

Objective

Q1: Fill in the blanks

1. Cluster
2. Size
3. MPI_Ssend, recv
4. Simple instruction, multiple data
5. Local Area Multicomputer
6. _np
7. Graphics Processing Unit
8. MPI_Barrier
9. Parallel Random Access Machine
10. Don't know
11. EREW exclusive read exclusive write
12. $P(n) \times T(n)$
13. MIMD multiple instruction multiple data
14. Cuda, Libra
15. Don't know
16. MISD multiple instruction single data

Q2: True/False

1. F
2. F
3. F
4. F
5. T
6. T
7. T
8. F
9. F
10. F
11. F
12. F
13. T
14. T
15. T

4 11 12 are not confirmed

Q3: choose the best answer

1. MPI_Ssend, MPI_SRecv
2. MIMD
3. SIMD
4. Gather
5. Scatter
6. MIMD
7. SIMD
8. SISD
9. Don't know
10. Don't know

Subjective

Q1: what is parallel computing and why we need it ?

Parallel computing is a form of **computation** in which many calculations are carried out **simultaneously**, operating on the principle that **large problems can often be divided into smaller ones**, which are then **solved concurrently ("in parallel")**. We need parallel computing because we can **use multiple processors in parallel** to solve problems more quickly than with a single processor.

Q2: Briefly describe the working of the following with the help of diagram.

ARRAY PROCESSOR: Array processor is a synchronous parallel computer with **multiple ALU** called **processing elements (PE)** that can operate in parallel in lock step fashion. It is composed of N identical PE under the control of a single control unit and many memory modules. Array processor also frequently use a form of parallel computation called **pipelining** where an operation is divided into smaller steps and the steps are performed simultaneously.

VECTOR PROCESSOR: A vector processor is a central processing unit that **can work on an entire vector in one instruction**. The instruction to processor is in the form of one computer vector instead of its elements. It is also known as an **array processor**. It exhibits SIMD behavior by having operations that are applied to all elements on vector.

GPU: A graphics processing unit (GPU) is a **single drip processor** primarily used to manage and boost the performance of video and graphics. It is in PCs on a video card or mother board as well as mobile phones, display adapters, work stations and game consoles.

Q3(a): Differentiate the following:

CLUSTER COMPUTING

1. It is **homogenous** network. **Similar hardware component running a similar operating system are connected together in a cluster.**
2. They are **within the same location** or complex.
3. The resources of all the nodes in a cluster are centrally **managed by a resource manager.**

GRID COMPUTING

1. It is a **heterogenous** network. **Different computer hardware running various kinds of operating systems are connected together in a grid.**
2. They are **distributed over a LAN, MAN or WAN.** They can be geographically separated.
3. Each entity (node) in a grid **behaves like an independent entity.** This means it **manages its resources by itself.**

Multicomputer

1. A computer **made up of several computers.** similar to **parallel computing.**
2. Distributed computing deals with hardware and software systems containing more than one processing element, multiple programs, running under a loosely or tightly controlled regime.
3. multicomputer **have one physical address space per CPU.**
4. It can **run faster**
5. A multi-computer is **multiple computers, each of which can have multiple processors.**
Used for **true parallel processing.**

Multiprocessors

1. A multiprocessor system is simply a computer that has **more than one CPU on its motherboard.**
2. Multiprocessing is the use of **two or more central processing units (CPUs)** within a single computer system.
3. Multiprocessors have a **single physical address space (memory) shared by all the CPUs**
4. A multiprocessor would **run slower**, because it would be in ONE computer.
5. A multi-processor is a single system with multiple CPU's.

Q3(b): Mark where applicable.

	TCP	UDP	MPI
UNICASTING	yes	yes	yes
MULTICASTING	no	yes	yes
BROADCASTING	no	yes	yes
MANY TO ONE	no	no	yes
MANY TO MANY	no	no	yes

Q4: Write an MPI program to generate following cartesian topology.

ODD EVEN TOPOLOGY

```
#include "mpi.h"
void main(int argc, char *argv[])
{
    int nrow, mcol, root, Iam, ndim, p, rank;
    int dims[2], coords[2], cyclic, reorder;
    MPI_Comm comm, comm1, ceven, codd;
    MPI_Group e_group, o_group;

    MPI_Init(&argc, &argv);          /* starts MPI */
    MPI_Comm_rank(MPI_COMM_WORLD, &Iam); /* get current process id */
    MPI_Comm_size(MPI_COMM_WORLD, &p);  /* get number of processes */

    nrow = 4; mcol = 2; ndim = 2;
    root = 0; cyclic = 1; reorder = 1;

    dims[0] = nrow;    /* rows */
    dims[1] = mcol;    /* columns */

    for(int i=0; i<p/2; i++)
    {
        ranks[i] = i*2;
    }

    MPI_Group world_group;
    MPI_Comm_group(MPI_COMM_WORLD, &world_group);

    MPI_Group_incl(world_group, n/2, ranks, &e_group);
    MPI_Group_excl(world_group, n/2, ranks, &o_group);

    MPI_Comm_create_group(MPI_COMM_WORLD, o_group, 0, &comm);
    MPI_Comm_create_group(MPI_COMM_WORLD, e_group, 0, &comm1);

    MPI_Cart_create(comm, ndim, dims, cyclic, reorder, &codd);
    MPI_Cart_create(comm1, ndim, dims, cyclic, reorder, &ceven);

    MPI_Finalize();
}
```

Q5:

Matrix addition

```
#include "mpi.h"
#define row_size = 5
#define col_size = 3

void main(int argc, char *argv[])
{
    int matrix1[row_size * col_size];
    int matrix2[row_size * col_size];
    int matrix3[row_size * col_size];

    MPI_Init(&argc, &argv);          /* starts MPI */
    MPI_Comm_rank(MPI_COMM_WORLD, &Iam); /* get current process id */
    MPI_Comm_size(MPI_COMM_WORLD, &p); /* get number of processes */

    int row1[col_size];
    int row2[col_size];
    int row3[col_size];

    //matrix1 , col_size, MPI_INT -- for master
    //row1, col_size, MPI_INT -- for other

    //for first matrix
    MPI_Scatter(matrix1 , col_size, MPI_INT, row1, col_size, MPI_INT, 0,
MPI_COMM_WORLD);
    //for second matrix
    MPI_Scatter(matrix2 , col_size, MPI_INT, row2, col_size, MPI_INT, 0,
MPI_COMM_WORLD);

    sum(row1, row2, &row3);

    MPI_Gather(row3, col_size, MPI_INT, matrix3, col_size*row_size, MPI_INT, 0,
MPI_COMM_WORLD);
    if( Iam== 0){
        for(int i=0; i< row_size*col_size; i++){
            printf("%d", matrix3[i]);
        }
    }

    MPI_Finalize();
}
```

Q6:

SUM OF NUMBERS

```
#include "mpi.h"
#define SIZE = 1000

void main(int argc, char *argv[])
{
    int data[SIZE];
    MPI_Init(&argc, &argv);          /* starts MPI */
    MPI_Comm_rank(MPI_COMM_WORLD, &Iam); /* get current process id */
    MPI_Comm_size(MPI_COMM_WORLD, &p);  /* get number of processes */

    MPI_Bcast(data, SIZE, MPI_INT, 0, MPI_COMM_WORLD);

    int s=sum(data, Iam*SIZE/p, SIZE/p);
    int asum=0;

    MPI_Reduce(&s, &asum, 1, MPI_INT, MPI_SUM, 0, MPI_COMM_WORLD);

    if(Iam == 0){
        printf("%d", asum);
    }
    MPI_Finalize();

}

void sum(int[] data, int start, int range)
{
    int ans=0;
    for(int i=start; i<start + range; ++i)
    {
        ans += data[i];
    }
    return ans;
}
```

Q7:

Algorithm Broadcast_EREW

Processor P_1

y (in P_1 's private memory) $\leftarrow x$

$L[1] \leftarrow y$

For $i=0$ to $\log p - 1$ do

Forall P_j , where $2^i + 1 \leq j \leq 2^{i+1}$ do in parallel

y (in P_j 's private memory) $\leftarrow L[j - 2^i]$

$L[j] \leftarrow y$

endfor

endfor

Complexity Analysis

Run Time $\rightarrow T(n)$

Number of Processors $\rightarrow P(n)$

Cost $\rightarrow C(n) = T(n) * P(n)$

Q8:

Algorithm Broadcast_EREW

For $i=0$ to $\log n$ do

Forall P_j , where $2^{i-1} + 1 \leq j \leq n$ do in parallel

$A[j] \leftarrow A[j] + A[j - 2^{i-1}]$

endfor

endfor

Complexity Analysis

Run Time $\rightarrow T(n)$

Number of Processors $\rightarrow P(n)$

Cost $\rightarrow C(n) = T(n) * P(n)$