



Quantum Chemistry on GPGPU

Marcin Makowski,^{1,2} Grzegorz Mazur,^{1,2} Piotr Kuźniarowicz,^{1,3}
Roman Łazarski,^{1,4} Alina Mreńca,⁵ Krzysztof Kolasiński⁵

¹Faculty of Chemistry, Jagiellonian University

²Academic Computer Centre CYFRONET AGH

³Department of Molecular and Material Sciences, Kyushu University

⁴Institute of Material Science and Material Technology, F. Schiller University

⁵Faculty of Physics and Computer Science, AGH University of Science and
Technology

KU KDM 2013



Plan

Introduction

GPGPU

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POWIEW

HPC Infrastructure for Grand Challenges of Science and Engineering

- ▶ ICM, CYFRONET, PCSS
- ▶ Acquisition of modern HPC architectures allowing for large-scale calculations
- ▶ Research groups working optimal computational models for HPC



POWIEW in CYFRONET

- ▶ Hardware
 - ▶ GPGPU
 - ▶ vSMP
 - ▶ cluster of fat nodes
- ▶ Software
- ▶ Software development
- ▶ Workshops



Computational Chemistry

ZEUS

- ▶ 34% of grid workload
- ▶ 56% of local workload





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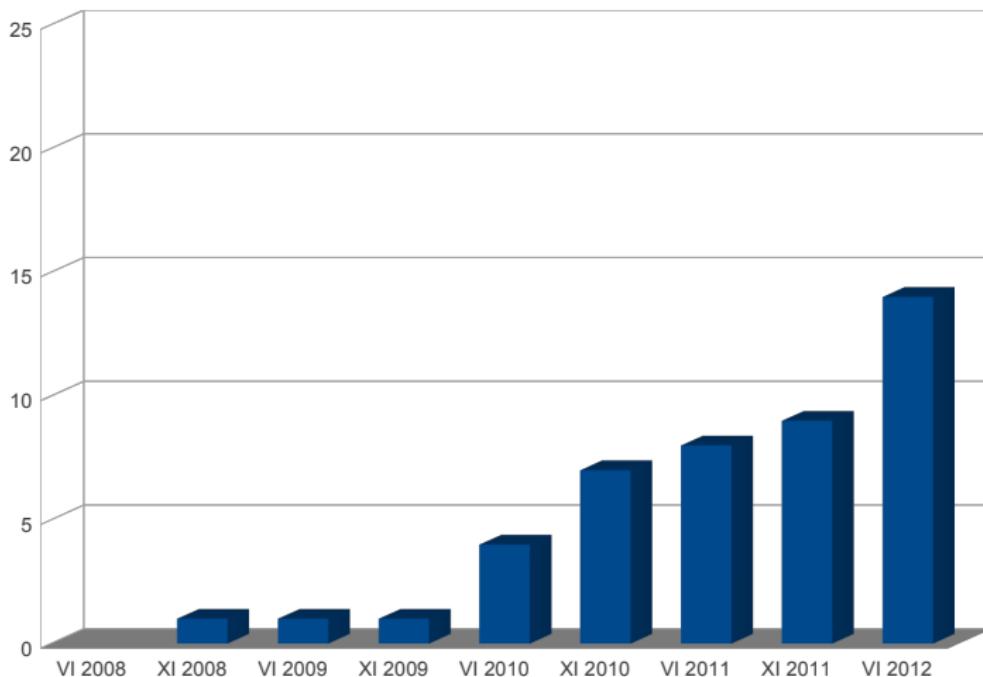
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GPGPU



GPGPU in TOP100





GPGPU in CYFRONET

- ▶ Hardware
 - ▶ 48 NVIDIA Tesla M2050 GPUs (24 servers)
 - ▶ 160 NVIDIA Tesla M2090 GPUs (20 servers)
 - ▶ almost 140 TFlops
- ▶ Software
 - ▶ TeraChem (quantum chemistry)
 - ▶ NAMD (molecular mechanics)
- ▶ Software development



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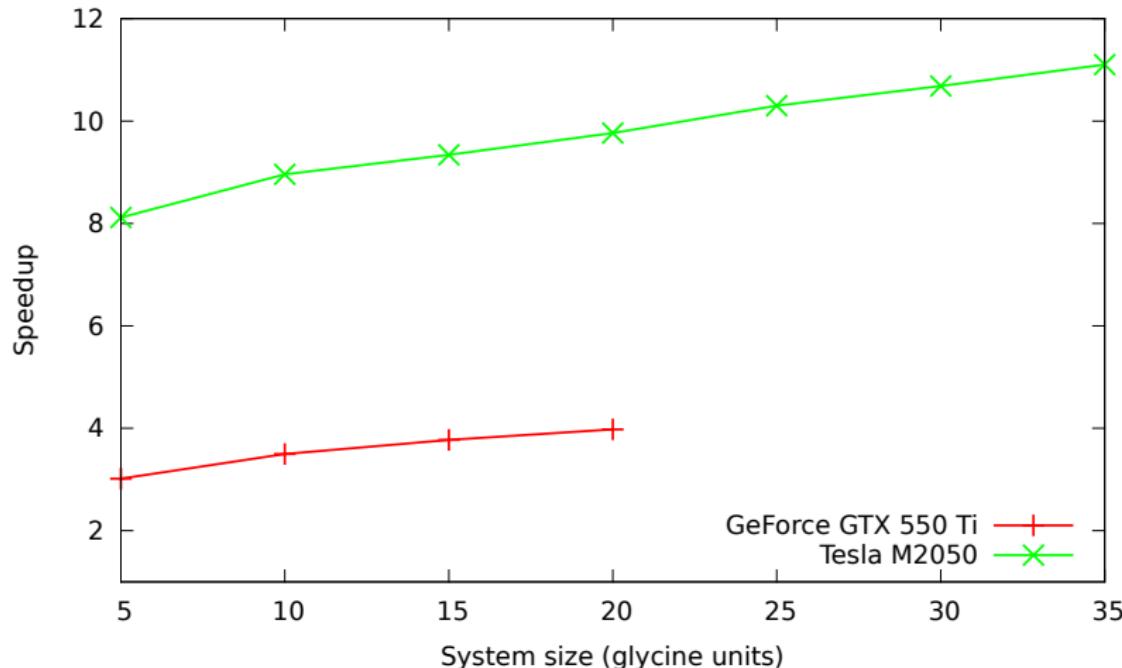
- ▶ General-purpose computational quantum-chemistry package
- ▶ Created and developed at Faculty of Chemistry (Jagiellonian University) in collaboration with CYFRONET
- ▶ Optimal implementation of existing methods
- ▶ Development of original, efficient methodologies
- ▶ Aiming at large systems
 - ▶ advanced hybrid parallelization
 - ▶ GPGPU computing



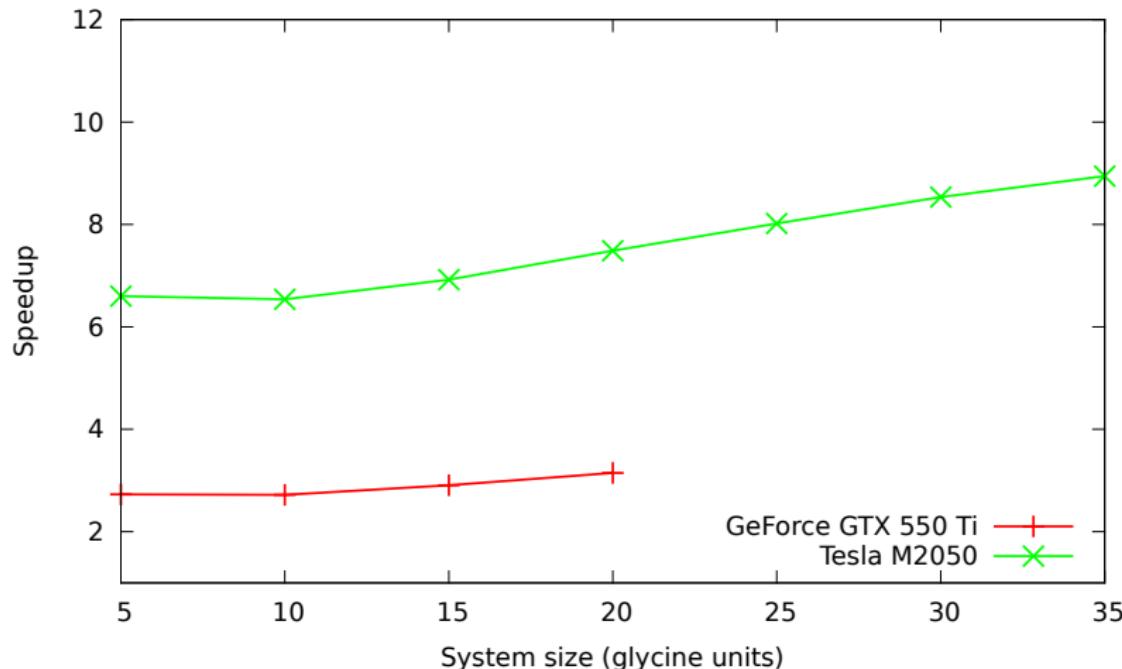
Testbed

- ▶ Hardware
- ▶ CPU: Intel Xeon X5670
- ▶ GPUs: Tesla M2050, GeForce 550 Ti
- ▶ Software
 - ▶ niedoida
 - ▶ TeraChem (ver. 1.50K)
- ▶ Molecular systems
 - ▶ polyglycine chains
 - ▶ 6-31G basis set
 - ▶ RHF and B3LYP calculations

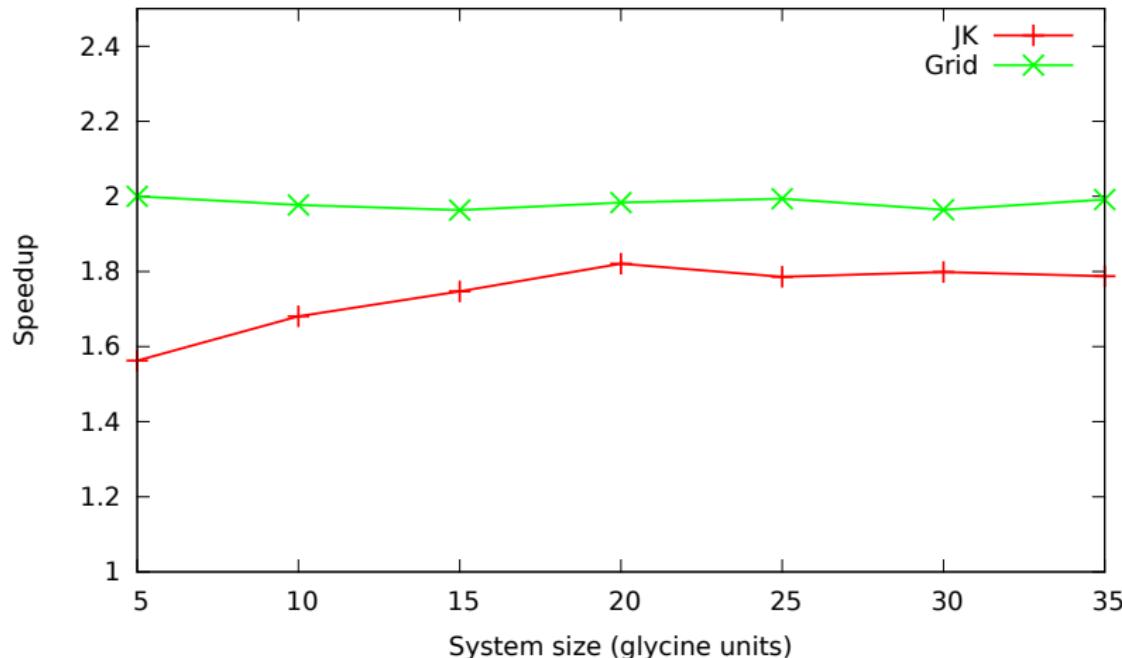
GPU vs CPU: Hartree-Fock



GPU vs CPU: B3LYP



Multi-GPU efficiency





Method	Speedup	
	vs CPU	vs TeraChem
HF	11.10	0.30
B3LYP	8.94	0.37
grid integration	2.33	0.69



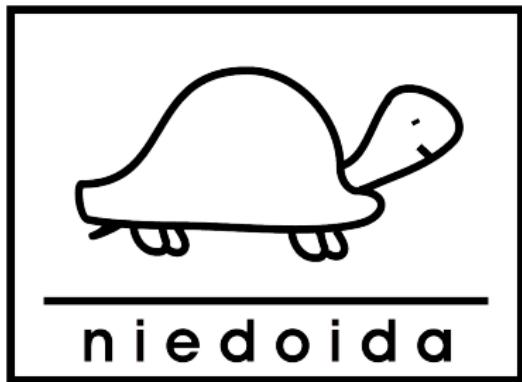
Future

Projects in progress/planned:

- ▶ integration for d- and f-type basis functions
- ▶ RI-DFT and RI-MP2
- ▶ CIS and TD-DFT excited states calculations
- ▶ gradients and second derivatives:
 - ▶ geometry optimization
 - ▶ Born-Oppenheimer molecular dynamics
 - ▶ vibrations
- ▶ hybrid parallelization



To Be Continued...



This work was carried out thanks to the
"POWIEW" POIG.02.03.00-00-018/08-00 project.
This project is co-funded by the European Regional Development Fund
(ERDF) as a part of the Innovative Economy program.