# Tradeoffs between synchronization, communication, and work in parallel schedules

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### Outline

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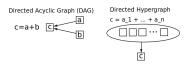
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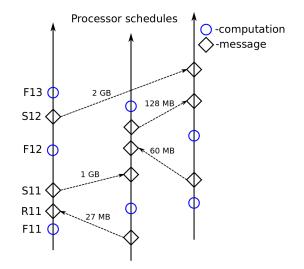
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  - $M_{ij}$  is the set of values sent by processor *i* at timestep *j*

### Parallel schedule example



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 The values ⋃<sub>j</sub> F<sub>ij</sub> = C<sub>i</sub> ⊂ V correspond to the vertices of dependency graph G computed by processor i

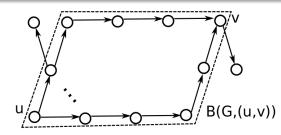
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- All dependencies must be satisfied by the schedule
- Dependent values must be communicated or computed previously
- For all non-local dependency paths in *G*, there must exist a sequence of messages in the schedule

#### Definition (Dependency bubble)

Given two vertices u, v in a directed acyclic graph G = (V, E), the dependency bubble B(G, (u, v)) is the union of all paths in G from u to v.



#### Definition (( $\epsilon, \sigma$ )-path-expander)

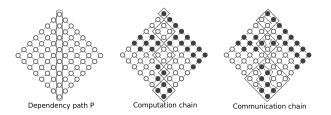
Graph G = (V, E) is a  $(\epsilon, \sigma)$ -**path-expander** if there exists a path  $(u_1, \ldots u_n) \subset V$ , such that the dependency bubble  $B(G, (u_i, u_{i+b}))$  has size  $\Omega(\sigma(b))$  and a minimum cut of size  $\Omega(\epsilon(b))$ .

#### Theorem (Path-expander communication lower bound)

Any parallel schedule of an algorithm, with a  $(\epsilon, \sigma)$ -path-expander dependency graph G = (V, E) about a path of length n incurs the computation (F), bandwidth (W), and latency (S) costs

 $F = \Omega \left( \sigma(b) \cdot n/b \right), \quad W = \Omega \left( \epsilon(b) \cdot n/b \right), \quad S = \Omega \left( n/b \right).$ 

## An example $(b, b^2)$ -path-expander



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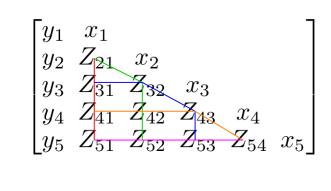
For lower triangular dense L, solve

$$\mathbf{L} \cdot \mathbf{x} = \mathbf{y},$$

i.e., 
$$\sum_{j=1}^{i} L_{ij} \cdot x_j = y_i$$
, for  $i \in \{1, ..., n\}$ .

 $\begin{aligned} \mathbf{x} &= \mathrm{TRSV}(\mathbf{L}, \mathbf{y}, n) \\ 1 \quad & \text{for } i = 1 \text{ to } n \\ 2 & \text{for } j = 1 \text{ to } i - 1 \\ 3 & Z_{ij} = L_{ij} \cdot x_j \\ 4 & x_i = \left(y_i - \sum_{j=1}^{i-1} Z_{ij}\right) / L_{ii} \end{aligned}$ 

### Dependency Hypergraph: Triangular solve



#### Theorem

Any parallelization of any dependency graph  $G_{\rm TRSV}(n)$  where two processors compute  $\lfloor n^2/p \rfloor$  elements of **Z** must incur a communication cost of

$$W_{\rm TRSV} = \Omega\left(n/\sqrt{p}\right).$$

#### Proof.

Proof by application of lower bound on 2D lattice Hypergraph cut.

#### Theorem

Any parallelization of any dependency graph  $G_{\text{TRSV}}(n)$  incurs the following computation (F), bandwidth (W), and latency (S) costs, for some  $b \in [1, n]$ ,

 $F_{\mathrm{TRSV}} = \Omega(n \cdot b), \qquad W_{\mathrm{TRSV}} = \Omega(n), \qquad S_{\mathrm{TRSV}} = \Omega(n/b),$ 

and furthermore,  $F_{\mathrm{TRSV}} \cdot S_{\mathrm{TRSV}} = \Omega\left(n^2\right)$ .

#### Proof.

Proof by application of path-based tradeoffs since  $G_{\text{TRSV}}(n)$  is a  $(b, b^2)$ -path-expander dependency graph.

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Diamond DAG lower bounds were also given by

- Papadimitriou and Ullman [P.U. 1987]
- Tiskin [T. 1998]

Efficient algorithms for TRSV attain above lower bounds

- wavefront algorithms (Heath 1988)
- also algorithms given by [P.U 1987] and [T. 1998]

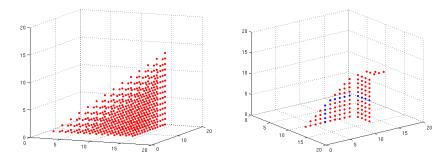
The Cholesky factorization of a symmetric positive definite matrix  ${\boldsymbol{\mathsf{A}}}$  is

 $\mathbf{A} = \mathbf{L} \cdot \mathbf{L}^{\mathcal{T}},$ 

for a lower-triangular matrix  $\boldsymbol{\mathsf{L}}.$ 

$$L = CHOLESKY(\mathbf{A}, n)$$
1 for  $j = 1$  to  $n$ 
2  $L_{jj} = \sqrt{A_{ij} - \sum_{k=1}^{j-1} L_{jk} \cdot L_{jk}}$ 
3 for  $i = j + 1$  to  $n$ 
4 for  $k = 1$  to  $j - 1$ 
5  $Z_{ijk} = L_{ik} \cdot L_{jk}$ 
6  $L_{ij} = (A_{ij} - \sum_{k=1}^{j-1} Z_{ijk})/L_{jj}$ 

### Cholesky dependency hypergraph



These diagrams show (a) the vertices  $Z_{ijk}$  in  $V_{GE}$  with n = 16 and (b) the hyperplane  $x_{12}$  and hyperedge  $e_{12,6}$  on  $H_{GE}$ .

#### Theorem

Any p-processor parallelization of the dependency graph  $G_{\rm GE}(n)$  must incur a communication of

$$W_{\rm GE} = \Omega\left(n^2/p^{2/3}
ight).$$

#### Proof.

Employs 3D lattice hypergraph cut lower bound and assumes some work balance.  $\hfill\square$ 

#### Theorem

Any parallelization of any dependency graph  $G_{GE}(n)$  incurs the following computation (F), bandwidth (W), and latency (S) costs, for some  $b \in [1, n]$ ,

$$\mathcal{F}_{\mathrm{GE}} = \Omega\left(n\cdot b^2
ight), \qquad \mathcal{W}_{\mathrm{GE}} = \Omega\left(n\cdot b
ight), \qquad \mathcal{S}_{\mathrm{GE}} = \Omega\left(n/b
ight),$$

and furthermore, 
$$F_{ ext{GE}} \cdot S_{ ext{GE}}^2 = \Omega\left(n^3
ight), \quad W_{ ext{GE}} \cdot S_{ ext{GE}} = \Omega\left(n^2
ight).$$

#### Proof.

Proof by showing that  $G_{GE}(n)$  is a  $(b^2, b^3)$ -path-expander about the path corresponding to the calculation of the diagonal elements of **L**.

The lower bounds are attainable for Cholesky and similar costs are achievable for QR and the symmetric eigenproblem

- Tiskin's non-pivoted recursive LU and pairwise-pivoted BSP algorithms
- 2.5D LU algorithm
- $W_{
  m GE}=n^2/\sqrt{cp}$  bandwidth cost  $S_{
  m GE}=\sqrt{cp}$  synchronization cost

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We consider the s-step Krylov subspace basis computation

$$\mathbf{x}^{(l)} = \mathbf{A} \cdot \mathbf{x}^{(l-1)},$$

for  $l \in \{1, ..., s\}$  where the graph of the symmetric sparse matrix **A** is a  $(2m+1)^d$ -point stencil.

#### Theorem

Any parallel execution of an s-step Krylov subspace basis computation for a  $(2m + 1)^d$ -point stencil, requires the following computational, bandwidth, and latency costs for some  $b \in \{1, \ldots s\}$ ,

$$F_{\mathrm{Kr}} = \Omega\left(m^d \cdot b^d \cdot s\right), W_{\mathrm{Kr}} = \Omega\left(m^d \cdot b^{d-1} \cdot s\right), S_{\mathrm{Kr}} = \Omega\left(s/b\right).$$

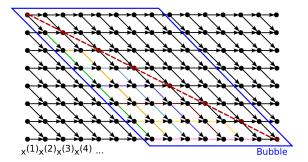
and furthermore,

$$F_{\mathrm{Kr}} \cdot S^d_{\mathrm{Kr}} = \Omega\left(m^d \cdot s^{d+1}\right), \quad W_{\mathrm{Kr}} \cdot S^{d-1}_{\mathrm{Kr}} = \Omega\left(m^d \cdot s^d\right).$$

#### Proof of tradeoffs for Krylov subspace methods

#### Proof.

Done by showing that the dependency graph of a *s*-step  $(2m+1)^d$ -point stencil is a  $(m^d b^d, m^d b^{d+1})$ -path-expander.



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The lower bounds may be attained via communication-avoiding *s*-step algorithms (PA1 in Demmel, Hoemmen, Mohiyuddin, and Yelick 2007)

$$F_{\mathrm{Kr}} = O\left(m^d \cdot b^d \cdot s\right), W_{\mathrm{Kr}} = O\left(m^d \cdot b^{d-1} \cdot s\right), S_{\mathrm{Kr}} = O\left(s/b\right),$$

under the assumption  $n/p^{1/d} = O(bm)$ .

### All-pairs shortest-paths problem

Given a weighted graph G = (V, E) with *n* vertices and a corresponding adjacency matrix **A**, we seek to find the shortest paths between all pairs of vertices in *G* 

- seek the closure, A\*, of A over the tropical semiring
  - $c = c \oplus a \otimes b$  on the tropical semiring implies  $c = \min(c, a + b)$
  - the identity matrix I on the tropical semiring is 0 on the diagonal and  $\infty$  everywhere else

$$\mathbf{A}^* = \mathbf{I} \oplus \mathbf{A} \oplus \mathbf{A}^2 \oplus \ldots \oplus \mathbf{A}^n = (\mathbf{I} \oplus \mathbf{A})^n$$

• numerical computation on the sum-product semiring can be computed by Gauss-Jordan Elimination

$$\mathbf{A}^* = (\mathbf{I} - \mathbf{A})^{-1}$$

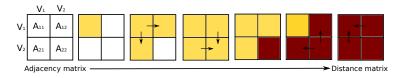
 on the tropical semiring it is commonly computed by the Floyd-Warshall algorithm Compute shortest paths between each pair of vertices using intermediate nodes  $\{1, 2, \dots k\}$ ,

```
D = FLOYD-WARSHALL(\mathbf{A}, n)
\mathbf{D} = \mathbf{A}
for k = 1 to n
for i = 1 to n
for j = 1 to n
d_{ij} = \min(d_{ij}, d_{ik} + d_{kj})
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### Gauss-Jordan elimination (Floyd Warshall algorithm)

$$\mathbf{A} = \begin{bmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{bmatrix} \rightarrow \begin{bmatrix} A_{11}^* & A_{12} \\ A_{21} & A_{22} \end{bmatrix} \rightarrow \begin{bmatrix} A_{11}^* & A_{11}^* A_{12} \\ A_{21}A_{11}^* & A_{22} \end{bmatrix}$$
$$\rightarrow \begin{bmatrix} A_{11}^* & A_{11}^* A_{12} \\ A_{21}A_{11}^* & A_{22} \oplus A_{21}A_{11}^* A_{12} \end{bmatrix} = \mathbf{B}$$
$$\begin{bmatrix} B_{11} & B_{12} \\ B_{21} & B_{22} \end{bmatrix} \rightarrow \begin{bmatrix} B_{11} \oplus B_{12}B_{22}^* B_{21} & B_{12}B_{22}^* \\ B_{22}^* B_{21} & B_{22}^* \end{bmatrix} = \mathbf{A}^*$$



The floating point cost of Gauss-Jordan elimination is  $F = \Theta(n^3/p)$ . Our lower bounds may be applied since the computation has the same structure as Gaussian Elimination, so

$$F \cdot S^2 = \Omega(n^3), \quad W \cdot S = \Omega(n^2).$$

These costs are achieved for  $W = O(n^2/p^{2/3})$  by schedules in

- Aggarwal, Chandra, and Snir 1990
- Tiskin 2007
- Solomonik, Buluc, and Demmel 2012

We can compute the tropical semiring closure

$$\mathbf{A}^* = \mathbf{I} \oplus \mathbf{A} \oplus \mathbf{A}^2 \oplus \ldots \oplus \mathbf{A}^n = (\mathbf{I} \oplus \mathbf{A})^n,$$

directly via repeated squaring (path-doubling)

$$(\mathbf{I}\oplus\mathbf{A})^{2k}=(\mathbf{I}\oplus\mathbf{A})^k\otimes(\mathbf{I}\oplus\mathbf{A})^k$$

with a total of log(n) matrix-matrix multiplications, with

$$F = O(n^3 \log(n)/p)$$

operations and  $O(\log(n))$  synchronizations, which can be less than the  $O(p^{1/2})$  required by Floyd-Warshall.

# Tiskin's path doubling algorithm

Tiskin gives a way to do path-doubling in  $F = O(n^3/p)$  operations. We can partition each  $\mathbf{A}^k$  by path size (number of edges)

$$\mathbf{A}^k = \mathbf{I} \oplus \mathbf{A}^k(1) \oplus \mathbf{A}^k(2) \oplus \ldots \oplus \mathbf{A}^k(k)$$

where each  $\mathbf{A}^{k}(I)$  contains the shortest paths of up to  $k \ge I$  edges, which have exactly I edges. We can see that

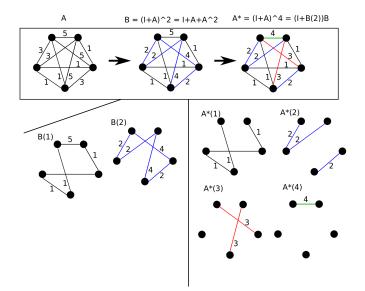
$$\mathbf{A}^{l}(l) \leq \mathbf{A}^{l+1}(l) \leq \ldots \leq \mathbf{A}^{n}(l) = \mathbf{A}^{*}(l),$$

in particular  $\mathbf{A}^*(I)$  corresponds to a sparse subset of  $\mathbf{A}^I(I)$ . The algorithm works by picking  $I \in [k/2, k]$  and computing

$$(\mathbf{I} \oplus \mathbf{A})^{3k/2} \leq (\mathbf{I} \oplus \mathbf{A}^k(I)) \otimes \mathbf{A}^k,$$

which finds all paths of size up to 3k/2 by taking all paths of size exactly  $l \ge k/2$  followed by all paths of size up to k.

## Path-doubling (Tiskin's algorithm)



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Earlier caveat:

$$(\mathbf{I} \oplus \mathbf{A})^{3k/2} \leq (\mathbf{I} \oplus \mathbf{A}^k(I)) \otimes \mathbf{A}^k,$$

does not hold in general. The fundamental property used by the algorithm is really

$$\mathbf{A}^*(l)\otimes \mathbf{A}^*(k)=\mathbf{A}^*(l+k).$$

All shortest paths of up to any length are composable (factorizable), but not paths up to a limited length. However, the algorithm is correct because  $\mathbf{A}^{l} \leq \mathbf{A}^{k}(l) \leq \mathbf{A}^{*}(k)$ .

Since the decomposition by path size is disjoint, one can pick  $\mathbf{A}^k(I)$  for  $I \in [k/2, k]$  to have size

$$|\mathbf{A}^k(I)| \geq 2n^2/k.$$

Each round of path doubling becomes cheaper than the previous, so the cost is dominated by the first matrix multiplication,

$$F = O(n^3/p)$$
  $W = O(n^2/p^{2/3})$   $S = O(\log(n)),$ 

solving the APSP problem with no  $F \cdot S^2$  or  $W \cdot S$  tradeoff and optimal flops.

Tiskin gives a way to lower the synchronization from  $S = O(\log(n))$  to  $O(\log(p))$ . For nonnegative edge lengths it is straightforward

- compute A<sup>p</sup> via path-doubling
- pick a small  $\mathbf{A}^p(I)$  for  $I \in [p/2, p]$
- replicate A<sup>p</sup>(l) and compute Dijkstra's algorithm for n/p nodes with each process, obtaining (A<sup>p</sup>(l))\*
- compute by matrix multiplication

$$\mathbf{A}^* = (\mathbf{A}^p(I))^* \otimes \mathbf{A}^p$$

since all shortest paths are composed of a path of size that is a multiple of  $l \le p$ , followed by a shortest path of size up to p

- obtained synchronization cost lower bound for any parallel schedule of Gaussian elimination
- same technique yields cost tradeoffs for Krylov subspace methods
- on the tropical semiring these are shortest-path graph algorithms, Floyd-Warshall and Bellman-Ford
- it is possible to use a different algorithm to circumvent the tradeoffs for the all-pairs shortest-paths problem
- Open question: can one circumvent the tradeoffs in an algorithm that obtain the closure of a numerical matrix?