

# Introduction to MPI Part 1

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# Overview

- Definition and rationale
- Important considerations
- Getting started
- Essential functions
- Exercises

# What is Message Passing Interface (MPI)?

- A communication library for parallel computing over a *distributed* system.
- Distributed=loosely coupled computers—they do not share a clock or memory.
- Programmer writes one program in C or Fortran.
- The one program is loaded on one or more processors and called a *process*.

# Major Parallel Computing Models

- Shared memory
  - Much like pthreads (multithreading library)
  - Processes (lightweight processes or threads) all share memory (i.e. globals)
  - Threads have some private memory (stack)
  - Locking is needed and handled through monitors and semaphores

# Major Parallel Computing Models

- Message Passing
  - Each process has its own memory
  - Programmer handles any sharing of data by passing messages
  - MPI falls into this category

# Parallel Programming Types

- Data Parallel: Single Instruction Multiple Data (SIMD) – instructions are same but work on different data
- Task Parallel: Multiple Instructions Multiple Data (MIMD) – instructions differ as does the data
- Single Program Multiple Data (SPMD): *program* (think source) is the same, but different parts of it execute depending on process rank (equiv. to MIMD)
- *Programming with MPI is typically SPMD.*

# SPMD Diagram

## Process 0

Code:

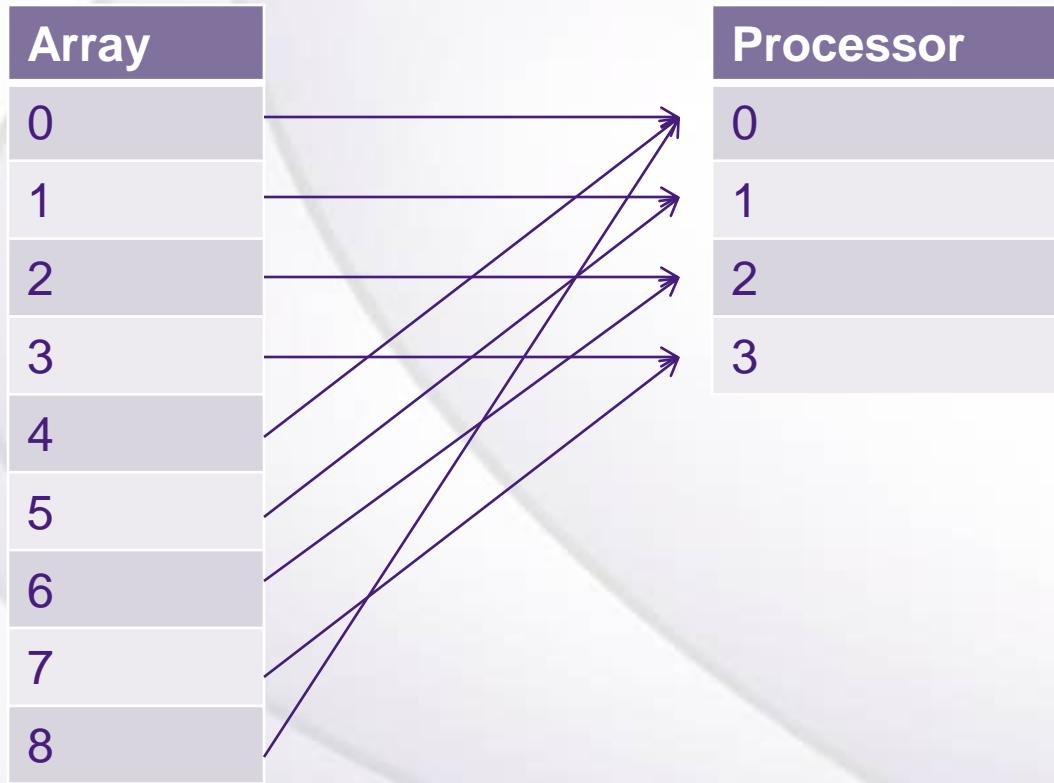
```
MPI_Init...
MPI_Comm_rank(MPI_C
OMM_WORLD, &rank);
if(rank==0) {
...
} else {
...
}
```

## Process 1

Code:

```
MPI_Init...
MPI_Comm_rank(MPI_C
OMM_WORLD, &rank);
if(rank==0) {
...
} else {
...
}
```

# Multiple Data



# Why Use MPI?

- It allows one to parallelize his/her own programs.
- It is a solid communication library for distributed computing.

# Process

- Each *process* is really just the one program, but usually executes differently on each *processor (core)*.
- Each process has a *rank* from 0 .. #CPUs, which is accessible from the running process.
- The programmer generally writes conditional blocks depending on this *rank*.

# Rank

- Depending on the rank, the process may:
  - work on a different part of the problem space (*slave*),
  - split up a problem space and distribute the work (*master*),
  - combine results and save/display them (*master*), or
  - anything else the programmer wishes to do.

# Writing a Program

- MPI has many functions in its library.
  - Point-to-point
  - Collective
- With only syntactic differences, the functions are the same in C and Fortran.

# Compiling a Program

- Compiling code only involves:
  - Including a C or Fortran header file
  - Compiling using included special compiler wrapper
- Running code is as easy as using mpirun or placing in a batch script for use with qsub.

# Basic MPI Program Structure

Call to Initialize MPI library.

If rank  $\neq 0$ : //Here, the master is 0.

- Read input data (from master or a file).

- Do necessary operations.

- Send result to process 0.

Else:

- Send all/part of input data to slave processes if necessary.

- For each process i:

  - Receive result from process i.

  - Combine result with final result.

  - Output final result.

Call to Finalize MPI library.

# Basic MPI Program in C

```
#include <stdio.h>
#include <mpi.h>
main(int argc, char **argv) {
    int np, rank;
    MPI_Init(&argc, &argv);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    MPI_Comm_size(MPI_COMM_WORLD, &np);
    printf("Hello World on process %d of %d.\n", rank,
    np);
    MPI_Finalize();
}
```

# Compiling & Running Program

```
$ mpicc -o 00basic 00basic.c  
$ qsub -I -l nodes=1:ppn=4 -l walltime=00:01:00 -q workq  
...  
$ mpirun -np 4 00basic
```

script:

```
#!/bin/bash  
#PBS -q workq  
#PBS -A loni_ccavatrain  
#PBS -l nodes=1:ppn=4  
#PBS -l walltime=00:00:01  
#PBS -o 00basic.out  
#PBS -j oe  
#PBS -N basic  
export WORK_DIR=/work/ccava/mpi  
cd $WORK_DIR  
export NPROCS=`wc -l $PBS_NODEFILE |gawk '{print $1}'`  
mpirun -machinefile $PBS_NODEFILE -np $NPROCS  
/home/ccava/mpi/00basic
```

```
$ qsub script
```

# Using Rank to Affect Flow

```
MPI_Comm_rank(MPI_COMM_WORLD, &rank);
MPI_Comm_size(MPI_COMM_WORLD, &np);
if( rank == 0 ) { /* process 0 */ }
else if( rank == 1 ) { /* process 1 */ }
else if( rank == 2 ) { /* process 2 */ }
else { /* other process */ }
```

# Sending Messages: MPI\_Send

```
int MPI_Send( void *buf, int count, MPI_Datatype  
datatype, int dest, int tag, MPI_Comm comm )
```

buf - initial address of send buffer (choice)

count - number of elements in send buffer (nonnegative integer)

datatype - datatype of each send buffer element (handle)

dest - rank of destination (integer)

tag - message tag (integer)

comm - communicator (handle)

# Receiving Messages: MPI\_Recv

```
int MPI_Recv( void *buf, int count, MPI_Datatype  
datatype, int source, int tag, MPI_Comm comm,  
MPI_Status *status )
```

buf - initial address of receive buffer (choice)

status - status object (Status)

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)

# MPI\_Send & MPI\_Recv Example

```
#include "mpi.h"
#include <stdio.h>
int main(argc,argv) int argc; char *argv[]; {
    int numtasks, rank, dest, source, rc, count, tag=1;
    char inmsg, outmsg='x';
    MPI_Status Stat;
    MPI_Init(&argc,&argv);
    MPI_Comm_size(MPI_COMM_WORLD, &numtasks);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    if (rank == 0) {
        dest = 1;
        source = 1;
        rc = MPI_Send(&outmsg, 1, MPI_CHAR, dest, tag, MPI_COMM_WORLD);
        rc = MPI_Recv(&inmsg, 1, MPI_CHAR, source, tag, MPI_COMM_WORLD, &Stat);
    } else if (rank == 1) {
        dest = 0;
        source = 0;
        rc = MPI_Recv(&inmsg, 1, MPI_CHAR, source, tag, MPI_COMM_WORLD, &Stat);
        rc = MPI_Send(&outmsg, 1, MPI_CHAR, dest, tag, MPI_COMM_WORLD);
    }
    rc = MPI_Get_count(&Stat, MPI_CHAR, &count);
    printf("Task %d: Received %d char(s) from task %d with tag %d \n", rank, count,
    Stat.MPI_SOURCE, Stat.MPI_TAG);
    MPI_Finalize();
}
```

# Data Types

MPI\_CHAR

MPI\_SHORT

MPI\_INT

MPI\_LONG

MPI\_UNSIGNED\_CHAR

MPI\_UNSIGNED\_SHORT

MPI\_UNSIGNED\_LONG

MPI\_UNSIGNED

MPI\_FLOAT

MPI\_DOUBLE

MPI\_LONG\_DOUBLE

MPI\_BYTE

MPI\_PACKED

# Distributing Data: MPI\_Bcast

```
int MPI_Bcast ( void *buffer, int count, MPI_Datatype  
datatype, int root, MPI_Comm comm )
```

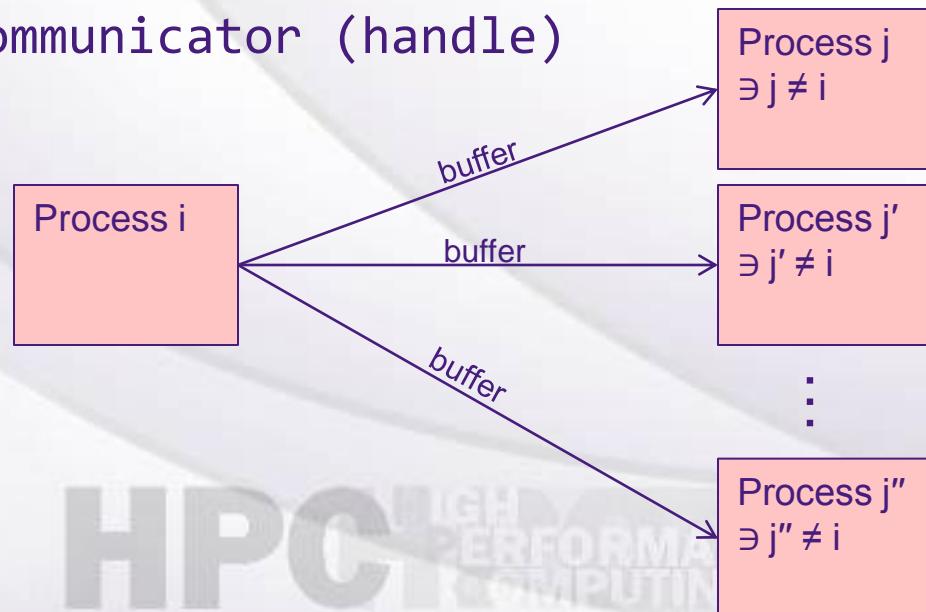
buffer - starting address of buffer (choice)

count - number of entries in buffer (integer)

datatype - data type of buffer (handle)

root - rank of broadcast root (integer)

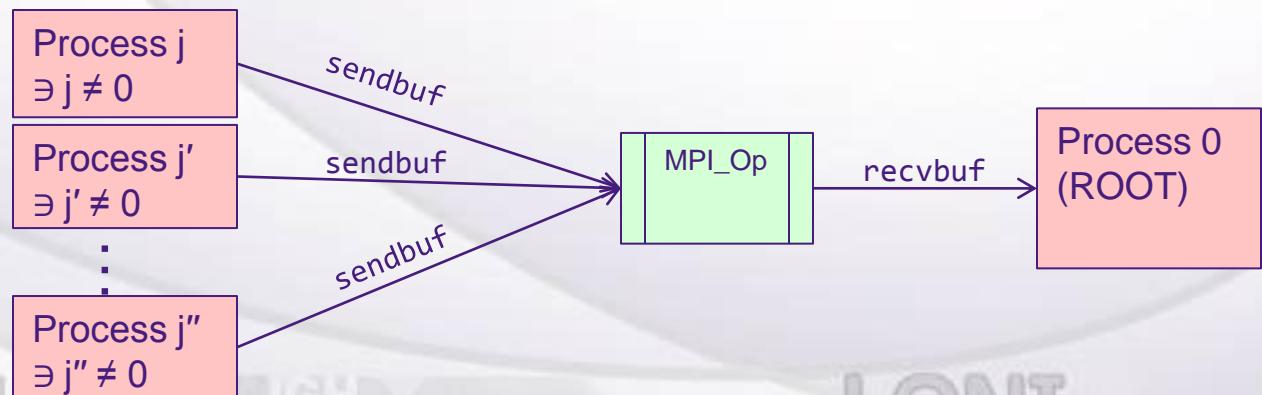
comm - communicator (handle)



# Collecting & Calculating with Data: MPI\_Reduce

```
int MPI_Reduce ( void *sendbuf, void *recvbuf, int count,
MPI_Datatype datatype, MPI_Op op, int root, MPI_Comm comm )
```

sendbuf - address of send buffer (choice)  
count - number of elements in send buffer (integer)  
datatype - data type of elements of send buffer (handle)  
op - reduce operation (handle)  
root - rank of root process (integer)  
comm - communicator (handle)  
recvbuf - address of receive buffer (choice, significant  
only at root )



# MPI\_Op Values

MPI\_MAX maximum

MPI\_MIN minimum

MPI\_SUM sum

MPI\_PROD product

MPI\_BAND logical and

MPI\_BAND bit-wise and

MPI\_LOR logical or

MPI\_BOR bit-wise or

MPI\_LXOR logical xor

MPI\_BXOR bit-wise xor

MPI\_MAXLOC max value and location

MPI\_MINLOC min value and location

Programmer may define own operation (using MPI\_Op\_create)

# MPI\_Bcast & MPI\_Reduce Example

```
#include "mpi.h"
#include <math.h>
int main(argc,argv) int argc; char *argv[]; {
    int done = 0, n, myid, numprocs, i, rc;
    double PI25DT = 3.141592653589793238462643;
    double mypi, pi, h, sum, x, a;
    MPI_Init(&argc,&argv);
    MPI_Comm_size(MPI_COMM_WORLD,&numprocs);
    MPI_Comm_rank(MPI_COMM_WORLD,&myid);
    while (!done){
        if (myid == 0) {
            printf("Enter the number of intervals: (0 quits) ");
            scanf("%d",&n); }
        MPI_Bcast(&n, 1, MPI_INT, 0, MPI_COMM_WORLD);
        if (n == 0) break;
        h = 1.0 / (double) n;
        sum = 0.0;
        for (i = myid + 1; i <= n; i += numprocs) {
            x = h * ((double)i - 0.5);
            sum += 4.0 / (1.0 + x*x); }
        mypi = h * sum;
        MPI_Reduce(&mypi, &pi, 1, MPI_DOUBLE, MPI_SUM, 0, MPI_COMM_WORLD);
        if (myid == 0)
            printf("pi is approximately %.16f, Error is %.16f\n", pi, fabs(pi - PI25DT)); }
    MPI_Finalize(); }
```

# MPI\_Barrier

Synchronizes all processes by blocking until all reach this function.

```
int MPI_Barrier ( MPI_Comm comm )  
comm - communicator (handle)
```

# Using MPI\_Barrier to Compute Time

```
double start,end;  
...  
MPI_Barrier(MPI_COMM_World);  
if(rank==0) start = MPI_Wtime(); //seconds  
...  
MPI_Barrier(MPI_COMM_World);  
if(rank==0) end = MPI_Wtime();  
...
```

# Exercises