Parallel algorithms for Dense Matrix Problems

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Logistics

Assignments

- A1 grades released
- ▶ Then Mini-Exam 1 grading have commenced
- ▶ Goal to have all grading done by Mon 10/11
- A2 is delayed: apologies

Reading

Grama Ch 8 on Dense Matrix Algorithms

- Naive Matrix Multiply
- Cannon's Algorithm
- LU Decomposition

Today

- Matrix algorithms
- ► Facebook?

Recall Matrix Transpose

- ightharpoonup Common operation on matrices is a **transpose** notated A^T
- ► Interchanges rows/columns of A: $a_{ij} \rightarrow a_{ji}$
- ▶ Diagonal elements stay the same
- Algorithms that perform operations on A can often be performed on A^T without re-arranging A how? Hint: consider summing rows of A vs summing rows of A^T

Original matrix A

0	5	10	15
20) 25	30	35
40	45	5 50	55

transpose(A)

0	20	40	60
5	25	45	65
10	30	50	70
15	35	55	75

Exercise: Matrix Partitioning Across Processors

Row Partition				
00	01	02	03	
10	11	12	13	
20	21	22	23	
30	31	32	33	

Column Partition				
00	01	02	03	
10	11	12	13	
20	21	22	23	
30	31	32	33	

Block Partition				
01	02	03		
11	12	13		
21	22	23		
31	32	33		
	11 21	11 12 21 22		

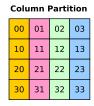
Proc Location		
	P0 / P00	
	P1 / P01	
	P2 / P10	
	P3 / P11	

- Recall several ways to partition matrices across processors
- Diagram shows these
 - Entry ij may be an individual element OR...
 - ► Entry ij may be a **Block**: ex. Block (2,3) is the submatrix from rows 200-299 and cols 300-399
- ► Assume **square** matrices : #rows = #cols
- ▶ Common to multiply to compute product: $C = A \times B$
- ▶ Ideal partitioning for A and B in matrix multiply?
- ▶ Ideal partitioning for $C = A^T \times B$
- $\blacktriangleright \ \ \text{Ideal partitioning for} \ C = A \times B^T$

4

Answers: Matrix Partitioning Across Processors

Row Partition				
00	01	02	03	
10	11	12	13	
20	21	22	23	
30	31	32	33	



Block Partition				
00	01	02	03	
10	11	12	13	
20	21	22	23	
30	31	32	33	

Proc Location		
	P0 / P00	
	P1 / P01	
	P2 / P10	
	P3 / P11	

- $ightharpoonup C = A \times B$
 - lacktriangle Ideally A is row-partitioned, B is column partitioned
 - ► Then block-partitioned C could be computed w/o communication
- $ightharpoonup C = A^T \times B$
 - ightharpoonup Ideally A and B column-partitioned
- $ightharpoonup C = A \times B^T$
 - ightharpoonup Ideally A and B row-partitioned
- ightharpoonup Block-partitioning often used: not ideal for any version but less communication required when both A and A^T will b used

Naive Parallel Dense Multiplication: Overview

Block Partitioning Appears Frequently

- Specific applications may be able to select a favorable partitioning (e.g. Row Partition for Page Rank)
- ▶ Many applications use both A and A^T so employ block-partitioned matrices: middle-way approach which does not favor rows or columns
- Parallel Libraries often use block partitions by default

Matrix Multiply with Blocks

- ➤ To compute Matrix-Matrix multiply, procs must (eventually) multiply full rows by full columns to compute an output block
- Naive method: each Proc stores full rows/columns needed for it to independently compute output block which it stores

Naive Parallel Dense Multiplication: Demo

P01

1. Initial data layout: each Proc holds a block of A. B. and C respectively. Processors are arragned in a logical grid that reflects their initial data.

data is bolded.



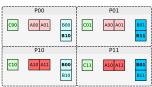
P00

B00

A00

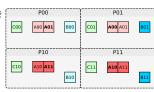
3. Each proc participates in an All-to-All sharing of data for the Column it is

This leaves each row with complete columns as well.

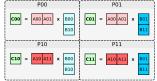


2. Each proc participates in an All-to-All sharing of data for the Row it is in.

This leaves each proc with entire rows of A.



4 Fach Proc now has a unique set of complete Rows and Columns and can independently compute a block of output matrix C through block multiplication.



Exercise: Analysis of Naive Dense Mult.

Assumptions

- ▶ Matrices A and B are size $N \times N$ so N^2 elements
- ▶ P processors in a $\sqrt{P} \times \sqrt{P}$ grid (P is a perfect square)
- ▶ Each Proc has block with N^2/P elements of A,B as a $(N/\sqrt{P}) \times (N/\sqrt{P})$ submatrix
- Simplified communication cost for All-to-All on a Ring with p #procs in ring, t_s comm startup time, t_w per word transfer rate, M message size:

$$t_{comm} = (p-1)(t_s + t_w M)$$

Questions

- 1. What is communication cost of this algorithm?
- 2. How much time does the final block matrix multiply take?
- 3. What is the memory requirement for each proc?
- 4. Biggest disadvantage for this algorithm?

Answers: Analysis of Naive Dense Mult.

- 1. What is communication cost of this algorithm?
 - ▶ #Procs in rows/cols is \sqrt{P} ~ ring size
 - $ightharpoonup M=N^2/P$: message size is num elements on each proc
 - ▶ 2 All-to-All shares : 1 for rows, 1 for cols

$$t_{comm} = 2(\sqrt{P} - 1) \times (t_s + t_w(N^2/P))$$

- 2. What is the memory requirement for each proc? E.g. how many submatrices of A,B are on each proc?
 - ► Full rows/cols on each proc
 - Requires \sqrt{P} submatrices for each Proc
- 3. How much time does the final block matrix multiply take?
 - Each proc has \sqrt{P} submats of A,B to multiply
 - ▶ MatMult is for size s is $O(s^3)$; submat size $s = N/\sqrt{P}$

$$t_{mult} = O((\sqrt{P}) \times ((N/\sqrt{P})^3)) = O(N^3/P)$$

- 4. Biggest disadvantage for this algorithm?
 - ▶ Major: The need to store \sqrt{P} sub matrices on all procs may be prohibitive: $2\sqrt{P} \times N^2/P$ space on each proc
 - Minor: Not much chance to overlap communication / computation in the algorithm

Cannon's Algorithm

- Proposed in Lynn Elliot Carter's 1969 thesis
- Target was very small parallel machines implementing a Kalman Filter algorithm in hardware
- "Communication" happening between small Procs with data in registers
- Scales nicely to large machines and overcomes the large memory requirement of the Naive Mat-Mult Algorithm

A CELLULAR COMPUTER TO IMPLEMENT THE KALMAN FILTER ALCORITHM

by

LYNN ELLIOT CANNON

By the conventional definition of matrix product, if A is multiplied by B, the result, call it C, is given by

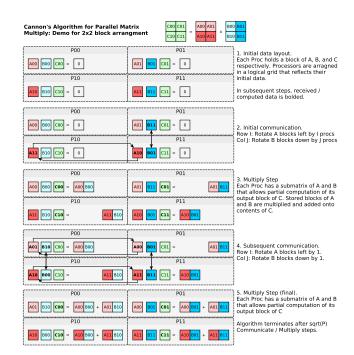
The symmetry of this product can be seen by comparing the ijth element with the jith element and noticing that one is obtained from

-24-

- A. 1. The first row of A is left alone.
 - 2. The second row of A is shifted left one column.
 - The third row of A is shifted left two columns.
 (Note, in general the ith row of A is shifted left i-l columns for i = 1, ..., n).
- B. 1. The first column of B is left alone.
 - 2. The second column of B is shifted up one row
 - The third column of B is shifted up two rows.
 (Note, in general the jth column of B is shifted up j-1 rows for j = 1, ..., n)

Once the registers have been shifted the multiplication pr

Demo



Cannon's Algorithm Pseudocode

```
Cannon_MM(i, j, Q){
  PE(i,j) has blocks A1=A(i,j) and B1=B(i,j)
  Q is the Block Dimension : A is Q*Q blocks
  Allocate space A2, B2, Cij sized as A1
  doboth send A1 to PE(i, j-i+Q % Q)
        recv A2 from PE(i, j+i+Q % Q)
  doboth send B1 to PE(i-j+Q % Q, j)
         recv B2 from PE(i+j+Q % Q, j)
  for(k=1 to 0){
    copy A2 into A1, B2 into B1
    Cij += A1 * B1
    doboth send A1 to PE(i, j-1+Q % Q)
           recv A2 from PE(i, j+1+Q % Q)
    doboth send B1 to PE(i-1+Q % Q, j)
           recv B2 from PE(i+1+Q % Q, j)
    // optionally skip last comm
  Cij now contains output block of C(i,j)
```

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

B _{0,0}	B _{0,1}	B _{0,2}	B _{0,3}
B _{1,0}	B _{1,1}	B _{1,2}	B _{1,3}
B _{2.0}	B _{2,1 &}	ў В _{2,2}	B _{2,3}
B _{3,0}	B _{3,1}	B _{3,2}	B _{3,3}

	1	4	1	4
ş -	A _{0,0} ⋖	A _{0.1} ~	A _{0,2} ~	A _{0,3} ~
	B _{0,0}	B _{1,1}	B _{2,2}	B _{3,3}
	A _{1,1} ~	A _{1,2} ~	A _{1,3} ~	A _{1,0}
	B _{1,0}	B _{2,1}	B _{3,2}	B _{0,3}
	A _{2,2}	A23	A _{2,0}	A _{2,1}

(a) Initial alignment of A

$B_{0,0}$	B _{0,1}	B _{0,2}	B _{0,3}
B _{1,0}	B _{1,1}	B _{1,2}	в _{1,3}
B _{2,0}	В _{2,1 в}	ў В _{2,2}	в _{2,3}
B _{3,0}	В _{3,1}	B _{3,2}	В _{3,3}

(b) In	nitial alig	nment of	В
1	1	1	1
~ A _{0.1} ~	A _{0,2} ~	A _{0.3} ~	A _{0,0} ~
B _{1,0}	B _{2,1}	B _{3,2}	$B_{0,3}$
A _{1,2} ~	A _{1,3} ~	A _{1,0} ~	A _{1,1}
B _{2,0}	B _{3,1}	B _{0,2}	B _{1,3}
- A _{2,3} -	A _{2,0} ~	A _{2,1} ~	A _{2,2}
B _{3,0}	$B_{0,1}$	B _{1,2}	B _{2,3}
A _{3,0} ~	A _{3,1} ~	A _{3,2} <	A _{3,3} <
B _{0,0}	B _{1,1}	B _{2,2}	B _{3,3}
1	1	1	

ſ	A _{0,3}	A _{0,0}	A _{0,1}	A _{0,2}
	B _{3,0}	B _{0,1}	B _{1,2}	B _{2,3}

4	A _{0,2} ~	A _{0,3} <	A _{0.0} ~	A _{0.1} ~
	B _{2,0}	B _{3,1}	B _{0,2}	B _{1,3}
4	A _{1,3} <	A _{1,0} <	A _{1,1} <	A _{1,2} ~
	B _{3,0}	B _{0,1}	B _{1,2}	B _{2,3}
ų.	A _{2,0} ~	A _{2,1} ~	A _{2,2} ~	A _{2,3} ~
	B _{0,0}	$B_{1,1}$	B _{2,2}	B _{3,3}
٠.	A _{3,1} <	A _{3,2} ~	A _{3,3} <	A _{3,0} ~
	B _{1,0}	B _{2,1}	B _{3,2}	B _{0,3}

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

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

Figure 8.3 The communication steps in Cannon's algorithm on 16 processes.

Exercise: Analysis of Cannon's Algorithm

Assumptions

- ▶ Matrices A and B are size $N \times N$ so N^2 elements
- ▶ P processors with block partitioning: initially N^2/P elements of A, B on each proc (assume P is a perfect square)
- Simplified communication cost for send/recv on a Ring:

$$t_{comm} = t_s + t_w M$$

with p #procs in ring, t_s comm startup time, t_w per word transfer rate, M message size.

Questions

- 1. What is communication cost of this Cannon's algorithm?
- 2. Is this any better/worse/same as the Naive algorithm?
- 3. What is the memory requirement for each proc?
- 4. Is this any better/worse/same as the Naive algorithm?

Answers: Analysis of Cannon's Algorithm

- 1. What is communication cost of this Cannon's algorithm?
 - ▶ In each step, each proc performs 2 send/recv ops
 - ▶ Each send/recv is a block of size N^2/P
 - ▶ Block Dim $Q = \sqrt{P}$ for square 2D Torus
 - \blacktriangleright Total \sqrt{P} steps : can skip last comm step

$$t_{comm} = 2(\sqrt{P} - 1) \times (t_s + t_w(N^2/P))$$

- 2. Is this any better/worse/same as the Naive algorithm?
 - Same communication cost as Naive algorithm
- 3. What is the memory requirement for each proc?
 - $ightharpoonup O(N^2/P)$: 5 blocks as stated in pseudcode,
 - ▶ 3 blocks for A_{ij}, B_{ij}, C_{ij}
 - ▶ 2 "workspaces" to allow send/recv of blocks:
 - Eliminate workspace blocks in a refinement
- 4. Is this any better/worse/same as the Naive algorithm?
 - ► Cannon's $O(N^2/P)$ vs Naive $O(\sqrt{P} \times N^2/P)$
 - Memory overhead is much better: constant number of blocks rather than the need to store entire rows/cols on single procs

Lessons from Cannon's Algorithm

- Illustrates "pipelining": blocks used to compute partial results then fed forward other processors
- Benefits greatly from a 2D Grid / Torus network which facilitates local communications that arise in the algorithm
- lacktriangle While not as ideal as row/col partitioning for A,B, realistic and relatively efficient
- Variants of central idea exist in some libraries such as Scalapack which has a parallel xGEMM() using many similar ideas
- Could really use some code support for
 - 2D Coordinates for processors rather than linear rank...
 - Sending/receiving in a ring...

MPI Tricks for Rings

Sendrecv in a Ring

MPI_Sendrecv() allows ring-link partnering

Sendrecv with Replacement

MPI_Sendrecv_replace() allows send/recv in the same buffer

In Cannon's Alg, no longer need A1 / A2: can send/receive block of A with a single buffer.

MPI Tricks for Grids: MPI_Cart_create()

MPI has special support for Grid/Torus network configs; allows creation of a MPI_Comm that maps processors to a N-D grid

▶ 2D Torus for Cannon's Alg

```
// cartesian comm.c
int dim_len = 2; // Set up the Cartesian topology
int dims[2] = {sqrt(npes), sqrt(npes)}; // # rows/cols
int periods[2] = {1, 1};  // wrap-around rows/cols
// Create the Cartesian topology, with rank reordering
MPI Comm comm 2d;
MPI_Cart_create(MPI_COMM_WORLD, // original comm
               dim_len, dims, periods, // cartesian comm props

    // re-order linear rank if beneficial

               &comm_2d); // new communicator with 2D coords
// Get the rank and coordinates with respect to the new topology
int my2drank = -1; // may be differ from world rank
MPI Comm_rank(comm_2d, &my2drank);
int mycoords[2] = \{-1, -1\}; // (i,j) coords
MPI_Cart_coords(comm_2d, my2drank, 2, mycoords);
printf("Proc %2d (%s): my2drank %3d mycoords (%3d, %3d)\n",
       myrank,processor_name,
       my2drank,mycoords[0],mycoords[1]);
```

MPI Tricks for Shifting

Shifts are eased by the MPI_Cart_shift() function

- Calculates linear rank of source/dest procs for shift operations in a Cartesian grid of procs.
- Data exchange via MPI_Sendrecv() is then direct

Cannon's Algorithm in MPI

- Grama Program 6.2 is Cannon's Matrix Multiply algorithm implemented via MPI
- Uses the tricks mentioned on the past 2 slides to ease implementation burden
- See cannon_grama.c for a source code version of it

Note: I haven't tested this code but everything from textbooks always works out the box, right?

Linear Equations

$$a_{11}x_1 + a_{12}x_2 + \dots + a_{1n}x_n = b_1$$

$$a_{21}x_1 + a_{22}x_2 + \dots + a_{2n}x_n = b_2$$

$$\vdots$$

$$a_{m1}x_1 + a_{m2}x_2 + \dots + a_{mn}x_n = b_m$$

Summarized in matrix form as

$$A\mathbf{x} = \mathbf{b}$$

$$A = \begin{bmatrix} a_{11} & a_{12} & \cdots & a_{1n} \\ a_{21} & a_{22} & \cdots & a_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{m1} & a_{m2} & \cdots & a_{mn} \end{bmatrix}, \quad \mathbf{x} = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix}, \quad \mathbf{b} = \begin{bmatrix} b_1 \\ b_2 \\ \vdots \\ b_m \end{bmatrix}$$

Usually given A, b, must find x. An inordinate amount of CPU cycles are spent on this problem.

Solving Triangular Systems

Easier than a general system via back substitution process

```
2 3 0 | 14 ]
                                [0560|28]
[0567|56]
[0 0 8 9 | 60]
                                [0 0 8 0 | 24]
                                                      60-9*4
                                [0 \ 0 \ 0 \ 1 \ | \ 4] \ x(3) == 4
[0 \ 0 \ 0 \ 1 \ | \ 4] \ x(3) == 4
    2 3 0 | 14 ]
                                           0 | 5 ]
                                                     14-3*3
[0560|28]
                                 [050010]
                                                      28-6*3
   0 \ 1 \ 0 \ | \ 3 \ ] \ x(2) == 3
                                [0 \ 0 \ 1 \ 0 \ | \ 3] \ x(2) == 3
                                 [0\ 0\ 0\ 1\ |\ 4\ ]\ x(3) == 4
[0 \ 0 \ 0 \ 1 \ | \ 4 \ ] \ x(3) == 4
   2 0 0 | 5 ]
[0 \ 1 \ 0 \ 0 \ | \ 2] \ x(1) == 2
                                 [ 0 1 0 0 | 2 ] x(1) == 2
   0 \ 1 \ 0 \ | \ 3 \ ] \ x(2) == 3
                                 [0\ 0\ 1\ 0\ |\ 3\ ]\ x(2) == 3
[0 \ 0 \ 0 \ 1 \ | \ 4] \ x(3) == 4
                                 [ 0 0 0 1 I
                                                 4] x(3) == 4
         0 \mid 1 \mid x(0) == 1
[0 \ 1 \ 0 \ 0 \ | \ 2] \ x(1) == 2
[0 \ 0 \ 1 \ 0 \ ] \ x(2) == 3
[0 \ 0 \ 0 \ 1 \ | \ 4 \ ] \ x(3) == 4
```

Standard Code for Back Substitution

```
BACK SUBSTITUTE(A,b,x){
  N = nrows(A)
  for(j=N-1 downto 0) {
    x[j] = b[j] / A[j,j]
    for[ i=j-1 downto 0 ] {
      b[i] = b[i] - A[i,j] *x[j]
      A[i,j] = 0 // OPTIONAL
  x \( \) now contains solutions
  b∏ has been modified
  A[] has been modified if OPTIONAL is executed
```

Computational complexity for square matrix of size N?

Getting a Triangular Matrix via Gaussian Elimination

- ▶ Standard solution algorithm to find x in Ax = b
- ightharpoonup Converts A to U which is upper triangular

```
[ 1 2 3 -4 | -14 ]
[ 2 7 21 10 | 38 ] -2 row0
[ 4 13 43 2 | 24 ] -4 row0
[ -2 -2 7 15 | 60 ] +2 row0
 1 2 3 -4 | -14 ]
0 3 15 18 | 66 ]
0 5 31 18 | 80 ]
0 2 13 7 1
                                                    -5/3 row1
                                                  -2/3 row1
   1 2 3 -4 | -14 ]
0 3 15 18 | 66 ]
0 0 6 -12 | -30 ]
0 0 3 -5 | -12 ]
                                                        -1/2 row3
   1 2 3 -4 | -14 ] [ 1 0 0
0 3 15 18 | 66 ] [ 2 1 0
0 0 6 -12 | -30 ] [ 4 5/3 1
0 0 0 1 | 3 ] [ -2 2/3 1/2
                                                                                           L is formed from negative
                                                                                           coefficients found via
                                                                                0] Gaussian elimination with
                                                                                           unit main diagonal.
                                          b'
```

LU: The Lower Upper Decomposition

- ightharpoonup By tracking the coefficients used during the Gaussian elimination, one gets a matrix L which is lower triangular
- lacktriangle Modifications to A become an upper triangular matrix U
- ▶ One can verify that A = LU

octave> L =	rats	s(L)			octave> U U =
1		0	0	0	1 2 3 -4
2		1	0	0	0 3 15 18
4	5,	/3	1	0	0 0 6 -12
-2	2/	/3	1/2	1	0 0 0 1
octave>	• L *	U			octave> L * U - A
ans =					ans =
1	2	3	-4		0 0 0 0
2	7	21	10		0 0 0 0
4	13	43	2		0 0 0 0
-2	-2	7	15		0 0 0 0

Exercise: LU Factorization Pseudocode

```
LU_FACTORS(A[] : square matrix){
      N = nrows(A)
      Allocate L as N*N identity mat
      Allocate U as copy of A
      for (d=0 \text{ to } N-1) {
                                             // leading row d
        for(i=d+1 to N-1){
                                             // remaining rows i
          scale = U[i,d] / U[d,d]
                                             // scale for this row
          L[i.d] = scale
                                             // record scale in L
          for(j=d to N-1){
                                            // iterate over this row j
10
            U[i,j] = U[i,j] - scale*U[d,j] // subtract off scaled leading row
11
12
1.3
14
      7
15
      return
16
        L: a lower triangle matrix with factors and unit diagonal
17
        U: an upper triangle matrix, obeys L*U = A
18
```

- Computational Complexity?
- Could anything go sideways numerically?

Answers: LU Factorization Pseudocode

```
LU_FACTORS(A[] : square matrix){
      N = nrows(A)
      Allocate L as N*N identity mat
      Allocate U as copy of A
6
      for (d=0 \text{ to } N-1) {
                                             // leading row d
        for(i=d+1 to N-1){
                                            // remaining rows i
          scale = U[i,d] / U[d,d]
                                            // scale for this row
          L[i,d] = scale
                                            // record scale in L
10
          for(j=r to N-1){
                                       // iterate over this row j
            U[i,j] = U[i,j] - scale*U[d,j] // subtract off scaled leading row
11
12
13
14
15
      return
        L: a lower triangle matrix with factors and unit diagonal
16
        U: an upper triangle matrix, obeys L*U = A
17
18 }
```

- ▶ Computational Complexity?: $O(N^3)$ 3 nested loops
- Could anything go sideways numerically? Division by 0 at line 8
 - To fix this requires pivoting
 - ► Robust versions permute rows so the row with the largest U[:,d] element used at iteration d

Utility of LU Decomposition

General Process

- 1. Want x in Ax = b
- 2. Compute LU = A via Gaussian elimination
- 3. Use forward-substitution to find y in Ly = b
- 4. Use back-substitution to find x in Ux = y

Solving in this fashion exploits the following identities

$$Ly = b = Ax$$
 so $Ax = Ly$
 $LU = A$ so $LUx = Ly$

$$L^{-1}LUx = L^{-1}Ly \quad \text{so} \qquad Ux = y$$

vs Gaussian Elimination

- ► LU factorization costs little more than Gaussian Elim
- Saving the LU Factorization allows solving for a new b with only passes of back/forward substitution

$$Ax_1 = b_1, Ax_2 = b_2, Ax_3 = b_3$$

1 LU decomposition, 3 rounds of back/forward substitution

- ▶ LU Decomp is $O(N^3)$
- ▶ Back/Forward Sub is (N^2)

Variants

- ► To save space, overwrite L,U in A
 - Upper triangle of A becomes U including main diagonal
 - ► Lower triangle of A would have been 0's, store L there, implied 1 diagonal
- lackbox Grama's variant makes main diagonal of U all 1's: saves some ops in back/forward substitution
- We are ignoring the need to **pivot** and permute the matrix rows for numerical stability: doing so yields the LUP decomposition with permutation matrix P

Exercise: Now, about Parallelizing...

```
LU_FACTORS_INPLACE(A[] : square matrix){
      N = nrows(A)
      // Will overwrite A with its L*U factors, no allocation of L or U
 5
      for (d=0 \text{ to } N-1)
                                             // leading row d
        for(i=d+1 to N-1){
                                            // remaining rows i
          scale = A[i,d] / A[d,d]
                                             // scale for this row
          A[i,d] = scale
                                            // record scale in L
          for(j=r to N-1){
                                       // iterate over this row j
            A[i,j] = A[i,j] - scale * A[d,j] // subtract off scaled leading row
10
11
12
13
      }
14
      return; // A now has its L,U factors in its lower/upper triangles
15
```

Assuming an in-place variant how would one go about parallelizing this?

- Decomposition / distribution of A?
- Communication at which steps?

Pitch some ideas

Answers: Now, about Parallelizing...

- Block decomposition means that some processors idle
- ► Row decomposition also leads to some idling, is described in Grama 8.3
- A cyclic decomposition leads to better balance
 - ► 100 x 100 matrix, 4 Procs, row cyclic
 - P0: rows 4*i+0 = 0,4,8,12,...
 - P1: rows 4*i+1 = 1,5,9,13,...
 - etc.
- Broadcast leading row from owning proc to all others

P_0	1	(0,1)	(0,2)	(0,3)	(0,4)	(0,5)	(0,6)	(0,7)
P_1	0	1	(1,2)	(1,3)	(1,4)	(1,5)	(1,6)	(1,7)
P_2	0	0	1	(2,3)	(2,4)	(2,5)	(2,6)	(2,7)
P ₃	0	0	0	(3,3)	(3,4)	(3,5)	(3,6)	(3,7)
P_4	0	0	0	(4,3)	(4,4)	(4,5)	(4,6)	(4,7)
P ₅	0	0	0	(5,3)	(5,4)	(5,5)	(5,6)	(5,7)
P ₆	0	0	0	(6,3)	(6,4)	(6,5)	(6,6)	(6,7)
P_7	0	0	0	(7,3)	(7,4)	(7,5)	(7,6)	(7,7)

P_0	1	(0,1)	(0,2)	(0,3) (0,4) (0,5) (0,6) (0,7)
P	0	1	(1,2)	(1,3) (1,4) (1,5) (1,6) (1,7)
P ₂	0	0	1	(2,3) (2,4) (2,5) (2,6) (2,7)
P ₃	0	0	0	1 (3,4) (3,5) (3,6) (3,7)
P ₄	0	0	0	(4,3)¥(4,4)¥(4,5)¥(4,6)¥(4,7)
P ₅	0	0	0	(5,3)¥(5,4)¥(5,5)¥(5,6)¥(5,7)
P ₆	0	0	0	(6,3)¥(6,4)¥(6,5)¥(6,6)¥(6,7)
P ₇	0	0	0	(7,3)\$\dagge(7,4)\$\dagge(7,5)\$\dagge(7,6)\$\dagge(7,7)\$
	_			

P_0	1	(0,1)	(0,2)	(0,3)	(0,4)	(0,5)	(0,6)	(0,7)
P ₁	0	1	(1,2)	(1,3)	(1,4)	(1,5)	(1,6)	(1,7)
P ₂	0	0	1	(2,3)	(2,4)	(2,5)	(2,6)	(2,7)
P ₃	0	0	0	1	(3,4)	(3,5)	(3,6)	(3,7)
P_4	0	0	0	(4,3)	(4,4)	(4,5)	(4,6)	(4,7)
P ₅	0	0	0	(5,3)	(5,4)	(5,5)	(5,6)	(5,7)
P ₆	0	0	0	(6,3)	(6,4)	(6,5)	(6,6)	(6,7)
P_7	0	0	0	(7,3)	(7,4)	(7,5)	(7,6)	(7,7)

- (a) Computation:
 - (i) A[k,j] := A[k,j]/A[k,k] for $k \le j \le j$
 - (ii) A[k,k] := 1

(b) Communication:

One-to-all broadcast of row A[k,*]

(c) Computation:

- (i) A[i,j] := A[i,j] A[i,k] × A[k,j]
 for k ≤ i ≤ n and k ≤ i ≤ n
- (ii) A[i,k] := 0 for $k \le i \le n$

Figure 8.6 Gaussian elimination steps during the iteration corresponding to k = 3 for an 8 \times 8 matrix partitioned rowwise among eight processes.

Analysis of LU Decomposition

- ightharpoonup Serial algorithm runs in $O(N^3)$
- Parallel approaches use
 - 1. N iterations of each row as the leading row
 - 2. Broadcast of leading row d to all P procs : N broadcasts
 - 3. Parallel modification of N-d lower block of A[] to store L,U factors in it $O(N^2/P)$
- lacktriangle For a ring of P procs to broadcast length N row

$$t_{broadcast} = (\log_2 P) \times t_s + t_w \times N \times P$$

leading to overall complexity of

$$T = N \times (N^2/P + (\log_2 P) \times t_s + t_w \times N \times P)$$

= $N^3/P + N \times \log_2 P \times t_s + N \times t_w \times N \times P$
= $O(N^3/P)$

- Main overhead is the need to broadcast a row at each step,
- Pipelined Broadcast improves on this: good implementations of MPI_Bcast() has a node pass on messages, begin computation again ASAP, not idle while broadcast completes