

# CMSC 714

## Lecture 23

### Applications

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## Notes

- Group projects due tomorrow, 6PM
  - Course evaluation web site open  
<http://www.CourseEvalUM.umd.edu>

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## Anton

- Special purpose machine built for molecular dynamics simulations, built at D.E. Shaw Research
  - to simulate biological processes that occur on very small time scales ( $10^{-15}$  sec), such as protein folding, interaction between proteins, etc.
  - and simulate those processes for a long time
- Molecular dynamics
  - force calculation followed by integration step to move particles
  - biomolecular forces have 3 parts
    - bonded forces – small atomic groups with covalent bonds
    - van der Waals forces – all pairs of atoms, but fall off quickly with distance (so only need close ones)
    - electrostatic forces – all pairs of atoms, fall off slowly with distance – divide into 2 parts to avoid all pairs computation

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## Anton

- Anton machine
  - up to 512 nodes in 8x8x8 torus
  - each node has 2 parts on 1 chip
    - high throughput interaction subsystem (HTIS) for range-limited interactions, using 32 hardwired pairwise point interaction pipelines (PPIPs)
    - flexible subsystem with 8 programmable geometry cores (GCs) for less structured part of MD computation, 4 Tensilica processors, 4 data transfer engines
    - plus DRAM controllers, 6 network interfaces, and host interface for I/O
- Most of computational time mapped to PPIPs, which run those computations maybe 100x faster than standard microprocessor core
- And computations spatially decomposed across nodes, with some twists to deal with communication as particles move between spatial domains
- Uses fixed-point arithmetic, with various bit widths, for several reasons:
  - performance – fixed-point hardware fast and small
  - fixed point arithmetic is truly associative
  - gain determinism – run same simulation again get exact same results bit-for-bit (doesn't really help, since MD is a chaotic system, so need ensemble)
  - computations are reversible

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## Anton

- Performance results show that can run a large chemical system at much higher rates than any previous system
  - can run multiple microseconds of simulation time per day of wall clock time
    - maybe 500 times faster than 512 node Intel Xeon cluster
    - and have run simulated systems up to over 1000 microseconds, which showed interesting behavior of the molecules
  - and results are validated very well
    - both against “known” results and using statistical error tests

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## Virtual Microscope

- Software emulation of a light microscope, to view and manipulate very large slide images, built at UMD and Johns Hopkins med school
  - for viewing and processing images captured from standard pathology specimens (need special purpose hardware for high throughput data capture)
  - problem is very large data sizes
    - a slide is maybe 30K pixels on a side at high resolution, so one focal plane is maybe 10GB uncompressed
    - and need multiple focal planes for some samples
    - and JHU hospital produces >400K slides per year
- Client/server system design
  - client runs on user desktop machine – Java GUI
  - server stores, retrieves, processes slide image data on parallel machine or workstation cluster
    - implemented both with Active Data Repository OO framework and with DataCutter component framework

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## Virtual Microscope

- Client provides drag/zoom interface to browse through a slide
  - use thumbnail to keep track of where you are on a slide
  - standalone client can cache image data for improved response time – using both memory and disk on client machine
- Server basic computation is map-reduce
  - map one or more input pixels at highest resolution to desired output resolution, and aggregate if multiple pixels map to same output pixel
- Active Data Repository
  - user defined functions used for map and reduce, framework orchestrates parallel execution across data stored on multiple nodes of a cluster or parallel machine
  - data blocks distributed across disks for parallel access and are indexed for fast retrieval (more important for more complex map functions)
  - images also need to be decompressed from stored JPEG form before map and reduce steps, and clipped to query window
  - experiments show that ADR implementation scales well, to handle multiple clients, with low overhead

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## Virtual Microscope

- DataCutter
  - component framework for processing large datasets in a distributed environment
  - filter-stream programming model (think *disklets*)
    - each filter is a component, and filters connected via streams, which deliver data buffers between filters
  - supports flexible placement of filters, filter replication for load balancing (transparent copies)
  - VM filter pipeline is: read-data, decompress, clip, zoom, view
- Performance results show that DataCutter implementation deals better than ADR with load balance issues, but ADR can process large queries faster from parallel execution of a single query
  - for DataCutter, filter placement matters – communication between filters adds latency if on different hosts
- Overall performance results for VM show that can achieve interactive response times for real slide data, on not-too-large server system configurations

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