Parallel Systems Course: Chapter VI

Dense Matrix Algorithms

KUMAR Chapter 8

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Overview

- 1. Matrix-vector Multiplication
- 2. Matrix-matrix Multiplication
- 3. Shared-memory

Utility of Matrix Algorithms

Applied in several numerical and nonnumerical contexts:

- 3D image calculations
- Solving (linear) equations
- Simulations of physical systems

E.g.: makes the basis of LINPACK, a software library for performing numerical linear algebra, by using the BLAS (Basic Linear Algebra Subprograms) libraries for performing basic vector and matrix operations

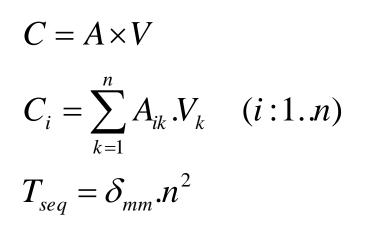
Dense versus Sparse Matrices

Dense matrices: have no or few known zero entries

Sparse matrices: are populated primarily with zeros

- often appear in science or engineering when solving partial differential equations
- easily compressed
- very large sparse matrices are impossible to manipulate with the standard algorithms, due to memory limitations
- special versions of the algorithms are necessary and are more efficient

Matrix - Vector Multiplication



n: number of elements in the vector, number of rows and columns of the matrix

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

```
A

A11 A12 A13 .. .. A1n

A21 A22 .. .. A2n

...

Ai1 Ai2 Ai3 .. .. Ain

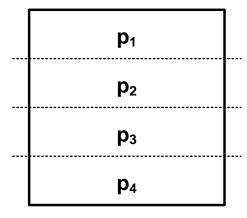
...

Ci

...

An1 An2 An3 .. .. Ann
```

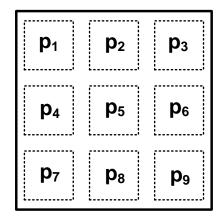
Matrix/Vector Partitioning?

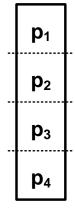


p₁ p₂ p₃ p₄

Row-wise block striping

Column-wise block striping





Checkerboard partitioning

Vector partitioning

Data & results distributed

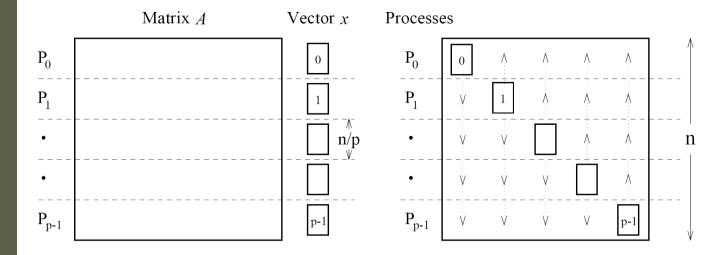
Now: distributed memory solutions

 In parallel applications, data remains distributed while operations (such as vectormatrix operations) are performed on them.

 In the following we assume that the data is already distributed among the processors

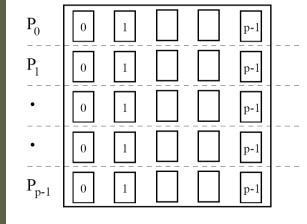
Parallel M x V Multiplication Version 1

$$(n=p)$$



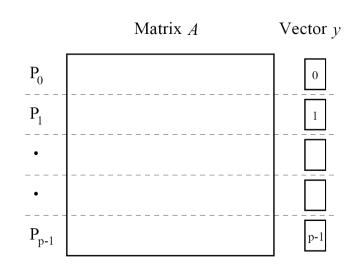
(a) Initial partitioning of the matrix and the starting vector *x*

Row-wise partitioning



(c) Entire vector distributed to each process after the broadcast

(b) Distribution of the full vector among all the processes by all-to-all broadcast



(d) Final distribution of the matrix and the result vector *y*

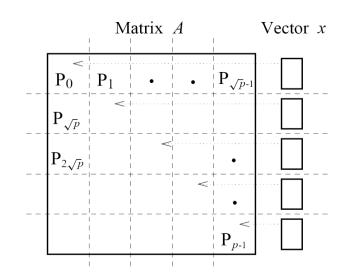
Figure 8.1 Multiplication of an $n \times n$ matrix with an $n \times 1$ vector using rowwise block 1-D partitioning. For the one-row-per-process case, p = n.

Parallel M x V Multiplication Version 1 p < n

n/p rows of matrix and n/p elements of vector per processor

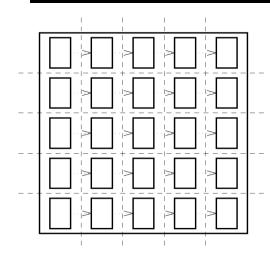
Parallel M x V Multi-plication Version 2

$$(n^2=p)$$

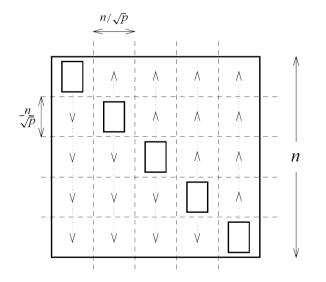


(a) Initial data distribution and communication steps to align the vector along the diagonal

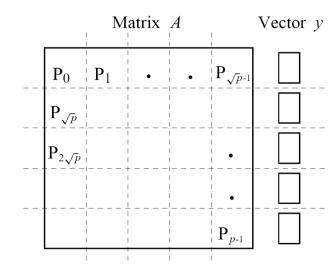
checkerboard partitioning



(c) All-to-one reduction of partial results



(b) One-to-all broadcast of portions of the vector along process columns



(d) Final distribution of the result vector

Figure 8.2 Matrix-vector multiplication with block 2-D partitioning. For the one-element-per-process case, $p = n^2$ if the matrix size is $n \times n$.

Parallel M x V Multiplication Version 2 p < n²

 (n/\sqrt{p}) x (n/\sqrt{p}) blocks of matrix and n/\sqrt{p} elements of vector per processor



Matrix Multiplication

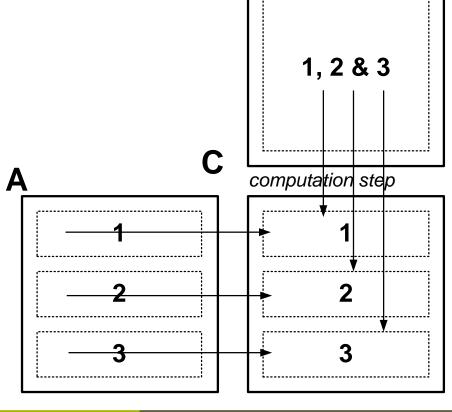
```
C = A \times B
                                                                                  B11 B12 . . B1j . . . . B1n
       C_{ij} = \sum_{i}^{n} A_{ik} . B_{kj} \quad (i, j:1.n)
                                                                                   B21 B22 .. B2j
       T_{s} = \delta_{mm}.n^{3}
                                                                                   Bn1 Bn2 .. Bnj .. Bnn
for (i=0; i< n; i++) {
                                                         A11 A12 A13 .. .. A1n
                                                         A21 A22 .. .. A2n
    for (j=0; j< n; j++) {
          C[i,j]=0;
                                                         Ai1 Ai2 Ai3 ...
          for (k=0; k< n; k++) {
              C[i,j]+=A[i,k]*B[k,j];
                                                         An1 An2 An3 .. .. Ann
```

MxM: one-step version

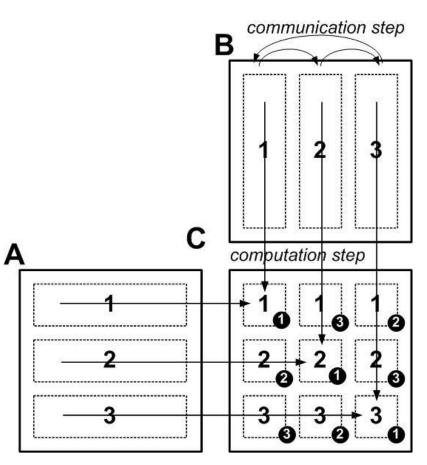
B is sent to all processors

Computation in 1 step

Same amount of communication

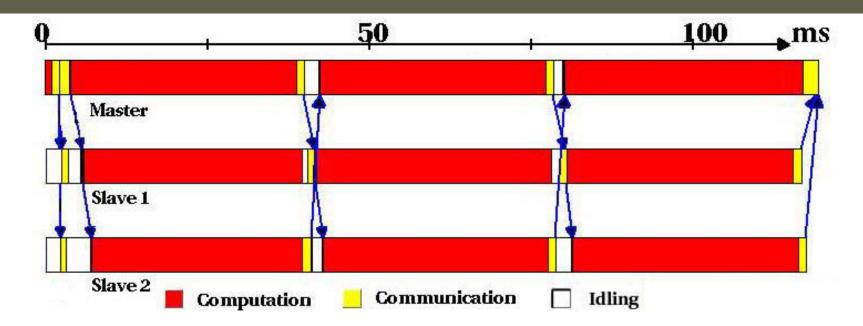


Alternate shift-compute version



- Algorithm alternates p computation and communication steps
- Computation step: each processor multiplies its A submatrix with its B submatrix, resulting in a submatrix of C. The black circles indicate the step in which each submatrix is computed.
- After multiplication: processor sends it B submatrix to next processor and receives one from the preceding processor. The communication forms a circular shift operation.

Parallel MxM: Execution Profile



Speedup=2.55 Efficiency = 85%

Note that the communication after the first step can be hidden behind the computation (Mehdi Moghaddmfar 2017).

MPI-2 supports non-blocking collective communication operations.

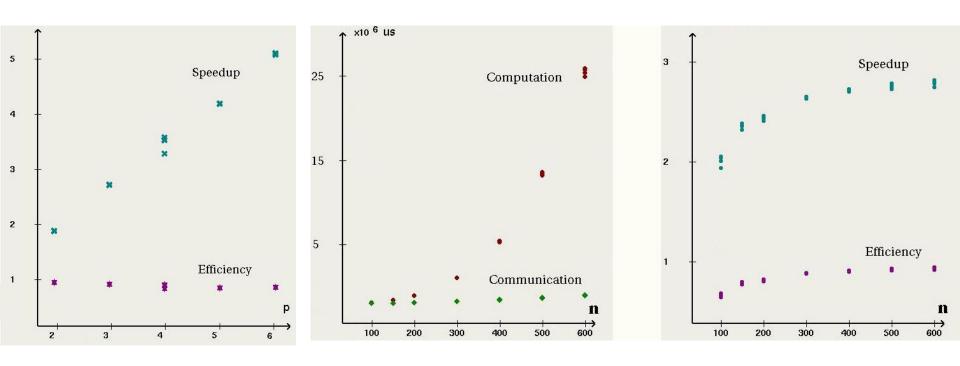
Theoretical Analysis of Matrix Multiplication

- Computation time (slave) $T_{work}^i = \frac{n^3}{p}.\delta_{mm}$
- Communication time (slave) including data distribution (initially 2x1/p, each round 1/p [p-1 rounds] + 1/p result)

$$T_{comm}^{i} = (p+1).t_s + (1+\frac{2}{p}).n^2.t_w$$

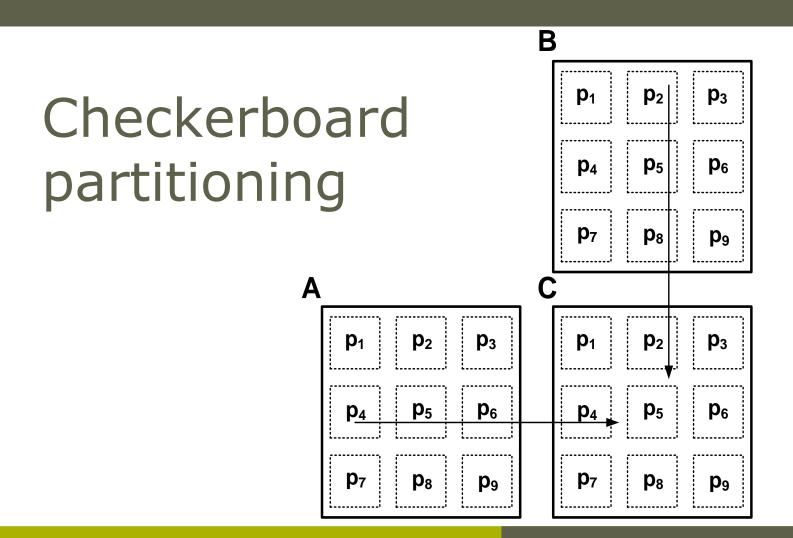
- \rightarrow Overhead ratio=overhead/computation = O(n²)/O(n³/p) = O(p/n)
- →Scalable since efficiency remains constant if we increase both p and n

Parameter Dependence of Matrix Multiplication



n: work size, here: matrix size

V3: Cannon's algorithm

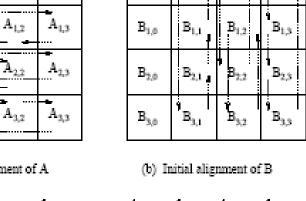


Cannon's parallel MxM

Communication steps on 16 processes

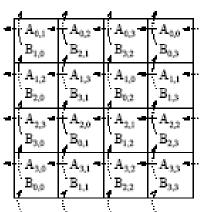
$A_{0,0}$	$\mathbf{A}_{0,1}$	$A_{0,2}$	$A_{0,3}$
$\mathbf{A}_{1,0}$	$\mathbf{A}_{1,1}$	A _{1,2}	A _{1,3}
A _{2,0}	A _{2,1}	A _{2,2}	A _{2,3}
$A_{3,0}$	$A_{3,1}$	A _{3,2}	A _{3,3}

(a) Initial alignment of A



 $B_{n,n}$

(c) A and B after initial alignment



 $B_{0,2}$

 $\mathbb{B}_{0,1}$

 $\mathbb{B}_{0,3}$

(d) Submatrix locations after first shift

	4	/ 1	4
$-1 - A_{0,2} -$	$A_{0,3}$	$-A_{0,0}$	A _{0,1} ==
$B_{2,0}$	$B_{3,1}$	$B_{0,2}$	$B_{1,3}$
A _{1,3}	$A_{1,0} =$	$-A_{1,1} -$	A _{1,2}
B _{3,0}	$\mathbf{B}_{0,1}$	B _{1,2}	B _{2,3}
- A _{2,0} -	$A_{2,1}$	$-A_{2,2}$	A _{2,3}
$B_{0,0}$	$B_{1,1}$	B _{2,2}	B _{3,3}
$-A_{3,1}$	$A_{3,2}$	+A _{3,3} -	A _{3,0}
B _{1,0}	$B_{2,1}$	B _{3,2}	$B_{0,3}$
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 $A_{0,3}$ $A_{0,0}$ $A_{0,1}$ $A_{0,2}$ $B_{n,n}$ $\mathbf{B}_{0.1}$ $B_{1,2}$ $B_{2,3}$ $A_{1,0}$ $A_{1,2}$ $A_{2,1}$ $A_{2,2}$ $A_{2.3}$ $A_{2.0}$ $\mathbb{B}_{2,1}$ $\mathbf{B}_{1,0}$ $B_{3,2}$ $B_{0.3}$ $A_{3,0}$ $A_{3,2}$ $A_{3,3}$ $A_{3.1}$ $\mathbb{B}_{1.3}$

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(e) Submatrix locations after second shift (f) Submatrix locations after third shift

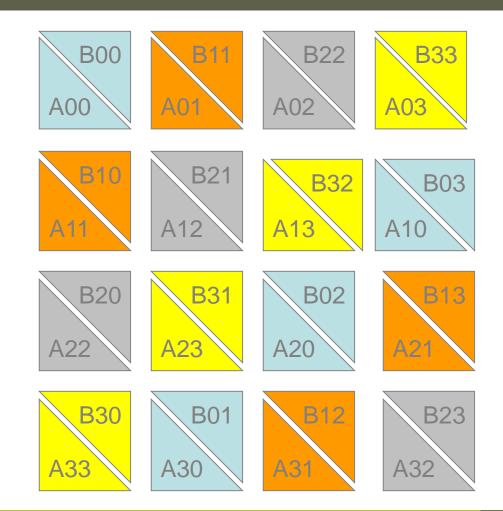
Initially, blocks Need to Be Aligned

Each triangle represents a matrix block

Only same-color triangles should be multiplied

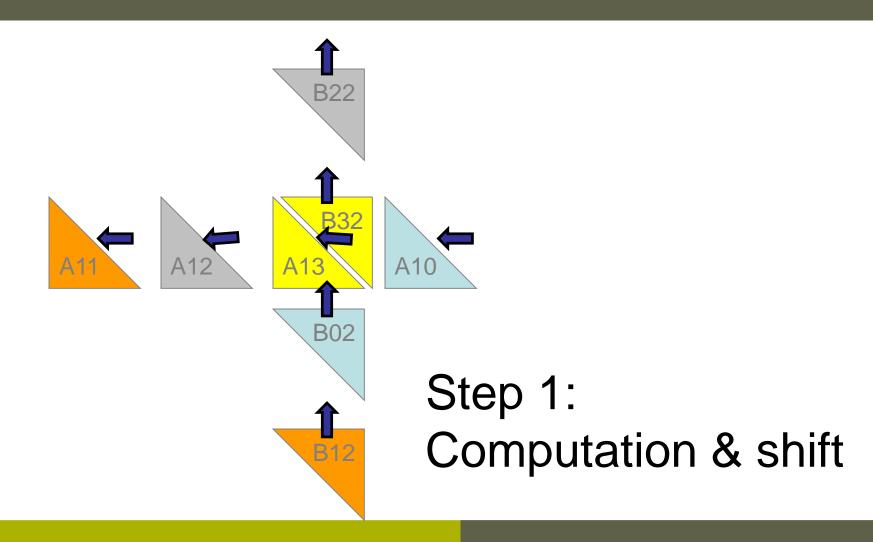


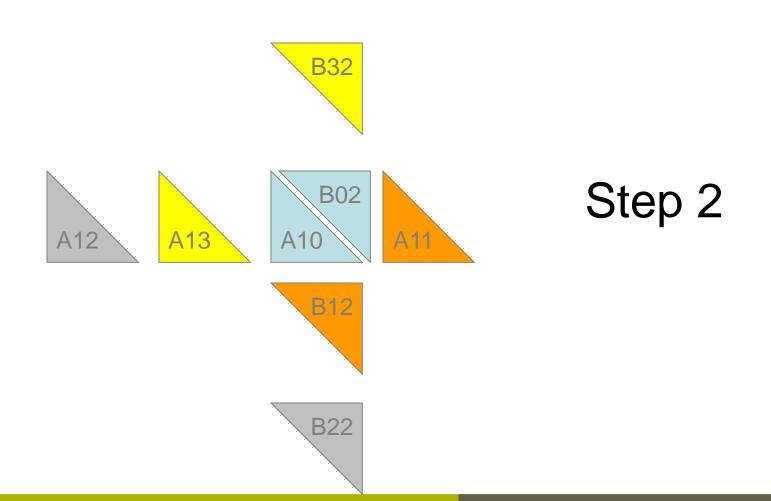
Rearrange Blocks

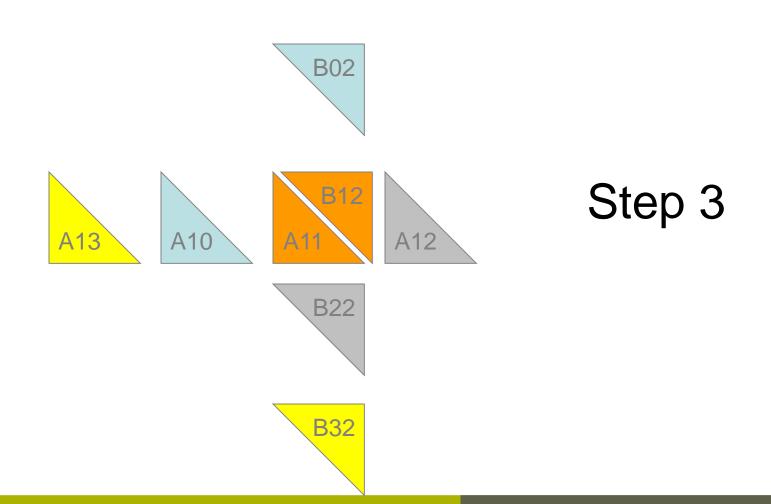


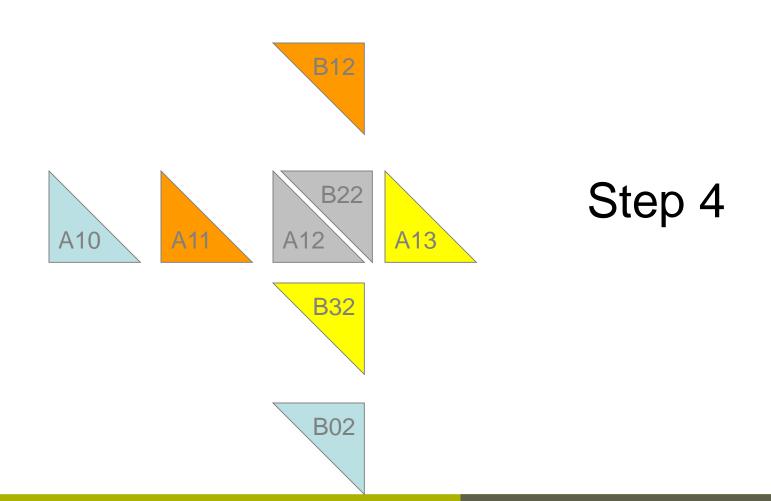
Block A_{ij} cycles left *i* positions

Block B*ij* cycles up *j* positions

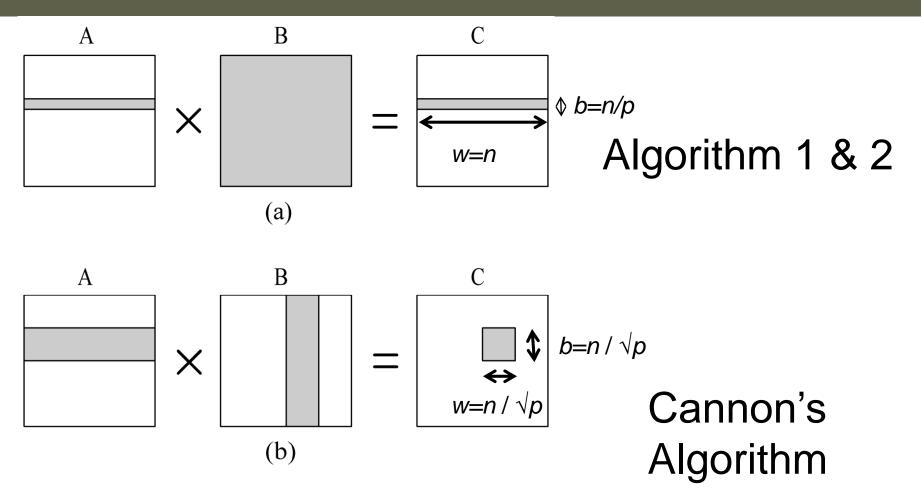








Elements of A and B Needed to Compute the process's portion of C



Why Cannon's requires less communication

- Amount of computations per process:
 n³/p=breadth.width.n/p ~ surface
- Amount of communication per process:
 breadth.n+width.n=(breadth+width).n ~ circumference
 - Version 1 & 2: n^2/p (A-matrix) + n^2 (B-matrix) = (n+n/p).n
 - Cannon: n^2/\sqrt{p} (A-matrix) + n^2/\sqrt{p} (B-matrix) = $(2n/\sqrt{p}).n$
- Granularity = computations / communication
 - = surface / circumference
 - should be minimized
 - for a given surface, circumference is minimal for a

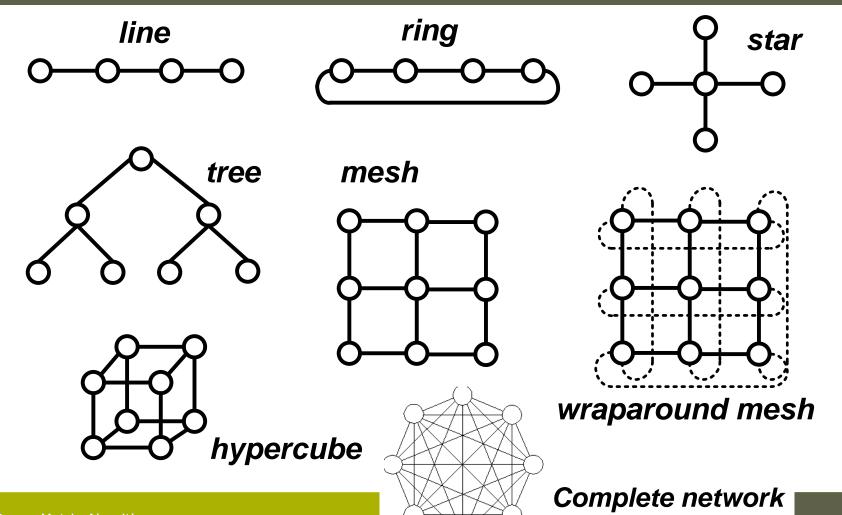
square

=> Cannon (square) is optimal and better than rectangle (v1&2)

Complexity Analysis

- o Algorithm has \sqrt{p} iterations
- During each iteration process multiplies two (n / \sqrt{p}) × (n / \sqrt{p}) matrices: $\Theta(n^3 / p^{3/2})$
- o Computational complexity: $\Theta(n^3/p)$ [the same]
- o During each iteration process sends and receives two blocks of size $(n / \sqrt{p}) \times (n / \sqrt{p})$
- Communication complexity: $\Theta(n^2/\sqrt{p})$ [lower!]
- ⇒ Overhead ratio= $O(n^2 / \sqrt{p}) / O(n^3/p) = O(\sqrt{p/n})$
- → Super-scalable since efficiency drops if we increase both p and n

Efficient Interconnection Networks for Cannon's MxM?



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Memory need of each processor

As a function of **n**, **p** and **b** (the number of bytes per element)

- Sequential algorithm:
- One-step algorithm:
- Alternate algorithm:
- Cannon's algorithm:

Special MPI functions

- MxV version 2: Reduce operation
- Cannon: Shift operation through a SendRecv_replace call
- Submatrix sending:
 - Consist of equally spaced blocks
 - DataType.Vector(int count, int blocklength, int stride, <u>Datatype</u> oldtype)



Shared-memory systems

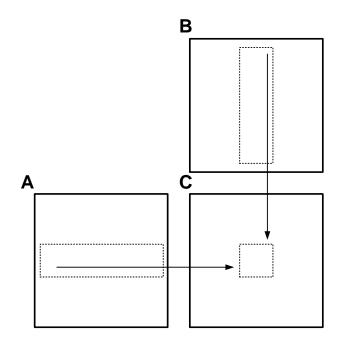
- Computation of each element is independent of the others
 - Can be done by a thread without the need of synchronization (see chapter of shared memory)
- Data is accessible by all threads
- However...

MxM on GPU

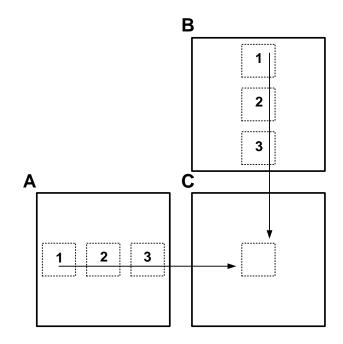
Example: GPUmat toolbox in Matlab

- Initially, matrices are copied to GPU
 - If they are not still in memory from previous matrix operations, keep pointers in CPU to the data on the GPU
- Every thread computes 1 element of C.
- Not enough memory to put all data in shared memory (16K)
- On one multiprocessor, 1 block of threads computes
 1 block of the C matrix
- Iteratively copy A row blocks and B column blocks to shared memory.
- Result: 200x speedup, 50x if compared to quadcore

MxM on GPU



Each multiprocessor computes 1 block

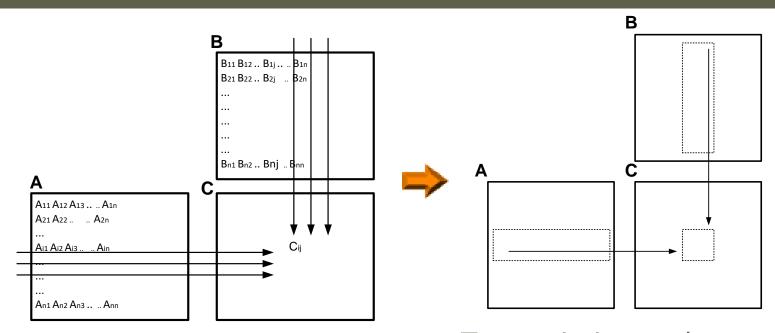


If rows and columns do not fit in local memory: compute blockby-block

Cache efficiency?

- Multicores or GPUs with cache can directly access RAM/global memory
- Data is cached automatically
- However: is the data in cache reused optimally?
- Better is to organize the computations as with the GPU implementation
 - Multicores: each thread calculates the multiplication of 2 submatrices of A and B

Cache efficiency



- Row of A is reused in computing row of C
- Column of B is reused in computing column of C

To maximize cache usage, compute block by block (cf Cannon which minimizes communication)



If A rows and B columns do not fit in cache: alternate over A and B blocks like GPU version

Experimental results

