execution of code until all processes have signaled that they have reached a barrier
- Many ways to implement a barrier
  - We may discuss these during the course
- Only use a Barrier if you need to
  - Hurts performance due to idle processes

# **Barrier!**

#### **MPI\_Barrier**

Blocks until all processes in the communicator have reached this routine. **Synopsis** 

int MPI\_Barrier( MPI\_Comm comm )

**Input Parameter** 

comm

communicator (handle)

## **Code Example**

### Let's see how to use a Barrier

```
MPI Comm rank (comm, &myrank);
if (my rank == 0) {
//p0 does something in stage 1
 MPI Barrier(comm);
 //p0 does something in stage 2
 MPI Barrier(comm);
else if (my rank == 1) {
 //p1 does something in stage 1
 MPI Barrier(comm);
//p1 does something in stage 2
 MPI_Barrier(comm);
}
```

# **Advanced MPI**

- Other types of Send/Recv
  - Buffered: Copies data to another buffer
  - MPI\_Ssend: Won't return until recv is completed
  - MPI\_Rsend: May only be used if matching recv is already active
  - MPI\_SendRecv: Combines Send and Receive into a single command

# **Collective Operations**

- An alternative to point to point operations.
- Involve communication and synchronization between many processes.
- The two most common are:
  - MPI\_Bcast(..., root, ...):
    - All processes call the same function.
    - All processes receive data from process root.
  - MPI\_Reduce( ..., root, ... )
    - A reduction operation is done with data from each process and the result is given to the root process.

# **Collective Example: Intuitive**

- TA normally wants to inform you we have a huge project.
  - So TA "Broadcast" the information in a mass email.
- TA will also collect the "quiz" after the midterm exam
  - a reduction is performed where all of you put your quizzes in the basket of the root process(TA).

# **Dot Product**

}

```
#include "mpi.h"
main(int argc, char* argv[]) {
//Step 1:initialize vector a and b
float loc dot=0.0f;
float dot=0.0f;
float a[N],b[N],loc dots[N];
for (i = 0;i<N;i++) {
   a[i] = i;
   b[i] = i+1;
//Step2 initialize MPI
MPI Init(&argc, &argv);
MPI Comm rank(MPI COMM WORLD,&my rank);
MPI_Comm_size(MPI_COMM_WORLD,&p);
//Step3: Each processor computes a local dot product
   loc n = N/p;
   bn = (my rank)*loc n;
   en = bn + loc n;
   loc dot = 0.0;
   for (i = bn;i <en; i++) {
    loc dot = loc dot + a[i]*b[i];
```

#### // Step4: collect result

MPI Reduce(&loc\_dot, dot, 1, MPI\_FLOAT, MPI SUM, 0, MPI COMM WORLD);

```
if(my rank==0)
 printf( "dot product = %f",dot);
```

```
/* mpi is terminated. */
   MPI Finalize();
```

# **Thoughts on Advanced MPI**

- Be very careful
- Using collectives may kills performance
  - If only because they are blocking
- There may be special cases where the specialized send and recv are useful
- But unless you are in HPC, use whatever is most intuitive
## Can I use MPI with...

- Fortran?
  - Yes
- C++?
  - Yes, and it is actually a very intuitive interface that I really like
- Java?
  - Not easily
- Python?
  - Yes
- Matlab?
  - Not easily

## References

- The CSP paper
  - <u>http://portal.acm.org/citation.cfm?id=359576.35</u>
    <u>9585</u>
- A Reference for MPI
  - <u>http://www.mcs.anl.gov/research/projects/mpi/</u> <u>www/www3/</u>