

## Yeshwantrao Chavan College of Engineering

(An Autonomous Institution affiliated to Rashtrasant Tukadoji Maharaj Nagpur University)
Hingna Road, Wanadongri, Nagpur - 441 110







### Department of Artificial Intelligence & Data Science

Vision of the Department

To be a well-known centre for pursuing computer education through innovative pedagogy, value-based education and industry collaboration.

Mission of the Department

To establish learning ambience for ushering in computer engineering professionals in core and multidisciplinary area by developing Problem-solving skills through emerging technologies.

### Session 2025-2026

Vision: Dream of where you want.	Mission: Means to achieve Vision

**Program Educational Objectives of the program (PEO):** (broad statements that describe the professional and career accomplishments)

PEO1	Preparation	P: Preparation	Pep-CL abbreviation
PEO2	<b>Core Competence</b>	E: Environment	pronounce as Pep-si-lL
		(Learning Environment)	easy to recall
PEO3	Breadth	P: Professionalism	
PEO4	Professionalism	C: Core Competence	
PEO5	Learning	L: Breadth (Learning in	
	Environment	diverse areas)	

**Program Outcomes (PO):** (statements that describe what a student should be able to do and know by the end of a program)

### **Keywords of POs:**

Engineering knowledge, Problem analysis, Design/development of solutions, Conduct Investigations of Complex Problems, Engineering Tool Usage, The Engineer and The World, Ethics, Individual and Collaborative Team work, Communication, Project Management and Finance, Life-Long Learning

PSO Keywords: Cutting edge technologies, Research

"I am an engineer, and I know how to apply engineering knowledge to investigate, analyse and design solutions to complex problems using tools for entire world following all ethics in a collaborative way with proper management skills throughout my life." *to contribute to the development of cutting-edge technologies and Research*.

**Integrity:** I will adhere to the Laboratory Code of Conduct and ethics in its entirety.

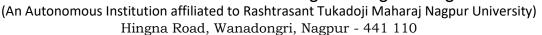
Sanskruti. Paunikar 02/09/2025

Name and Signature of Student and Date

(Signature and Date in Handwritten)









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Session	2025-26 (ODD)	Course Name	High Performance Computing Lab
Semester	7 AIDS	Course Code	22ADS702
Roll No	21	Name of Student	Sanskruti. Paunikar

Practical Number	5
Course Outcome	CO1:-Understand and Apply Parallel Programming Concepts CO1:-Analyze and Improve Program Performance. CO3:-Demonstrate Practical Skills in HPC Tools and Environments.
Aim	Basics of MPI Programming
Theory (100 words)	Message Passing Interface (MPI) is a standardized library employed for parallel programming on distributed memory architectures. It allows several processes on distinct nodes to exchange and synchronize information by passing messages. MPI is extensively applied in High-Performance Computing (HPC) to address large-scale computational issues. Programs are authored in C, C++, or Fortran, compiled by mpicc, and run with mpirun or mpiexec. Fundamental operations are MPI_Init, MPI_Comm_size, MPI_Comm_rank, MPI_Send, MPI_Recv, and MPI_Finalize. MPI ensures scalability, portability, and optimized communication for parallel processes on Linux clusters.
Procedure and Execution	Steps of Implementation:- 1. Install MPI library (e.g., OpenMPI or MPICH) on Linux.
(100 Words)	<ol> <li>Write MPI program in C/C++ using functions like MPI Init, MPI Comm_rank, and MPI Send/Recv.</li> <li>Compile using mpicc program.c -o program.</li> <li>Run with multiple processes: mpirun -np 4 ./program.</li> <li>Observe outputs from each process (rank IDs)</li> </ol>
	sudo apt-get install openmpi-bin openmpi-common libopenmpi-dev brew install open-mpi mpicc hello.c -o hello mpirun -np 4 ./hello mpirun -np 4 ./program_name





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```
GNU nano 5.6.1
include <stdio.
                                                              broadcast.c
                                                                                                                       Modified
 include <mpi.h>
  nt main(int argc, char* argv[]) {
  int rank, size, data;
      MPI_Init(&argc, &argv);
MPI_Comm_rank(MPI_COMM_WORLD, &rank);
MPI_Comm_size(MPI_COMM_WORLD, &size);
       if (rank == 0) {
            data = 42; // root process sets the data printf("Process 0 is broadcasting data %d\n", data);
      // Broadcast data from process 0 to all processe
MPI_Bcast(&data, 1, MPI_INT, 0, MPI_COMM_WORLD);
      printf("Process %d received data %d\n", rank, data);
                     ^O Write Out ^W Where Is ^K Cut
^R Read File ^\ Replace ^U Paste
                                                                                         ^T Execute
^J Justify
                                                                                                               ^C Location
^_ Go To Line
    Exit
    oadcast.c:14:41: error: stray '\' in program
14 | printf("Process %d received data %d\n", rank, data);
prterun was unable to launch the specified application as it lacked permissions to execute an executable:
Executable: ./broadcast Node: localhost
 while attempting to start process rank 0.
[labl@localhost openmpi-5.0.5]$ nano broadcast.c -o broadcast
[labl@localhost openmpi-5.0.5]$ mpicc broadcast.c -o broadcast
[labl@localhost openmpi-5.0.5]$ mpirun -np 4 ./broadcast
Process 0 is broadcasting data 42
Process 0 received data 42
Process 2 received data 42
Process 1 received data 42
Process 3 received data 42
Process 3 received data 42
 lab1@localhost openmpi-5.0.5]$
                                                                                                                         Modified
  GNU nano 5.6.1
                                                                   reduce.c
 include <stdio.h
  nt main(int argc, char* argv[]) {
int rank, size;
int value, sum;
     MPI_Init(&argc, &argv);
MPI_Comm_rank(MPI_COMM_WORLD, &rank);
MPI_Comm_size(MPI_COMM_WORLD, &size);
      value = rank;
      // Reduce operation: sum of all values, result stored in root (process 6) MPI_Reduce(&value, &sum, 1, MPI_INT, MPI_SUM, 0, MPI_COMM_WORLD);
             printf("Sum of all ranks = %d\n", sum);
                      ^O Write Out ^W Where Is ^K Cut
^R Read File ^\ Replace ^U Past
                                                                                          ^T Execute
                                                                                                                ^C Location
     Help
```

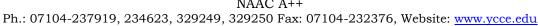




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```
compilation terminated
[lab1@localhost openmpi-5.0.5]$ nano hello_mpi.c
[lab1@localhost openmpi-5.0.5]$ mpicc hello_mpi.c -o hello_mpi
[lab1@localhost openmpi-5.0.5]$ mpirun -np 4 ./hello_mpi
Hello from process 1 of 4
Hello from process 3 of 4
Hello from process 0 of 4
Hello from process 2 of 4
[labl@localhost openmpi-5.0.5]$ nano send_recv.c
[labl@localhost openmpi-5.0.5]$ mpicc send_recv.c -o send_recv
mpirun -np 2 ./send_recv
Process 0 sent data 100 to process 1
Process 1 received data 100 from process 0
[lab1@localhost openmpi-5.0.5]$ nano broadcast.c
[lab1@localhost openmpi-5.0.5]$ mpicc broadcast.c -o broadcast
npirun -np 4 ./broadcast
  nclude <stdio.h&gt;
#include <mpi.h&gt;
int main(int argc, char* argv[]) {
int rank, size, data;
MPI_Init(&argc, &argv);
MPI_Comm_rank(MPI_COMM_WORLD, &rank);
MPI_Comm_size(MPI_COMM_WORLD, &size);
 f (rank == 0) {
data = 42; // Root proces
// Broadcast data from process 0 to all other processes
MPI_Bcast(&data, 1, MPI_INT, 0, MPI_COMM_WORLD);
printf("Process %d received data %d\n", rank, data);
MPI_Finalize();
     rn 0;
                                                                              ^C Location
                G Help
                                                               ^T Execute
 [lab1@localhost openmpi-5.0.5]$ nano hello_mpi.c
[lab1@localhost openmpi-5.0.5]$ mpicc hello_mpi.c -o hello_mpi
[lab1@localhost openmpi-5.0.5]$ mpirun -np 4 ./hello_mpi
Hello from process 1 of 4
Hello from process 3 of 4
Hello from process 0 of 4
Hello from process 2 of 4
[lab1@localhost openmpi-5.0.5]$ nano send_recv.c
cc1:
                        hello_mpi.c: No such file or directory
compilation terminated.
[lab1@localhost openmpi-5.0.5]$ nano hello_mpi.c
[lab1@localhost openmpi-5.0.5]$ mpicc hello_mpi.c -o hello_mpi
[lab1@localhost openmpi-5.0.5]$ mpirun -np 4 ./hello_mpi
Hello from process 1 of 4
Hello from process 3 of 4
Hello from process 0 of 4
Hello from process 2 of 4
[lab1@localhost openmpi-5.0.5]$ nano send_recv.c
[lab1@localhost openmpi-5.0.5]$ mpicc send_recv.c -o send_recv
mpirun -np 2 ./send_recv
Process 0 sent data 100 to process 1
Process 1 received data 100 from process 0
[lab1@localhost openmpi-5.0.5]$ nano broadcast.c
```



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```
lab1@localhost openmpi-5.0.5]$ nano hello_mpi.c
[lab1@localhost openmpi=5.0.5]$ mpicc hello_mpi.c -o hello_mpi
[lab1@localhost openmpi-5.0.5]$ mpirun -np 4 ./hello_mpi
Hello from process 1 of 4
Hello from process 3 of 4
Hello from process 0 of
Hello from process 2 of 4
[lab1@localhost openmpi-5.0.5]$ nano send_recv.c
[lab1@localhost openmpi-5.0.5]$ mpicc send_recv.c -o send_recv
mpirun -np 2 ./send_recv
                                                                                                              Modified
 ≠include <stdio.h
≠include <mpi.h>
  nt main(int argc, char* argv[]) {
   int rank, size, data;
     MPI_Init(&argc, &argv);
MPI_Comm_rank(MPI_COMM_WORLD, &rank);
MPI Comm size(MPI_COMM_WORLD, &size);
       if (rank == 0) {
            data = 100;
            WPI_Send(&data, 1, MPI_INT, 1, 0, MPI_COMM_WORLD);
printf("Process 0 sent data %d to process 1\n", data);
       velse if (rank == 1) {
    MPI_Recv(&data, 1, MPI_INT, 0, 0, MPI_COMM_WORLD, MPI_STATUS_
    printf("Process 1 received data %d from process 6\n", data);
                     ^O Write Out ^W Where Is ^K Cut
^R Read File ^\ Replace ^U Pass
                                                                                                      ^C Location
    Help
                                                                                  ^T Execute
                                                                                                            Q
                                           lab1@localhost:~/openmpi-5.0.5
                                                                                                                     ≡
make[2]: Leaving directory '/home/lab1/openmpi-5.0.5'
make[1]: Leaving directory '/home/lab1/openmpi-5.0.5'
[lab1@localhost openmpi-5.0.5]$ export PATH=$HOME/.openmpi/bin:$PATH
export LD_LIBRARY_PATH=$HOME/.openmpi/lib:$LD_LIBRARY_PATH
[lab1@localhost openmpi-5.0.5]$ mpicc --version
gcc (GCC) 11.5.0 20240719 (Red Hat 11.5.0-5)
Copyright (C) 2021 Free Software Foundation, Inc.
This is free software; see the source for copying conditions. There is NO
warranty; not even for MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE.
[labl@localhost openmpi-5.0.5]$ gcc -fopenmp -o hello_omp hello_omp.ccl: fatal error: hello_omp.c: No such file or directory
compilation terminated.
[lab1@localhost openmpi-5.0.5]$ mpicc hello_mpi.c -o hello_mpi
                            hello_mpi.c: No such file or directory
 compilation terminated.
[lab1@localhost openmpi-5.0.5]$ nano hello_mpi.c
[lab1@localhost openmpi-5.0.5]$ mpicc hello_mpi.c -o hello_mpi
 [lab1@localhost openmpi-5.0.5]$ mpirun -np 4 ./hello_mpi
Hello from process 1 of 4
Hello from process 3 of 4
Hello from process 0 of 4
 Hello from process 2 of 4
   lab1@localhost openmpi-5.0.5]$
```



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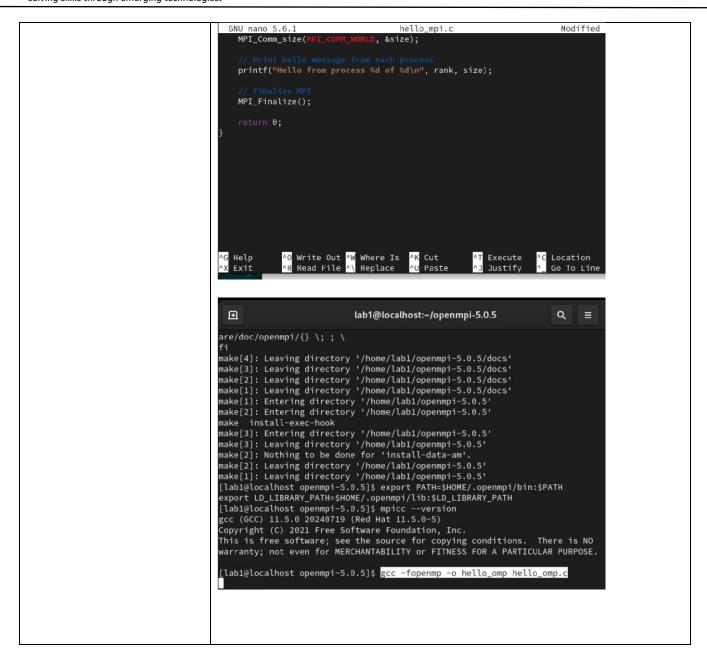
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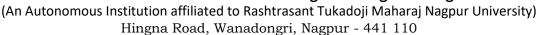
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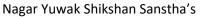
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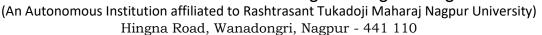
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	• lab1@localhost:~/openmpi-5.0.5
	make install-exec-hook make[3]: Entering directory '/home/labl/openmpi-5.0.5' make[3]: Leaving directory '/home/labl/openmpi-5.0.5' make[2]: Nothing to be done for 'install-data-am'. make[2]: Leaving directory '/home/labl/openmpi-5.0.5' make[1]: Leaving directory '/home/labl/openmpi-5.0.5' make[1]: Leaving directory '/home/labl/openmpi-5.0.5' [labl@localhost openmpi-5.0.5]\$ export PATH=\$HOME/.openmpi/bin:\$PATH export LD_LIBRARY_PATH=\$HOME/.openmpi/lib:\$LD_LIBRARY_PATH [labl@localhost openmpi-5.0.5]\$ mpiccversion gcc (GCC) 11.5.0 20240719 (Red Hat 11.5.0-5) Copyright (C) 2021 Free Software Foundation, Inc. This is free software; see the source for copying conditions. There is NO warranty; not even for MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE.  [labl@localhost openmpi-5.0.5]\$ gcc -fopenmp -o hello_omp hello_omp.c ccl: fatal error: hello_omp.c: No such file or directory compilation terminated.  [labl@localhost openmpi-5.0.5]\$ mpicc hello_mpi.c -o hello_mpi ccl: fatal error: hello_mpi.c: No such file or directory compilation terminated.  [labl@localhost openmpi-5.0.5]\$ nano hello_mpi.c [labl@localhost openmpi-5.0.5]\$ mpicc hello_mpi.c -o hello_mpi [labl@localhost openmpi-5.0.5]\$ mpirun -np 4 ./hello_mpi
Output Analysis	When executed with multiple processes, each process prints its rank and the total number of processes. Communication via MPI_Send and MPI_Recv demonstrates message passing between processes. The output confirms correct distribution of tasks across processes, proving parallel execution on Linux.
Github link	https://github.com/sanskruti-1234/HPC.git
Conclusion	MPI allows efficient communication between processes for parallel programming. It provides scalability and portability for distributed computing on Linux systems













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