THE ARCTIC UNIVERSITY OF NORWAY

Lecture 3b: Message-Passing Computing - MPI

Parallell Programming (INF-3201)

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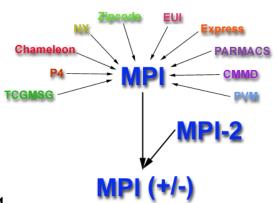
MPI (Message Passing Interface)

What?

 a specification for the developers and users of message passing libraries (not a library).

Why use MPI?

- Standardization MPI is the only message passing library interface which can be considered a standard.
- Portability There is no need to modify your source code when you port your application to a different platform.
- Performance Vendor implementations can exploit native hardware features to optimize performance.
- Functionality Over 115 routines are defined in MPI-1 alone.
- Availability A variety of implementations are available.



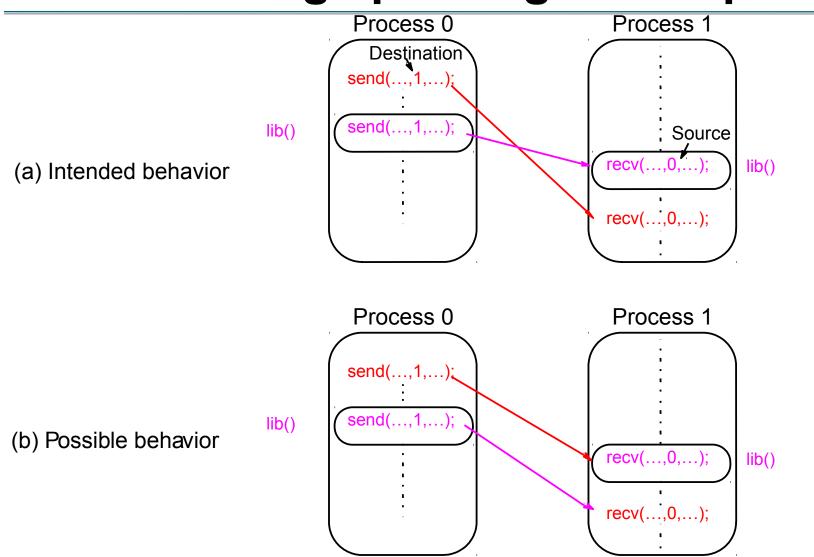
MPI Process Creation and Execution

- Purposely not defined Will depend upon implementation.
- MPI version 1
 - Only static process creation supported.
 - All processes must be defined prior to execution and started together.
- MPI version 2
 - Dynamic process creation supported (MPI_Comm_spawn)
- Originally SPMD model of computation.
 - MPMD also possible with static creation

Using SPMD Computational Model

where master() and slave() are to be executed by master process and slave process, respectively.

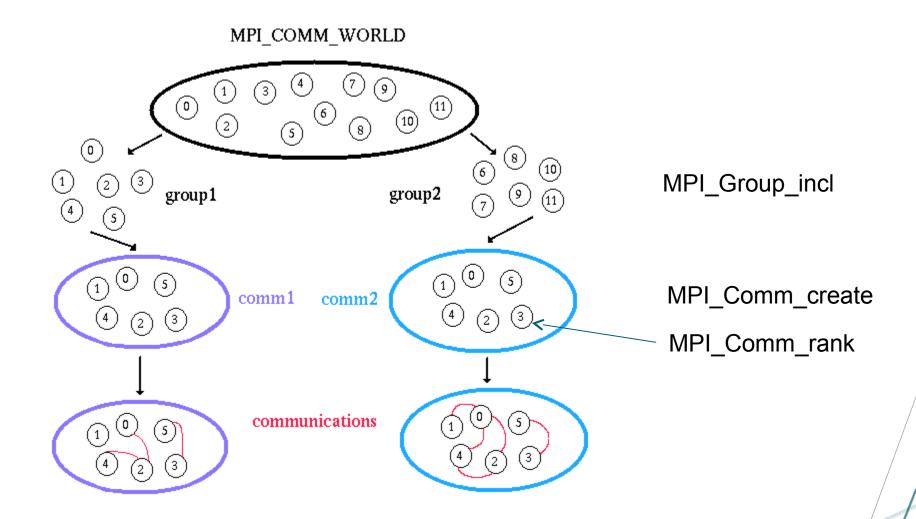
Unsafe message passing - Example



MPI Solution: "Communicators"

- Defines a communication domain a set of processes that are allowed to communicate between themselves.
- Processes have ranks associated with communicator (0 to n-1)
- Initially, all processes enrolled in a "universe" called MPI_COMM_WORLD
- Other communicators can be established for groups of processes.
- Communication domains of libraries can be separated from that of a user program.
- Used in all point-to-point and collective MPI message-passing communications.

Communicator Example



MPI Point-to-Point Communication

- Involves one source process and one destination process
- Uses send and receive routines with message tags (and communicator).
- Wild card message tags available

MPI Blocking Routines

- Return when locally complete
 - location used to hold message can be used again or altered without affecting message being sent.
- When blocking send returns
 - process free to move on without adversely affecting message.
 - does not mean that message has been received,

Parameters of blocking send

MPI_Send(buf, count, datatype, dest, tag, comm)

Address of Datatype of each item

Number of items Rank of destination Communicator to send process

Parameters of blocking receive

MPI_Recv(buf, count, datatype, src, tag, comm, status)

Address of receive buffer

Maximum number of items to receive process

Address of Patatype of

Example

To send an integer x from process 0 to process 1,

```
MPI_Comm_rank(MPI_COMM_WORLD,&myrank); /* find rank */
if (myrank == 0) {
   int x;
   MPI_Send(&x, 1, MPI_INT, 1, msgtag, MPI_COMM_WORLD);
} else if (myrank == 1) {
   int x;
   MPI_Recv(&x, 1, MPI_INT, 0, msgtag, MPI_COMM_WORLD, status);
}
```

MPI Nonblocking Routines

 Nonblocking send - MPI_Isend() - will return immediately even before source location is safe to be altered.

 Nonblocking receive - MPI_Irecv() - will return even if no message to accept.

Nonblocking Routine Formats

```
MPI Isend (buf, count, datatype, dest, tag, comm, request);
MPI Irecv (buf, count, datatype, source, tag, comm, request);
Completion detected by MPI Wait() and MPI Test().
MPI Wait (request, ...)
  waits until operation completed and returns then.
MPI Test (request, ...)
  returns with flag (set) indicating whether operation completed at that
   time.
```

Example

To send an integer x from process 0 to process 1 and allow process 0 to continue,

```
MPI_Comm_rank(MPI_COMM_WORLD, &myrank); /* find rank */
if (myrank == 0) {
    int x;
    MPI_Isend(&x,1,MPI_INT, 1, msgtag, MPI_COMM_WORLD, req1);
    compute();
    MPI_Wait(req1, status);
} else if (myrank == 1) {
    int x;
    MPI_Recv(&x,1,MPI_INT,0,msgtag, MPI_COMM_WORLD, status);
}
```

Send Communication Modes

- Buffered Mode Send may start and return before a matching receive. Necessary to specify buffer space via routine MPI_Buffer_attach().
- Synchronous Mode Send and receive can start before each other but can only complete together.
- Standard Mode Not assumed that corresponding receive routine has started. If buffering provided, send could complete before receive reached.
- Ready Mode Send can only start if matching receive already reached, otherwise error. Use with care.

- Each of the four modes can be applied to both blocking and nonblocking send routines.
- Any type of send routine can be used with any type of receive routine.

Collective Communication

All or None

- Must involves all processes in the scope of an intra-communicator.
- It is the programmer's responsibility to insure that all processes within a communicator participate in any collective routines.

Programming Considerations and Restrictions

- Routines are blocking.
- Routines do not take message tag arguments.
- Can only be used with MPI predefined datatypes. Solution?
 - Convert user-defined datatypes to an array of integer (cf. cast operator in C)

Principal collective routines:

- MPI_Bcast() Broadcast from root to all other processes
- MPI Gather () Gather values for group of processes
- MPI_Scatter() Scatters buffer in parts to group of processes
- MPI_Alltoall() Sends data from all processes to all processes
- MPI_Reduce ()
 Combine values on all processes to single value
- MPI_Reduce_scatter() Combine values and scatter results
- MPI_Scan () Compute prefix reductions of data on processes

MPI_Bcast

MPI_Bcast Broadcasts a message to all other processes of that group count = 1;source = 1; broadcast originates in task 1 MPI_Bcast(&msg, count, MPI_INT, source, MPI_COMM_WORLD); task 0 task 1 task 2 task 3 7 msg (before) msg (after)

MPI_Gather

MPI_Gather

Gathers together values from a group of processes

```
sendcnt = 1;
recvent = 1;
src = 1;
                   messages will be gathered in task 1
MPI_Gather(sendbuf, sendcnt, MPI_INT,
             recvbuf, recvent, MPI INT,
             src, MPI COMM WŌRLD);
task 0
             task 1
                         task 2
                                      task 3
               2
                            3
                                                     sendbuf (before)
  1
                                        4
               1
               2
                                                      recybuf (after)
               3
               4
```

MPI_Scatter

MPI_Scatter

Sends data from one task to all other tasks in a group

```
sendcnt = 1;
recvent = 1;
src = 1;
                      task 1 contains the message to be scattered
MPI_Scatter(sendbuf, sendcnt, MPI_INT, recvbuf, recvcnt, MPI_INT,
               src, MPÍ COMM WŌRLD);
task 0
               task 1
                             task 2
                                            task 3
                 1
                 2
                                                              sendbuf (before)
                 3
                 4
                                3
                                                             recybuf (after)
                 2
```

MPI_Alltoall

MPI_Alltoall

Sends data from all to all processes. Each process performs a scatter operation.

```
sendcnt = 1;
recvent = 1;
MPI_Alltoall(sendbuf, sendcnt, MPI_INT, recvbuf, recvcnt, MPI_INT,
             MPI_CÓMM_WORLD);
task 0
             task 1
                           task 2
                                        task 3
                                          13
                5
                             9
  1
  2
                6
                             10
                                          14
                                                         sendbuf (before)
                            11
                                          15
  3
                7
   4
                8
                            12
                                          16
  1
                2
                             3
                                           4
  5
                6
                             7
                                           8
                                                       - recybuf (after)
  9
               10
                            11
                                          12
  13
                            15
                                          16
               14
```

MPI_Reduce

MPI_Reduce

Perform and associate reduction operation across all tasks in the group and place the result in one task

```
count = 1;
                         result will be placed in task 1
dest = 1;
MPI_Reduce(sendbuf, recvbuf, count, MPI_INT, MPI_SUM,
             dest, MPI_COMM_WÓRLD);
                          task 2
                                      task 3
task 0
             task 1
                                                      sendbuf (before)
                            3
               2
  1
                                         4
                                                      recybuf (after)
               10
```

MPI_Reduce_scatter

MPI_Reduce_scatter

Perform reduction operation on vector elements across all tasks in the group, then distribute segments of result vector to tasks

recvcount = 1;

MPI_Reduce_scatter(sendbuf, recvbuf, recvcount, MPI_INT, MPI_SUM, MPI_COMM_WORLD);

task 0	task 1	task 2	task 3	
0	0	0	0	sendbuf (before)
1	1	1	1	
2	2	2	2	
3	3	3	3	
0	4	8	12	recvbuf (after)

MPI_Scan

MPI_Scan

Computes the scan (partial reductions) of data on a collection of processes

	task 3	task 2	task 1	task 0
sendbuf (before)	4	3	2	1
- recybuf (after)	10	6	3	1

Example

To gather items from group of processes into process 0, using dynamically allocated memory in root process:

gathers from all processes, including root.

Barrier routine

- A means of synchronizing processes by stopping each one until they all have reached a specific "barrier" call.
 - MPI_Barrier (comm)

Sample MPI program

```
#include "mpi.h"
#include <stdio.h>
#include <math.h>
#define MAXSIZE 1000
void main(int argc, char *argv)
    int myid, numprocs;
    int data[MAXSIZE], i, x, low, high, myresult, result;
    char fn[255];
    char *fp;
   MPI_Init(&argc,&argv);
   MPI Comm size (MPI COMM WORLD, &numprocs);
   MPI Comm rank (MPI COMM WORLD, &myid);
    if (myid == 0) { /* Open input file and initialize data */
       strcpy(fn,getenv("HOME"));
       strcat(fn,"/MPI/rand data.txt");
       if ((fp = fopen(fn,"r")) == NULL) {
           printf("Can't open the input file: %s\n\n", fn);
           exit(1);
       for(i = 0; i < MAXSIZE; i++) fscanf(fp,"%d", &data[i]);</pre>
    MPI Bcast(data, MAXSIZE, MPI INT, 0, MPI COMM WORLD); /* broadcast data */
    x = n/nproc; /* Add my portion Of data */
   low = myid * x;
   high = low + x;
    for (i = low; i < high; i++)
        myresult += data[i];
    printf("I got %d from %d\n", myresult, myid); /* Compute global sum */
   MPI Reduce (&myresult, &result, 1, MPI INT, MPI SUM, 0, MPI COMM WORLD);
    if (myid == 0) printf("The sum is %d.\n", result);
    MPI Finalize();
```

Estimating speedup

Sequential execution time, $t_{s:}$ Estimate by counting computational steps of best sequential algorithm.

Parallel execution time, $t_{p:}$ In addition to number of computational steps, t_{comp} , need to estimate communication overhead, t_{comm} .

$$t_p = t_{\text{comp}} + t_{\text{comm}}$$

Computation Time

- Count number of computational steps.
- When more than one process executed simultaneously, count computational steps of most complex process.
- •Generally, function of number of data elements *n* and number of processors *p*, i.e.

$$t_{\text{comp}} = f(n, p)$$

Often break down computation time into parts. Then

$$t_{\text{comp}} = t_{\text{comp1}} + t_{\text{comp2}} + t_{\text{comp3}} + \dots$$

Analysis usually done assuming that all processors are same and operating at same speed.

Communication Time

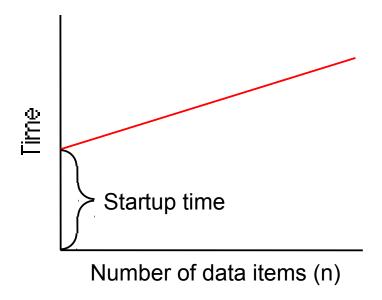
Many factors, including network structure and network contention. As a first approximation, use

$$t_{\text{comm1}} = t_{\text{startup}} + nt_{\text{data}}$$

 t_{startup} is startup time, essentially time to send a message with no data. Assumed to be constant.

 t_{data} is transmission time to send one data word, also assumed constant, and there are n data words.

Idealized Communication Time



Final communication time, t_{comm}

Sum of communication times of all sequential messages from a process, i.e.

$$t_{\text{comm}} = t_{\text{comm1}} + t_{\text{comm2}} + t_{\text{comm3}} + \dots$$

- Assumption: Communication patterns of all processes are the same and take place together
- \Rightarrow only one process need be considered.

Both t_{startup} and t_{data} , measured in units of one computational step, so that can add t_{comp} and t_{comm} together to obtain parallel execution time, t_{p} .

Benchmark Factors

With t_{s_i} t_{comp_i} and t_{comm} , can establish speedup factor and computation/communication ratio for a particular algorithm/implementation.

Speedup factor =
$$\frac{t_S}{t_p} = \frac{t_S}{t_{comp} + t_{comm}}$$

Computation/communication ratio =
$$\frac{t_{comp}}{t_{comm}}$$

Both functions of number of processors, p, and number of data elements, n.

Factors give indication of scalability of parallel solution with increasing number of processors and problem size.

Computation/communication ratio will highlight effect of communication with increasing problem size and system size.

References

- Barry Wilkinson & Michael Allen. Parallel Programming: Techniques and Applications Using Networked Workstations and Parallel Computers.
- Blaise Barney, "Message Passing Interface (MPI)", Livermore Computing.