Day1

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1. Agenda

- Recap
- Communicators
- MPI Communications

- MPI Programs
- Point to point communications

2. MPI Communicators

Communicators in MPI define a group of processes that can communicate with each other. The default communicator is $MPI_{COMMWORLD}$, which includes all the processes. Custom communicators can be created to define subgroups of processes for specific communication patterns.

3. Types of MPI Communications

MPI offers various communication mechanisms to facilitate different types of data exchanges between processes:

3.1. Point-to-Point Communication:

- **Blocking**: The sending and receiving operations wait until the message is delivered (e.g., `MPI_{Send}`, `MPI_{Recv}`).
- **Non-Blocking**: The operations return immediately, allowing computation and communication to overlap (e.g., `MPI_{Isend}`, `MPI_{Irecv}`).

3.2. Collective Communication:

These operations involve a group of processes and include:

- Broadcast: Send data from one process to all other processes (`MPI_{Bcast}`).
- Scatter: Distribute distinct chunks of data from one process to all processes (`MPI_{Scatter}`).
- Gather: Collect chunks of data from all processes to one process (`MPIGather`).
- All-to-All: Every process sends and receives distinct chunks of data (`MPI_{Alltoall}`).

Collectives can also include operations like reductions (`MPI_{Reduce}`, `MPI_{Allreduce}`) which perform computations on data from all processes and distribute the result.

4. Sample MPI code

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

int main(int argc, char **argv) {
    MPI_Init(&argc, &argv);
    int size;
    MPI_Comm_size(MPI_COMM_WORLD, &size);
    int rank;
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    printf("Hello from process %d of %d\n", rank, size);

    MPI_Finalize();
    return 0;
}
```

This example demonstrates a simple MPI program where each process prints its rank and the total number of processes.

5. Scripts

5.1. compile script

```
#!/bin/sh

#source /opt/ohpc/pub/apps/spack/share/spack/setup-env.sh
#spack load gcc/5i5y5cb
#spack load openmpi/c7kvqyq
source ~/git/spack/share/spack/setup-env.sh
spack load openmpi

inputFile=$1
outputFile="${1%.*}.out" # extract the name of the file without extension and adding extension .out
#cmd=`mpicc $inputFile -o $outputFile`
cmd="mpicc $inputFile -o $outputFile" # running code using MPI
```

```
echo "-----"
echo "Command executed: $cmd"
echo "-----"
$cmd

echo "Compilation successful. Check at $outputFile"
echo "-----"
```

5.2. run script

```
#!/bin/sh
#source /opt/ohpc/pub/apps/spack/share/spack/setup-env.sh
#spack load gcc/5i5y5cbc
source ~/git/spack/share/spack/setup-env.sh
spack load openmpi
cmd="mpirun -np $2 $1"
echo "------"
echo "Command executed: $cmd"
echo "-----
OUTPUT
echo
mpirun -np $2 $1
echo
echo "#########
```

6. Hello World in C

```
#include<stdio.h>
int main(){
    printf("Hello World\n");
    return 0;
}
```

Hello World

7. Hello World just by using MPI

```
#include<stdio.h>
int main(){
  printf("Hello World\n");
  return 0;
Hello World
bash compile.sh hello mpi.c
Command executed: mpicc hello mpi.c -o hello mpi.out
Compilation successful. Check at hello mpi.out
_ ·
bash run.sh ./hello mpi.out 6
Command executed: mpirun -np 6 ./hello mpi.out
OUTPUT
Hello World
Hello World
Hello World
Hello World
```

8. Hello World with MPI routines

```
bash compile.sh hello_mpi_processes.c
```

```
Command executed: mpicc hello_mpi_processes.c -o hello_mpi_processes.out

Compilation successful. Check at hello_mpi_processes.out
```

```
bash run.sh ./hello_mpi_processes.out 8
```

9. Hello from only even rank of processes

```
#include<stdio.h>
#includ
```

```
bash compile.sh hello_mpi_even_processes.c
```

```
Command executed: mpicc hello_mpi_even_processes.c -o hello_mpi_even_processes.out

Compilation successful. Check at hello_mpi_even_processes.out
```

```
bash run.sh ./hello_mpi_even_processes.out 10
```

10. MPI Initialization: MPI_{Init} vs. $MPI_{Initthread}$

MPI provides two main functions to initialize the MPI environment: $`MPI_{Init}`$ and $`MPI_{Initthread}`$. The primary difference is that $`MPI_{Initthread}`$ allows you to specify the desired level of thread support.

10.1. Levels of Thread Support

- `MPI_{THREADSINGLE}`: Only one thread will execute.
- \bullet `MPI_{THREADFUNNELED}`: The process may be multi-threaded, but only the main thread will make MPI calls.
- `MPI_{THREADSERIALIZED}`: Multiple threads may make MPI calls, but only one at a time.
- `MPI_{THREADMULTIPLE}`: Multiple threads may make MPI calls with no restrictions.

10.2. MPI_{Init} Example

This example uses `MPI $_{\mbox{\footnotesize{Init}}}$ to initialize the MPI environment.

```
#include <mpi.h>
#include <stdio.h>
int main(int argc, char** argv) {
```

```
// Initialize the MPI environment
MPI_Init(&argc, &argv);

// Get the number of processes
int world_size;
MPI_Comm_size(MPI_COMM_WORLD, &world_size);

// Get the rank of the process
int world_rank;
MPI_Comm_rank(MPI_COMM_WORLD, &world_rank);

// Print off a hello world message
printf("Hello world from processor %d out of %d processors\n", world_rank, world_size);

// Finalize the MPI environment.
MPI_Finalize();
return 0;
}
```

10.3. Compilation and Execution (MPI $_{Init}$)

• Compile the program:

```
Compilation successful. Check at mpi_init.out
```

• Run the program:

```
bash run.sh ./mpi_init.out 6

Command executed: mpirun -np 6 ./mpi_init.out
```

10.4. MPI_{Initthread} Example

This example uses `MPI_{Initthread}` to initialize the MPI environment with thread support.

```
#include <mpi.h>
#include <stdio.h>
int main(int argc, char** argv) {
    int provided;
    // Initialize the MPI environment with thread support
    MPI Init thread(&argc, &argv, MPI THREAD MULTIPLE, &provided);
    // Check the level of thread support provided
    if (provided < MPI THREAD MULTIPLE) {</pre>
        printf("MPI does not provide required thread support\n");
        MPI Abort(MPI COMM WORLD, 1);
    // Get the number of processes
    int world size;
    MPI Comm size(MPI COMM WORLD, &world size);
    // Get the rank of the process
    int world rank:
   MPI Comm rank(MPI COMM WORLD, &world rank);
    // Print off a hello world message
```

```
printf("Hello world from processor %d out of %d processors with thread support level %d\n", world_rank, world_size, provided);

// Finalize the MPI environment.

MPI_Finalize();

return 0;
}
```

10.5. Compilation and Execution (MPI_{Initthread})

• Compile the program:

```
bash compile.sh mpi_init_thread.c
Command executed: mpicc mpi_init_thread.c -o mpi_init_thread.out
Compilation successful. Check at mpi_init_thread.out
```

• Run the program:

```
bash run.sh ./mpi_init_thread.out 5
```

10.6. Summary

- `MPI_{Init}` is used for standard MPI initialization without considering threading.
- `MPI_{Initthread}` allows the program to specify and check the level of thread support.
 - o Important for applications that require multi-threading in conjunction with MPI.
 - o Ensures that the required thread support is available.

11. Point-to-point communication

```
#include"stdio.h"
#include"mpi.h"
int main(int argc, char **argv)
    int myid, size;
    int myval;
    MPI Status status;
    //Initialize MPI environment
    MPI Init(&argc,&argv);
    //Get total number of processes
    MPI Comm size(MPI COMM WORLD, &size);
    //Get my unique ID among all processes
    MPI Comm rank(MPI COMM WORLD, &myid);
    // Process with ID exactly equal to 0
    if(myid==0){
       //Initialize data to be sent
        myval = 100;
        //Print the data to be sent
       printf("\nmyid: %d \t myval = %d", myid, myval);
        //Send data
        MPI Send(&myval, 1, MPI INT, 1, 0, MPI COMM WORLD);
        printf("\nmyid: %d \t Data sent.\n", myid);
    else if(myid==1){ // Process with ID exactly equal to 1
```

```
//Initialize receive array to some other data
myval = 200;
MPI_Recv(&myval, 1, MPI_INT, 0, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
printf("\nmyid: %d \t Data received.", myid);
//Print received data
printf("\nmyid: %d \t myval = %d", myid, myval);
printf("\n\nProgram exit!\n");
}

//End MPI environment
MPI_Finalize();
}
```

```
bash compile.sh p2p_mpi.c

Command executed: mpicc p2p_mpi.c -o p2p_mpi.out

Compilation successful. Check at p2p_mpi.out
```

```
bash run.sh ./p2p_mpi.out 2
```

```
Command executed: mpirun -np 2 ./p2p mpi.out
OUTPUT
myid: 0
       mvval = 100
myid: 0
       Data sent.
myid: 1
       Data received.
myid: 1
       myval = 100
Program exit!
#########
              DONE
```

11.1. Sending array to process 1

```
#include"stdio.h"
#include"mpi.h"
#define N 100
int main()
    int myid, size;
    int myval;
    int arr[N];
    //Initialize MPI environment
    MPI Init(NULL, NULL);
    //Get total number of processes
    MPI Comm size(MPI COMM WORLD, &size);
    //Get my unique ID among all processes
    MPI Comm rank(MPI COMM WORLD, &myid);
    // Process with ID exactly equal to 0
    if(myid==0){
        //Initialize data to be sent
        for(int i = 0; i < N; i++) arr[i] = i + 1;
        //Send data
        MPI Send(arr, N, MPI INT, 1, 0, MPI COMM WORLD);
        printf("\nmyid: %d \t Data sent.\n", myid);
    else if(myid==1){  // Process with ID exactly equal to 1
        //Initialize receive array to some other data
        MPI Recv(arr, N, MPI INT, 0, 0, MPI COMM WORLD, MPI STATUS IGNORE);
        printf("\nmyid: %d \t Data received.\n", myid);
        //Print received data
        for(int i = 0; i < N; i++)
          printf("%d ", arr[i]);
    }
    //End MPI environment
    MPI Finalize();
}
```

bash compile.sh p2p mpi array.c

Command executed: mpicc p2p mpi array.c -o p2p mpi array.out Compilation successful. Check at p2p_mpi_array.out

bash run.sh ./p2p mpi array.out 2

Command executed: mpirun -np 2 ./p2p mpi array.out

OUTPUT

Data received. myid: 1

myid: 0 Data sent.

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48

########## DONE

Author: Abhishek Raj

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