CS 3006 Parallel and Distributed Computer

Fall 2022

- Learn about parallel and distributed computer architectures.(1)
- Implement different parallel and distributed programming paradigms and algorithms using Message-Passing Interface (MPI) and OpenMP.(4)
- Perform analytical modelling, dependence, and performance analysis of parallel algorithms and programs. (2)
- Use Hadoop or MapReduce programming model to write bigdata applications.(5)

Week # 9 – Lecture # 22, 23, 24

20th, 22nd, 23rd Rabi ul Awwal, 1444 17th, 19th, 20th October 2022

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Lecture # 22 – Topics (Lab # 9)

- Review of MPI_Send/Recv
- MPI_Get_Count to get more info. using MPI_Recv return status
- Parallizing the Trapezoidal Rule
- Collective Operation like MPI_Broadcast, MPI_Reduce, etc.
- MPI_Reduce and MPI_AllReduce
- Collective vs Point-to-Point communication
- Non-Blocking Send and Receive. MPI_Wait and MPI_Test
- Reading and Distributing a vector. MPI_Scatter, MPI_Gather, MPI_AllGather

```
int MPI Send(
   void*
               msg buf /* input */,
               msg size /* input */,
   int
   MPI Datatype msg type /* input */,
                dest /* input */,
   int
                tag /* input */,
   int
                     /* input */)
   MPI Comm
                comm
int MPI Recv(
   void*
               msg buf /* output */,
               msg size /* input */,
   int
   MPI Datatype msg type /* input */,
               source /* input */,
   int
   int
                       /* input */,
          tag
                    /* input */,
   MPI Comm comm
   MPI Status* status p /* output */)
```

- MPI_Send may behave differently with regard to buffer size, cutoffs and blocking.
- MPI_Recv always blocks until a matching message is received.

Data types

MPI datatype	C datatype
MPI_CHAR	signed char
MPI_SHORT	signed short int
MPI_INT	signed int
MPI_LONG	signed long int
MPI_LONG_LONG	signed long long int
MPI_UNSIGNED_CHAR	unsigned char
MPI_UNSIGNED_SHORT	unsigned short int
MPI_UNSIGNED	unsigned int
MPI_UNSIGNED_LONG	unsigned long int
MPI_FLOAT	float
MPI_DOUBLE	double
MPI_LONG_DOUBLE	long double
MPI_BYTE	
MPI_PACKED	

```
1 const int MAX NUMBERS = 100;
 2 int numbers[MAX NUMBERS];
   int number amount;
 4
 5 if (world rank == 0) {
 6 srand(time(NULL));
      number amount = (rand() / (float) RAND_MAX) * MAX_NUMBERS;
      MPI Send(numbers, number amount, MPI INT, 1, 0, MPI COMM WORLD);
      printf("0 sent %d numbers to 1\n", number_amount);
10 - } else if (world rank == 1) {
11
      MPI Status status;
    MPI Recv(numbers, MAX NUMBERS, MPI INT, 0, 0, MPI COMM WORLD, &status);
12
     MPI_Get_count( & status, MPI_INT, & number_amount); //check how many actually recv.
13
      printf("1 received %d numbers from 0. Message source = %d, tag = %d\n",
14 -
15
              number amount, status.MPI SOURCE, status.MPI TAG);
16 }
```

```
Get a, b, n;
      h = (b-a)/n;
      local_n = n/comm_sz;
      local_a = a + my_rank*local_n*h;
4
      local_b = local_a + local_n*h;
      local_integral = Trap(local_a, local_b, local_n, h);
6
      if (my_rank != 0)
         Send local_integral to process 0;
9
      else /* my_rank == 0 */
10
         total_integral = local_integral;
         for (proc = 1; proc < comm_sz; proc++)
11
12
            Receive local_integral from proc;
13
            total_integral += local_integral;
14
15
      if (my_rank == 0)
16
```

Parallelizing the Trapezoidal Rule

- 1. Partition problem solution into tasks.
- 2. Identify communication channels between tasks.
- 3. Aggregate tasks into composite tasks.
- 4. Map composite tasks to cores.

- All collective operations must be called by all processes in the communicator
- MPI_Bcast is called by both the sender (called the root process) and the processes that are to receive the broadcast
 - MPI_Bcast is not a "multi-send"

 $if (my_rank == 0) {$

"root" argument is the rank of the sender; this tells MPI which process originates the broadcast and which receive

printf("Enter a, b, and n\n");
 scanf("%lf %lf %d", a_p, b_p, n_p);
}

MPI_Bcast(a_p, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
MPI_Bcast(b_p, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
MPI_Bcast(n_p, 1, MPI_INT, 0, MPI_COMM_WORLD);

Seen by all processes master and slaves

MPI_Bcast MPI_Reduce 2) 3) MPI_SUM 0 • 1 • 2 • 3 • MPI_Allgather MPI_Scatter MPI_Reduce 2 3 7 8 MPI_SUM ① • 1 • 2 • 3 • 18 14 MPI_Allreduce MPI_Gather 2 3 7 8 MPI_SUM

MPI_Reduce

Operation Value	Meaning
MPI_MAX	Maximum
MPI_MIN	Minimum
MPI_SUM	Sum
MPI_PROD	Product
MPI_LAND	Logical and
MPI_BAND	Bitwise and
MPI_LOR	Logical or
MPI_BOR	Bitwise or
MPI_LXOR	Logical exclusive or
MPI_BXOR	Bitwise exclusive or
MPI_MAXLOC	Maximum and location of maximum
MPI_MINLOC	Minimum and location of minimum

Collective vs. Point-to-Point Communications

- <u>All</u> the processes in the communicator must call the same collective function.
- For example, a program that attempts to match a call to MPI_Reduce on one process with a call to MPI_Recv on another process is erroneous, and, in all likelihood, the program will hang or crash.

- The arguments passed by each process to an MPI collective communication must be "compatible."
- For example, if one process passes in 0 as the dest_process and another passes in 1, then the outcome of a call to MPI_Reduce is erroneous, and, once again, the program is likely to hang or crash.
- Point-to-point communications are matched on the basis of tags and communicators.
- Collective communications don't use tags.
- They're matched solely on the basis of the communicator and the order in which they're called.

MPI_Allreduce

 Useful in a situation in which all of the processes need the result of a global sum in order to complete some larger computation.

Non-blocking is not part of the syllabus. You are encourage to include it in your Semester Project code.

Non-blocking Send

```
int MPI Isend(
    void*
                 msg buf /* input */,
    int
                 msg size /* input */,
    MPI Datatype msg type /* input */,
    int
                 dest
                         /* input */,
                         /* input */,
    int
                 tag
    MPI Comm
                         /* input */,
                 comm
    MPI Request &request /* output */)
```

Identical to MPI_Send but added argument for getting handle to MPI_Request struct.

Non-blocking Test

```
int MPI_Test(
   MPI_Request request /* input */,
   int    &flag /* output */
   MPI_Status /* output */)
```

Flag returns 1 if completed, otherwise 0.

```
MPI_Issend(buf,count,datatype,dest,tag,comm,request);
do {
    MPI_Test(request,flag,status);
} while (flag != 1);
```

Non-blocking Receive

```
int MPI Irecv(
                 msg buf /* output */,
    void*
                 msq size /* input */,
    int
    MPI Datatype msg type /* input */,
                 source
                         /* input */,
    int
                         /* input */,
    int
                 tag
                         /* input */,
    MPI Comm
                 comm
    MPI Request &request /* output */)
```

Identical to MPI_Read but replaces MPI_Status argument with MPI_Request struct. Status is now retrieved in the MPI_Wait.

Blocking Wait

```
int MPI_Wait(
   MPI_Request request /* input */,
   MPI_Status &status /* output */)

   Same status that would have been returned from
   MPI_Recv, not as useful in MPI_Send.
```

```
MPI_Isend (buf, count, datatype, dest, tag, comm, request);
MPI_Wait (request, status);
if (status...) {
```

Scatter

 MPI_Scatter can be used in a function that reads in an entire vector on process 0 but only sends the needed components to each of the other processes.

Gather

 Collect all of the components of the vector onto process 0, and then process 0 can process all of the components.

```
int MPI_Gather(
     void*
                send_buf_p /* in */,
                send_count /* in */,
     int
                send_type /*in */,
    MPI_Datatype
    void*
                recv_buf_p /* out */,
        recv_count /*in */,
     int
    MPI_Datatype recv_type /*in */,
                dest_proc /* in */,
     int
                comm /* in */);
    MPI_Comm
```

Reading and distributing a vector

```
void Read vector(
                                                    double local_a[] /* out */,
                                                    int local n /* in */.
                                                    int n /* in */,
                                                    char vec_name[] /* in */,
                                                    int my_rank /* in */,
if (my_rank == 0) {
                                                    MPI_Comm comm /* in */)
   a = malloc(n*sizeof(double));
                                                  double * a = NULL:
   printf("Enter the vector %s\n", vec_name)
                                                  int i:
   for (i = 0; i < n; i++)
      scanf("%lf", &a[i]);
   MPI_Scatter(a, local_n, MPI_DOUBLE, local_a, local_n, MPI_DOUBLE,
         0, comm);
   free(a);
  else {
   MPI_Scatter(a, local_n, MPI_DOUBLE, local_a, local_n, MPI_DOUBLE,
         0, comm);
```

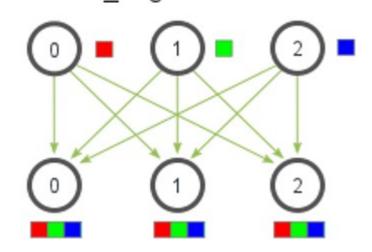
Print a distributed vector

```
void Print_vector(
     double
             local b[] /* in */.
     int
             local n
                       /* in */.
     int
                       /* in */.
     char
             title[]
                       /* in */.
     int
             my_rank /* in */,
     MPI_Comm comm /* in */)
  double * b = NULL;
  int i:
```

```
if (my_rank == 0) {
  b = malloc(n*sizeof(double));
  MPI_Gather(local_b, local_n, MPI_DOUBLE, b, local_n, MPI_DOUBLE,
        0, comm);
  printf("%s\n", title);
  for (i = 0; i < n; i++)
     printf("%f ", b[i]);
  printf("\n");
  free(b):
 else
  MPI_Gather(local_b, local_n, MPI_DOUBLE, b, local_n, MPI_DOUBLE,
         0, comm);
```

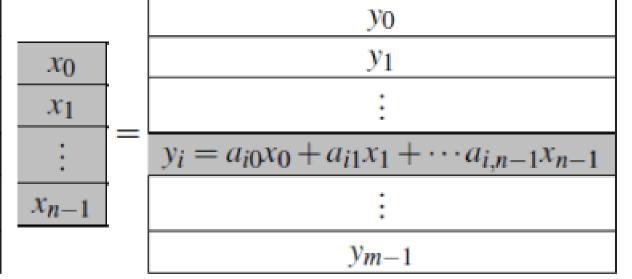
Allgather

- Concatenates the contents of each process' send_buf_p and stores this in each process' recv_buf_p.
- As usual, recv_count is the amount of data being received from each process.



Matrix-vector multiplication

<i>a</i> ₀₀	<i>a</i> ₀₁		$a_{0,n-1}$
a_{10}	a_{11}		$a_{1,n-1}$
:	:		:
a_{i0}	a_{i1}	• • •	$a_{i,n-1}$
<i>a_{i0}</i> :	<i>a</i> _{i1}	•••	$a_{i,n-1}$



```
void Mat_vect_mult(
      double local_A[] /* in */,
      double local_x[] /* in */,
                                                 MPI_Allgather
      double local_y[] /* out */,
      int local_m /* in */,
                      /* in */,
      int
          n
         local_n /* in */,
      int
     MPI_Comm comm /* in */) {
  double * x;
   int local_i, j;
   int local_ok = 1;
   x = malloc(n*sizeof(double));
   MPI_Allgather(local_x, local_n, MPI_DOUBLE,
        x, local_n, MPI_DOUBLE, comm);
   for (local_i = 0; local_i < local_m; local_i++) {
     local_y[local_i] = 0.0;
      for (j = 0; j < n; j++)
        local_y[local_i] += local_A[local_i*n+j]*x[j];
                                      y_i = a_{i0}x_0 + a_{i1}x_1 + \cdots + a_{i,n-1}x_{n-1}
   free(x);
   /* Mat_vect_mult */
```

Lecture # 23 – Topics

- Interaction among Tasks running of different processors
- Task Interaction Graph
 - Capturing data dependencies
- Task Interaction graph- An example
 - Sparse Matrix Vector Multiplication
 - Task interaction graph
- Processes and Mapping
- Criteria for Mapping
- Mapping data query to processes Example

Task Interaction Graphs

- Subtasks generally exchange data with others in a decomposition.
 - For example, even in the trivial decomposition of the dense matrix-vector product, if the vector is not replicated across all tasks, they will have to communicate elements of the vector.

- The graph of tasks (nodes) and their interactions/data exchange (edges) is referred to as a *task interaction graph*.
 - Task interaction graphs represent data dependencies,
 - Task dependency graphs represent control dependencies.

The pattern of interaction among tasks is captured by what is known as a **task-interaction graph**. The nodes in a task-interaction graph represent tasks and the edges connect tasks that interact with each other. The nodes and edges of a task-interaction graph can be assigned weights proportional to the amount of computation a task performs and the amount of interaction that occurs along an edge, if this information is known. The edges in a task-interaction graph are usually undirected, but directed edges can be used to indicate the direction of flow of data, if it is unidirectional. The edge-set of a task-interaction graph is usually a superset of the edge-set of the task-dependency graph. In the database query example discussed earlier, the task-interaction graph is the same as the task-dependency graph. We now give an example of a more interesting task-interaction graph that results from the problem of sparse matrix-vector multiplication.

Example 3.3 Sparse matrix-vector multiplication

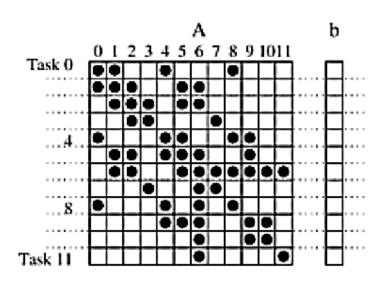
Consider the problem of computing the product y = Ab of a sparse $n \times n$ matrix A with a dense $n \times 1$ vector b. A matrix is considered sparse when a significant number of entries in it are zero and the locations of the non-zero entries do not conform to a predefined structure or pattern. Arithmetic operations involving sparse matrices can often be optimized significantly by avoiding computations involving the zeros.

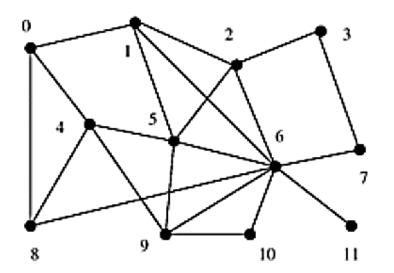
One possible way of decomposing this computation is to partition the output vector y and have each task compute an entry in it. Figure 3.6(a) illustrates this decomposition. In addition to assigning the computation of the element y[i] of the output vector to Task i, we also make it the "owner" of row A[i, *] of the matrix and the element b[i] of the input vector. Note that the computation of y[i] requires access to many elements of b that are owned by other tasks. So Task i must get these elements from the appropriate locations. In the message-passing paradigm, with the ownership of b[i], Task i also inherits the responsibility of sending b[i] to all the other tasks that need it for their computation. For example, Task 4 must send b[4] to Tasks 0, 5, 8, and 9 and must get b[0], b[5], b[8], and b[9] to perform its own computation. The resulting task-interaction graph is shown in Figure 3.6(b). \blacksquare

Task Interaction graph- An example

$$\begin{bmatrix} a & b & c & d \\ e & f & g & h \\ i & j & k & l \\ m & n & o & p \end{bmatrix} \times \begin{bmatrix} x \\ y \\ z \\ w \end{bmatrix} = \begin{bmatrix} ax + by + cz + dw \\ ex + fy + gz + hw \\ ix + jy + kz + lw \\ mx + ny + oz + pw \end{bmatrix}$$

Figure 3.6. A decomposition for sparse matrix-vector multiplication and the corresponding task-interaction graph. In the decomposition Task i computes $\sum_{0 \le j \le 11, A[i,j] \ne 0} A[i,j], b[j]$.





(a)

Processes and Mapping

- Mapping: the mechanism by which tasks are assigned to processes for execution.
- Process: a logic computing agent that performs tasks, which is an abstract entity that uses the code and data corresponding to a task to produce the output of that task.
- Why use processes rather than processors?
 - We rely on OS to map processes to physical processors.
 - We can aggregate tasks into a process

Criteria of Mapping

- Maximize the use of concurrency by mapping independent tasks onto different processes
- Minimize the total completion time by making sure that processes are available to execute the tasks on critical path as soon as such tasks become executable
- Minimize interaction among processes by mapping tasks with a high degree of mutual interaction onto the same process.

Basis for Choosing Mapping

Task-dependency graph

Makes sure the max. concurrency
Task-interaction graph

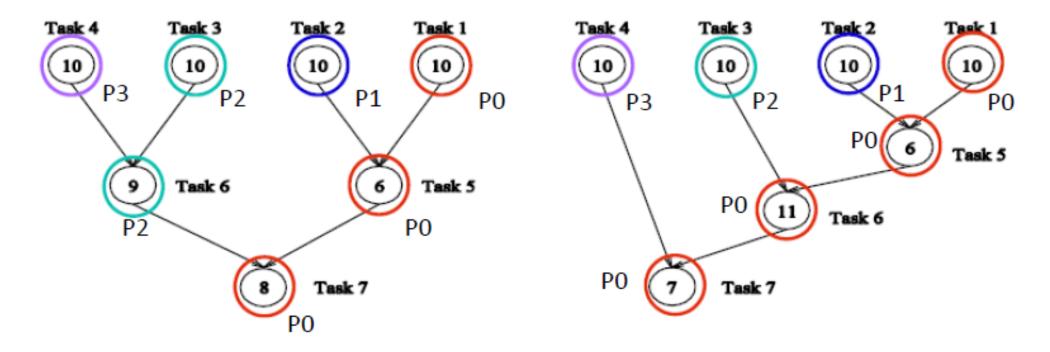
Minimum communication.

These criteria often conflict with each other.

For example, a decomposition into one task (or no decomposition at all) minimizes interaction but does not result in a speedup at all!

- Finding a balance that optimizes the overall parallel performance is the key to a successful parallel algorithm.
- Therefore, mapping of tasks onto processes plays an important role in determining how efficient the resulting parallel algorithm is.

Example: Mapping Database Query to Processes



No two nodes in a level have dependencies, therefore, single level tasks are assigned to different processes.

- 4 processes can be used in total since the max. concurrency is 4.
- Assign all tasks within a level to different processes.

Lecture # 24 – Topics

Decomposition Techniques

- Recursive Decomposition
- Data Decomposition
- Exploratory Decomposition
- Speculative Decomposition

Decomposition Techniques

So how does one decompose a task into various subtasks?

While there is no single recipe that works for all problems, we present a set of commonly used techniques that apply to broad classes of problems. These include:

- recursive decomposition
- data decomposition
- exploratory decomposition
- speculative decomposition

Recursive Decomposition

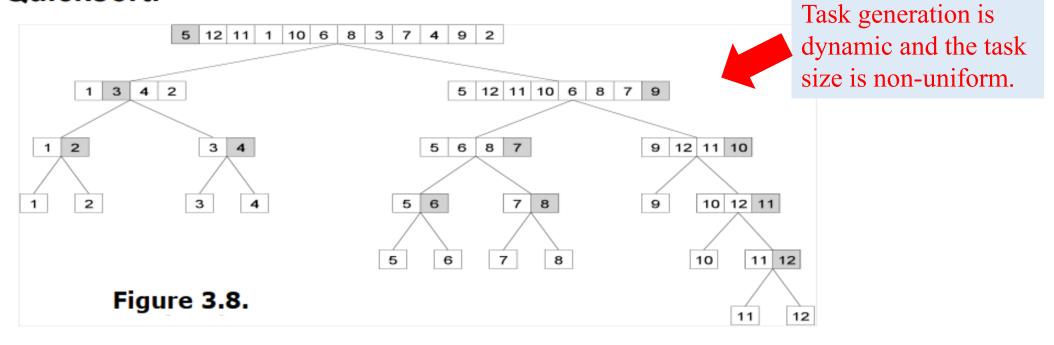
 Generally suited to problems that are solved using the divideand-conquer strategy.

 A given problem is first decomposed into a set of subproblems.

 These sub-problems are recursively decomposed further until a desired granularity is reached.

Recursive Decomposition: Example

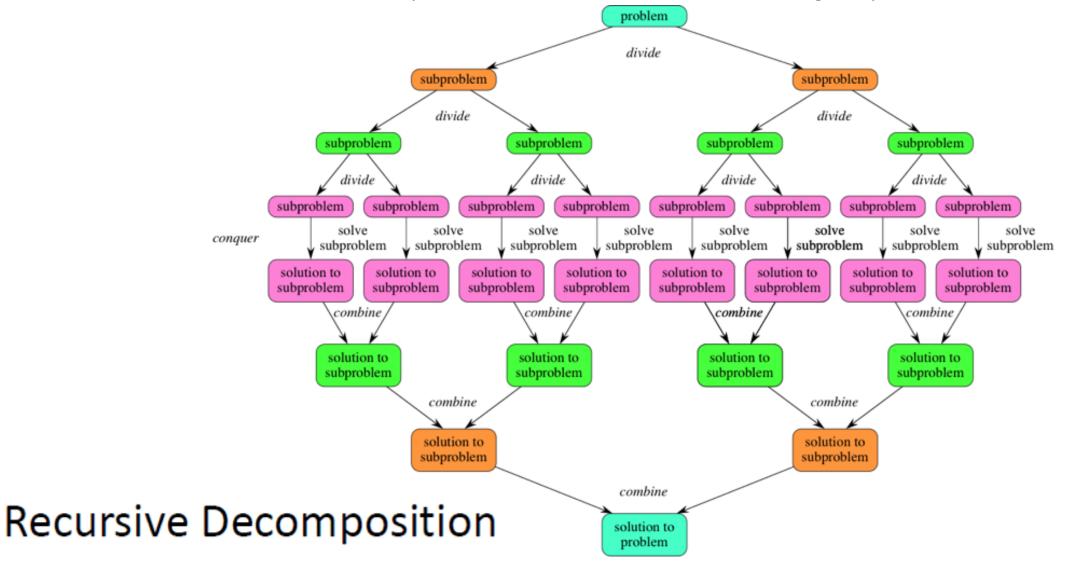
A classic example of a divide-and-conquer algorithm on which we can apply recursive decomposition is Quicksort.



In this example, a task represents the work of partitioning a (sub)array. Note that each subarray represents an independent subtask. This can be repeated recursively.

You should think of a divide-and-conquer algorithm as having three parts:

- **1.Divide** the problem into a number of subproblems that are smaller instances of the same problem.
- 2.Conquer the subproblems by solving them recursively. If they are small enough, solve as base cases.
- **3.Combine** the solutions to the subproblems into the solution for the original problem.



Algorithm 3.1 A serial program for finding the minimum in an array of

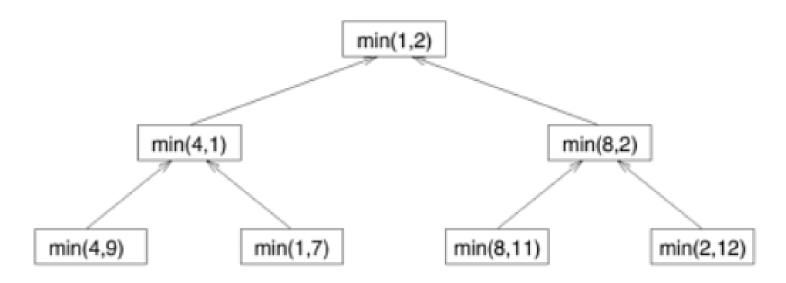
numbers A of length n.

Algorithm 3.2 A recursive program for finding the minimum in an array of numbers A of length n.

```
procedure RECURSIVE_MIN (A, n)
     begin
     if (n = 1) then
         min := A[0];
     else
         lmin := RECURSIVE MIN (A, n/2);
       rmin := RECURSIVE MIN (&(A[n/2]), n - n/2);
         if (lmin < rmin) then</pre>
             min := lmin;
10.
         else
11.
             min := rmin;
12.
         endelse;
     endelse;
13.
14.
     return min;
15.
     end RECURSIVE MIN
```

```
    procedure SERIAL_MIN (A, n)
    begin
    min = A[0];
    for i := 1 to n - 1 do
    if (A[i] < min) min := A[i];</li>
    endfor;
    return min;
    end SERIAL_MIN
```

Figure 3.9. The task-dependency graph for finding the minimum number in the sequence {4, 9, 1, 7, 8, 11, 2, 12}. Each node in the tree represents the task of finding the minimum of a pair of numbers.



Data Decomposition

 Ideal for problems that operate on large data structures

Steps

- The data on which the computations are performed are partitioned
- Data partition is used to induce a partitioning of the computations into tasks.

Data Partitioning

- Partition output data
- Partition input data
- Partition input + output data
- Partition intermediate data

Data Decomposition: Output Data Decomposition

- Often, each element of the output can be computed independently of others (but simply as a function of the input).
- A partition of the output across tasks decomposes the problem naturally.
- Input: if each output is described as a function of the input directly. Some combination of the individual results may be necessary.
- Output data decomposition: if it applies, it can result in less communication.
- Intermediate data decomposition more rare.
- Owner computes rules: the process that owns a part of the data performs all the computations related to it.

Consider the problem of multiplying two $n \times n$ matrices A and B to yield matrix C. The output matrix C can be partitioned into four tasks as follows:

$$\begin{pmatrix} A_{1,1} & A_{1,2} \\ A_{2,1} & A_{2,2} \end{pmatrix} \cdot \begin{pmatrix} B_{1,1} & B_{1,2} \\ B_{2,1} & B_{2,2} \end{pmatrix} \to \begin{pmatrix} C_{1,1} & C_{1,2} \\ C_{2,1} & C_{2,2} \end{pmatrix}$$

Task 1:
$$C_{1,1} = A_{1,1}B_{1,1} + A_{1,2}B_{2,1}$$

Task 2:
$$C_{1,2} = A_{1,1}B_{1,2} + A_{1,2}B_{2,2}$$

Task 3:
$$C_{2,1} = A_{2,1}B_{1,1} + A_{2,2}B_{2,1}$$

Task 4:
$$C_{2,2} = A_{2,1}B_{1,2} + A_{2,2}B_{2,2}$$

A given data decomposition does not result in a unique decomposition into tasks.

Figure 3.11. Two examples of decomposition of matrix multiplication into eight tasks.

Decomposition I	Decomposition II
Task 1: $C_{1,1} = A_{1,1}B_{1,1}$	Task 1: $C_{1,1} = A_{1,1}B_{1,1}$
Task 2: $C_{1,1} = C_{1,1} + A_{1,2}B_{2,1}$	Task 2: $C_{1,1} = C_{1,1} + A_{1,2}B_{2,1}$
Task 3: $C_{1,2} = A_{1,1}B_{1,2}$	Task 3: $C_{1,2} = A_{1,2}B_{2,2}$
Task 4: $C_{1,2} = C_{1,2} + A_{1,2}B_{2,2}$	Task 4: $C_{1,2} = C_{1,2} + A_{1,1}B_{1,2}$
Task 5: $C_{2,1} = A_{2,1}B_{1,1}$	Task 5: $C_{2,1} = A_{2,2}B_{2,1}$
Task 6: $C_{2,1} = C_{2,1} + A_{2,2}B_{2,1}$	Task 6: $C_{2,1} = C_{2,1} + A_{2,1}B_{1,1}$
Task 7: $C_{2,2} = A_{2,1}B_{1,2}$	Task 7: $C_{2,2} = A_{2,1}B_{1,2}$
Task 8: $C_{2,2} = C_{2,2} + A_{2,2}B_{2,2}$	Task 8: $C_{2,2} = C_{2,2} + A_{2,2}B_{2,2}$

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Task 2: $C_{1,1} = C_{1,1} + A_{1,2}B_{2,1}$	Task 2: $C_{1,1} = C_{1,1} + A_{1,2}B_{2,1}$
Task 3: $C_{1,2} = A_{1,1}B_{1,2}$	Task 3: $C_{1,2} = A_{1,2}B_{2,2}$
Task 4: $C_{1,2} = C_{1,2} + A_{1,2}B_{2,2}$	Task 4: $C_{1,2} = C_{1,2} + A_{1,1}B_{1,2}$
Task 5: $C_{2,1} = A_{2,1}B_{1,1}$	Task 5: $C_{2,1} = A_{2,2}B_{2,1}$
Task 6: $C_{2,1} = C_{2,1} + A_{2,2}B_{2,1}$	Task 6: $C_{2,1} = C_{2,1} + A_{2,1}B_{1,1}$
Task 7: $C_{2,2} = A_{2,1}B_{1,2}$	Task 7: $C_{2,2} = A_{2,1}B_{1,2}$
Task 8: $C_{2,2} = C_{2,2} + A_{2,2}B_{2,2}$	Task 8: $C_{2,2} = C_{2,2} + A_{2,2}B_{2,2}$

Consider the problem of counting the instances of given itemsets in a database of transactions. In this case, the output (itemset frequencies) can be partitioned across tasks.

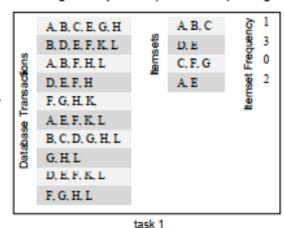
(a) Transactions (input), itemsets (input), and frequencies (output)

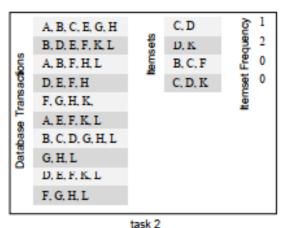
suc	A, B, C, E, G, H B, D, E, F, K, L A, B, F, H, L		A, B, C D, E C, F, G	1 60 0
Fransaction	D, E, F, H	msets	A, E	Ped 2
	F, G, H, K, A, E, F, K, L	ē	C, D D, K	₩ 1 ₩ 2
Catabase	B, C, D, G, H, L G, H, L		B, C, F C, D, K	= 0
	D, E, F, K, L			
	F, G, H, L			

Problem: Find the number of times that each itemset in *I* appears in all the transactions; i.e., the number of transactions of which each itemset is a subset of.

(b) Partitioning the frequencies (and itemsets) among the tasks

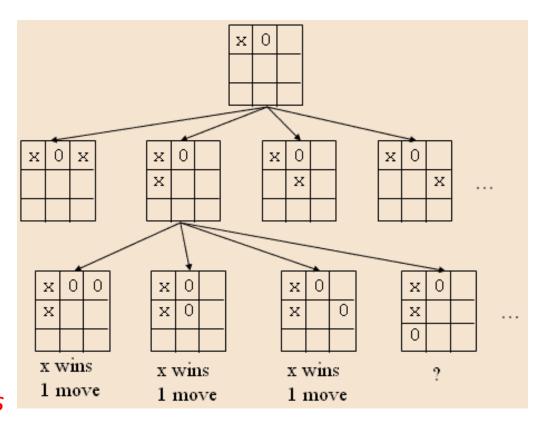
Part (b) shows how two tasks can achieve results by partitioning the output into two parts and having each task compute its half of the frequencies.





Exploratory Decomposition

- Example: looking for the best move in a game.
- Simple case: generate all possible configurations from the starting position.
- Send each of the configurations to a child process.
- Each process will look for possible best moves for the opponent recursively eventually using more processes.
- When it finds the result, it sends it back to the parent.
- The parent selects the best move from all of the results received from the child (eventually the worst move for the opponent).



Speculative Decomposition

- Switch statement in a program: We wait to know the value of the expression and execute only the corresponding case.
- In speculative decomposition we execute some or all of the cases in advance.
- When the value of the expression is know, we keep only the results from the computation to be executed in that case.
- The gain in performance comes from anticipating the possible computations.

```
Sequential version
                       Parallel version
                       Slave(i)
compute expr;
switch (expr) {
                         compute ai;
  case 1:
                         Wait (request);
    compute a1;
                         if (request)
    break:
                           Send(ai, 0);
  case 2:
                      Master()
    compute a2;
    break:
  case 3:
                         compute expr;
    compute a3;
                         swicth (expr) {
    break; ....
                           case 1:
                             Send(request, 1);
                             Receive(a1, i);
```

The difference with the exploratory decomposition is that we can compute the possible states before the next move is performed.