

# Adaptive Heterogeneous Computing with OpenCL: Harnessing hundreds of GPUs and CPUs

Simon McIntosh-Smith simonm@cs.bris.ac.uk
Head of Microelectronics Research
University of Bristol, UK







#### **K** A brief biography







ClearSpeed<sup>®</sup>





Graduated as Valedictorian in Computer Science from Cardiff University in 1991

Joined Inmos to work for David May as a microprocessor architect

Moved to Pixelfusion in 1999 – a high-tech startup designing the first GPGPU, a many-core general purpose graphics processor

Co-founded ClearSpeed in 2002 as Director of Architecture and Applications

Joined the CS department at the University of Bristol in April 2009 to focus on High Performance Computing and architectures

#### Recent HPC activities

- European HPC:
  - PRACE
    - Mont Blanc ARM-based supercomputer
    - EESI exascale software initiative

- UK HPC:
  - ARCHER project working group
    - Next UK national supercomputer
    - Several PFLOPS, goes live summer '13



#### **K** Collaborators

- Richard B. Sessions, Amaurys Avila Ibarra
  - University of Bristol, Biochemistry
  - Developers of the docking code BUDE
- James Price (port to OpenCL)
  - University of Bristol, Computer Science
- Tsuyoshi Hamada, Felipe Cruz (GPUs)
  - University of Nagasaki, Japan
  - Winners of the 2009 Gordon Bell price/ performance prize





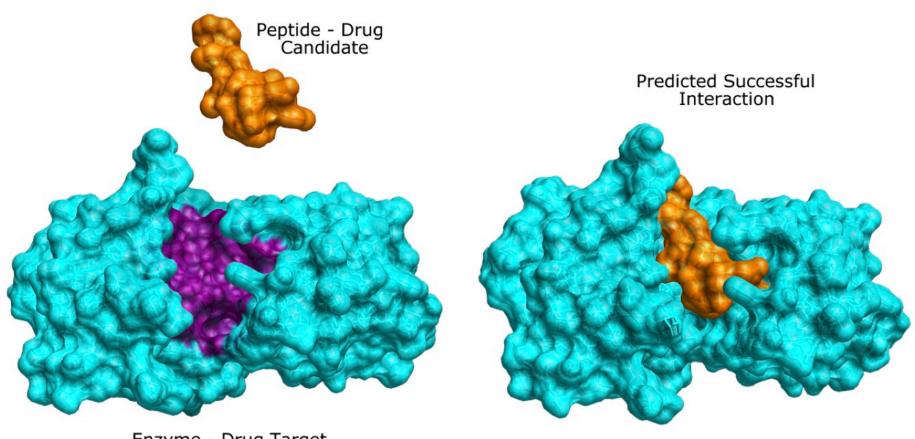
# Molecular docking







# Molecular docking

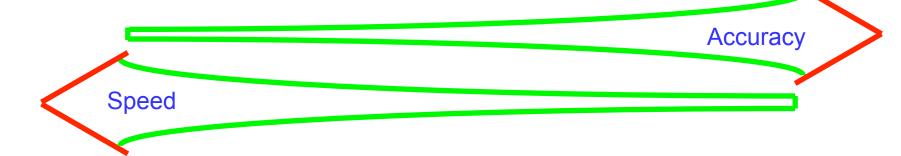






Proteins typically O(1,000) atoms Ligands typically O(100) atoms

# **BUDE**: Bristol University Docking Engine



Typical docking scoring **functions** 

No

**Empirical Free Energy Forcefield BUDE** 

Free Energy calculations  $MM^{1,2}$  QM/MM<sup>3</sup>

Yes

Yes

Yes

Yes

Yes

#### **Entropy:**

solvation No Yes configurational Approx **Approx Electrostatics Approx** All atom No Yes **Explicit solvent** 



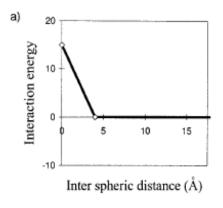
1. MD Tyka, AR Clarke, RB Sessions, J. Phys. Chem. B 110 17212-20 (2006)

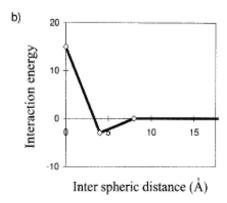
No

- 2. MD Tyka, RB Sessions, AR Clarke, J. Phys. Chem. B 111 9571-80 (2007)
- 3. CJ Woods, FR Manby, AJ Mulholland, J. Chem. Phys. 128 014109 (2008)

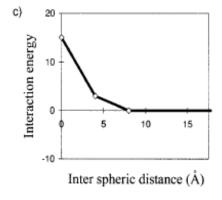
# Empirical Free Energy Function (atom-atom)

$$\Delta G_{\text{ligand binding}} = \sum_{i=1}^{N_{\text{protein}}} \sum_{j=1}^{N_{\text{ligand}}} f(x_i, x_j)$$





# Parameterised using experimental data<sup>†</sup>



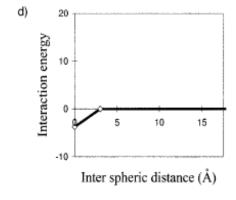
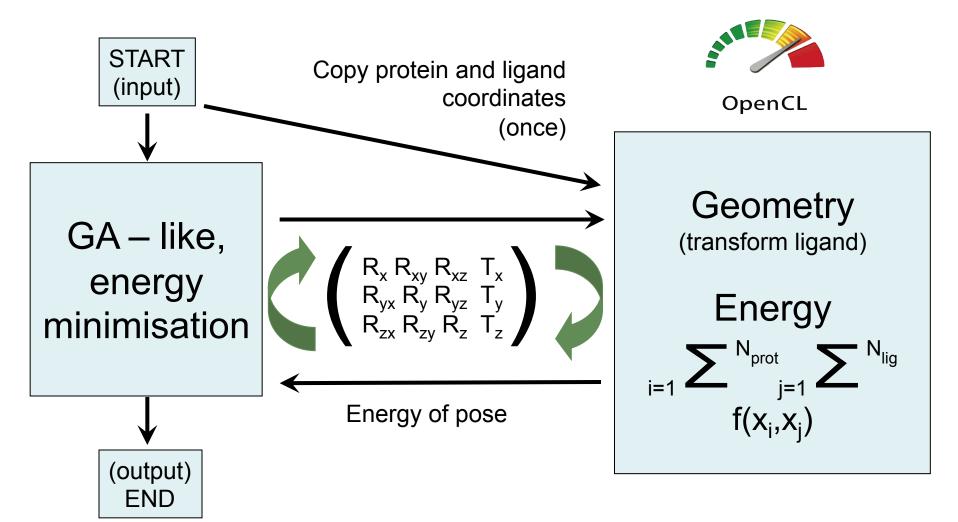


Fig. 1. Inter-residue sphere-sphere interaction energy functions of the force field, a: Between two polar spheres, or between a backbone sphere and any other non-hydrogen-bonding sphere, b: Between two non-polar spheres, c: Between a non-polar sphere and a polar sphere, d: Between a hydrogen bond donor sphere and a hydrogen bond acceptor sphere.



# **BUDE** Acceleration with OpenCL



# Why OpenCL?

- Open standard with fast growing support
  - Evolving rapidly: C++, HSA, OpenCL 2.0, ...

- Platform portability
  - Have used dozens of different GPUs & CPUs

- Naturally supports heterogeneous systems
  - Harness CPUs and GPUs simultaneously



# Multiple levels of parallelism

- O(10<sup>8</sup>) conformers from O(10<sup>7</sup>) ligands
- O(10<sup>5</sup>) poses per conformer (ligand)
- O(10<sup>3</sup>) atoms per protein
- O(10<sup>2</sup>) atoms per ligand (drug molecule)
- Conformers all independent
- Poses all independent, but there are benefits in grouping all poses of one conformer to one OpenCL device

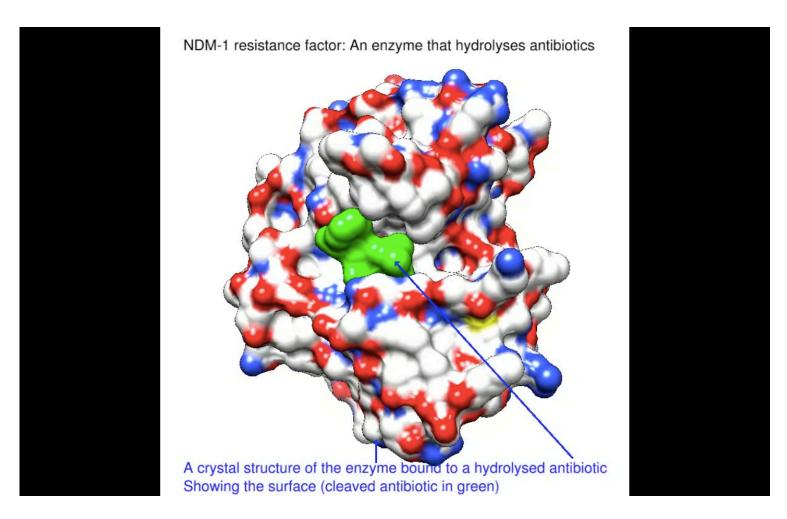


# Parallelism strategy

- Distribute ligands across nodes 10<sup>7</sup>-way parallelism
- All the poses of one conformer distributed across all the OpenCL devices in a node – 10<sup>3</sup>-way parallelism
- Each Work-Item (thread) performs an entire conformer-protein docking – 10<sup>5</sup> atom-atom force calculations



#### **K** How BUDE's EMC works







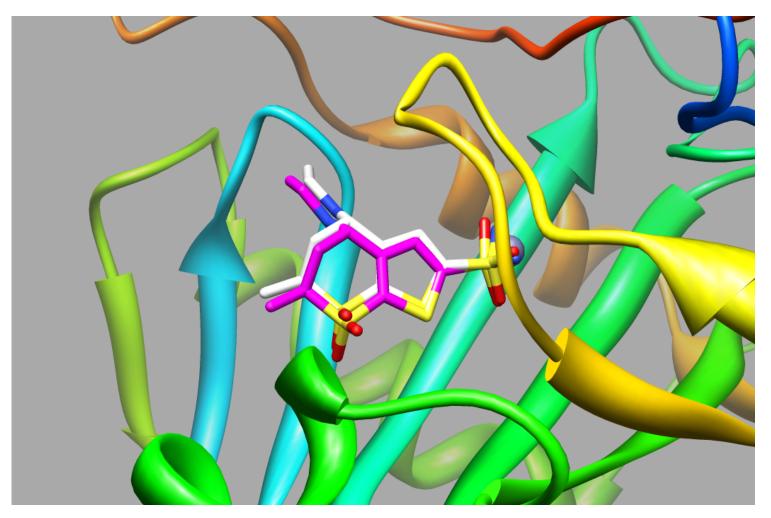
#### **Experimental results**





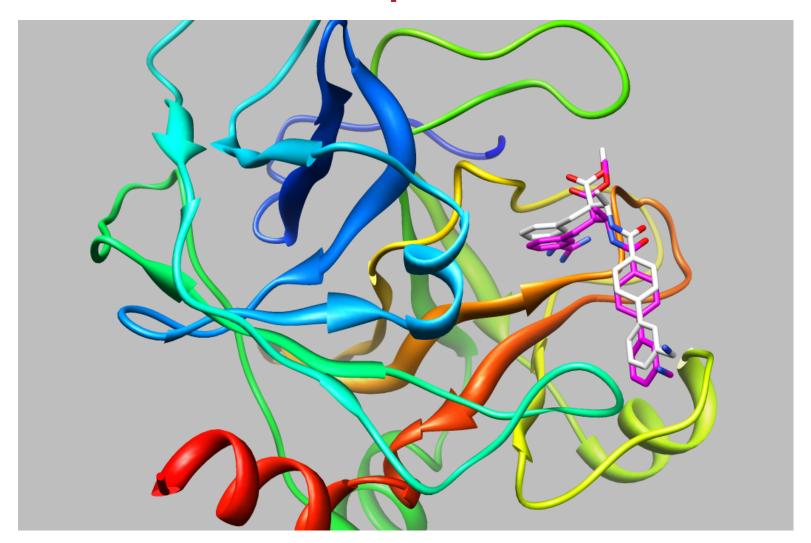


# Redocking into Xray Structure



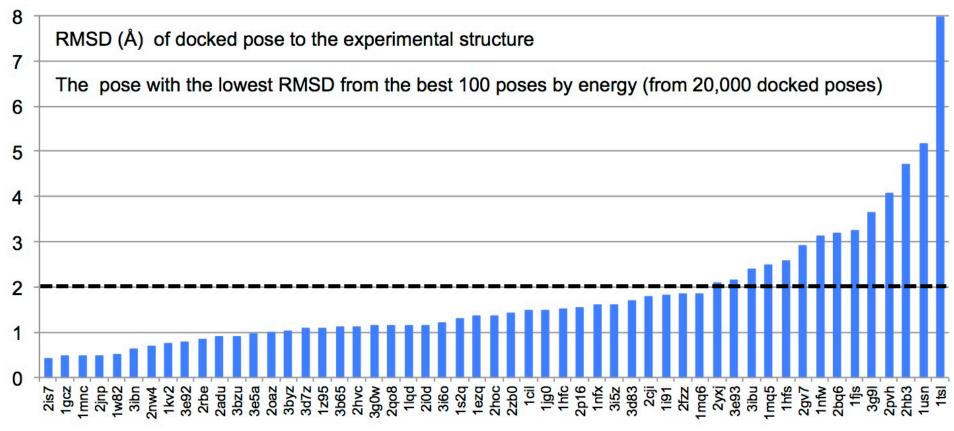


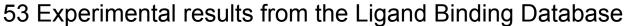
# Another example





# The science is working well









#### **Heterogeneous Systems**



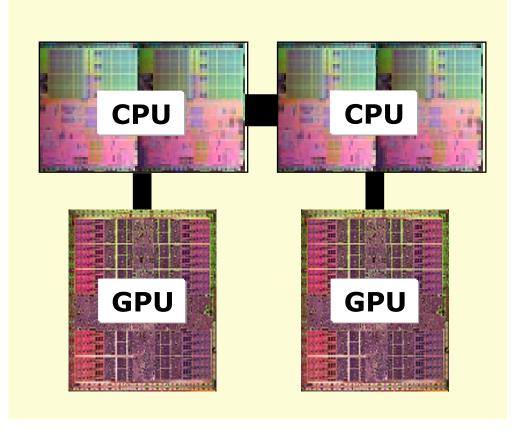




#### OpenCL for heterogeneous computing

A modern computer includes:

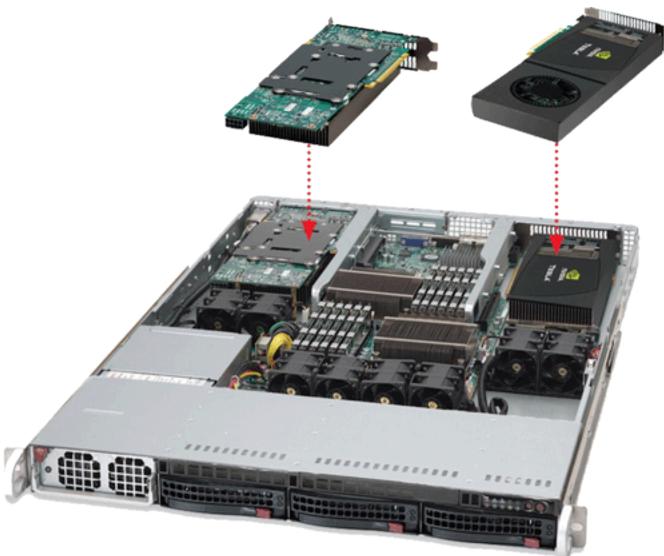
- One or more CPUs
- One or more GPUs



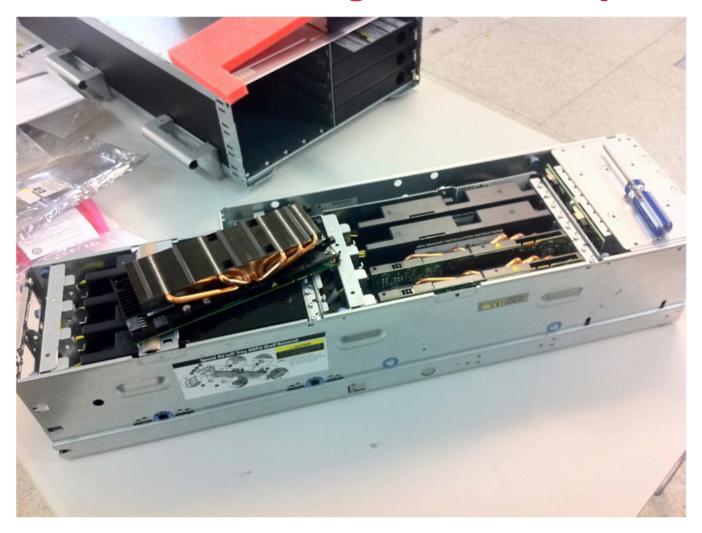
OpenCL (Open Compute Language) lets programmers write a single <u>portable</u> program that uses <u>ALL</u> resources in the heterogeneous platform



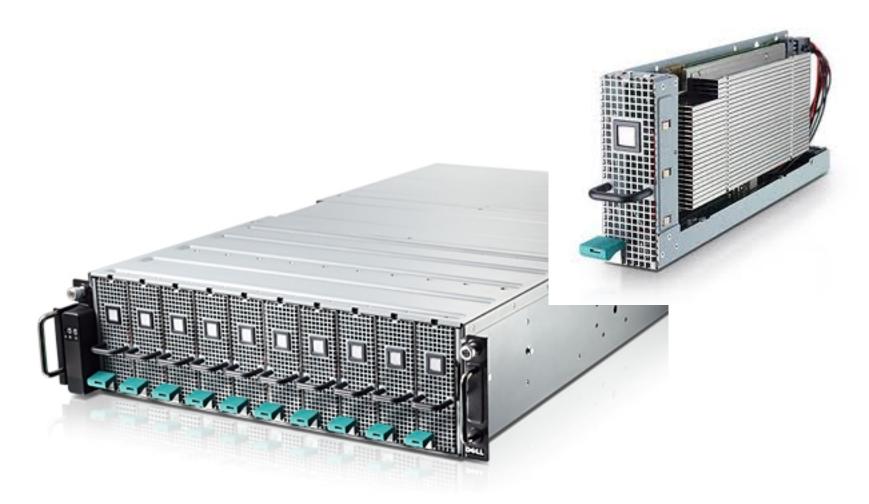






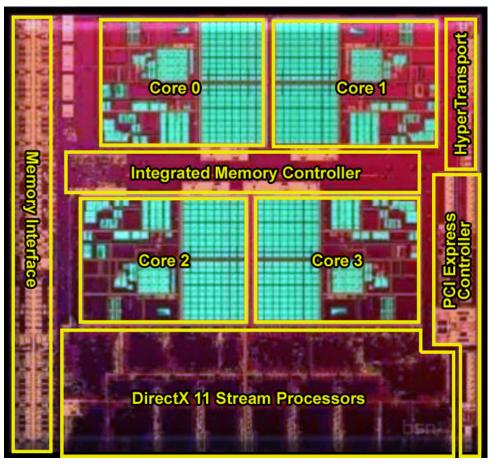














AMD Llano Fusion APUs



Intel Core2 Duo CPU P8600 @ 2.40GHz, NVIDIA GeForce 9400M integrated GPU, NVIDIA GeForce 9600M GT discrete GPU





#### **Benchmark results**







#### BUDE's heterogeneous approach

#### Within a node:

- Discover all OpenCL platforms/devices, including CPUs and GPUs
- 2. Run a <u>micro benchmark</u> on each device, actually a short piece of real work
- 3. Load balance using micro benchmark results
- 4. Re-run micro benchmark at regular intervals to adapt to load changes



#### BUDE's heterogeneous approach

#### Between nodes:

- Partition entire ligand database (~160M) into subsets that will take a few minutes to process on one node
- 2. Distribute initial subsets across nodes
- 3. Work stealing scheduler to load balance across nodes of different performance
- 4. Time-outs trigger reallocation of MIA subsets to working nodes

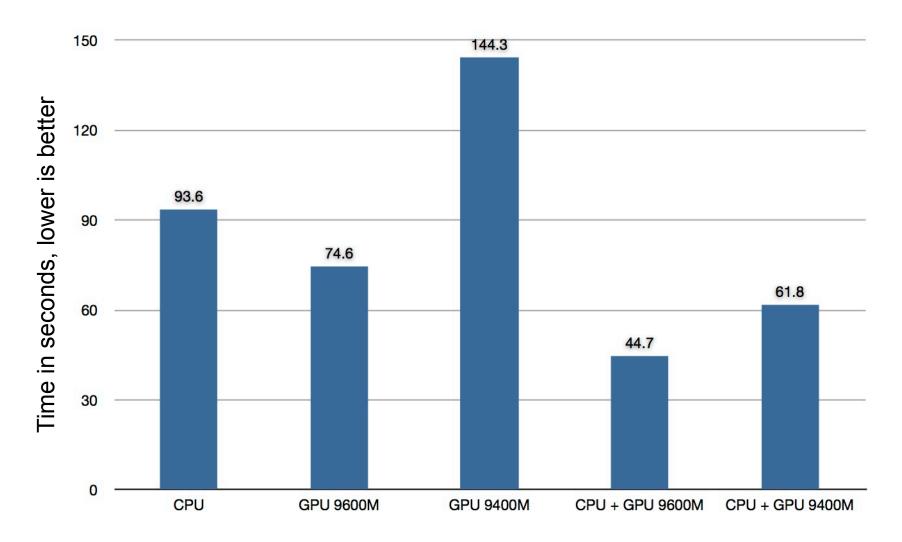


# Benchmarking methodology

- Use the same power measurement equipment for all the systems under test
- Watts Up? Pro meter
- Measures complete system power 'at the wall'
- Run as fast as possible on all available resources (i.e. all cores or all GPUs simultaneously)

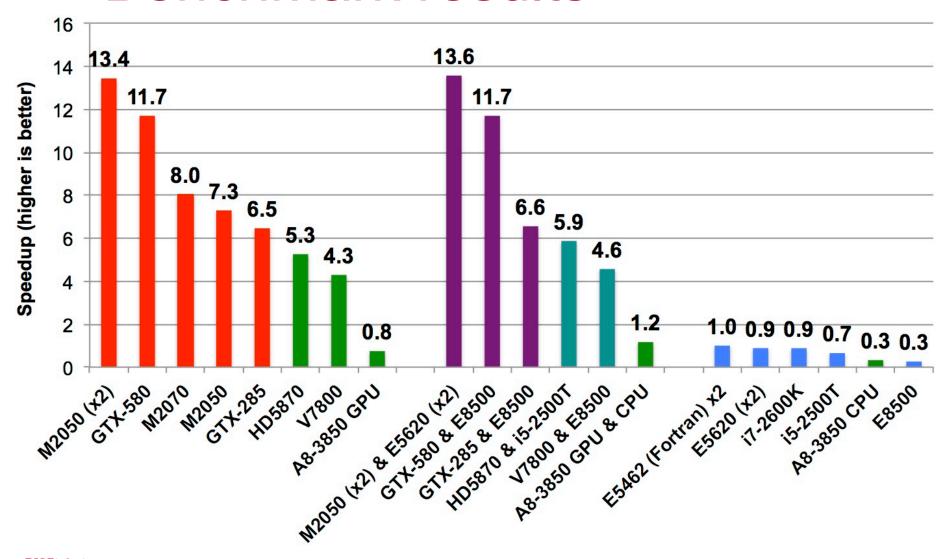


#### MacBook Pro 2009 results

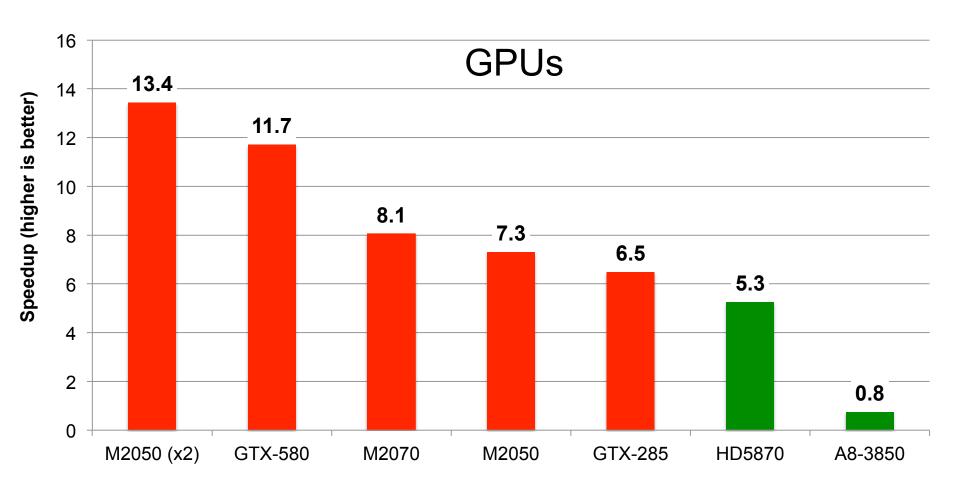




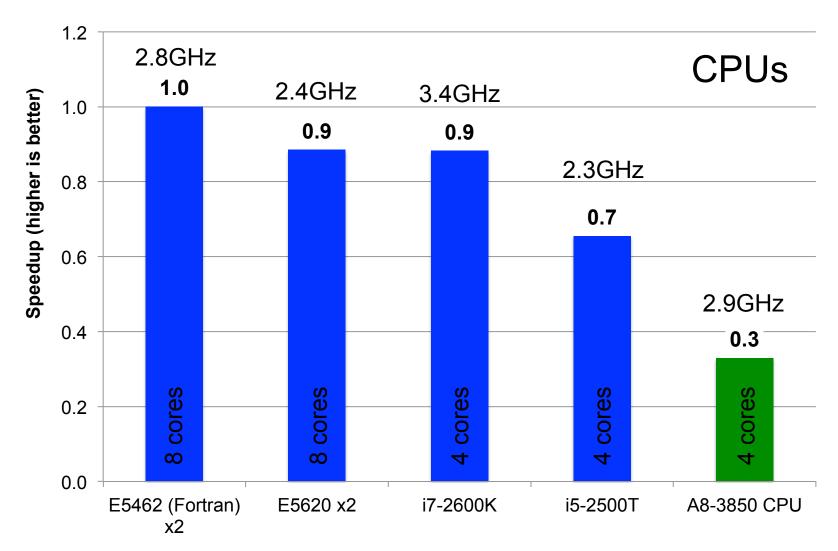
#### **Benchmark results**



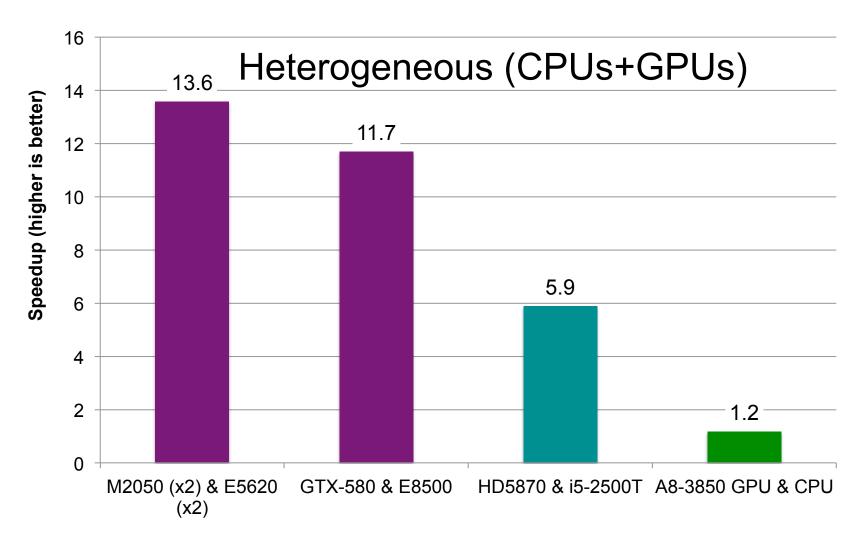




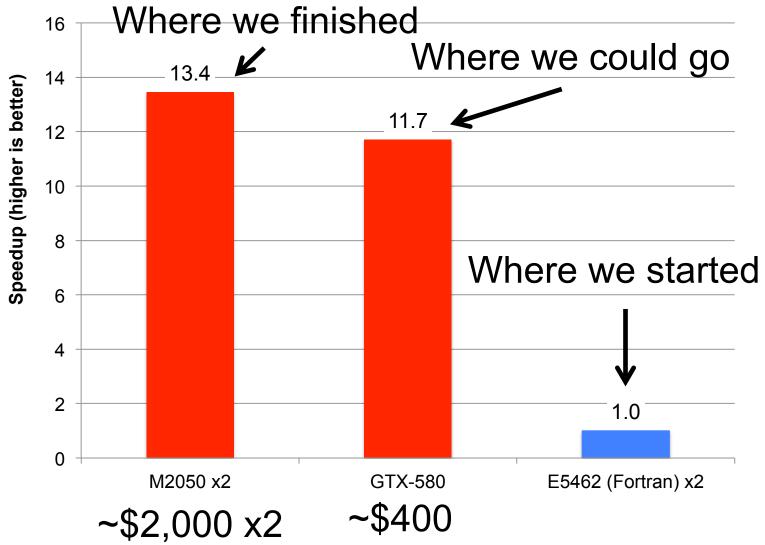






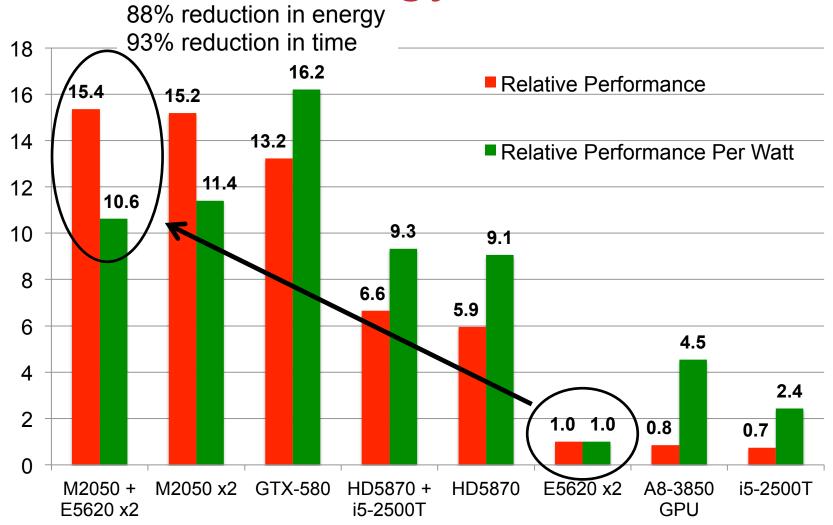








# Relative energy and run-time





Measurements are for a constant amount of work. Energy measurements are "at the wall" and include any idle components.



#### What does this let us do?







## Potentially save lives

Discovery News > Human News > Japan Detects Antibiotic-Resistant Superbug

New Delhi Metallo-betalactamase-1 (NDM-1) is an enzyme that makes bacteria resistant to antibiotics, giving rise to "superbugs"

http://news.discovery.com/ human/superbug-found-injapan.html

#### JAPAN DETECTS ANTIBIOTIC-**RESISTANT SUPERBUG**

The first case of an antibiotic-resistant "superbug" in South Asia has triggered a global health alert.

Tue Sep 7, 2010 01:03 PM ET Content provided by AFP (1) Comments | Leave a Comment



168 people like this. Be the first of vour friends.











The New Delhi metallo-lactamase-1 (NDM-1) gene has created a drug-resistant superbug. Click to enlarge this image. iStockphoto

#### THE GIST

- · A hospital north of Tokyo detected a drug-resistant
- . The bacterium has been identified as carrying the New Delhi metallo-lactamase-1 (NDM-1) gene.
- · In August a Belgian man became the first such known fatality.

Japan on Monday said it has detected its first case of an antibiotic-resistant "superbug" that surfaced in South Asia and has triggered a global health alert.

A hospital linked to the Dokkyo Medical University in Tochigi prefecture north of Tokyo detected a drugresistant "superbug," a bacterium carrying the New Delhi metallo-lactamase-1 (NDM-1) gene, in a patient last year, a hospital official told AFP.

The case follows a warning from the World Health Organization (WHO) last month calling on global health

authorities to monitor the drug-resistant superbug that is believed to have spread



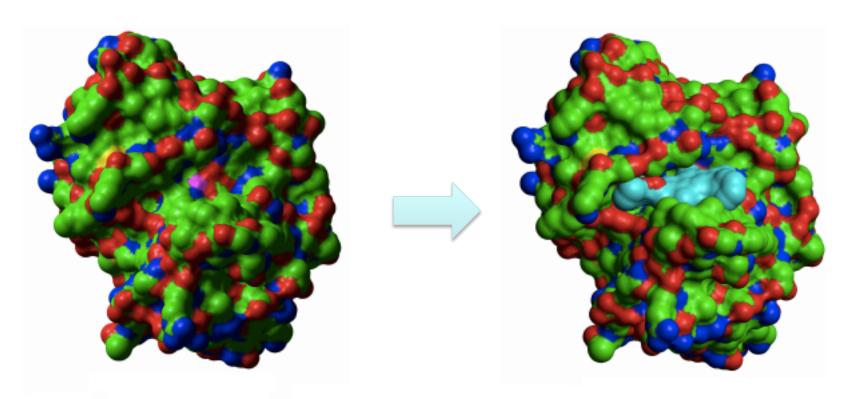
swine flu originated and what happens once it

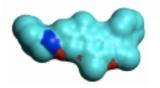
"A patient who came home from India, in his 50s, had fever symptoms while he was hospitalized in May last year, and after a blood test the hospital detected an antibiotics-resistant bacterium," the official said, adding that the patient fully

After examination doctors found that the bacterium contained the NDM-1 gene.



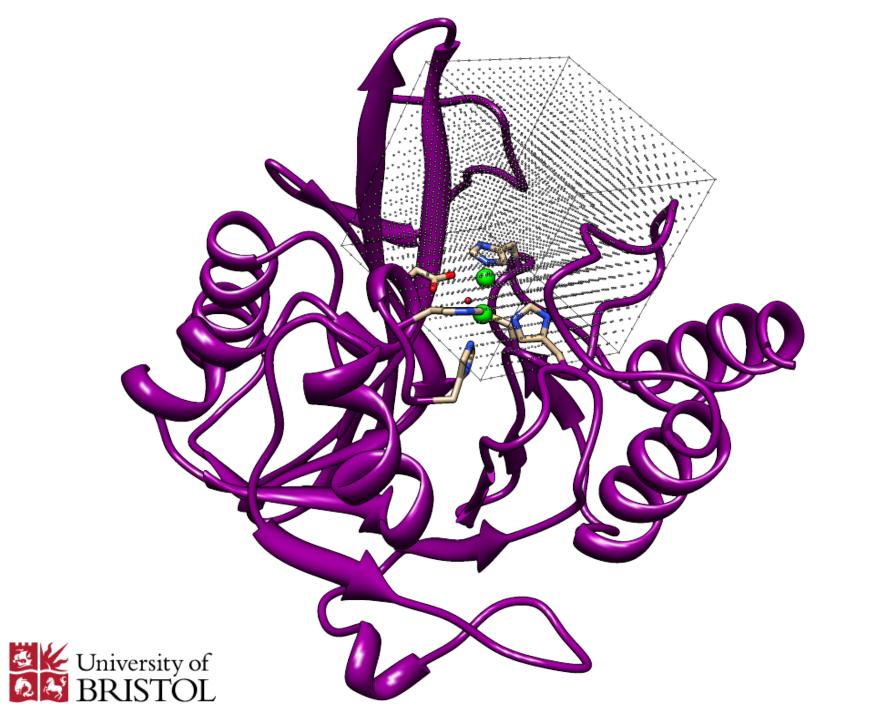
# 



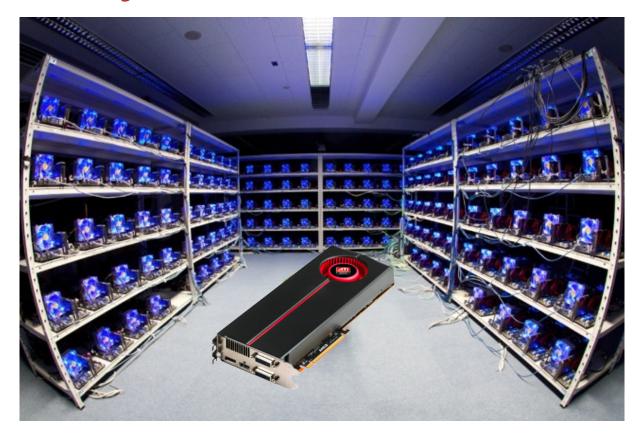


NDM-1 protein made up of 939 atoms





## **GPU-system DEGIMA**



- Used 222 GPUs in parallel for drug docking simulations
  - ATI Radeon HD5870 (2.72 TFLOPS) & Intel i5-2500T
- ~600 TFLOPS single precision peak performance
- Courtesy of Tsuyoshi Hamada and Felipe Cruz, Nagasaki



## **№ NDM-1** experiment

- 7.65 million candidate drug molecules,
   21.8 conformers each → 166.7x10<sup>6</sup>
   dockings
- 4.168 x 10<sup>12</sup> poses calculated
- ~98 hours actual wall-time
- Top 300 "hits" being analysed, down selecting to 10 compounds for further investigation in the lab





## **Future work**







#### **K** Future work

- Run on new, 372 GPU "Emerald" cluster
  - HP SL390 system, Nvidia M2090s GPUs
  - Also Gnodal fast Ethernet interconnect
- Further improve forcefield (underway)
  - RMSD already under 2Å, down from >4Å
- Port to emerging, OpenCL systems
  - E.g. ARM-based, Imagination, ...
- Benchmark on Kepler, Graphics Core Next
- Explore fault tolerance within a node



## **K** Conclusions

- OpenCL enables truly heterogeneous computing, harnessing all hardware resources in a system
- GPUs can yield significant savings in energy costs (and equipment costs)
- OpenCL can work well for multi-core CPUs as well as for GPUs

It's possible to screen libraries of millions of molecules against complex targets using highly accurate, computationally-expensive methods in hours using equipment costing O(£100K)



## **References**

- S. McIntosh-Smith, T. Wilson, A.A. Ibarra, J. Crisp and R.B. Sessions, "Benchmarking energy efficiency, power costs and carbon emissions on heterogeneous systems", The Computer Journal, September 12th 2011. DOI: 10.1093/comjnl/bxr091
- N. Gibbs, A.R. Clarke & R.B. Sessions, "Abinitio Protein Folding using Physicochemical Potentials and a Simplified Off-Lattice Model", Proteins 43:186-202,200



## **References**

- Kumarasamy, K.K., et al., Emergence of a new antibiotic resistance mechanism in India, Pakistan, and the UK: a molecular, biological, and epidemiological study. Lancet Infectious Diseases, 2010. 10(9): p. 597-602.
- Zhang, H.M. and Q. Hao, Crystal structure of NDM-1 reveals a common beta-lactam hydrolysis mechanism. Faseb Journal, 2011. 25(8): p. 2574-2582.
- Irwin, J.J. and B.K. Shoichet, ZINC A free database of commercially available compounds for virtual screening. Journal of Chemical Information and Modeling, 2005. 45(1): p. 177-182.



## **References**

 Wolfenden, R., et al., Affinities of Amino-Acid Side-Chains for Solvent Water.
 Biochemistry, 1981. 20(4): p. 849-855.

 Gibbs, N., A.R. Clarke, and R.B. Sessions, Ab initio protein structure prediction using physicochemical potentials and a simplified off-lattice model. Proteins, 2001. 43(2): p. 186-202.

