Parallel Algorithms and Programming

Parallel Algorithms in Distributed Memory Systems (Part 1)

Thomas Ropars

Email: thomas.ropars@univ-grenoble-alpes.fr

Website: tropars.github.io

In this lecture

- Performance of parallel distributed algorithms
- Parallel operations on vectors and matrices
 - Matrix-vector multiplication
 - Matrix-matrix multiplication

Introduction

Introduction

Assumptions

- A distributed memory system
 - Processes communicate by send and receiving messages
- A **virtual ring** topology

Main costs to consider

- The cost of running computation
 - floating point operations in our context
- The cost of moving data
 - from the memory to the processors
 - between nodes (over the interconnection network)

Cost of a distributed parallel algorithm

• A computational term:

$$nb of flops \times time per flop$$

• A bandwidth term:

$$\text{amount of data moved} \times \frac{1}{bandwidth}$$

• A latency term:

nb of messages \times *latency*

Cost of a distributed parallel algorithm

We should remember that (in general):

time per flop
$$\ll \frac{1}{bandwidth} \ll latency$$

Consequence

- Minimizing communication is important for performance
- It is also important for energy efficiency
 - Moving data from/to DRAM or over the interconnection network are the most energy consuming operations

About memory accesses

Cost of moving data between the memory and the processor

- Can often be ignored
 - Cost mostly hidden by hardware prefetchers

Prefetchers

- Hardware mechanisms to load data in the cache before it is needed
- Based on the memory access pattern of the program
- Works well for regular access patterns
 - It is the case for the programs studied in this lecture

Matrix-vector multiplication

Specification of the problem

We consider:

- A is a matrix of size $n \times n$
- ullet x, y are vectors of size n

We want to compute:

$$y = Ax$$

Specification of the problem

$$ext{Ax} = egin{bmatrix} a_{11} & a_{12} & \cdots & a_{1n} \ a_{21} & a_{22} & \cdots & a_{2n} \ dots & dots & dots & dots \ a_{m1} & a_{m2} & \cdots & a_{mn} \end{bmatrix} egin{bmatrix} x_1 \ x_2 \ dots \ x_n \end{bmatrix} = egin{bmatrix} a_{11}x_1 + a_{12}x_2 + \cdots + a_{1n}x_n \ a_{21}x_1 + a_{22}x_2 + \cdots + a_{2n}x_n \ dots \ a_{21}x_1 + a_{22}x_2 + \cdots + a_{2n}x_n \ dots \ a_{m1}x_1 + a_{m2}x_2 + \cdots + a_{mn}x_n \end{bmatrix}$$

Sequential implementation

```
for (i=0; i<n; i++) {
    y[i] = 0;
    for (j=0; j<n; j++) {
        y[i] = y[i] + A[i][j] * x[j]
    }
}</pre>
```

Observations

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- The computation of each value of y is the scalar product between a row of A and the vector x.
- Each scalar product is independent

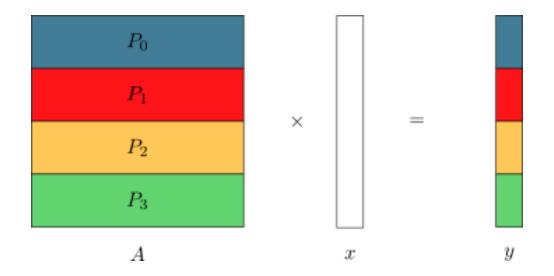
Parallel algorithm

Observations

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Parallel algorithm

- Distribute the rows of A over p processors
 - Each processor has a copy of x
- ullet Each processor computes $r=rac{n}{p}$ scalar products in parallel



- Distributed matrix-vector multiplication over 4 processors.
- All data with the same color are on the same processor

Improving performance and solving larger problems

Impact of distributing the data and the computation over several nodes

- Improve the execution time (hopefully)
- Solve larger problems
 - Matrix A might not fit in the memory of a single node
 - \blacksquare Distributing the rows of matrix ${\tt A}$ over ${\tt p}$ nodes allows storing a much larger matrix in memory

Making the algorithm more generic

Initial distribution of the blocks of x

- An execution scenario:
 - y = Ax followed by z = By
 - At the begining of the second computation, blocks of y are distributed over the nodes (see previous figure)

A generic matrix-vector multiplication function

- Function that implements y = Ax
 - With both A and x distibuted over the nodes

Basic steps

Basic steps

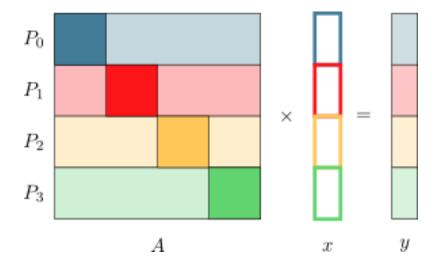
- Computation step:
 - ullet Each processor computes a partial result using the elements of x it has in its local memory
- 2. Communication step:
 - ullet Each processor sends its block of x to its successor and receives from its predecessor in the virtual ring

Basic steps

- Computation step:
 - ullet Each processor computes a partial result using the elements of x it has in its local memory
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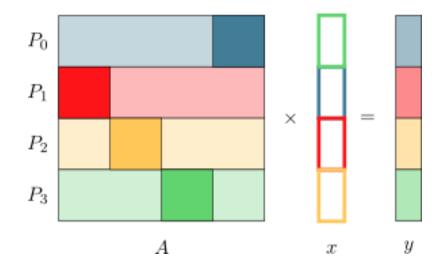
Number of iterations to terminate: *p*

Computation step 0



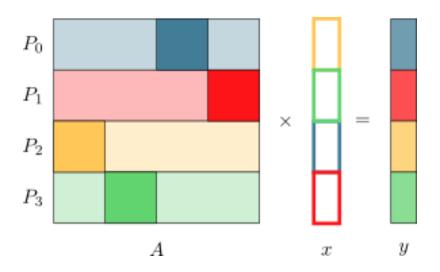
- Distributed matrix-vector multiplication over 4 processors
 - Each processor stores $\frac{n}{4}$ rows of A
 - Each processor stores $\frac{n}{4}$ values of x.
- Colors for blocks of vector x refer to the initial position of the blocks.

Computation step 1

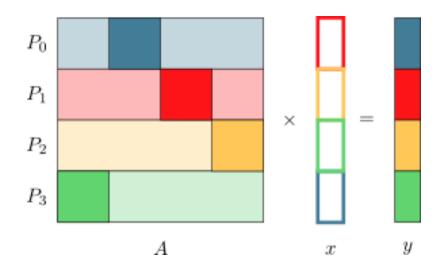


- Blocks of x have been *shifted* in the virtual ring
 - Focus on the color of the blocks

Computation step 2



Computation step 3



ullet The last round of communication will bring x back into its initial state

```
double A[r][n]; /* rows of A, assumed to be already initialized*/
double x[r]; /* values of x, assumed to be already initialized*/
double y[r];

int rank = my_rank();
int P = num_procs();

double tempS[r], tempR[r];
```

```
tempS <- x; /* copy of the values */
for(step = 0; step < P; step++){
    Send(tempS, (rank+1) % P);
    Recv(tempR, (rank-1) % P);
    int block = (rank-step) % P;
    for (i=0; i<r; i++) {</pre>
        for(j=0; j<r; j++){
          y[i] = y[i] + A[i, r * block + j)] * tempS[j]
    tempS <- tempR;</pre>
```

Computing the indexes of elements of A to use at each step

- On which block of A, processor K should compute at step S?
 - Vector x is divided in p blocks
 - At step 0, processor K computes using block K of x
 - At step 1, processor K computes using block K-1 of x
 - At step 2, processor K computes using block K-2 of x

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- Computation of the block index: $block = (rank step) \mod p$
- ullet Computation of the indexes in a row of A

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- Computation of the block index: $block = (rank step) \mod p$
- Computation of the indexes in a row of A
 - $r = \frac{n}{p}$ (p is the number of processors)
 - start index: $block \times r$
 - lacksquare end index: (block+1) imes r-1

About the execution in parallel

- Algorithm designed to allow the communication and the computation inside one step to occur in parallel (symbol | | |)
 - Requires to allocate an extra buffer (tempR) for the communication
 - The communication in the last iteration allows restoring the initial distribution of the blocks of x

Implementation with MPI

How to overlap communication and computation?

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Implementation with MPI

- How to overlap communication and computation?
 - Use of MPI_Isend() and MPI_Irecv(), plus MPI_Wait()

Parameters

- p processors
- matrix of size $n \times n$
- $r = \frac{n}{p}$

Model parameters

- L: the latency of a network communication
- B: the bandwidth of a network communication
- w: the computation time for one basic unit of work
 - = two floating point operations in this algorithm

- ullet Time for a communication: $T_{comm} = L + rac{r}{B}$
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- ullet Time for computation in one step: $T_{comp}=r^2 imes w$
 - lacksquare Computation on a block of size r imes r

Total time

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Total time

- *p* iterations
- Computation occur in parallel with communication

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Total time

- *p* iterations
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$$T_{MV}(p) = p imes \max(T_{comp}, T_{comm})$$

$$T_{MV}(p) = p imes \max(r^2 imes w, L + rac{r}{B}) = \max(rac{n^2}{p} imes w, p imes L + rac{n}{B})$$

Performance for large matrices

When n becomes large, the execution time is asymptotically equal to:

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Speedup of p with p processors

Distribution of the data

- Case 1: The rows of A are already distributed between the processors
 - A is the result of a previous computation
 - A is loaded from a file on network file system
 - Parallel IO libraries (e.g., MPI-IO) allow processors to load data in parellel
- Case 2: The data are initially on one node and should be aggregated on a single node at the end of the computation

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- Case 2: The data are initially on one node and should be aggregated on a single node at the end of the computation
 - Use of MPI_Scatter() and MPI_Gather() functions

Matrix-matrix multiplication

Specification of the problem

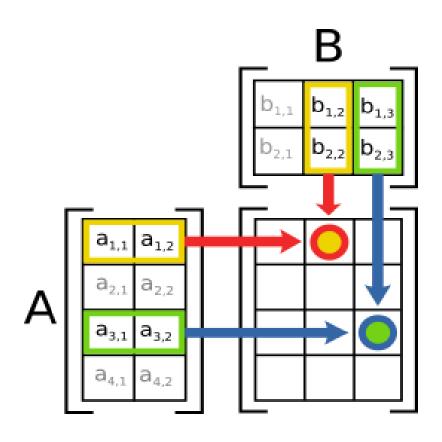
We consider:

• A, B, C are matrices of size $n \times n$

We want to compute:

$$C = A \times B$$

Specification of the problem



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

Source: Wikipedia (Bilou)

Sequential implementation

```
for (i=0; i<n; i++) {
    for (j=0; j<n; j++) {
        C[i][j] = 0;
        for (k=0; k<n; k++) {
              C[i][j] = C[i][j] + A[i][k] * B[k][j];
        }
    }
}</pre>
```

Parallel implementation

Same assumptions as before

- p processes
- A virtual ring topology

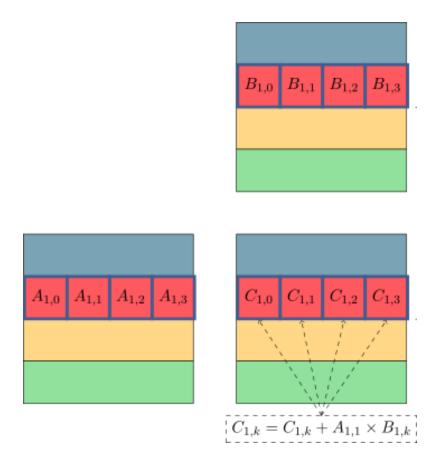
Basic idea of the algorithm

• Same idea as for matrix-vector multiplication

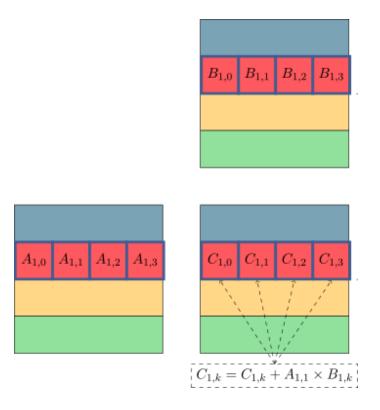
Description of the algorithm

- The 3 matrices are distributed over the processes based on the rows.
 - Each process stores $r=rac{n}{p}$ rows of each matrix.
- In each iteration:
 - 1. Each processor runs a partial matrix-matrix multiplication based on the data that are available locally.
 - 2. Each processor sends r rows of B to the next processor and receives r rows of B from the previous processor

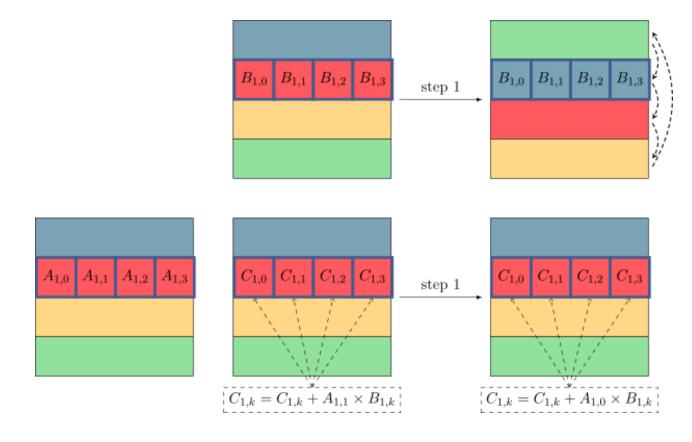
The algorithm requires p iterations to terminate.



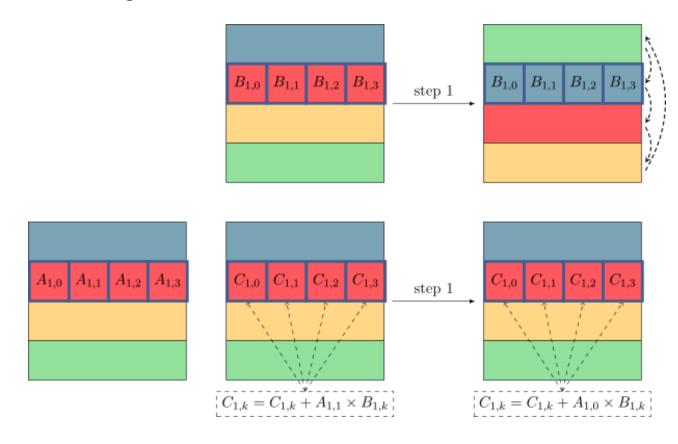
- The figure assumes 4 processors (4 colors)
 - The figure focuses on the computation run by processor P_1 .
- The notation $A_{l,k}$ refers to the **block** k in the part of matrix A stored by processor l



- ullet In step 0, processor P_k has rows [r imes k] to [r imes (k+1)-1] of B
 - P_1 has rows r to 2r-1 of B
- ullet P_k can only compute based on elements [r imes k] to [r imes (k+1)-1] of A in each row (block $A_{k,k}$)



- During step 0, processors exchange their rows of B in the virtual ring
- Processors keep the same rows of A during the whole algorithm



- ullet In step 1, processor P_k has rows [r imes (k-1)] to [r imes k-1] of B
 - $lacksquare P_1$ has rows 0 to r-1 of B

Algorithm in pseudo C code

```
double A[r][n]; /* rows of A, assumed to be already initialized*/
double B[r][n]; /* rows of B, assumed to be already initialized*/
double C[r][n];

int rank = my_rank();
int P = num_procs();

double tempS[r][n], tempR[r][n];
```

Algorithm in pseudo C code

```
tempS <- B; /* copy of the values */
for(step = 0; step < P; step++){
    Send(tempS, (rank+1) % P);
    Recv(tempR, (rank-1) % P);
    int block = (rank-step) % P;
    for (1=0; 1<P; 1++) {
        for(i=0; i<r; i++) {
            for(j=0; j<r; j++){
                for (k=0; k<r; k++) {
                     C[i][1 * r + j] = C[i][1 * r + j]
                     + A[i][block * r + k] * tempS[k][l * r + j]$;
    tempS <- tempR;</pre>
```

The additional index (l) is introduced to iterate over the blocks of B and C.

Parameters

- p processors
- matrix of size $n \times n$
- $r=rac{n}{p}$

Model parameters

- L: the latency of a network communication
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 - Processors send and receive r rows of B
- Time for computation in one step: $T_{comp} = n imes r^2 imes w$
 - lacktriangle Computation on a block of size n imes r with r operations for each value

Total time

- *p* iterations
- Computation occur in parallel with communication

$$T_{MM}(p) = p imes \max(T_{comp}, T_{comm})$$
 $T_{MM}(p) = p imes \max(n imes r^2 imes w, L + rac{n imes r}{B}) = \max(rac{n^3}{p} imes w, p imes L + rac{n^2}{B})$

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Speedup of p with p processors

About Bulk communication

Observation:

ullet The parallel matrix-matrix multiplication could be implemented as n matrix-vector multiplications (one for each column of B)

Total time

$$T_{MM-alt}(p) = n imes T_{MV}(p)$$
 $T_{MM-alt}(p) = \max(rac{n^3}{p} imes w, n imes p imes L + rac{n^2}{B})$

About Bulk communication

 T_{MM} vs T_{MM-alt}

$$T_{MM}(p) = \max(rac{n^3}{p} imes w, p imes L + rac{n^2}{B})$$
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About Bulk communication

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$$T_{MM}(p) = \max(rac{n^3}{p} imes w, p imes L + rac{n^2}{B})$$
 $T_{MM-alt}(p) = \max(rac{n^3}{p} imes w, n imes p imes L + rac{n^2}{B})$

- The asymptotic efficiency is the same
- The latency term is multiplied by n with the alternative version
 - Can have a big impact on performance in practice
 - Relates to the fact that instead of exchanging blocks of matrices, processors would exchange blocks of vectors

Sending data in bulk

- General technique to reduce the cost of communicating over the network in parallel algorithms
 - Reduces the impact of latency.

References

• Section 4.1 and 4.2 of the book "Parallel Algorithms" (by Casanova, Robert, and Legrand).