Parallel Computing (CMSC416 / CMSC616)



Parallel CSE Applications

Abhinav Bhatele, Department of Computer Science



Announcements



Student cluster competition

- CMSC416: If you are an undergrad interested in participating in International Student Cluster Competitions
- bhatele@cs.umd.edu

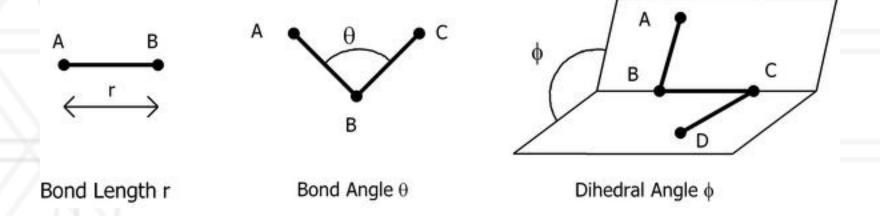


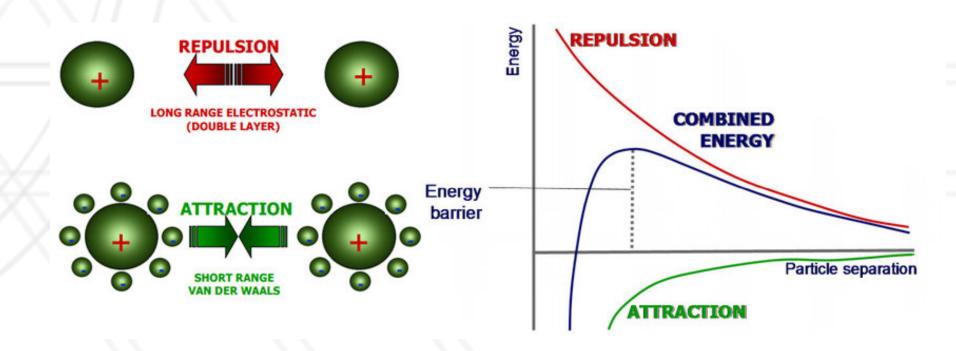


 Calculate trajectories of atoms and molecules by solving Newton's equations of motions

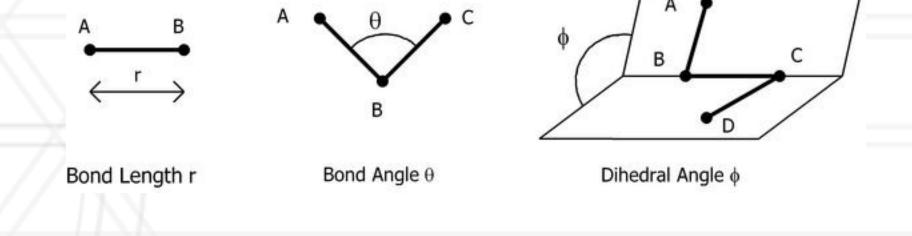


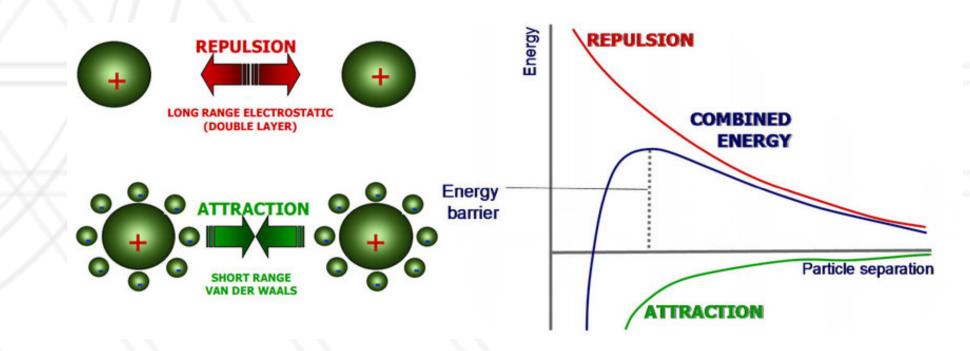
- Calculate trajectories of atoms and molecules by solving Newton's equations of motions
- Force calculations
 - Bonded interactions: bonds, angles, dihedrals
 - Non-bonded interactions: van der Waal's and electrostatic forces



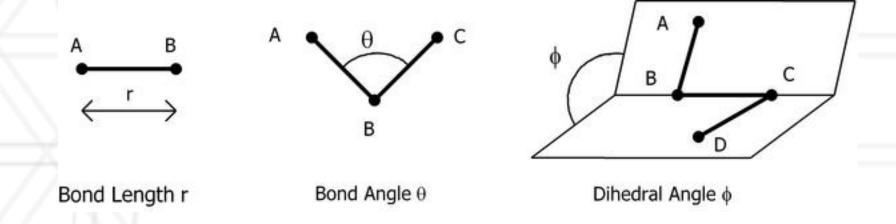


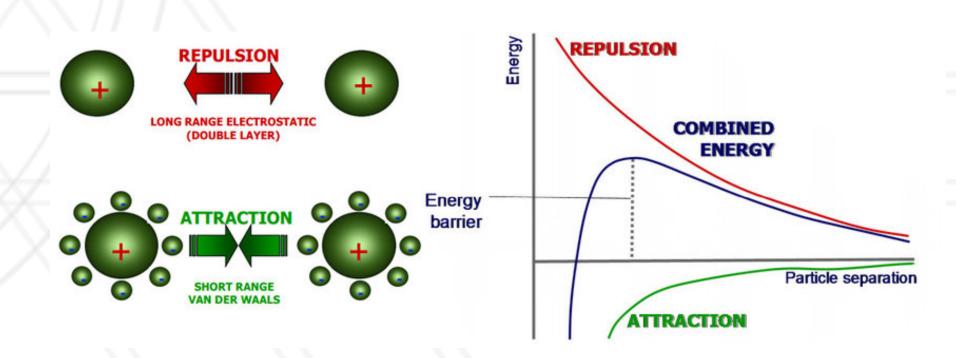
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- Number of atoms: thousands to millions



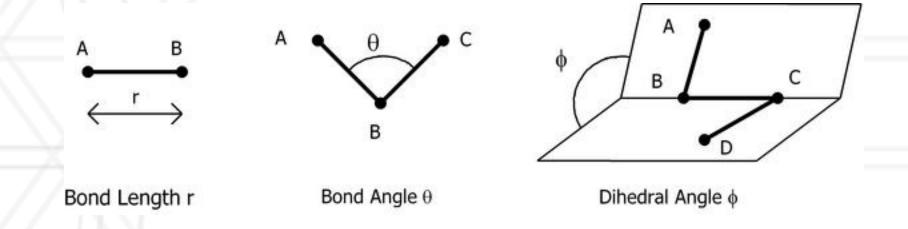


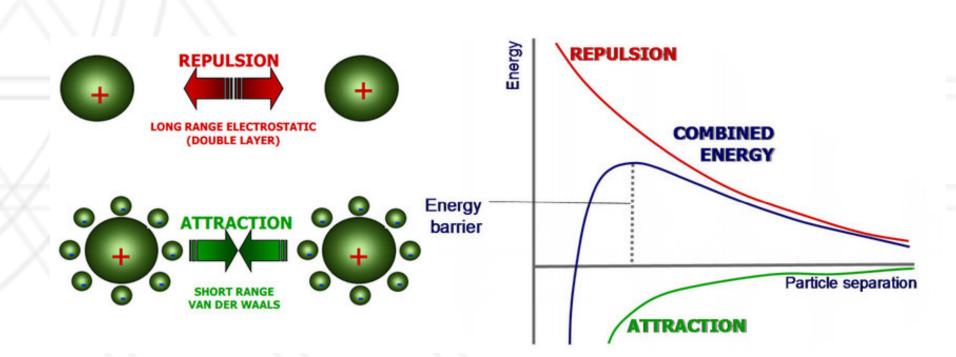
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- Number of atoms: thousands to millions
- Simulation step: ~I femtosecond (10-15 s)
- Used for drug design, materials design





Sequential Algorithm

- At every step, calculate forces on each atom
 - Calculate bonded and short-range forces every step
 - Calculate long-range non-bonded forces every few time steps (using PME or P3M etc.)
- Particle mesh Ewald (PME) summation:
 - Calculate long-range interactions in Fourier space
- Calculate velocities and new positions
- Repeat ...





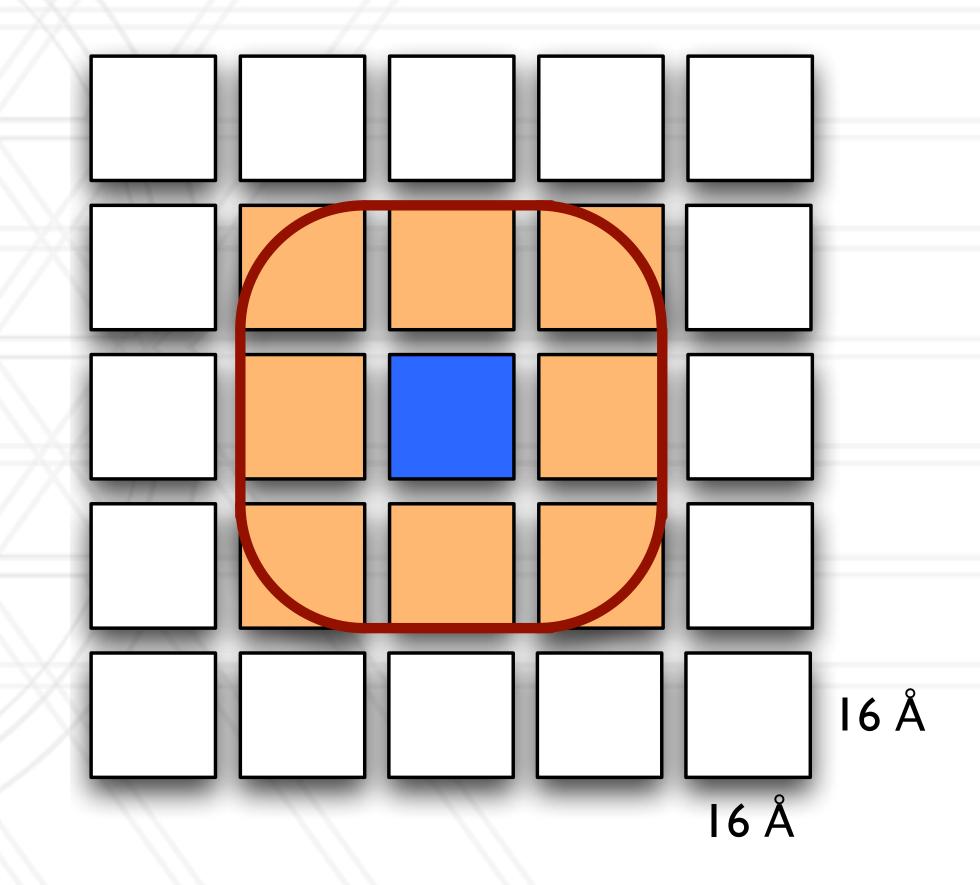
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- Force decomposition:
 - Distribute the force matrix to processes
 - Matrix is sparse and non-uniform



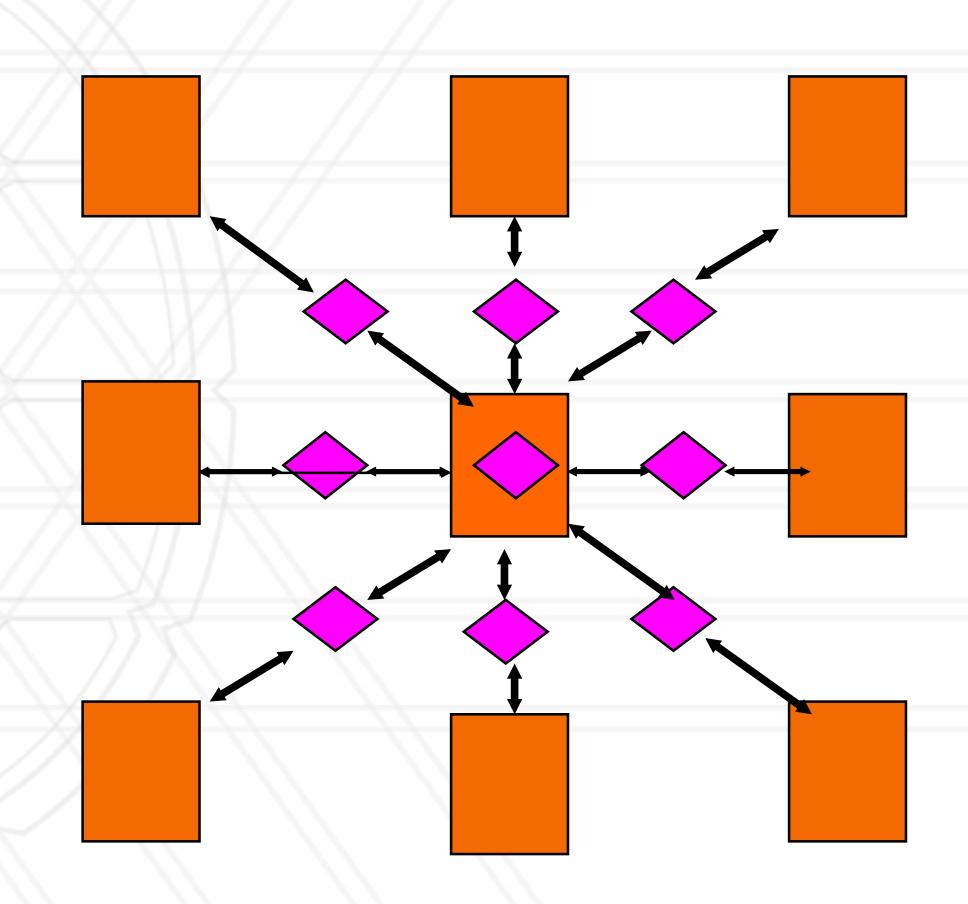
- Atom decomposition:
 - Partition the atoms across processes
- Force decomposition:
 - Distribute the force matrix to processes
 - Matrix is sparse and non-uniform
- Spatial decomposition:
 - Assign a region of the 3D simulation space to each process





Hybrid parallelization

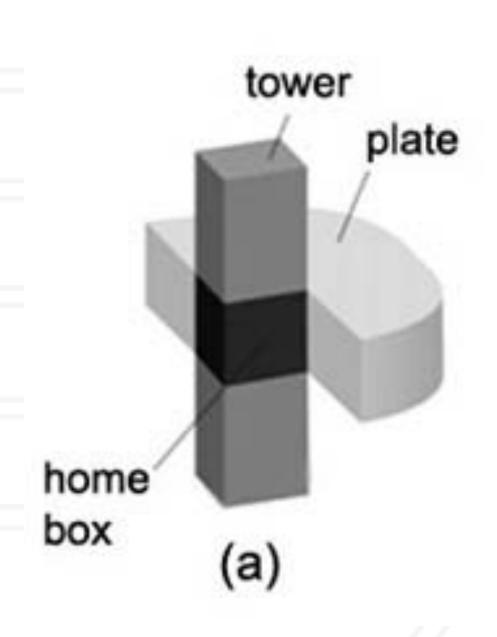
- Hybrid of spatial and force decomposition
- Decouple assignment of data and work to processes
- Distribute both atoms and the force calculations to different processes

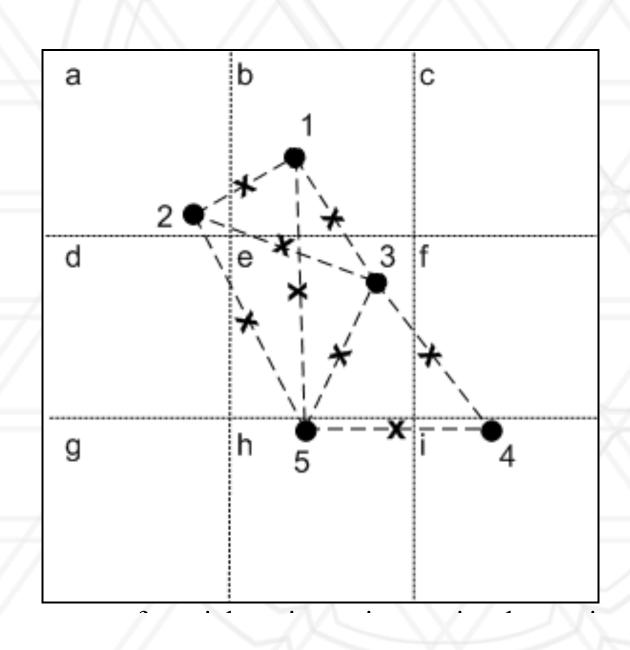




Neutral territory (NT) methods

Desmond's mid-point method





SC23 Test-of-time award:

Shaw DE, RO Dror, JK Salmon, et al. 2009. "Millisecond-scale molecular dynamics simulations on Anton," In *Proceedings of the Conference on High Performance Computing Networking, Storage and Analysis* (SC09), Portland, OR, USA, pp. 1-11, doi: 10.1145/1654059.1654126



Particle mesh Ewald

- Replace direct force calculations by:
 - Calculate short-range forces in real space
 - Calculate long-range forces in Fourier space
- Create a 3D mesh/grid representing charge densities of atoms
 - Compute a 3D Fast Fourier Transform (FFT)
- FFT computes the discrete Fourier transform (DFT) or inverse DFT
 - Reduces the complexity from $O(N^2)$ to $O(N \log N)$



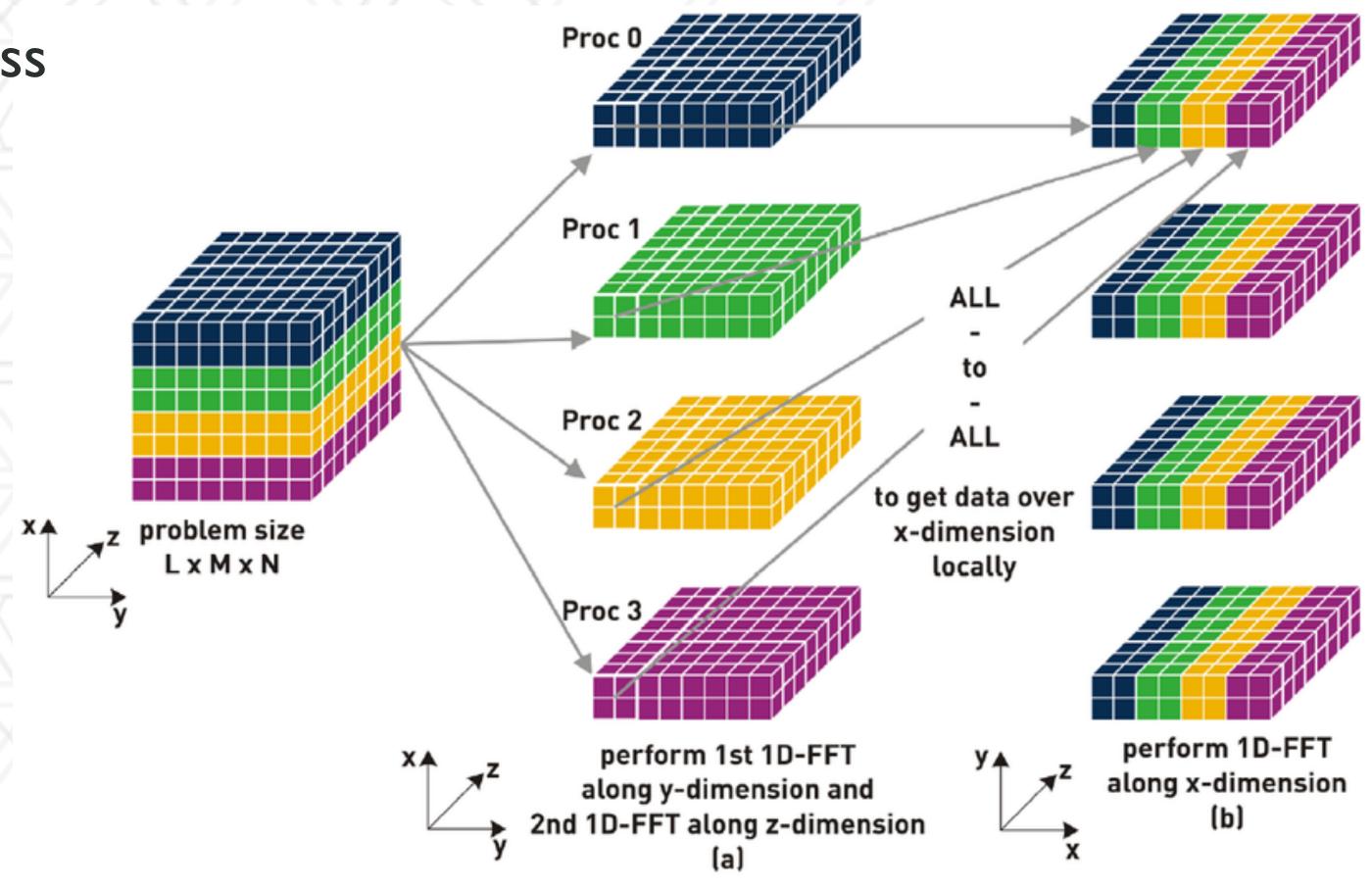


Bring all the data to one process



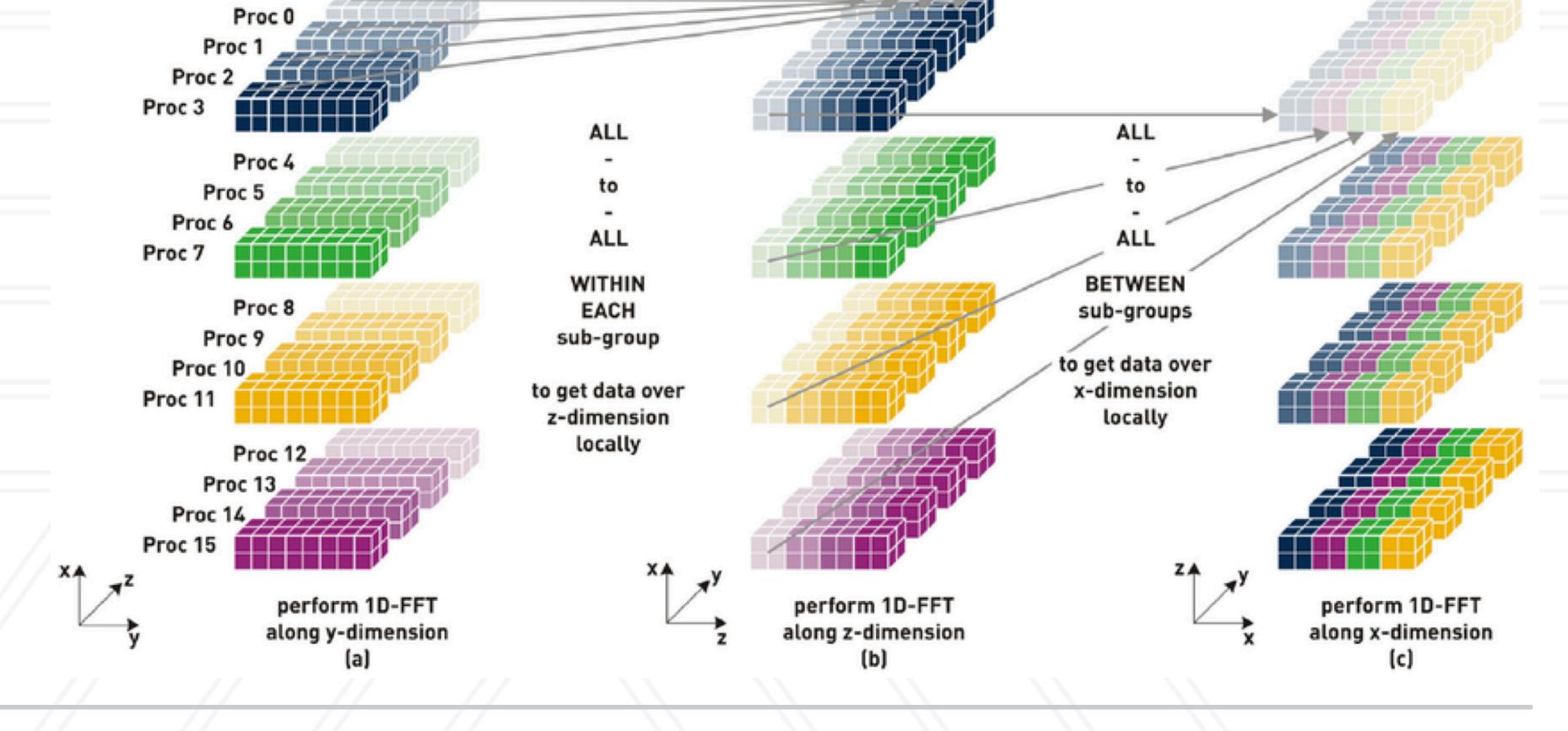
Bring all the data to one process

ID or slab decomposition





• 2D or pencil decomposition

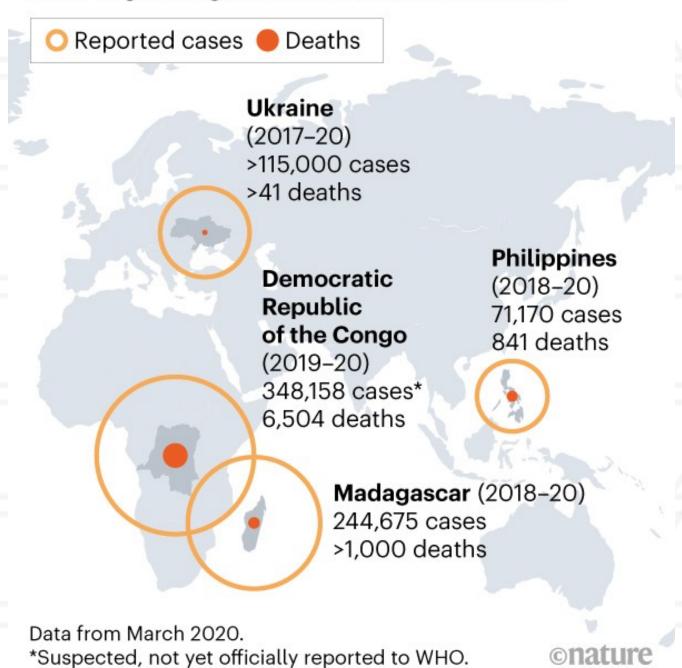




Measles killed 200,000 in 2020 alone!

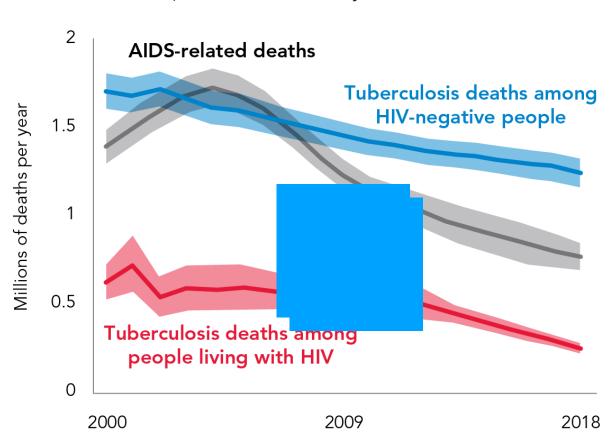
LARGE MEASLES OUTBREAKS

The epidemic in the Democratic Republic of the Congo is the largest single-nation outbreak for decades.



Global trends in the estimated number of deaths caused by TB and AIDS, 2000–2018

Shaded areas represent uncertainty intervals



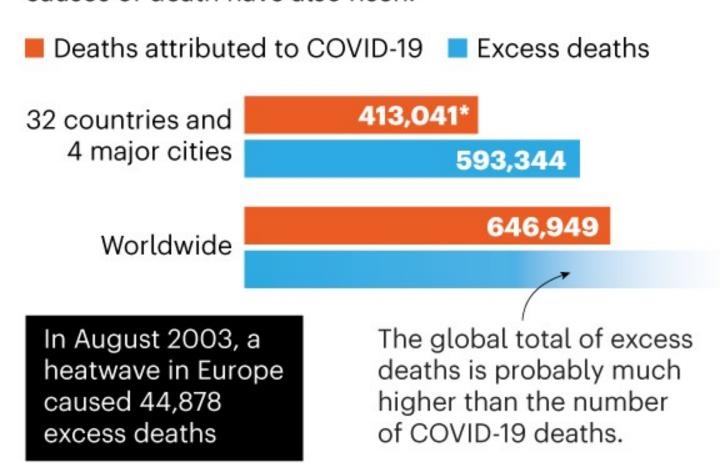
For AIDS, the latest estimates of the number of deaths in 2018 that have been published by UNAIDS are available at http://www.unaids.org/en/. For TB, the estimates for 2018 are those published in the Global Tuberculosis Report 2019. Deaths from TB among people living with HIV are officially classified as deaths caused by HIV/AIDS in the International Classification of Diseases.

Source: Global tuberculosis report 2019. Geneva: World Health Organization, 2019.

Predictions say that 1.66 million people died of tuberculosis in 2020

TERRIBLE TOLL

By the end of July 2020, there had been 646,949 COVID-19 deaths worldwide. In the 32 countries and 4 major cities with relevant data, there were more excess deaths than COVID-19 deaths, suggesting that some COVID-19 deaths are misclassified or that other causes of death have also risen.



*Cumulative deaths from outbreak onset to latest available data, as of 18 August 2020.

onature

https://www.nature.com/articles/d41586-020-01011-6

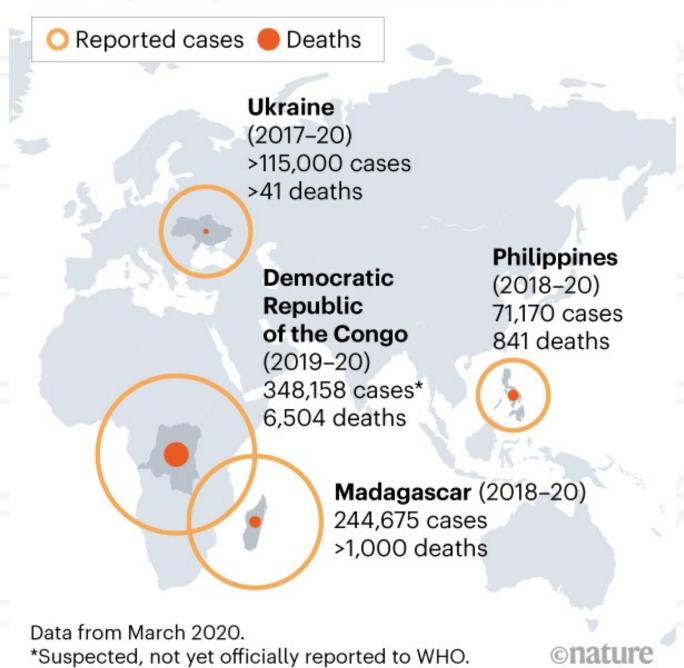
https://www.nature.com/articles/d41586-020-02497-w



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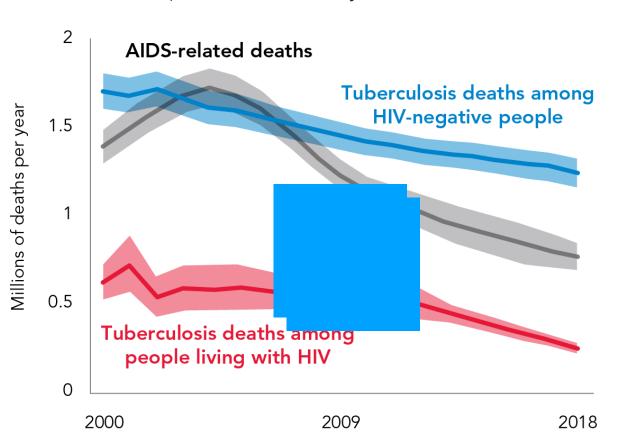
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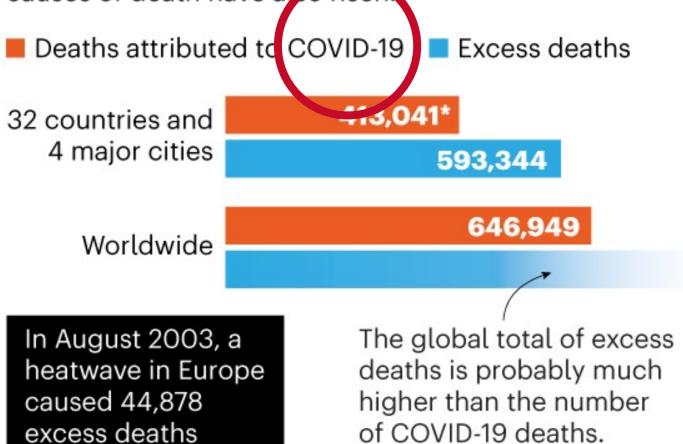
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Societal challenge

- Controlling the spread of infectious diseases is important
- Computational and mathematical modeling of epidemics important to assist governments in responding to outbreaks
- Made challenging due to:
 - increased and denser urbanization
 - increased local and global travel
 - increasingly immuno-comprised population



Approach: individual-based simulation

- Agent-based modeling to simulate epidemic diffusion
- Models agents (people) and interactions between them
- People interact when they visit the same location at the same time
- These "interactions" between pairs of people are represented as "visits" to locations
- Use a bi-partite graph of people and locations or a people-people interactivity graph



Serial algorithm

- At each timestep (typically a day):
 - Determine which people visit which locations
 - "Send" people to those locations
 - At each location "interactions" happen and transmission happens
 - Update people's states at the end of the day and continue
- Interventions (vaccinations, school closures) can be added on certain days to change people's susceptibility, movements etc.

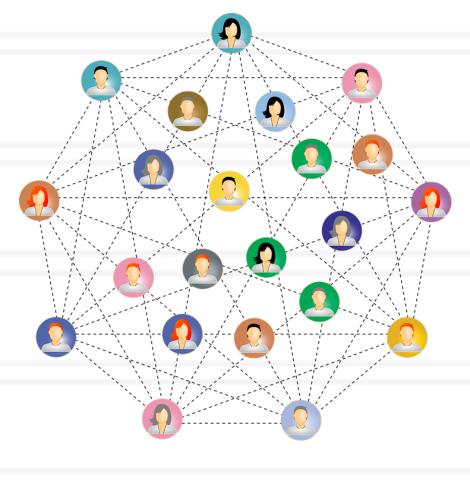


Hybrid time-stepped and discrete-event simulation



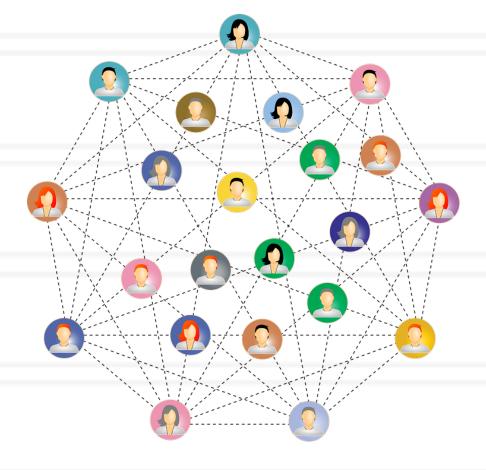
Hybrid time-stepped and discrete-event simulation

Contact model at each location

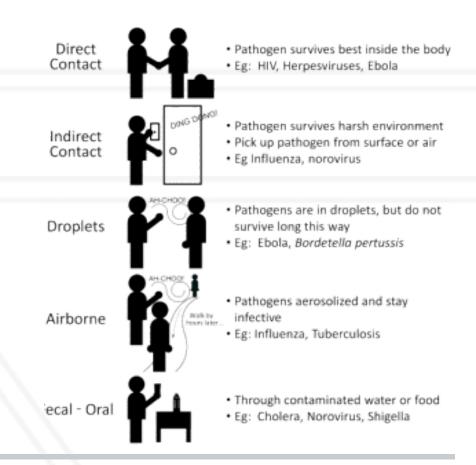


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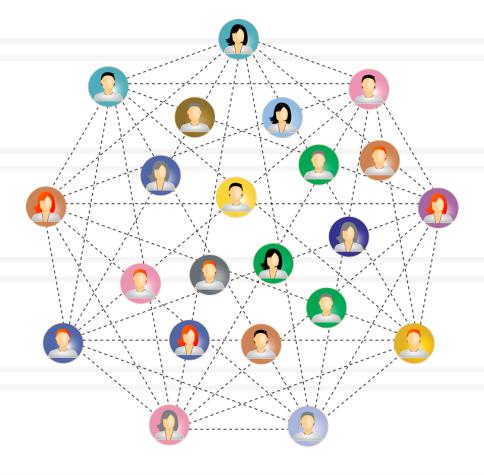


Model for human transmission

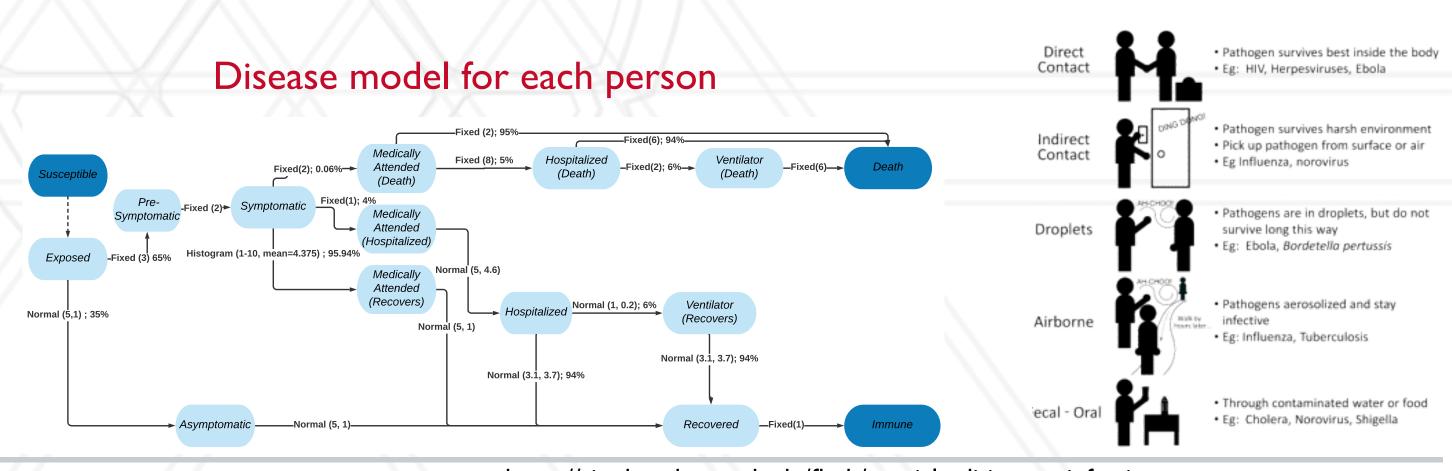


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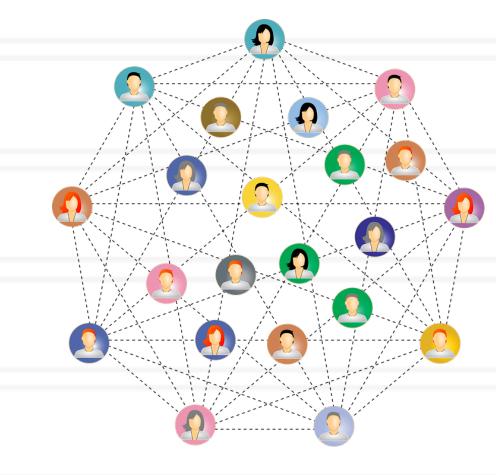
Hybrid time-stepped and discrete-event simulation

while d < num_days:
 for each person:
 Send visit messages to locations</pre>

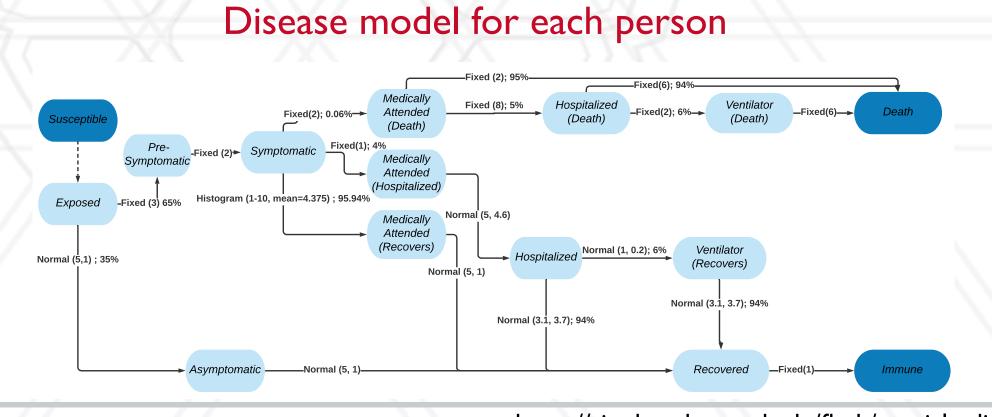
for each location:
Process all visit messages
Run discrete event simulation
Send interaction messages

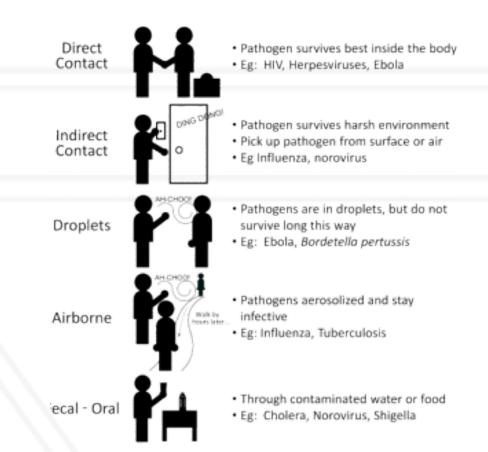
for each person:
Process interactions
Update disease state

Contact model at each location



Model for human transmission







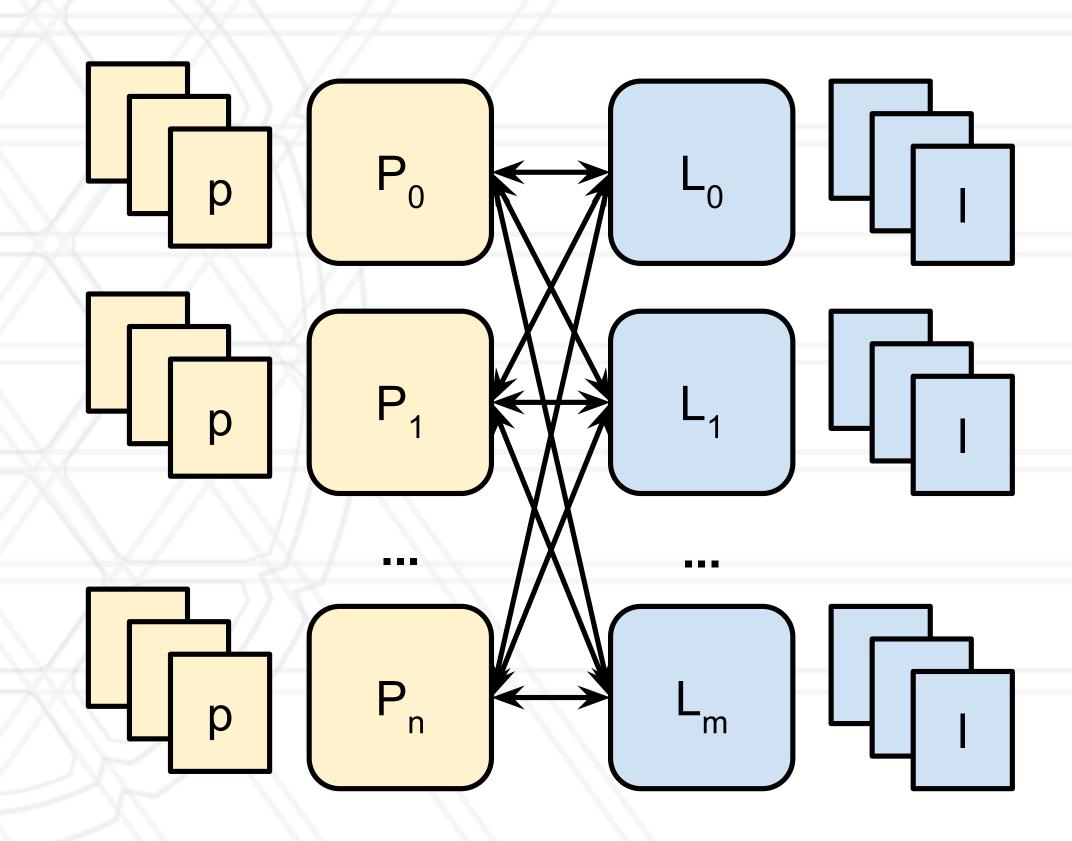
Parallel simulation is challenging

- Size and scale of the social contact network (6 billion agents for a global simulation)
 - Unstructured networks and complicated dependencies lead to high communication cost
- Individuals and their behaviors are not identical
- Co-evolving epidemics, public policies and agent behaviors make it impossible to apply standard model reduction techniques



Parallel implementation: Loimos

- All the people and locations are distributed among all processes
- DES computation can be done locally in parallel
- Communication when sending visit and infection messages
- Uses Charm++, a message-driven model





Application software stack

- Parallel programming model / runtime:
 - MPI, OpenMP, Charm++, CUDA, ...
- Libraries
 - Data and visualization libraries (mesh management, simulation output)
 - I/O libraries
 - Math/numerical libraries
 - Graph partitioning, load balancing ...

User program

Libraries

Parallel prog. runtime (MPI, Charm++ etc.)



Why use libraries?

- No need to reinvent the wheel
 - Libraries are highly optimized, have fewer bugs
- Avoids significant effort to write, optimize and maintain code
- Makes code more portable



Popular Libraries

- Data/visualization and I/O libraries
 - I/O: HDF5, pNetCDF, ADIOS
- Numerical libraries:
 - Fast Fourier transforms: FFTW
 - Dense linear algebra: BLAS, LAPACK, Intel MKL
 - Solvers for sparse systems: Hypre, PETSc, Trilinos
- Graph partitioning/load balancing:
 - METIS, Scotch, Zoltan, Chaco

https://events.prace-ri.eu/event/176/contributions/38/attachments/154/305/HPC_libraries.pdf



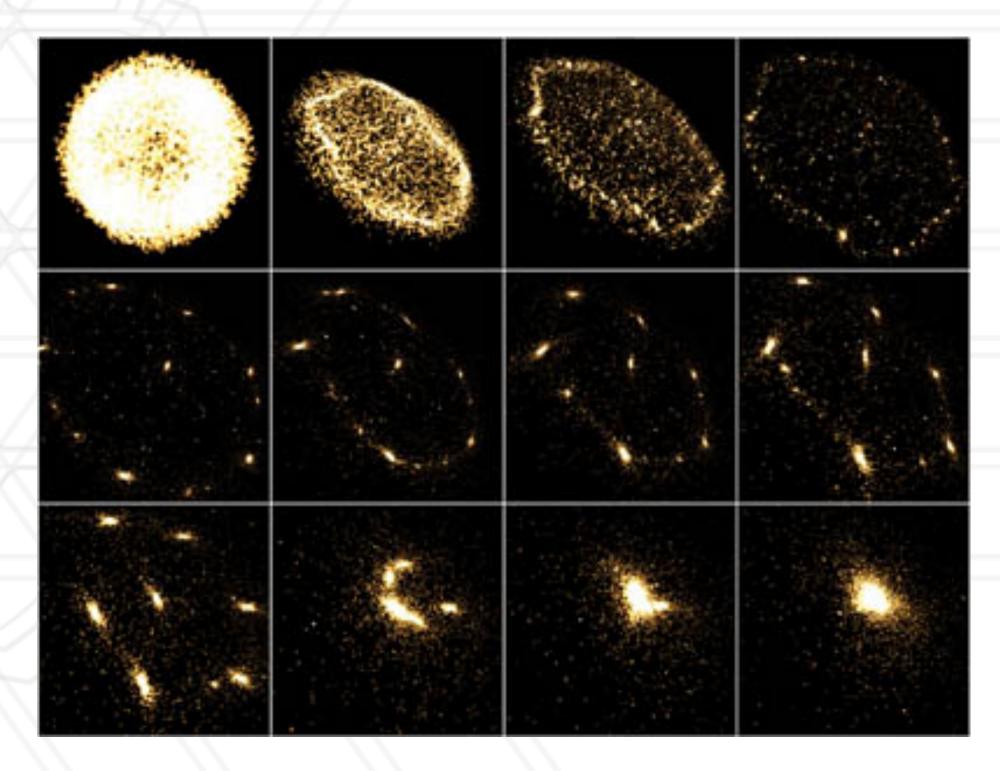
Domain-specific languages/frameworks

- Structured grids: SAMRAI, Chombo, AMREx
- Unstructured grids: MFEM, Quinoa



The n-body problem

- Simulate the motion of celestial objects interacting with one another due to gravitational forces
- Naive algorithm: $O(n^2)$
 - Every body calculates forces pair-wise with every other body (particle)



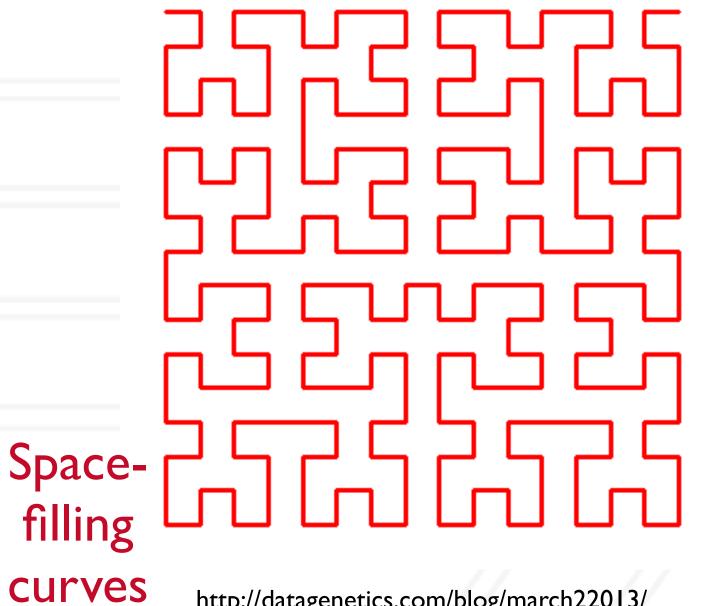
https://developer.nvidia.com/gpugems/gpugems3/part-v-physics-simulation/chapter-3 I -fast-n-body-simulation-cuda



- Naive approach: Assign n/p particles to each process
- Other approaches?



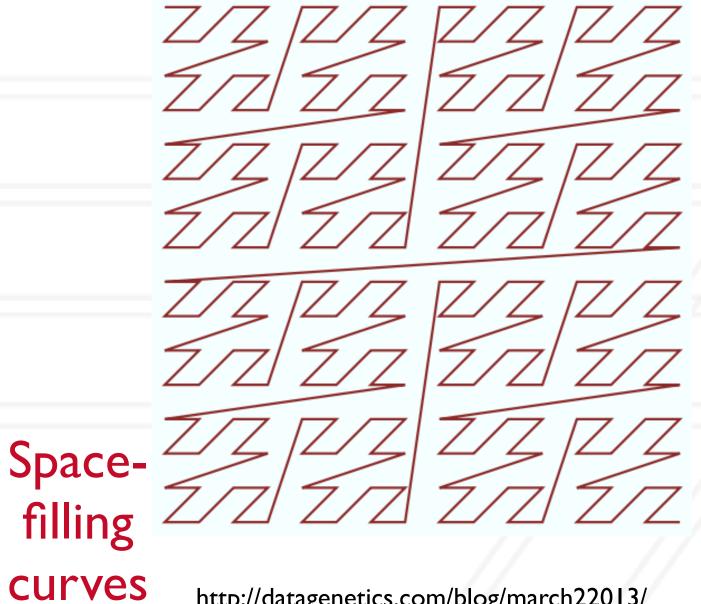
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http://datagenetics.com/blog/march22013/ https://en.wikipedia.org/wiki/Z-order_curve



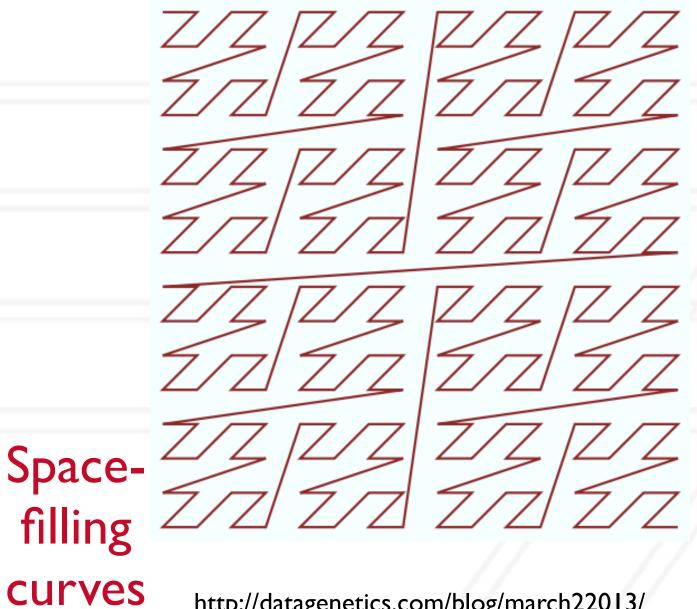
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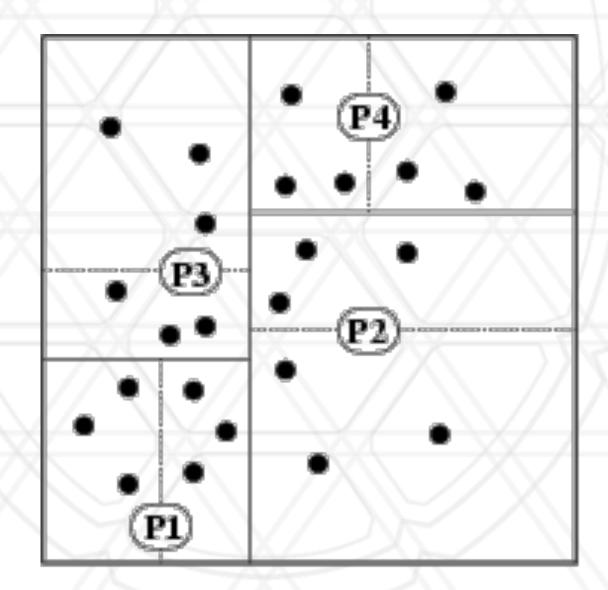
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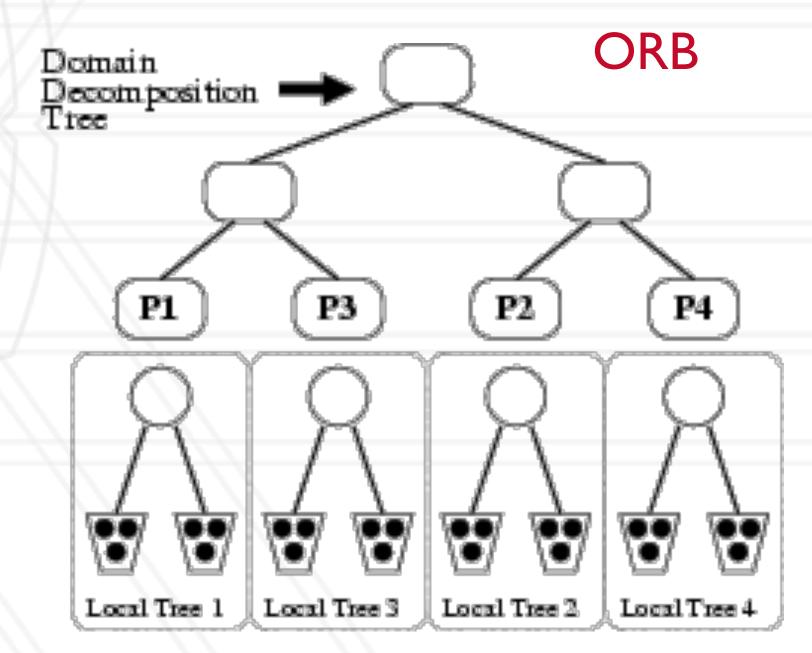


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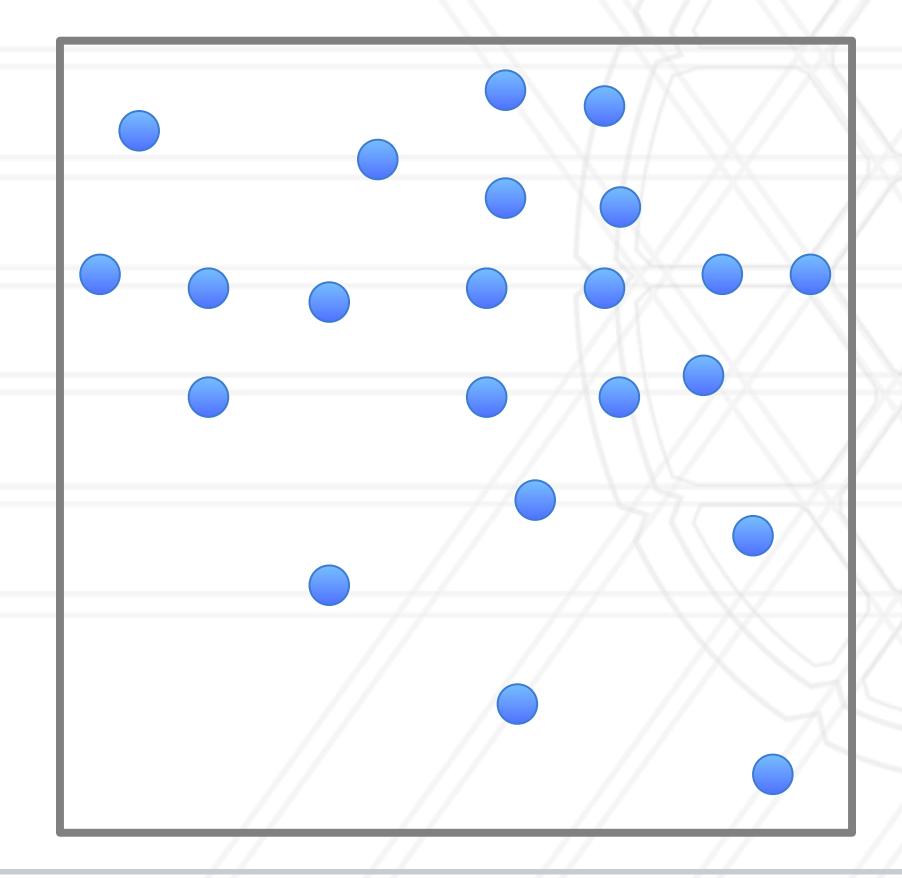




http://charm.cs.uiuc.edu/workshops/charmWorkshop2011/slides/CharmWorkshop2011_apps_ChaNGa.pdf

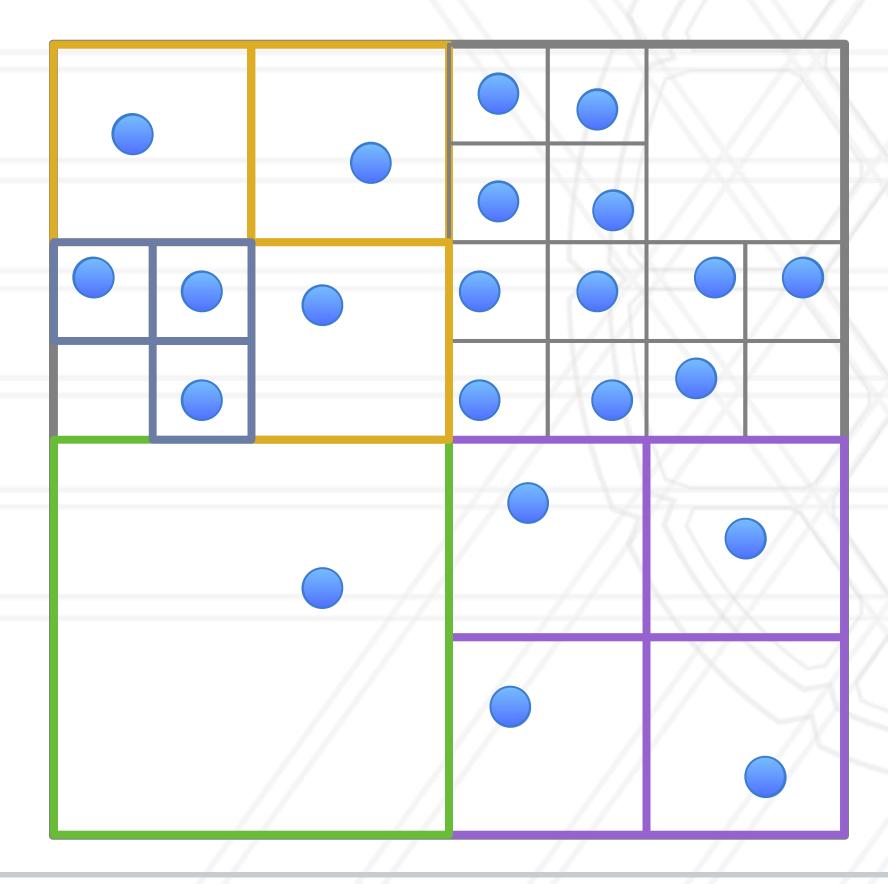


• Let us consider a two-dimensional space with bodies/particles in it



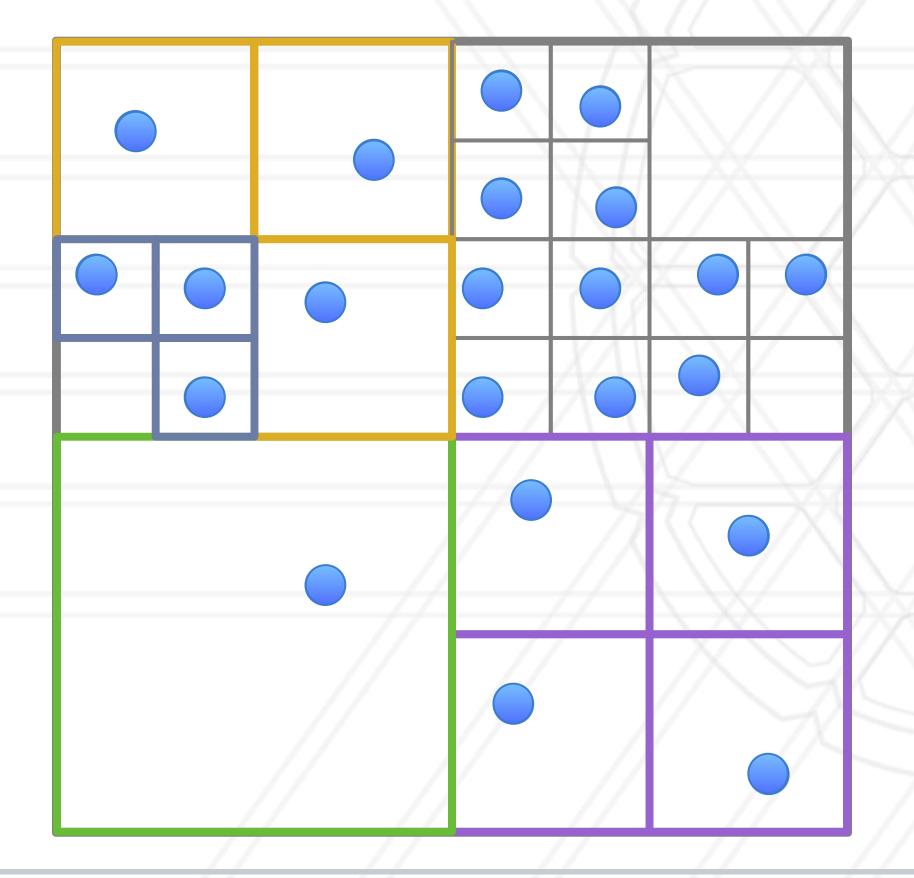


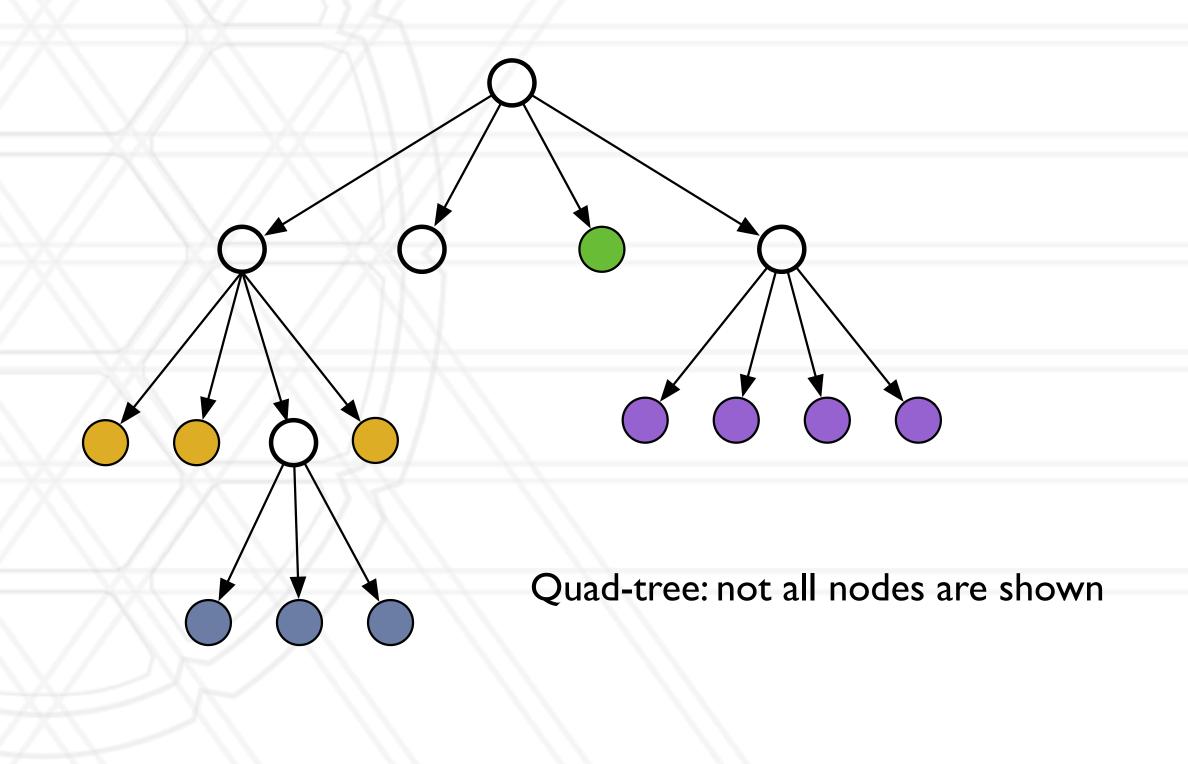
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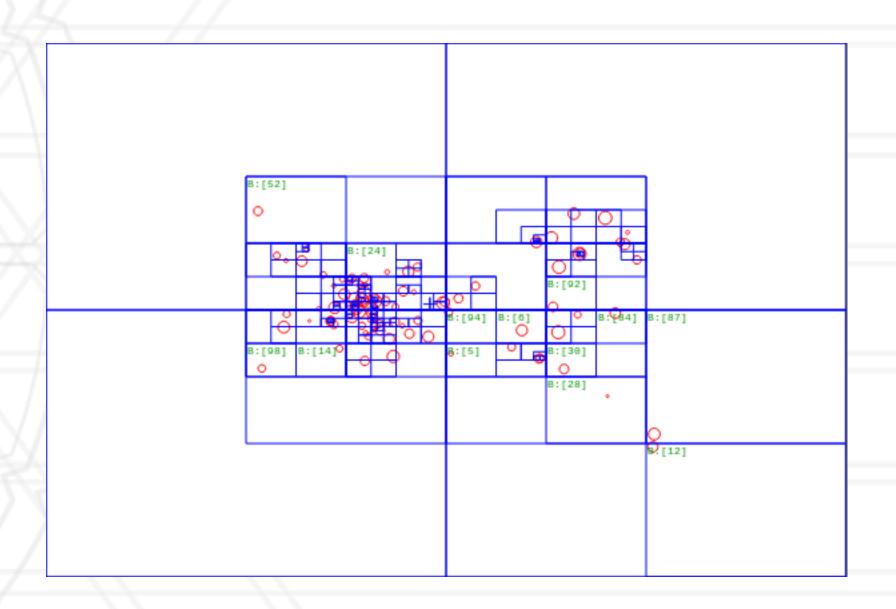
Different parallelization methods

- Tree codes: Barnes-Hut simulations
- Fast multipole methods (FMM): Greengard and Rokhlin
- Particle mesh methods
- Particle-particle particle-mesh (P³M) methods



Barnes-Hut simulation

- Represent the space containing the particles as an oct-tree
- Pairwise force calculations for nearby particles
- For tree nodes that are sufficiently far away,
 approximate the particles in the node by a single large particle at the center of mass
- O(N logN) algorithm



https://en.wikipedia.org/wiki/Barnes-Hut_simulation

Fast multipole methods

- Use multipole expansion for distant particles
- Takes advantage of the fact that for nearby particles, multipole-expanded forces from distant particles are similar
- Reduces the time complexity further to O(n)



Particle-particle particle-mesh methods

- Explicit calculation of forces on nearby particles
- Fourier-based Ewald summation for calculating potentials on a grid
- Smoothed particle hydrodynamics





Abhinav Bhatele

5218 Brendan Iribe Center (IRB) / College Park, MD 20742

phone: 301.405.4507 / e-mail: bhatele@cs.umd.edu