# CS 5220: Locality and parallelism in simulations II 

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## Basic styles of simulation

- Discrete event systems (continuous or discrete time)
- Game of life, logic-level circuit simulation
- Network simulation
- Particle systems
- Billiards, electrons, galaxies, ...
- Ants, cars, ...?
- Lumped parameter models (ODEs)
- Circuits (SPICE), structures, chemical kinetics
- Distributed parameter models (PDEs / integral equations)
- Heat, elasticity, electrostatics, ...

Often more than one type of simulation appropriate. Sometimes more than one at a time!

## Common ideas / issues

- Load balancing
- Imbalance may be from lack of parallelism, poor distribution
- Can be static or dynamic
- Locality
- Want big blocks with low surface-to-volume ratio
- Minimizes communication / computation ratio
- Can generalize ideas to graph setting
- Tensions and tradeoffs
- Irregular spatial decompositions for load balance at the cost of complexity, maybe extra communication
- Particle-mesh methods - can't manage moving particles and fixed meshes simultaneously without communicating


## Lumped parameter simulations

Examples include:

- SPICE-level circuit simulation
- nodal voltages vs. voltage distributions
- Structural simulation
- beam end displacements vs. continuum field
- Chemical concentrations in stirred tank reactor
- concentrations in tank vs. spatially varying concentrations

Typically involves ordinary differential equations (ODEs), or with constraints (differential-algebraic equations, or DAEs).

Often (not always) sparse.

## Sparsity

$$
1-2-3-4-5
$$

Matrix
Graph

Consider system of ODEs $x^{\prime}=f(x)$ (special case: $f(x)=A x$ )

- Dependency graph has edge $(i, j)$ if $f_{j}$ depends on $x_{i}$
- Sparsity means each $f_{j}$ depends on only a few $x_{i}$
- Often arises from physical or logical locality
- Corresponds to A being a sparse matrix (mostly zeros)


## Sparsity and partitioning



Matrix


Graph

Want to partition sparse graphs so that

- Subgraphs are same size (load balance)
- Cut size is minimal (minimize communication)

We'll talk more about this later.

## Types of analysis

Consider $x^{\prime}=f(x)$ (special case: $\left.f(x)=A x+b\right)$. Might want:

- Static analysis $\left(f\left(x_{*}\right)=0\right)$
- Boils down to $A x=b$ (e.g. for Newton-like steps)
- Can solve directly or iteratively
- Sparsity matters a lot!
- Dynamic analysis (compute $x(t)$ for many values of $t$ )
- Involves time stepping (explicit or implicit)
- Implicit methods involve linear/nonlinear solves
- Need to understand stiffness and stability issues
- Modal analysis (compute eigenvalues of $A$ or $f^{\prime}\left(x_{*}\right)$ )


## Explicit time stepping

- Example: forward Euler
- Next step depends only on earlier steps
- Simple algorithms
- May have stability/stiffness issues


## Implicit time stepping

- Example: backward Euler
- Next step depends on itself and on earlier steps
- Algorithms involve solves - complication, communication!
- Larger time steps, each step costs more


## A common kernel

In all these analyses, spend lots of time in sparse matvec:

- Iterative linear solvers: repeated sparse matvec
- Iterative eigensolvers: repeated sparse matvec
- Explicit time marching: matvecs at each step
- Implicit time marching: iterative solves (involving matvecs)

We need to figure out how to make matvec fast!

## An aside on sparse matrix storage

- Sparse matrix $\Longrightarrow$ mostly zero entries
- Can also have "data sparseness" - representation with less than $O\left(n^{2}\right)$ storage, even if most entries nonzero
- Could be implicit (e.g. directional differencing)
- Sometimes explicit representation is useful
- Easy to get lots of indirect indexing!
- Compressed sparse storage schemes help


## Example: Compressed sparse row storage



This can be even more compact:

- Could organize by blocks (block CSR)
- Could compress column index data (16-bit vs 64-bit)
- Various other optimizations - see OSKI


## Distributed parameter problems

Mostly PDEs:

| Type | Example | Time? | Space dependence? |
| :--- | :--- | :--- | :--- |
| Elliptic | electrostatics | steady | global |
| Hyperbolic | sound waves | yes | local |
| Parabolic | diffusion | yes | global |

Different types involve different communication:

- Global dependence $\Longrightarrow$ lots of communication (or tiny steps)
- Local dependence from finite wave speeds; limits communication


## Example: 1D heat equation



Consider flow (e.g. of heat) in a uniform rod

- Heat $(Q) \propto$ temperature $(u) \times \operatorname{mass}(\rho h)$
- Heat flow $\propto$ temperature gradient (Fourier's law)

$$
\begin{aligned}
\frac{\partial Q}{\partial t} \propto h \frac{\partial u}{\partial t} & \approx C\left[\left(\frac{u(x-h)-u(x)}{h}\right)+\left(\frac{u(x)-u(x+h)}{h}\right)\right] \\
\frac{\partial u}{\partial t} & \approx C\left[\frac{u(x-h)-2 u(x)+u(x+h)}{h^{2}}\right] \rightarrow C \frac{\partial^{2} u}{\partial x^{2}}
\end{aligned}
$$

## Spatial discretization

Heat equation with $u(0)=u(1)=0$

$$
\frac{\partial u}{\partial t}=C \frac{\partial^{2} u}{\partial x^{2}}
$$

Spatial semi-discretization:

$$
\frac{\partial^{2} u}{\partial x^{2}} \approx \frac{u(x-h)-2 u(x)+u(x+h)}{h^{2}}
$$

Yields a system of ODEs

$$
\frac{d u}{d t}=C h^{-2}(-T) u=-C h^{-2}\left[\begin{array}{ccccc}
2 & -1 & & & \\
-1 & 2 & -1 & & \\
& \ddots & \ddots & \ddots & \\
& & -1 & 2 & -1 \\
& & & -1 & 2
\end{array}\right]\left[\begin{array}{c}
u_{1} \\
u_{2} \\
\vdots \\
u_{n-2} \\
u_{n-1}
\end{array}\right]
$$

## Explicit time stepping

Approximate PDE by ODE system ("method of lines"):

$$
\frac{d u}{d t}=\mathrm{Ch}^{-2} \mathrm{Tu}
$$

Now need a time-stepping scheme for the ODE:

- Simplest scheme is Euler:

$$
u(t+\delta) \approx u(t)+u^{\prime}(t) \delta=\left(1-C \frac{\delta}{h^{2}} T\right) u(t)
$$

- Taking a time step $\equiv$ sparse matvec with $\left(1-C \frac{\delta}{h^{2}} T\right)$
- This may not end well...


## Explicit time stepping data dependence



Nearest neighbor interactions per step $\Longrightarrow$ finite rate of numerical information propagation

## Explicit time stepping in parallel


for $t=1$ to $N$
communicate boundary data ("ghost cell") take time steps locally end

## Overlapping communication with computation


for $t=1$ to $N$
start boundary data sendrecv
compute new interior values
finish sendrecv
compute new boundary values
end

## Batching time steps


for $t=1$ to $N$ by $B$
start boundary data sendrecv (B values)
compute new interior values
finish sendrecv (B values)
compute new boundary values
end

## Explicit pain



Unstable for $\delta>O\left(h^{2}\right)$ !

## Implicit time stepping

- Backward Euler uses backward difference for $d / d t$

$$
u(t+\delta) \approx u(t)+u^{\prime}(t+\delta t) \delta
$$

- Taking a time step $\equiv$ sparse matvec with $\left(I+C \frac{\delta}{h^{2}} T\right)^{-1}$
- No time step restriction for stability (good!)
- But each step involves linear solve (not so good!)
- Good if you like numerical linear algebra?


## Explicit and implicit

Explicit:

- Propagates information at finite rate
- Steps look like sparse matvec (in linear case)
- Stable step determined by fastest time scale
- Works fine for hyperbolic PDEs

Implicit:

- No need to resolve fastest time scales
- Steps can be long... but expensive
- Linear/nonlinear solves at each step
- Often these solves involve sparse matvecs
- Critical for parabolic PDEs


## Poisson problems

Consider 2D Poisson

$$
-\nabla^{2} u=\frac{\partial^{2} u}{\partial x^{2}}+\frac{\partial^{2} u}{\partial y^{2}}=f
$$

- Prototypical elliptic problem (steady state)
- Similar to a backward Euler step on heat equation


## Poisson problem discretization

$$
u_{i, j}=h^{-2}\left(4 u_{i, j}-u_{i-1, j}-u_{i+1, j}-u_{i, j-1}-u_{i, j+1}\right)
$$

$$
L=\left[\begin{array}{ccc|ccc|ccc}
4 & -1 & & -1 & & & & & \\
-1 & 4 & -1 & & -1 & & & & \\
& -1 & 4 & & & -1 & & & \\
\hline-1 & & & 4 & -1 & & -1 & & \\
& -1 & & -1 & 4 & -1 & & -1 & \\
& & -1 & & -1 & 4 & & & -1 \\
\hline & & & -1 & & & 4 & -1 & \\
& & & & -1 & & -1 & 4 & -1 \\
& & & & & -1 & & -1 & 4
\end{array}\right]
$$

## Poisson solvers in 2D/3D

$N=n^{d}=$ total unknowns

| Method | Time | Space |
| :--- | :--- | :--- |
| Dense LU | $N^{3}$ | $N^{2}$ |
| Band LU | $N^{2}\left(N^{7 / 3}\right)$ | $N^{3 / 2}\left(N^{5 / 3}\right)$ |
| Jacobi | $N^{2}$ | $N$ |
| Explicit inv | $N^{2}$ | $N^{2}$ |
| CG | $N^{3 / 2}$ | $N$ |
| Red-black SOR | $N^{3 / 2}$ | $N$ |
| Sparse LU | $N^{3 / 2}$ | $N \log N\left(N^{4 / 3}\right)$ |
| FFT | $N \log N$ | $N$ |
| Multigrid | $N$ | $N$ |

Ref: Demmel, Applied Numerical Linear Algebra, SIAM, 1997.

## General implicit picture

- Implicit solves or steady state $\Longrightarrow$ solving systems
- Nonlinear solvers generally linearize
- Linear solvers can be
- Direct (hard to scale)
- Iterative (often problem-specific)
- Iterative solves boil down to matvec!


## PDE solver summary

- Can be implicit or explicit (as with ODEs)
- Explicit (sparse matvec) - fast, but short steps?
- works fine for hyperbolic PDEs
- Implicit (sparse solve)
- Direct solvers are hard!
- Sparse solvers turn into matvec again
- Differential operators turn into local mesh stencils
- Matrix connectivity looks like mesh connectivity
- Can partition into subdomains that communicate only through boundary data
- More on graph partitioning later
- Not all nearest neighbor ops are equally efficient!
- Depends on mesh structure
- Also depends on flops/point

