

StarPU a runtime system for Scheduling Tasks, or How to get portable performance on accelerator-based platforms without the agonizing pain

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Once upon a time in computer architecture ...



Introduction

Programming Accelerator-based machines

- Prehistory (<2007)

- SMP machines (1960s !)
- NUMA architectures (1970s)
- Vector machines (1980s-90s)
- Multicore chips (2000s)

- GPGPU for masses ? (from 2007)

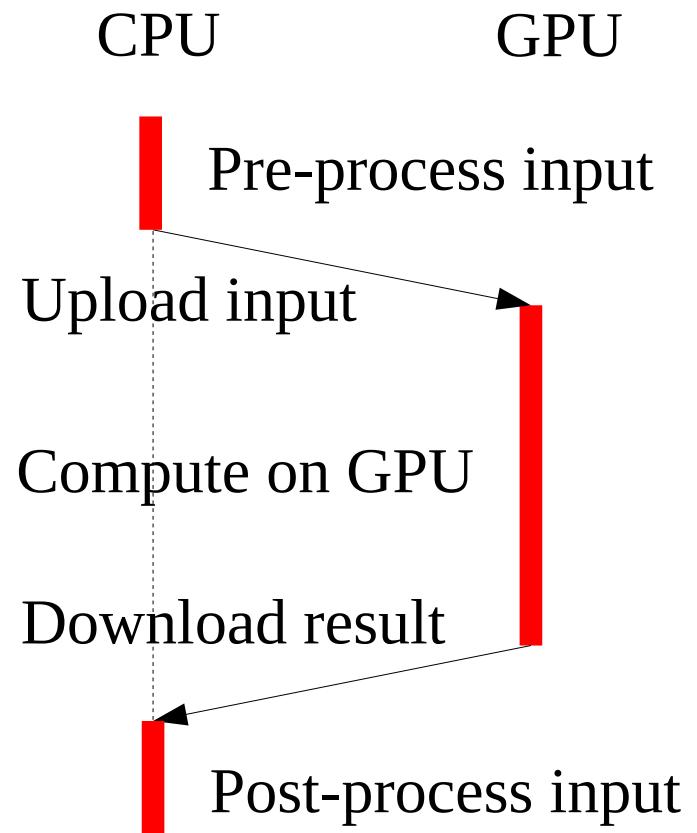
- CPUs are now deprecated ?
- Rewrite all codes for accelerators
- Pure offloading model



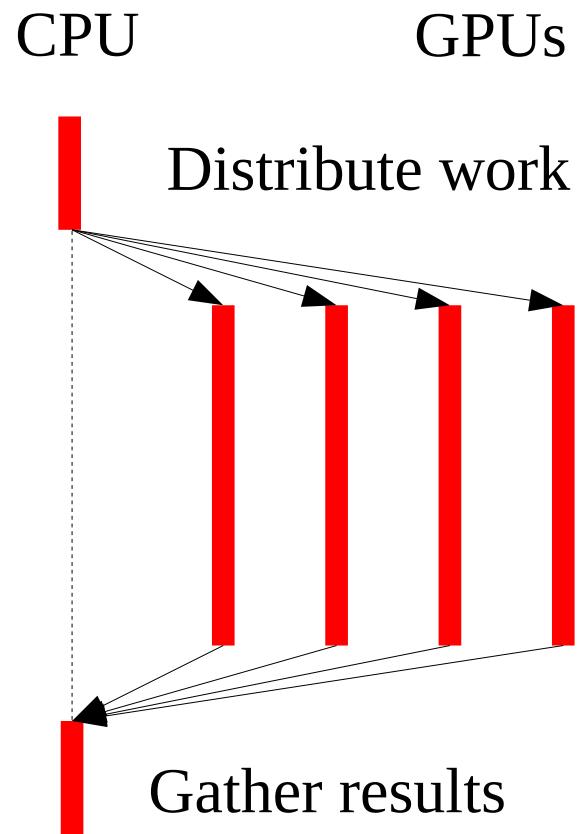
Introduction⁴

Programming Accelerator-based machines

- Pure offloading model
 - CUDA 1.x (2007 – 2008)
 - Synchronous cards
 - Ignore CPUs
 - Concentrate on efficient kernels
 - Complex memory accesses
 - CUDA heros (eg. V.Volkov)
 - Port standard libraries to CUDA
 - CUBLAS
 - CUFFT
 - GPUCV...



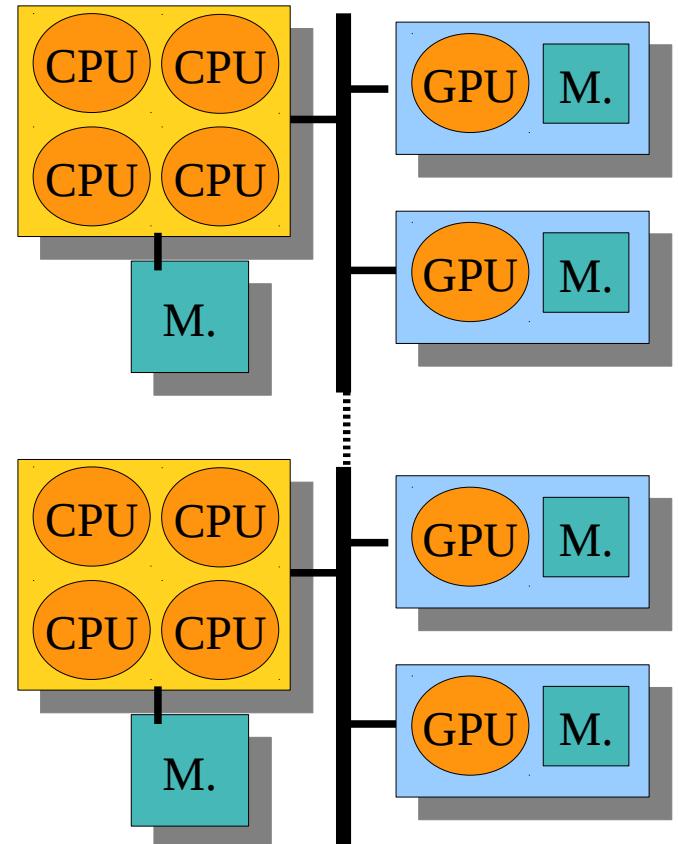
- Multi-GPU era
 - CUDA 2.x (2008 – 2009)
 - Asynchronous transfers
 - S1070 servers
 - Still Ignore CPUs (in general)
 - Suitable for regular applications
 - Massively parallel problems
 - Use previously written kernels
 - New problems
 - Parallel programming for real
 - PCI bus = bottleneck
 - Pre/Post-processing is costly



Introduction

Programming Accelerator-based machines

- GPU computing era
 - CUDA 3.x (2009 – ?)
 - End of GPGPU
 - Hybrid machines
 - Tightly coupled CPUs and GPUs
 - Take advantage of all resources
 - MUCH more complicated
 - Load balancing
 - Who does what ?
 - Heterogeneous capabilities
 - Data management
 - Numerous data transfers
 - Fully asynchronous model



Introduction

Challenging issues at all stages

- Applications

- Programming paradigm
- BLAS kernels, FFT, ...

HPC Applications

- Compilers

- Languages
- Code generation/optimization

Compiling environment

Specific libraries

- Runtime systems

- Resources management
- Task scheduling

Runtime system

Operating System

- Architecture

- Memory interconnect

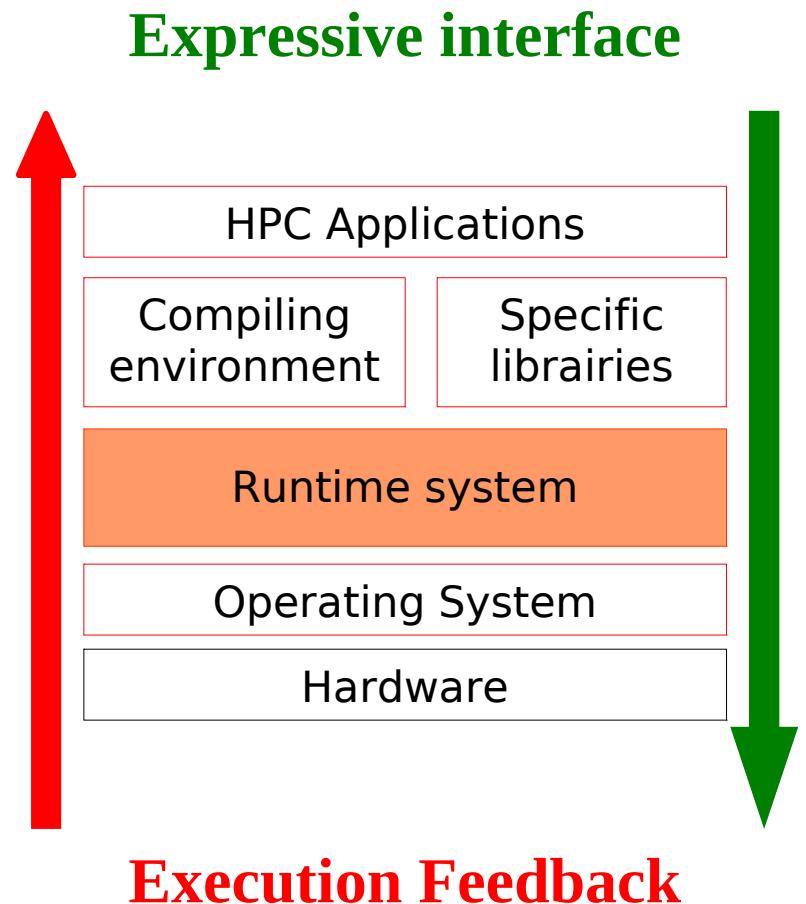
Hardware



Introduction

Challenging issues at all stages

- Applications
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 - BLAS kernels, FFT, ...
- Compilers
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Outline

- The StarPU runtime system
- Task Scheduling
 - Load balancing
 - Improving data locality
- Evaluation on dense linear algebra algorithms
 - Synthetic “LU” decomposition
 - Mixing PLASMA and MAGMA (Cholesky & QR)
- Scheduling parallel tasks
- Adding support for MPI in StarPU



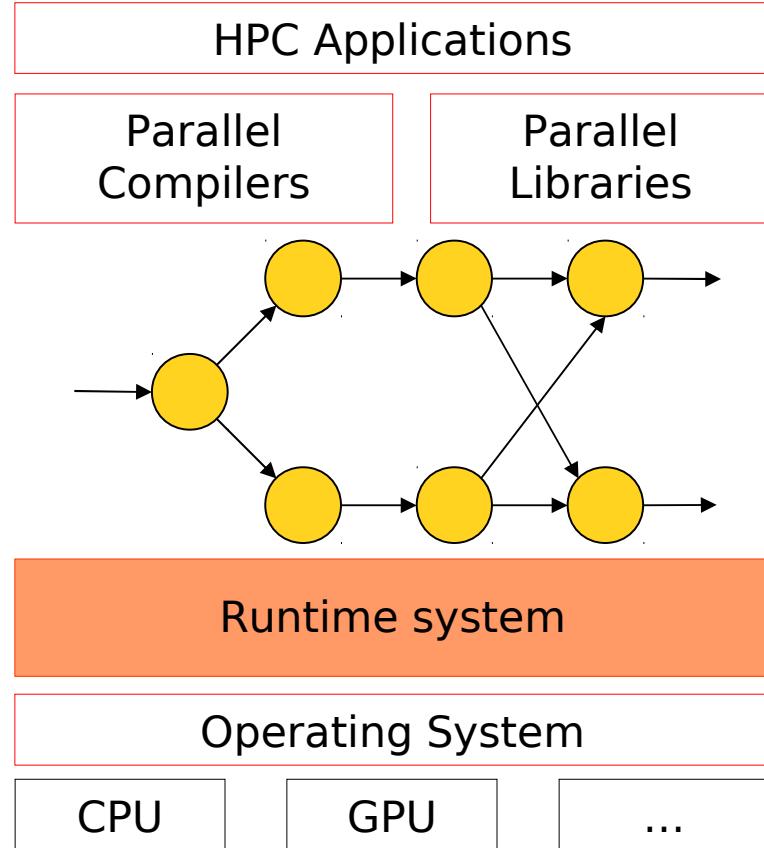
The StarPU runtime system



The StarPU runtime system

The need for runtime systems

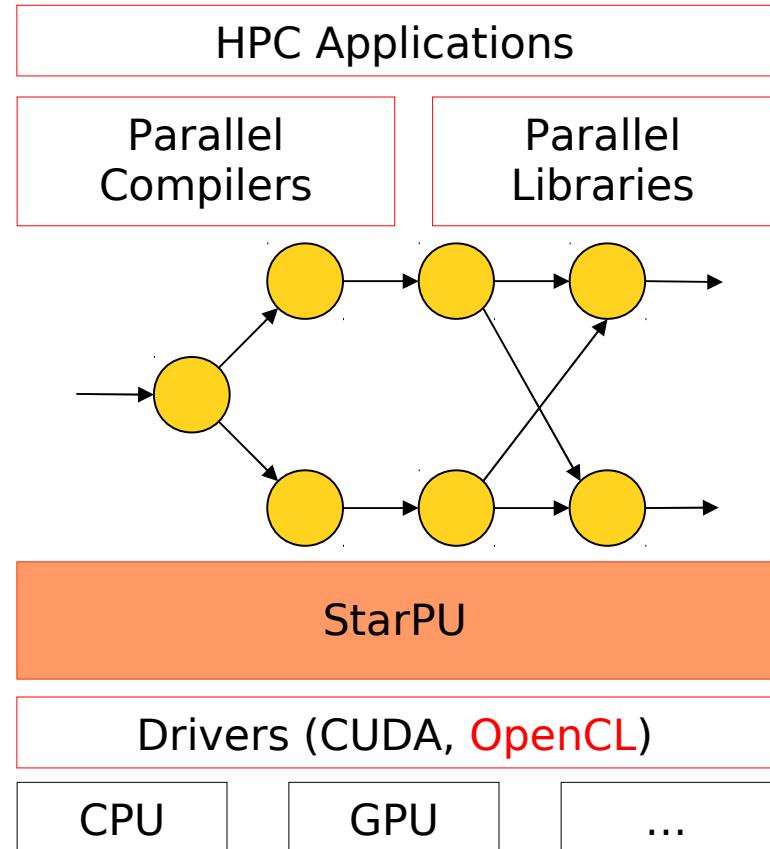
- “do dynamically what can’t be done statically anymore”
- Library that provides
 - Task scheduling
 - Memory management
- Compilers and libraries generate (graphs of) parallel tasks
 - Additional information is welcome!



The StarPU runtime system

Data management library

- StarPU provides a **Virtual Shared Memory** subsystem
 - Weak consistency
 - Replication
 - Single writer
 - High level API
 - Partitioning filters



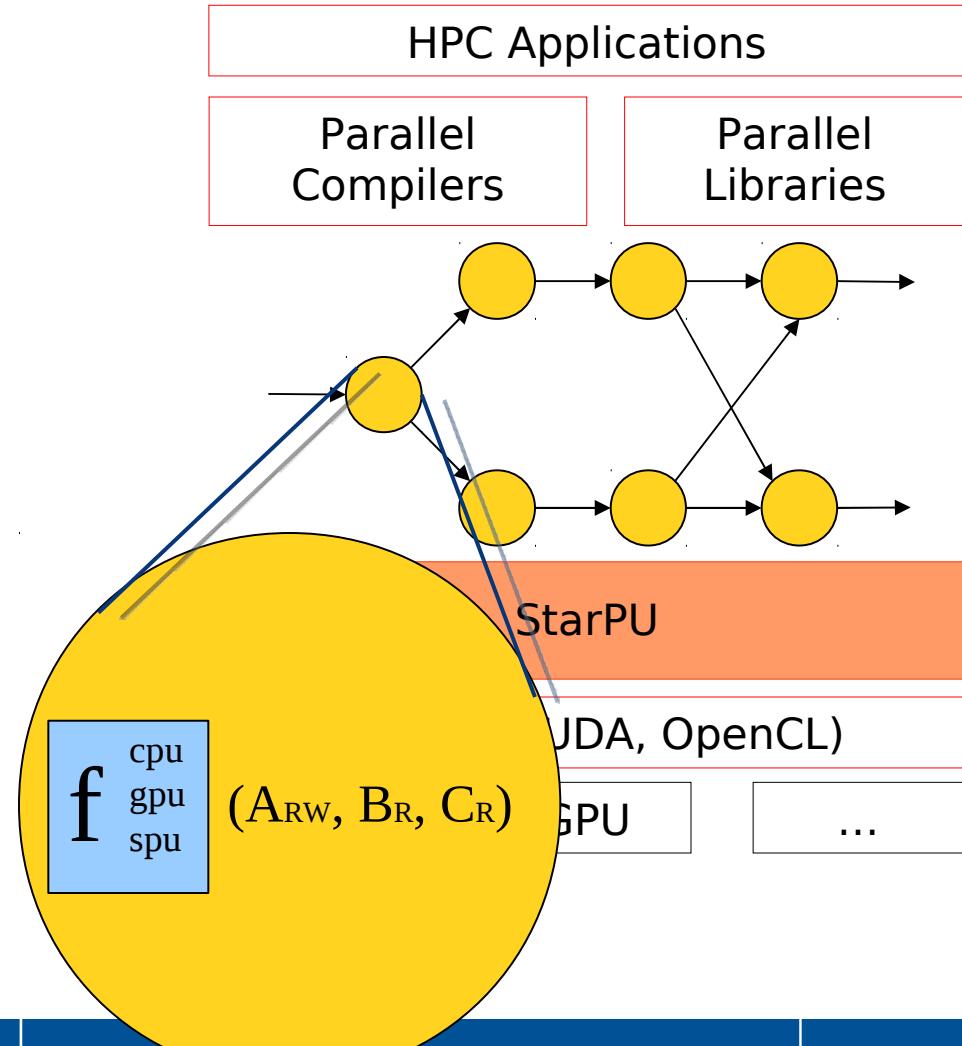
- Input & output of tasks = reference to VSM data



The StarPU runtime system

Task scheduling

- Tasks =
 - Data input & output
 - Reference to VSM data
 - Multiple implementations
 - E.g. CUDA + CPU implementation
 - Dependencies with other tasks
 - Scheduling hints



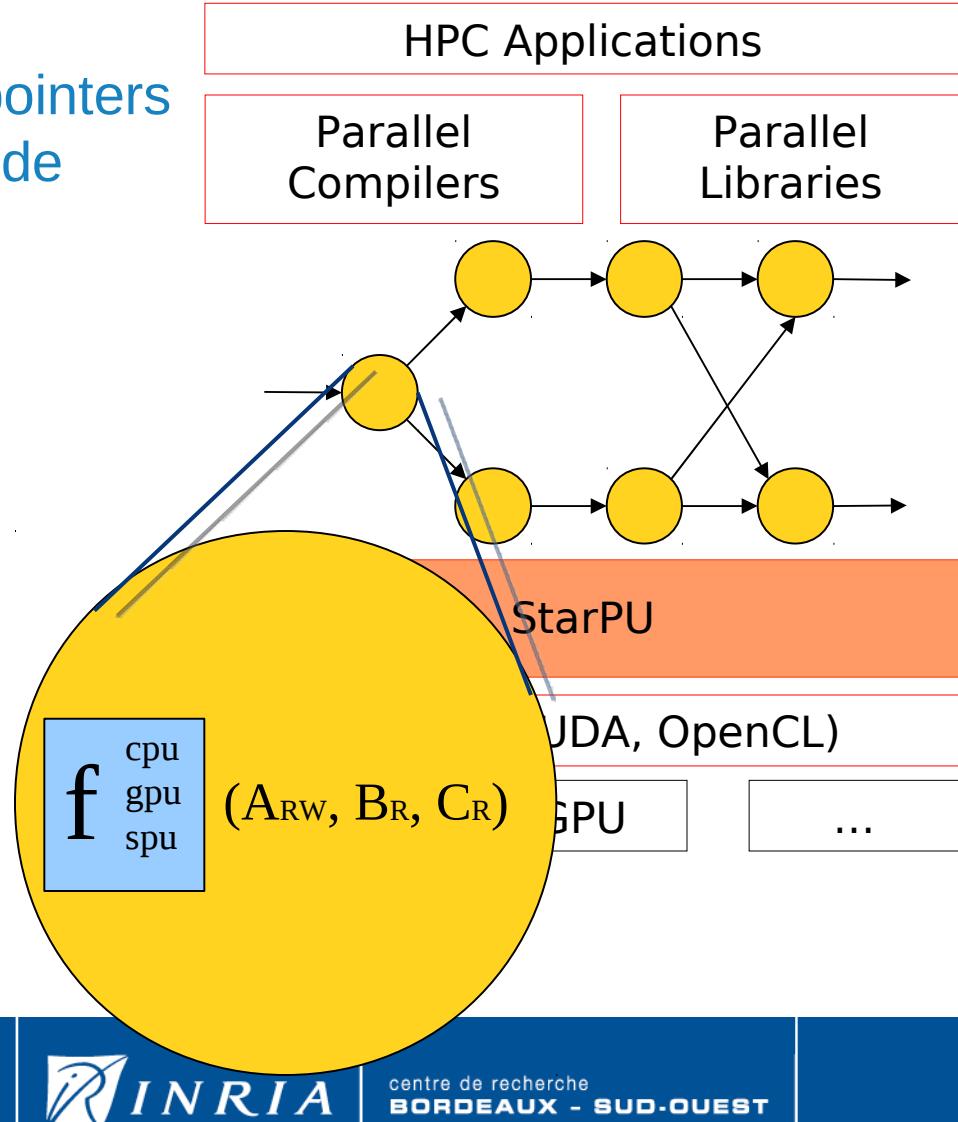
- StarPU provides an **Open Scheduling platform**
 - Scheduling algorithm = plug-ins



The StarPU runtime system

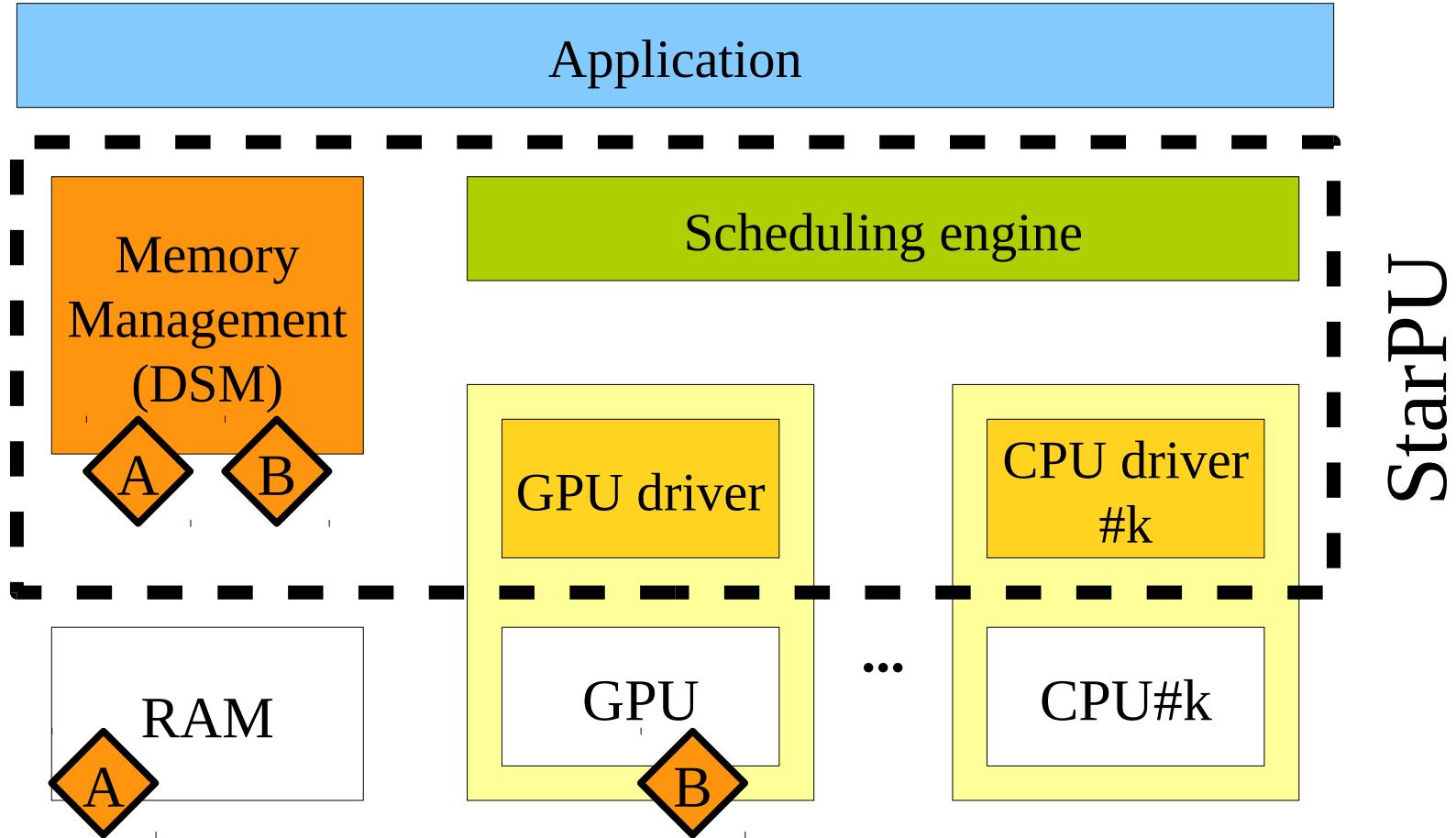
Task scheduling

- Who generates the code ?
 - StarPU Task = ~function pointers
 - StarPU don't generates code
- Programming heros ?
- Libraries era
 - PLASMA + MAGMA
 - FFTW + CUFFT...
- Rely on compilers
 - PGI accelerators
 - CAPS HMPP...



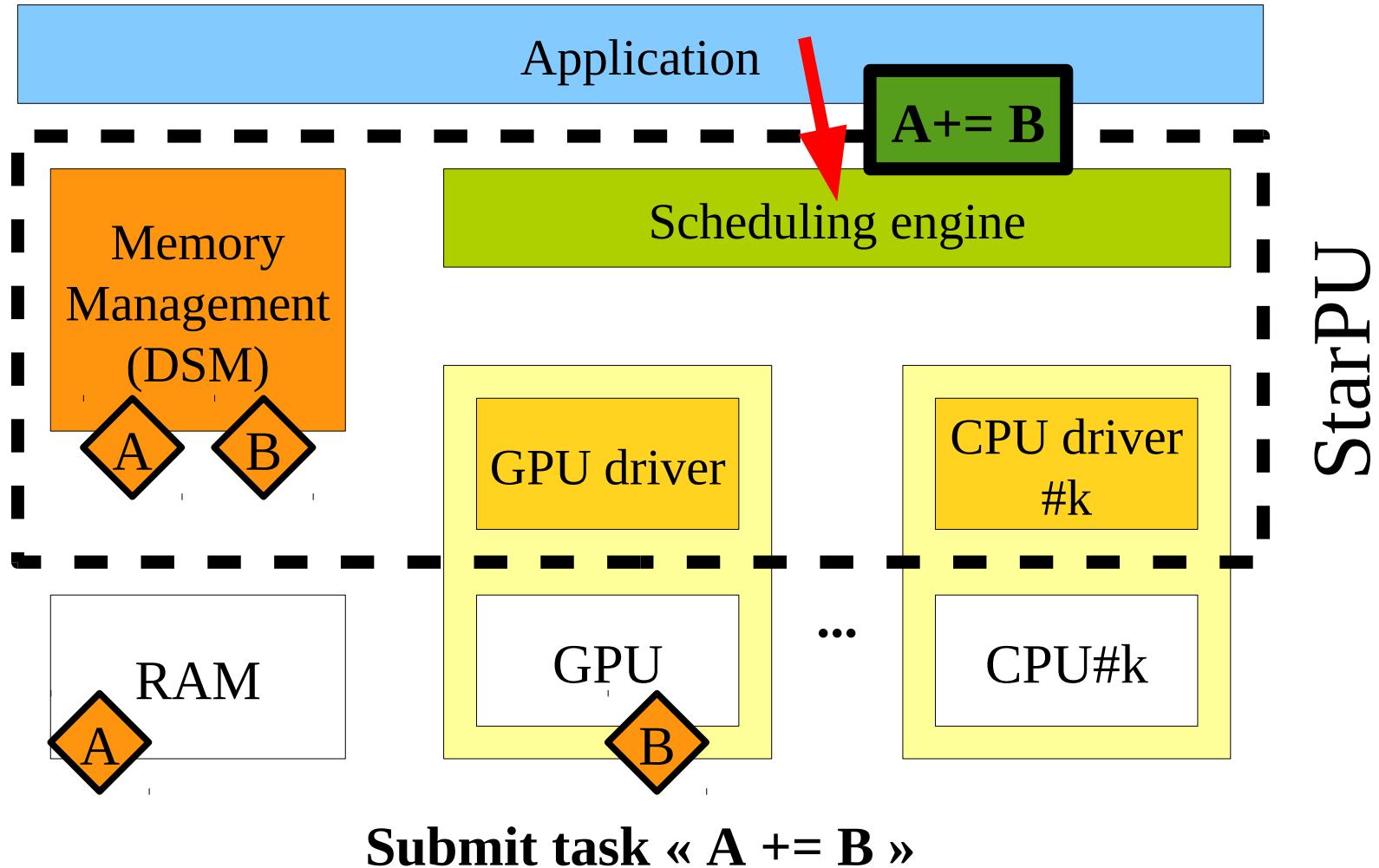
The StarPU runtime system

Execution model



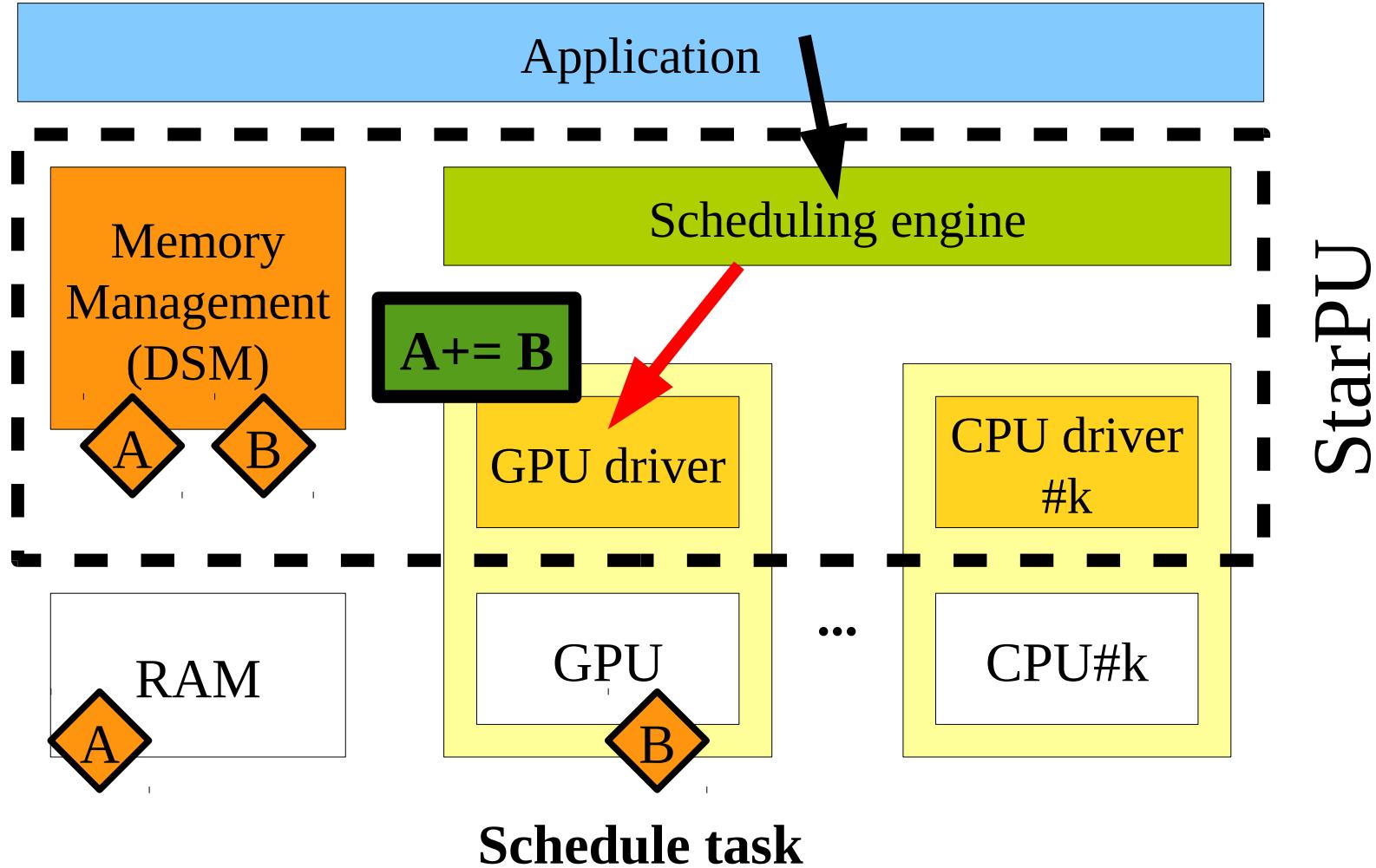
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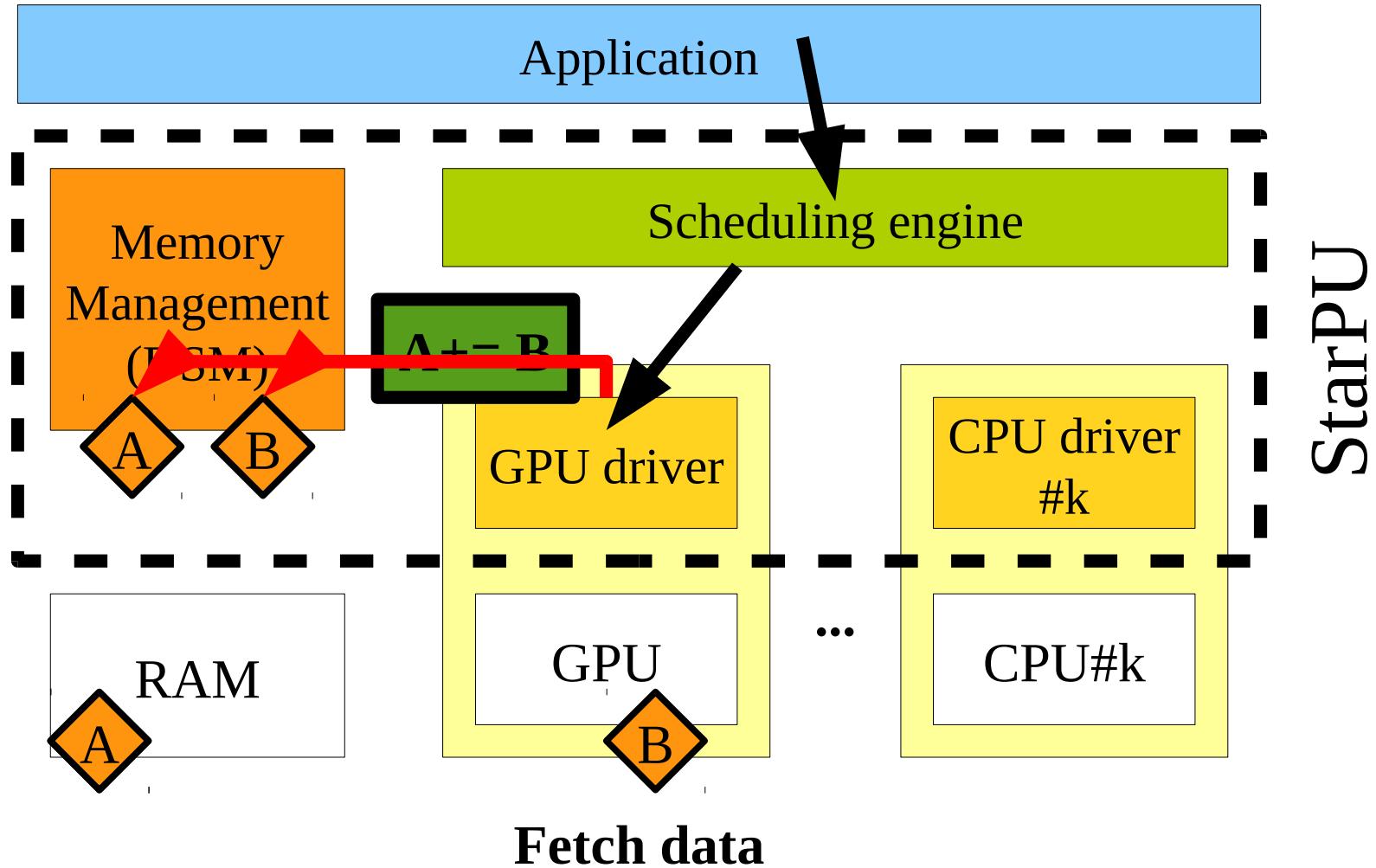
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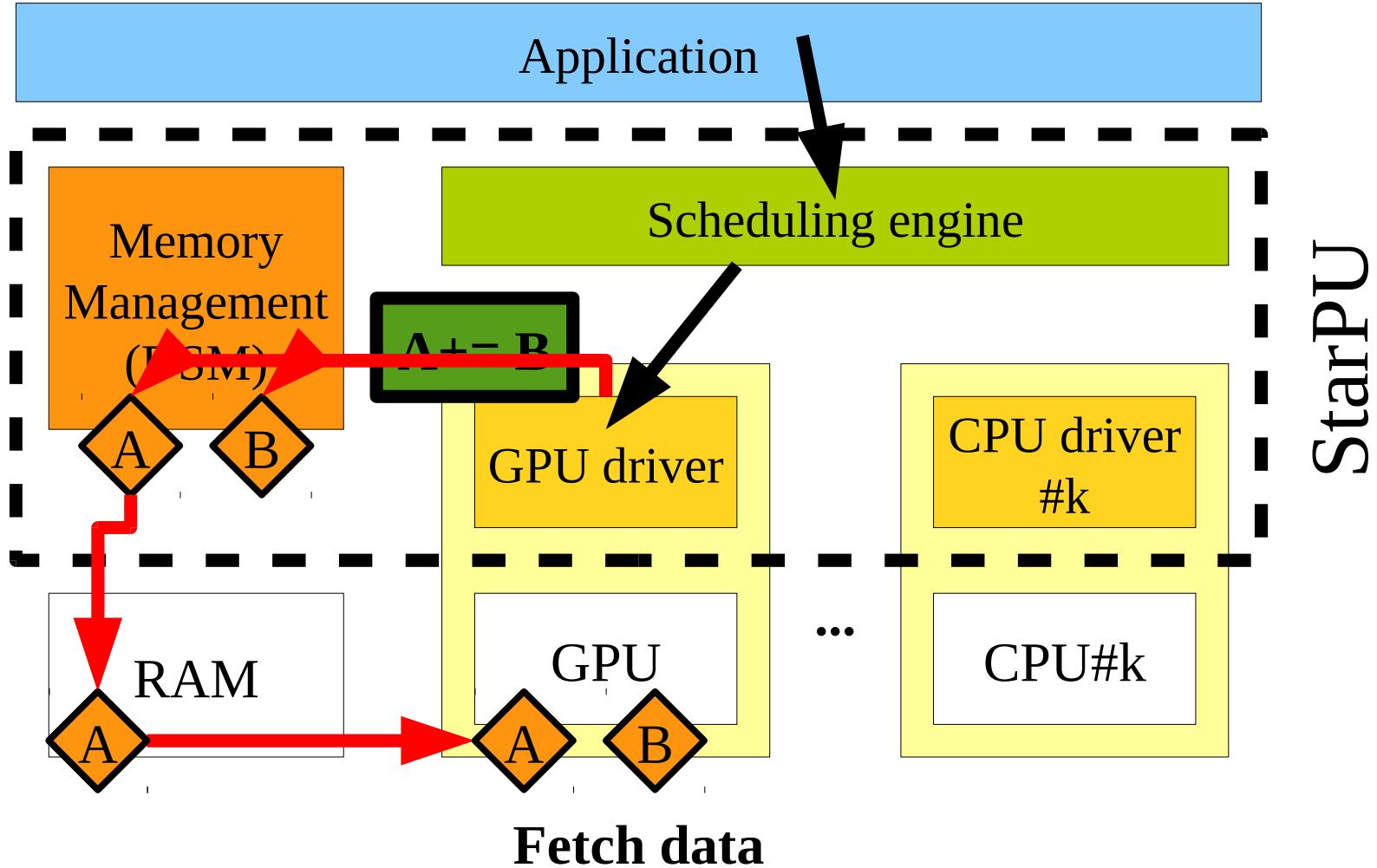
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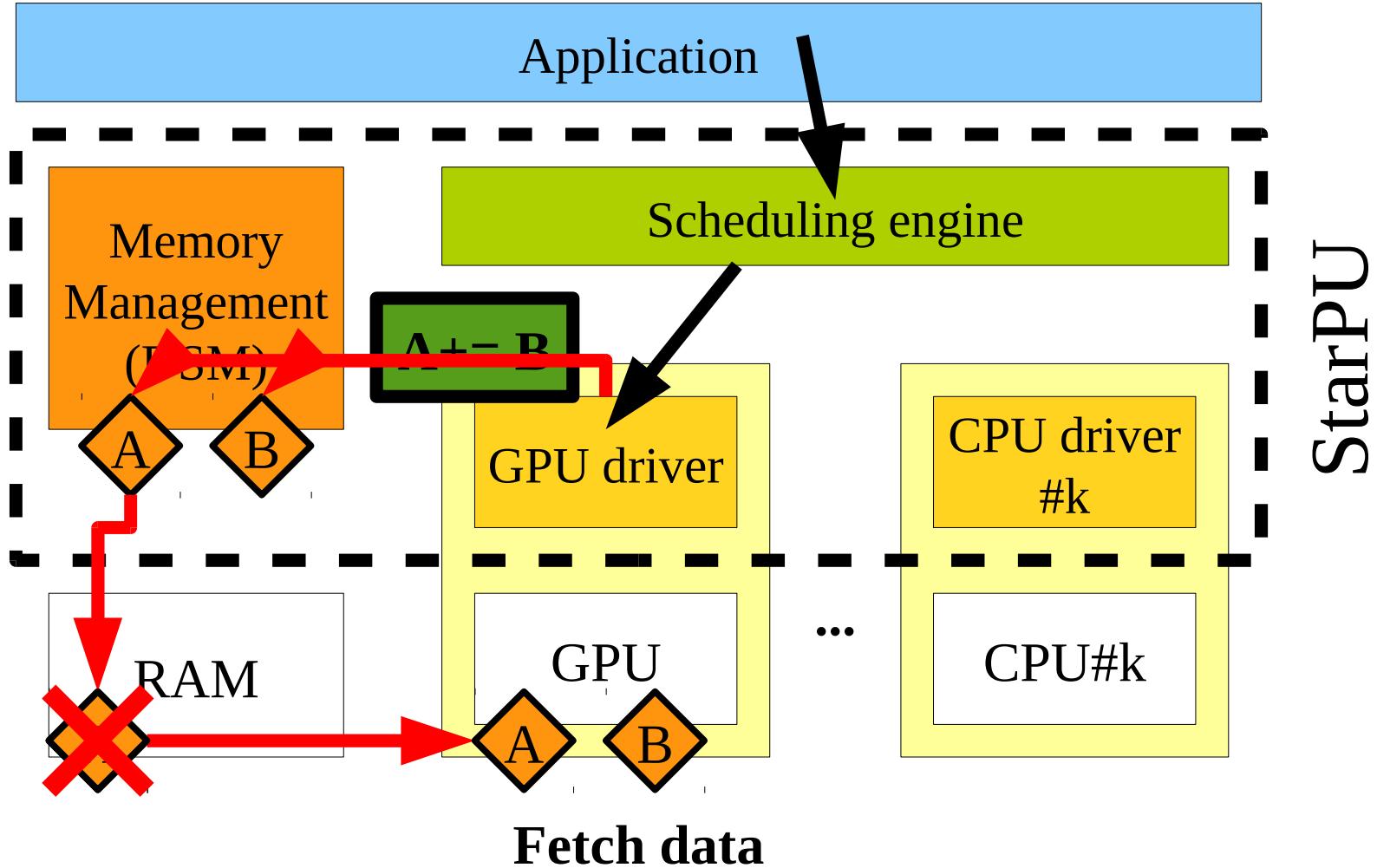
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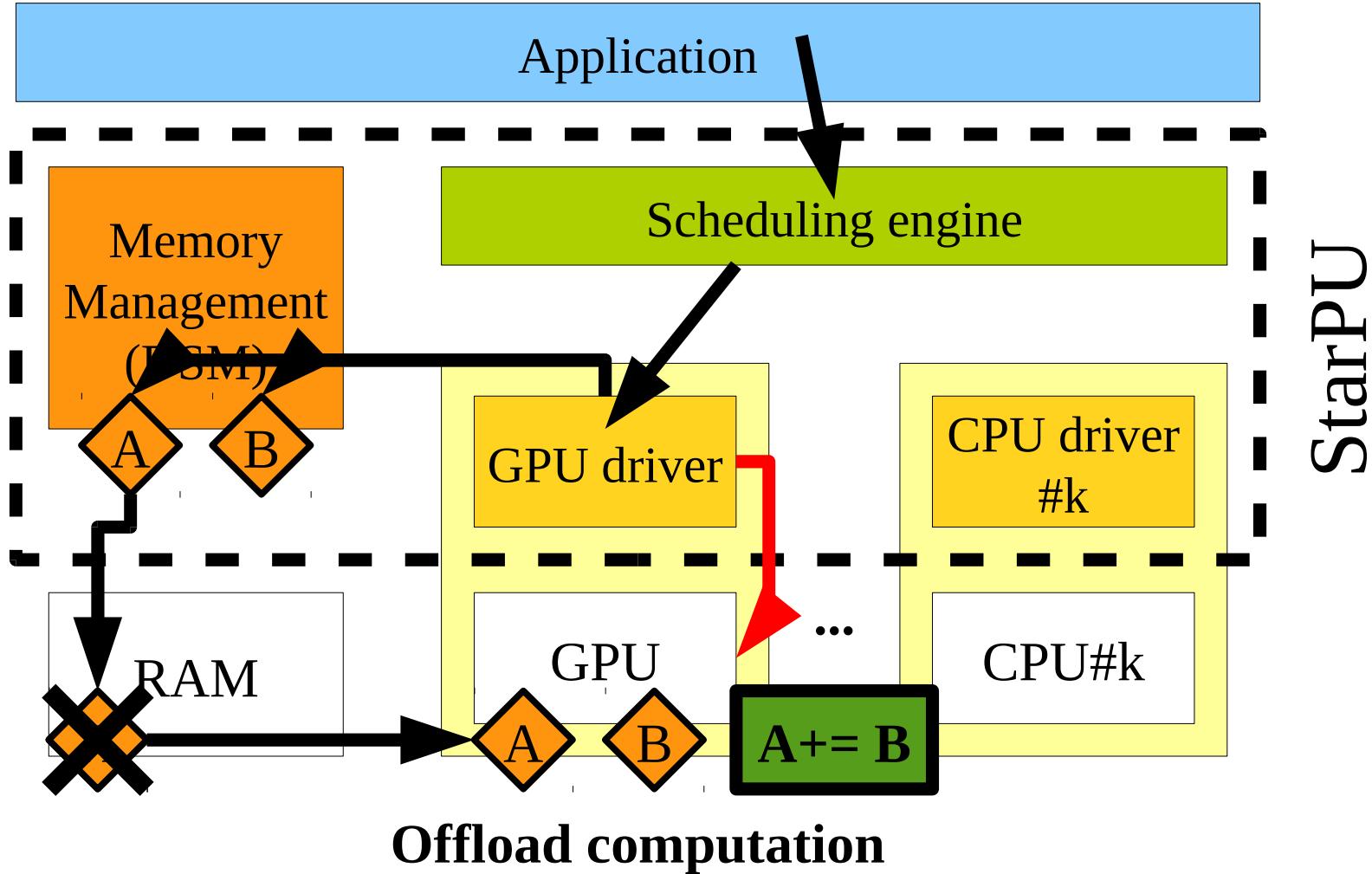
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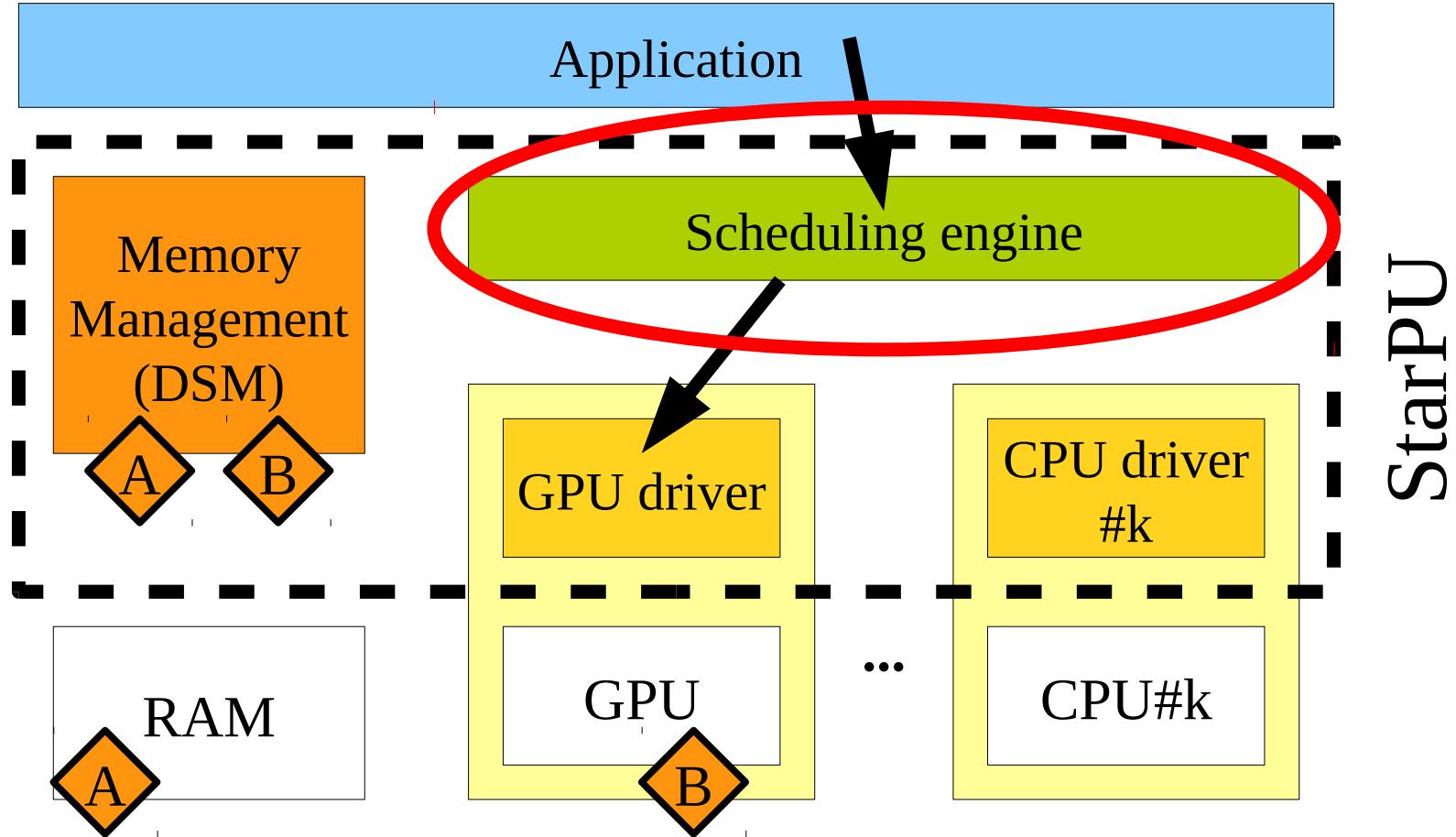
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The StarPU runtime system

Execution model



Task Scheduling



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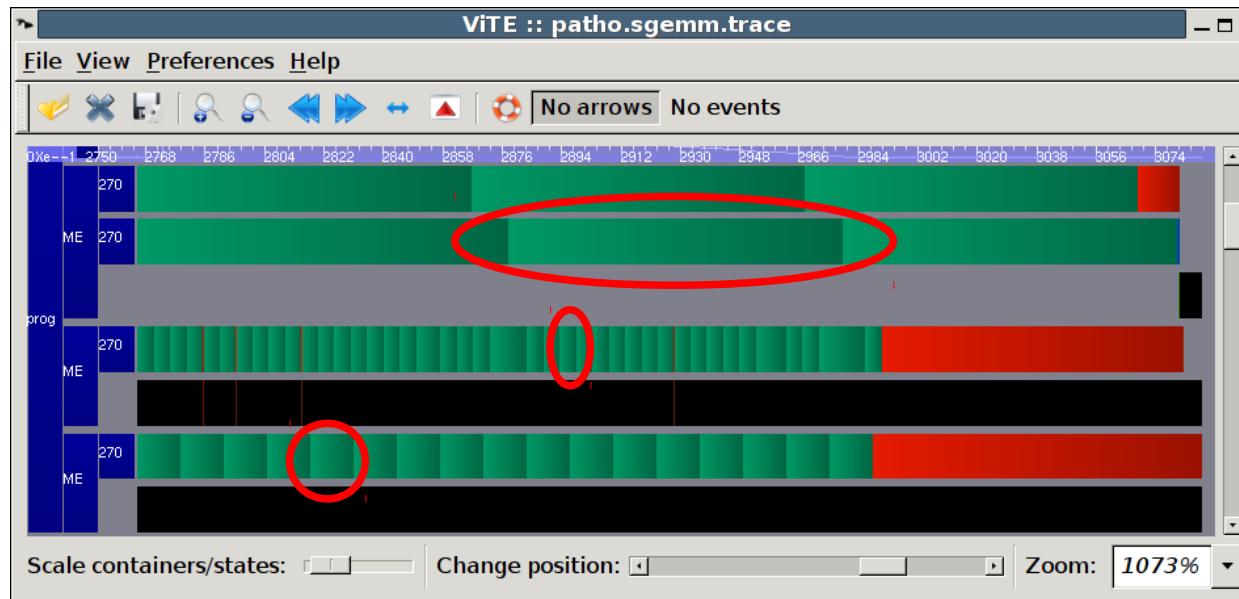
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Why do we need task scheduling ?

Blocked Matrix multiplication

Things can go (really) wrong even on trivial problems !

- Static mapping ?
 - Not portable, too hard for real-life problems
- Need Dynamic Task Scheduling
 - Performance models



2 Xeon cores

Quadro FX5800

Quadro FX4600



Predicting task duration

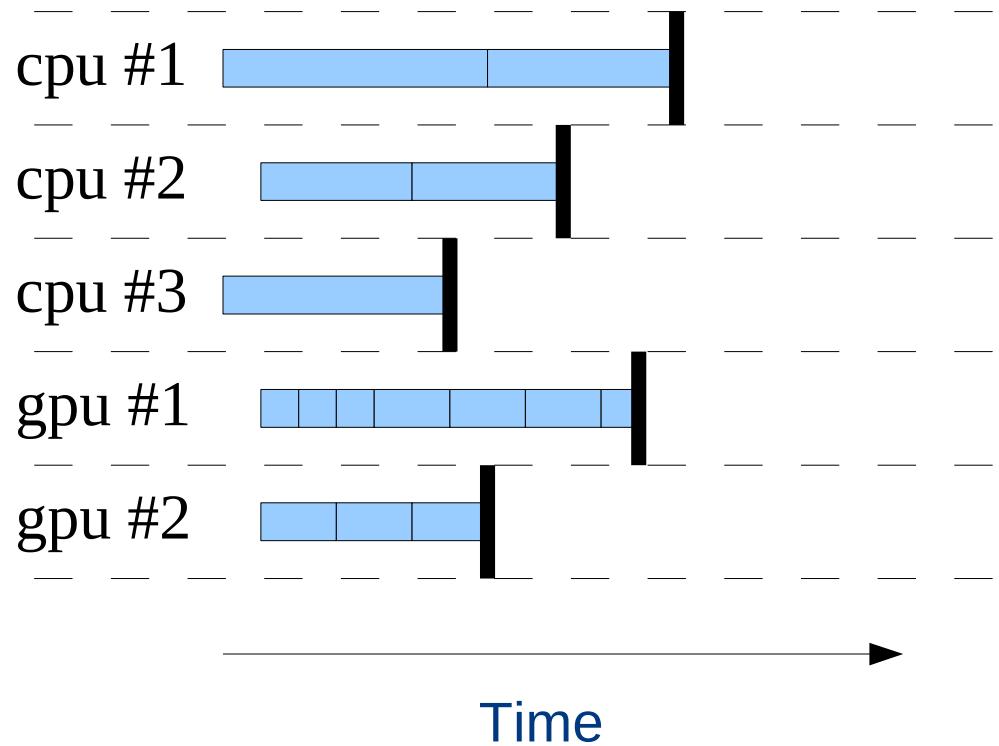
Load balancing

- Task completion time estimation

- History-based
- User-defined cost function
- Parametric cost model

- Can be used to improve scheduling

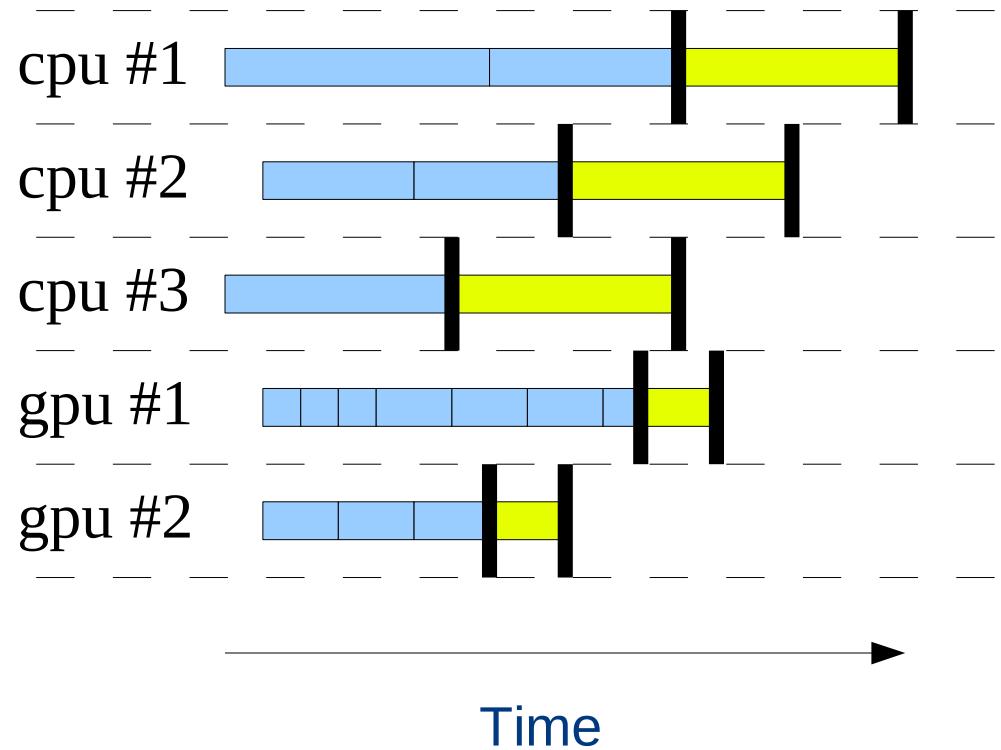
- E.g. Heterogeneous Earliest Finish Time



Predicting task duration

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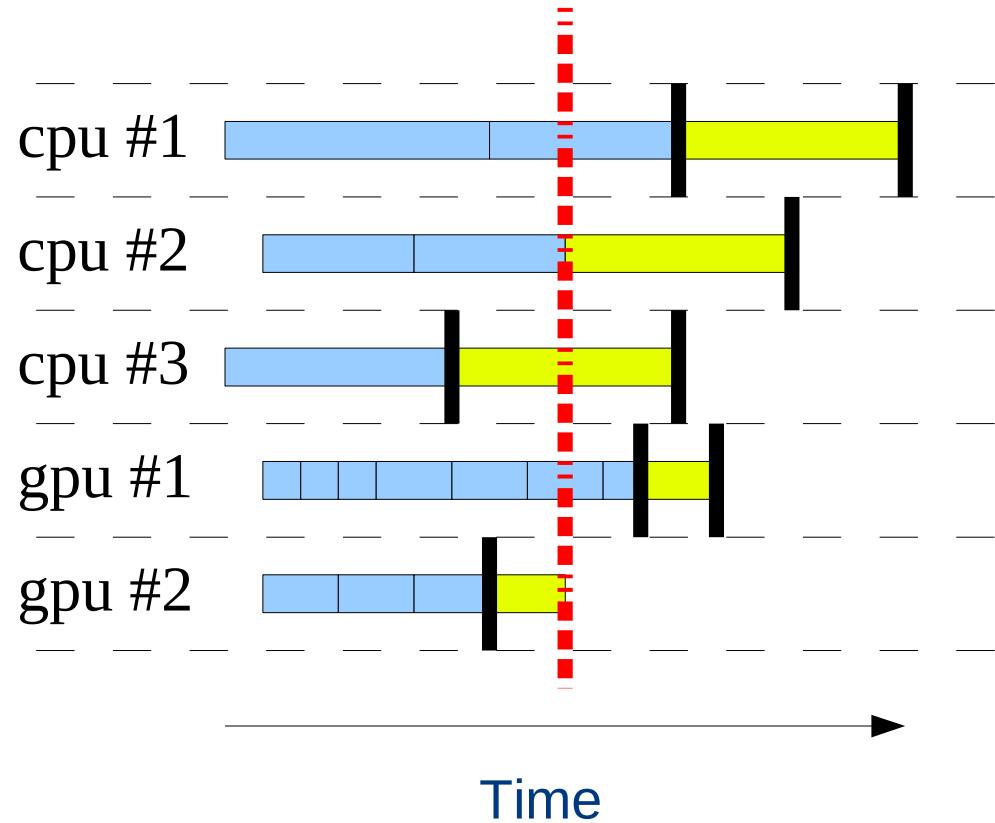
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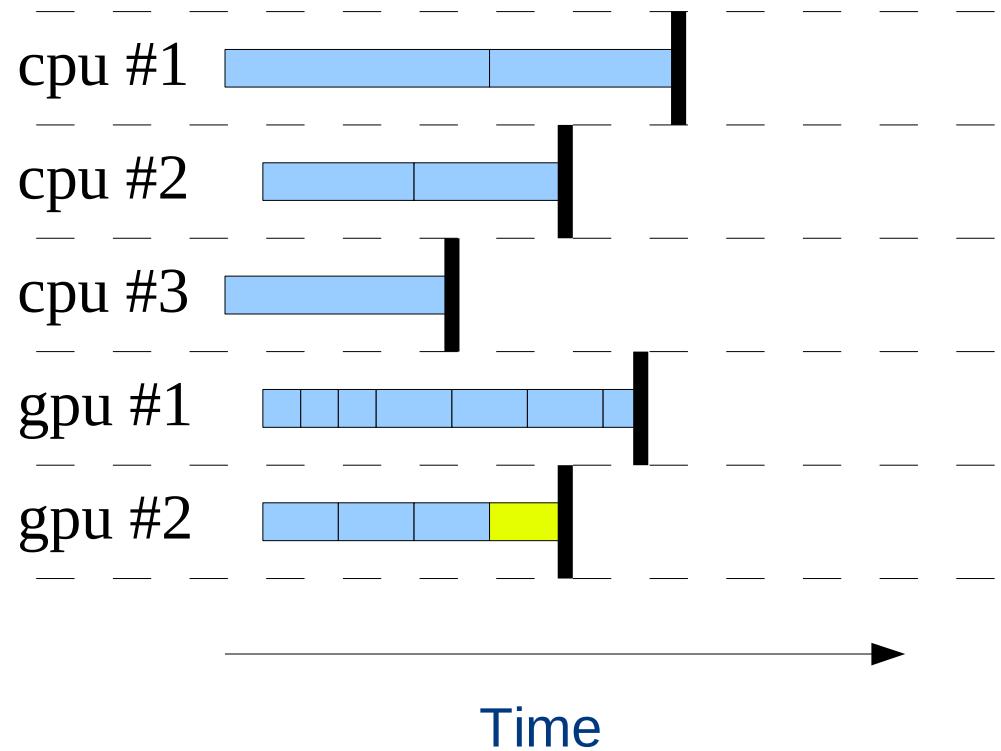
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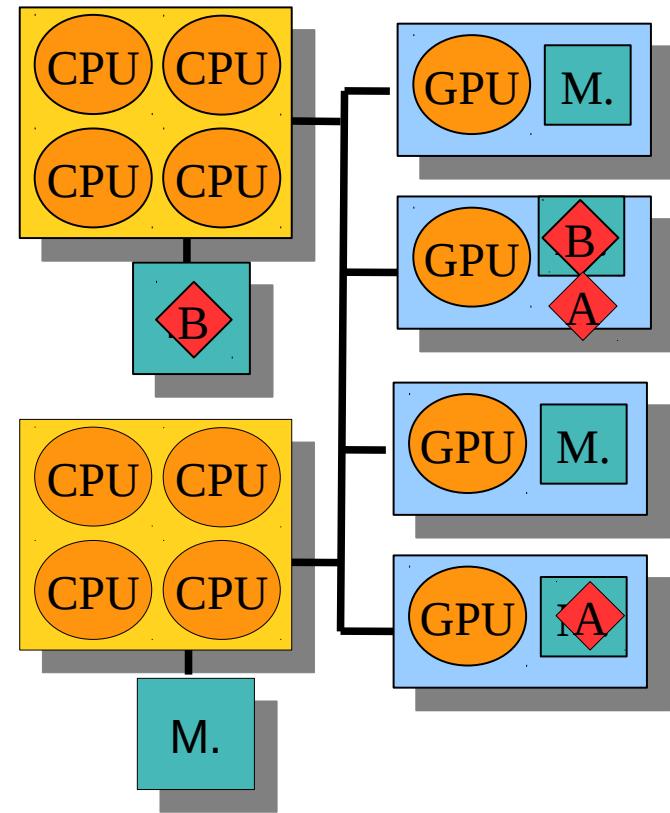
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Predicting data transfer overhead

Motivations

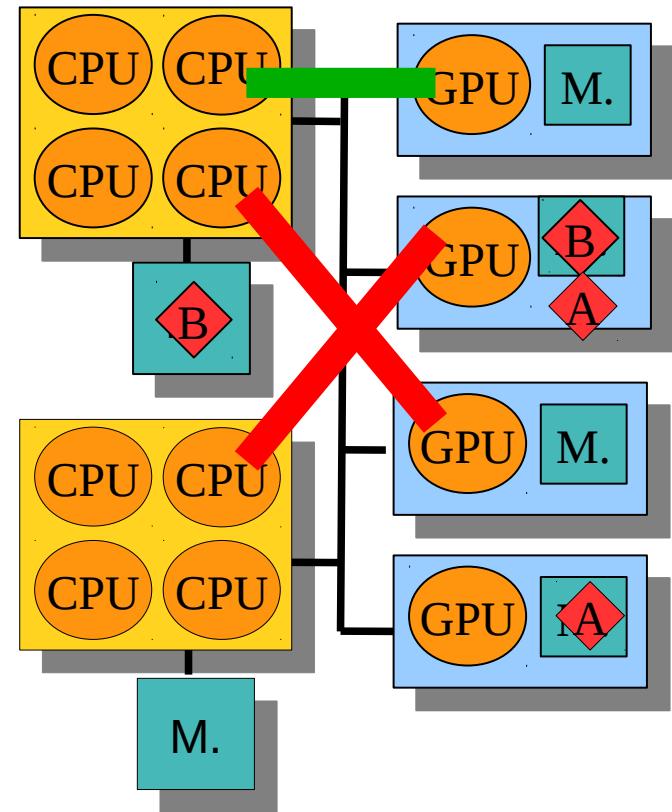
- Hybrid platforms
 - Multicore CPUs and GPUs
 - PCI-e bus is a precious ressource
- Data locality vs. Load balancing
 - Cannot avoid all data transfers
 - Minimize them
- StarPU keeps track of
 - data replicates
 - on-going data movements



Predicting data transfer overhead

Offline bus benchmarking

- Offline bus benchmarking
 - When StarPU is launched for the first time
 - Measure bandwidth and latency
 - Stored as files
 - Loaded when StarPU is initialized
- Detect CPU/GPU affinity
 - Control a GPU from the closest CPU
 - Significant impact on bus usage
- Straightforward cost prediction
 - Latency + size * bandwidth
 - Could be improved in many ways



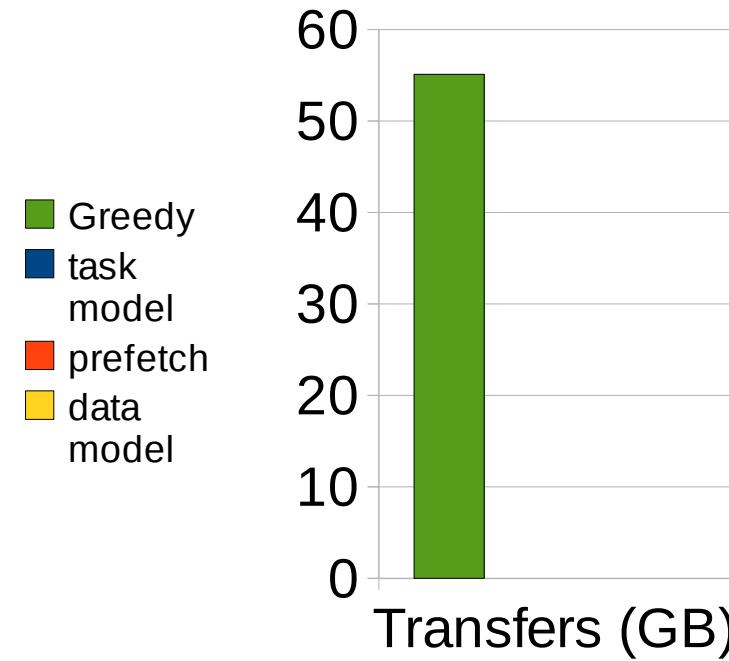
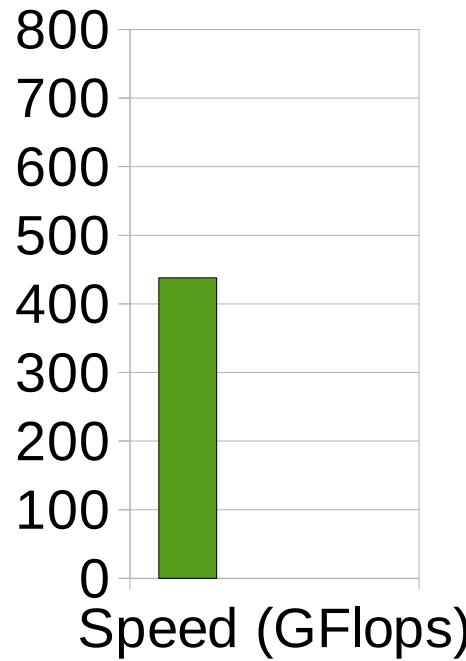
Impact of scheduling policy on a synthetic LU decomposition (without pivoting !)



Scheduling in a hybrid environment

Performance models

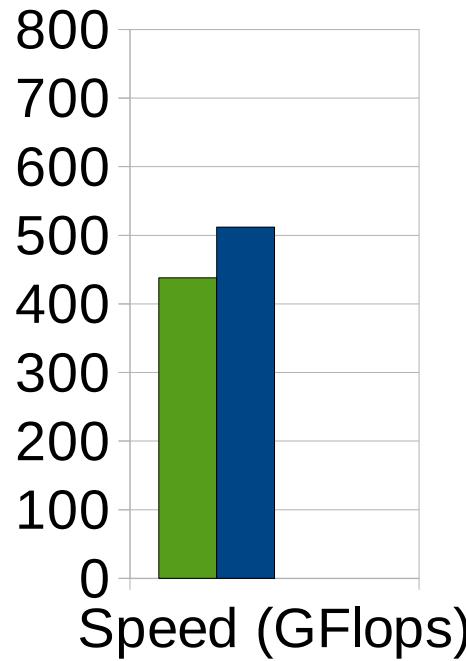
- LU without pivoting (16GB input matrix)
 - 8 CPUs (nehalem) + 3 GPUs (FX5800)



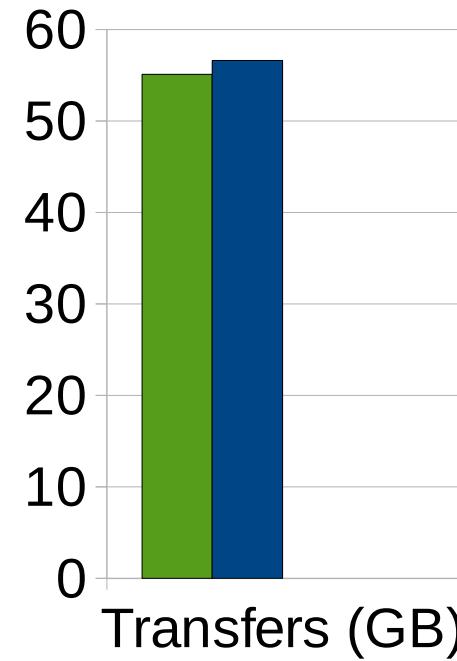
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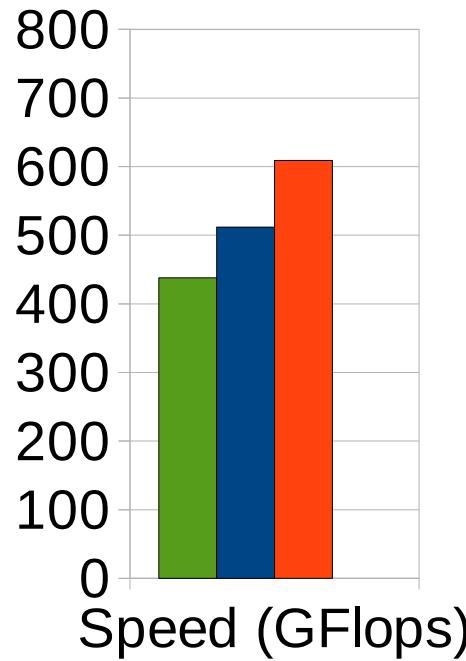
Greedy
task
model
prefetch
data
model



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Greedy
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Speed (GFlops)



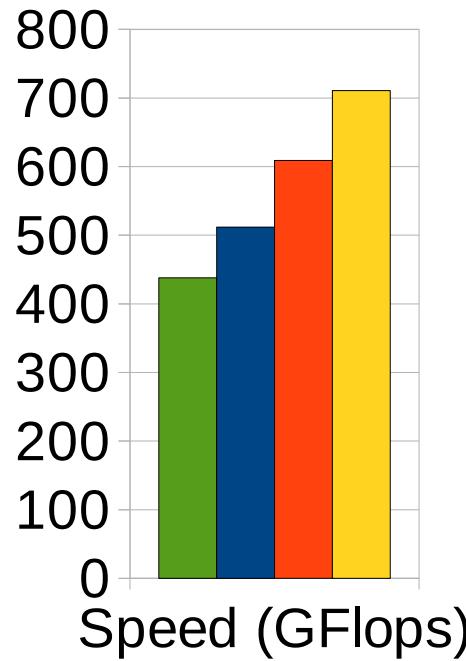
Transfers (GB)



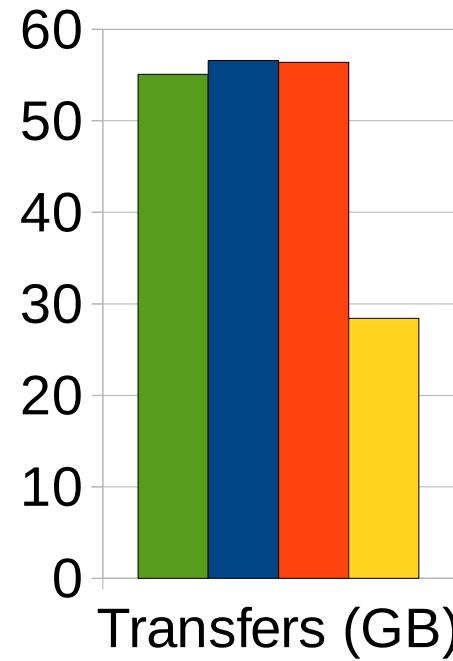
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Mixing PLASMA and MAGMA with StarPU

(in collaboration with UTK)

Cholesky & QR decompositions



Mixing PLASMA and MAGMA with StarPU

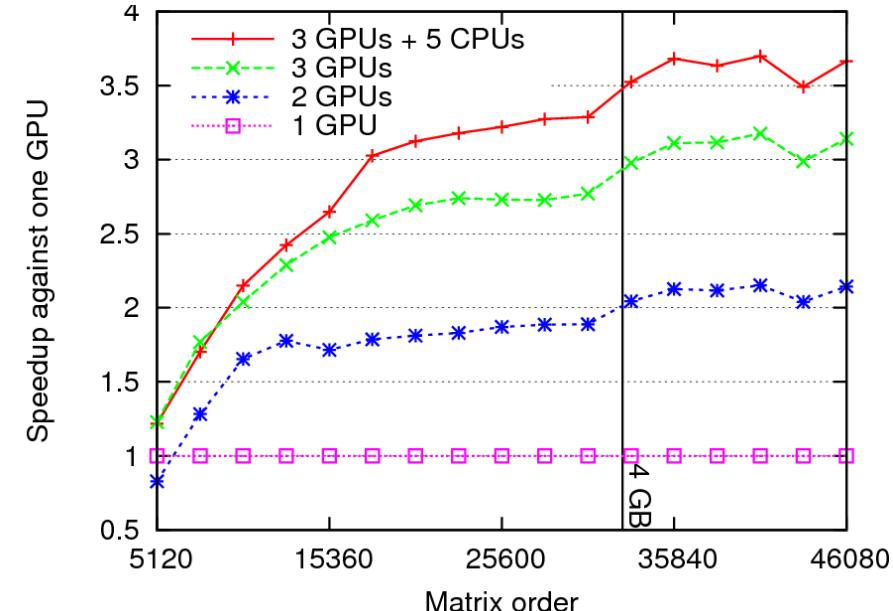
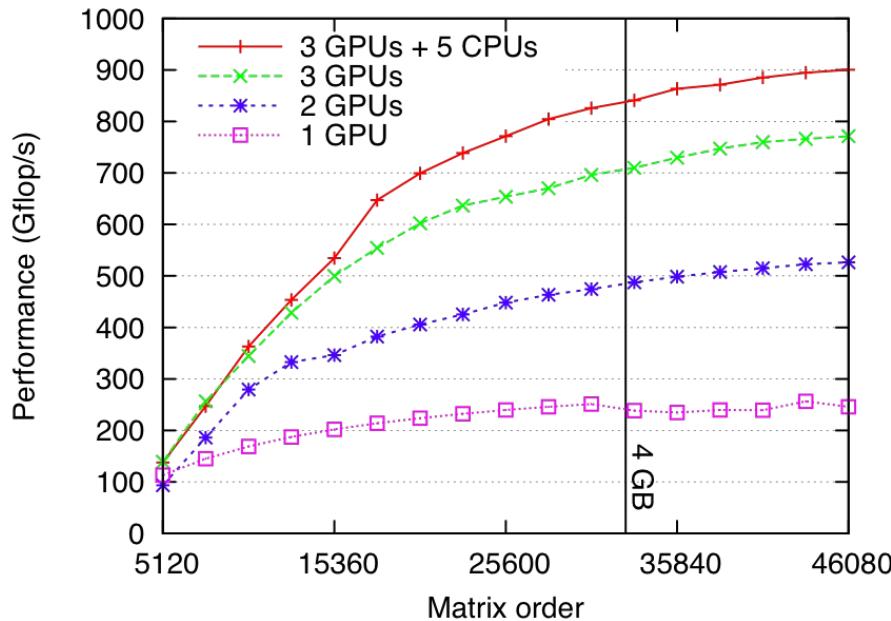
- State of the art algorithms
 - PLASMA (Multicore CPUs)
 - Dynamically scheduled with Quark
 - MAGMA (Multiple GPUs)
 - Hand-coded data transfers
 - Static task mapping
- General SLAGMA design
 - Use PLASMA algorithm with « magnum tiles »
 - PLASMA kernels on CPUs, MAGMA kernels on GPUs
 - Bypass the QUARK scheduler
- Programmability
 - Cholesky: ~half a week
 - QR : ~2 days of works
 - Quick algorithmic prototyping



Mixing PLASMA and MAGMA with StarPU

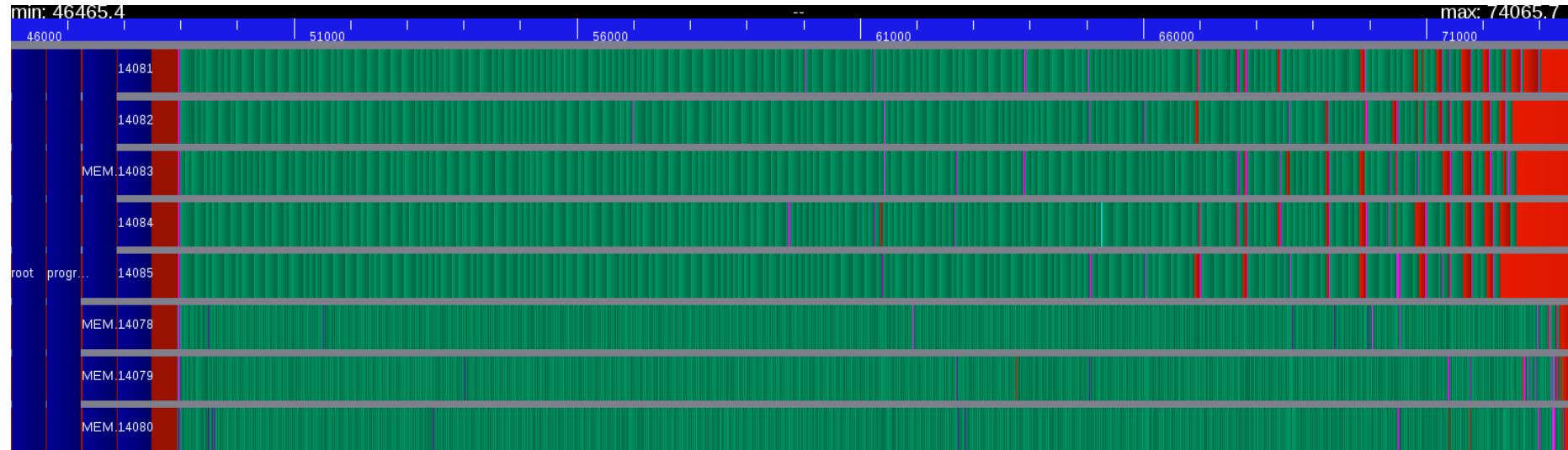
- Cholesky decomposition

- 5 CPUs (Nehalem) + 3 GPUs (FX5800)
- Efficiency > 100%



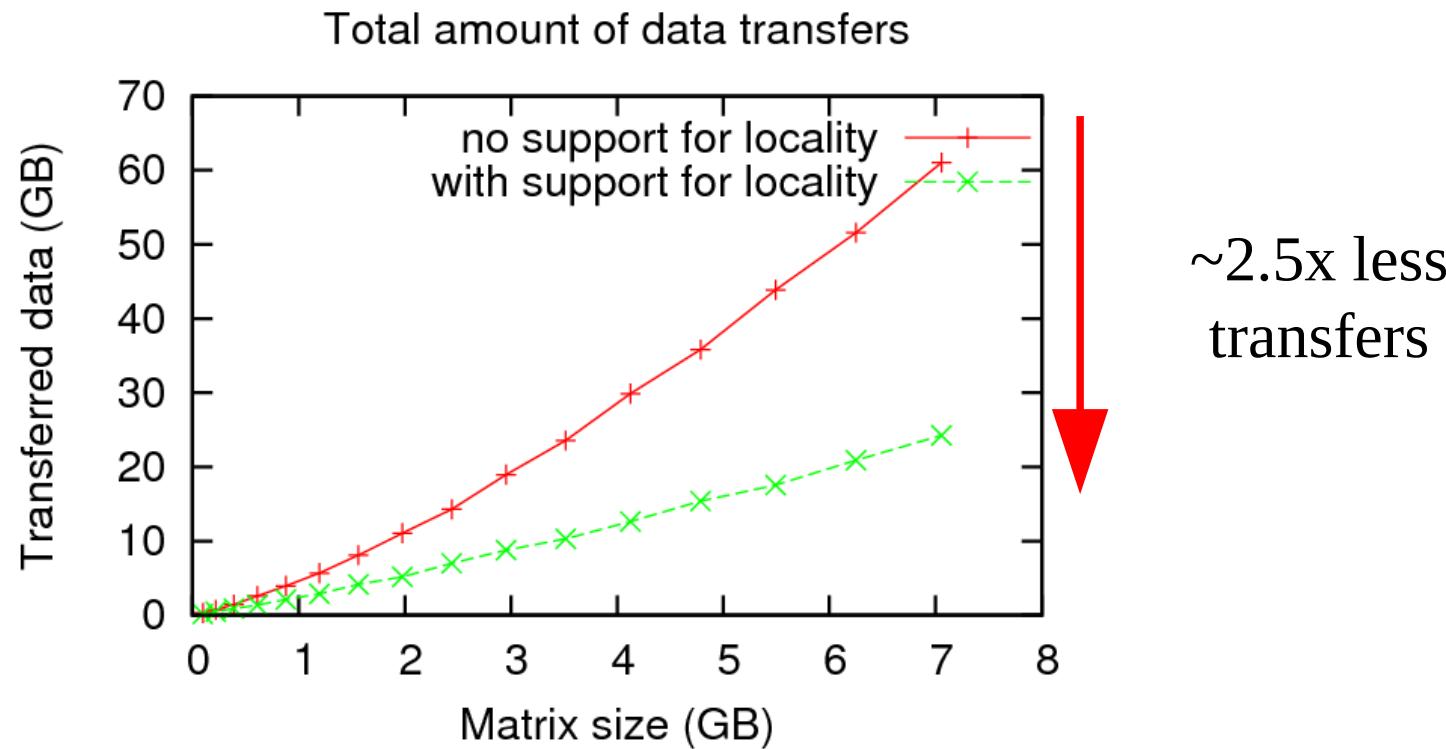
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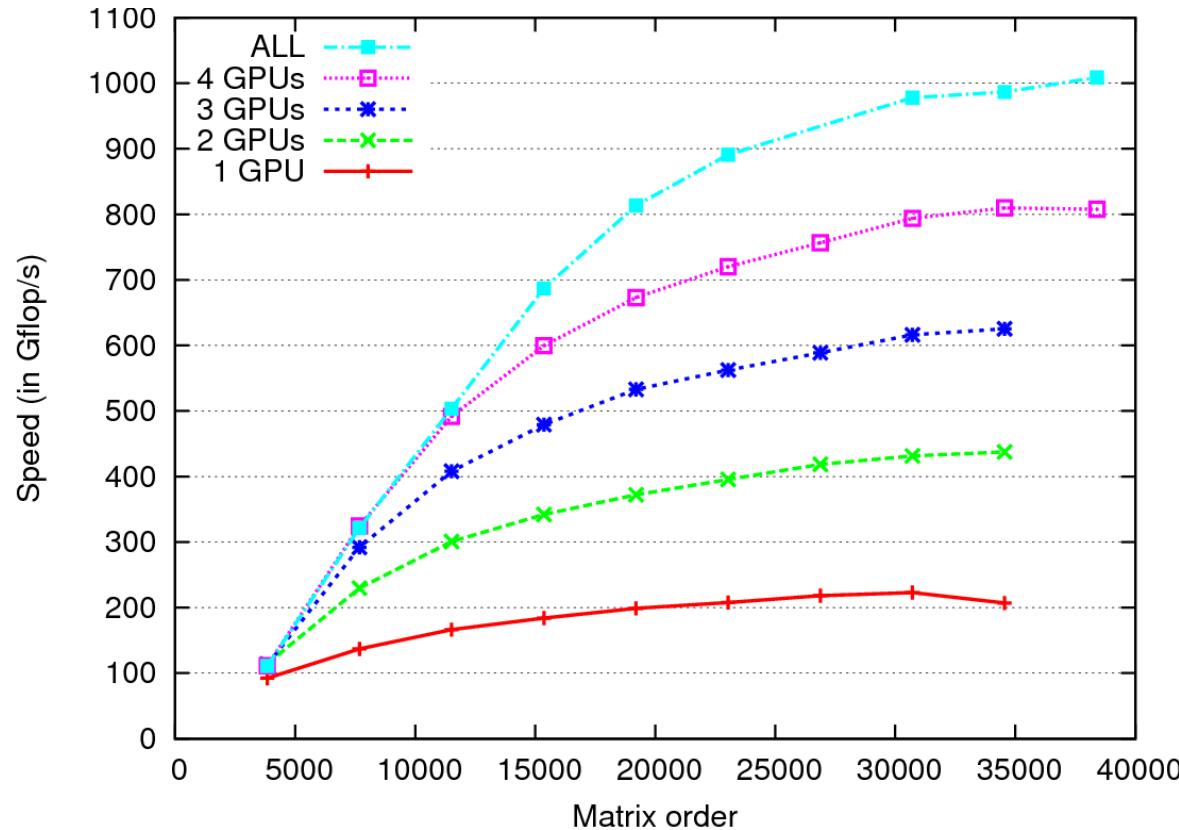
Mixing PLASMA and MAGMA with StarPU

- Memory transfers during Cholesky decomposition



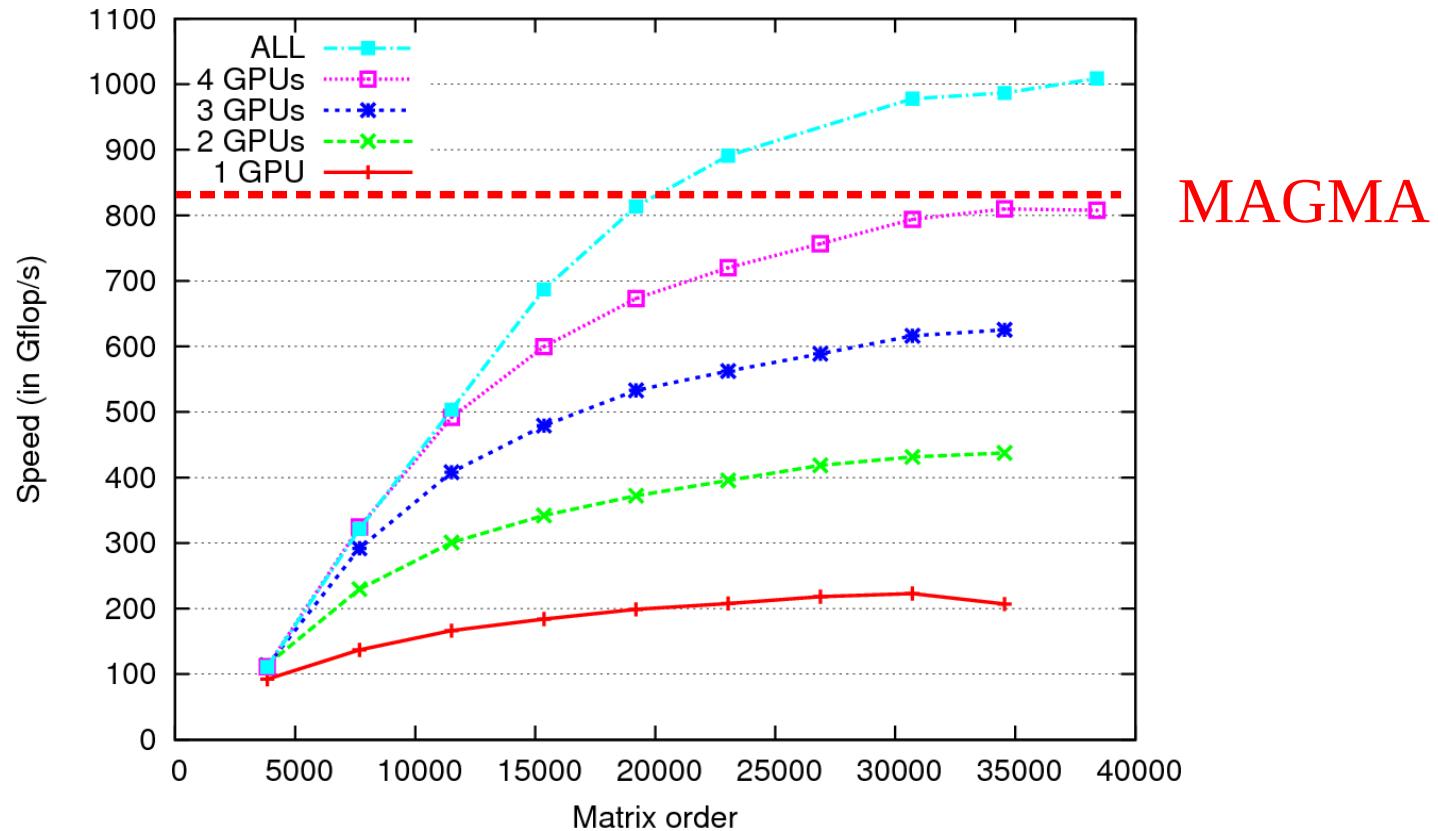
Mixing PLASMA and MAGMA with StarPU

- QR decomposition
 - Mordor8 (UTK) : 16 CPUs (AMD) + 4 GPUs (C1060)



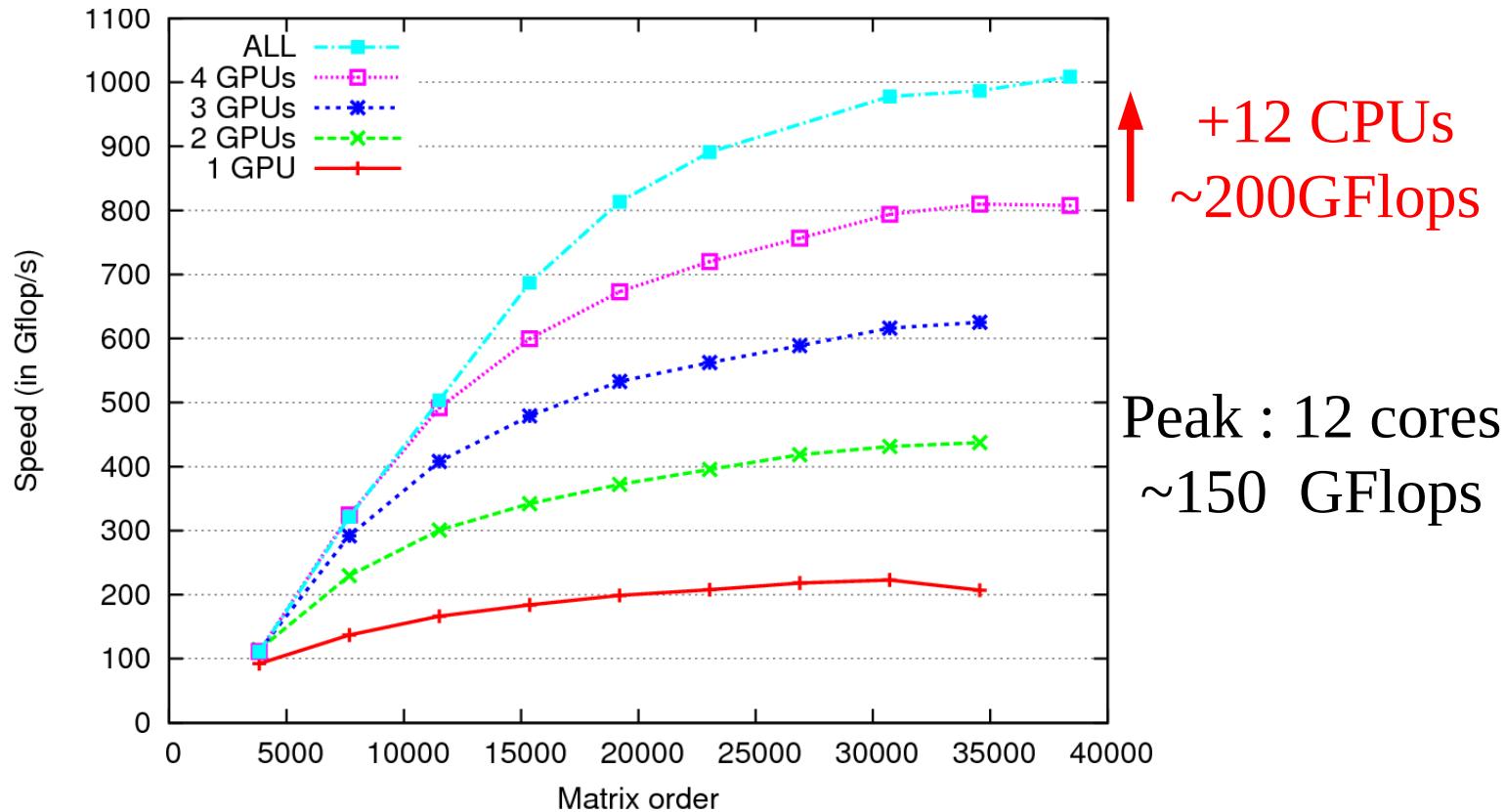
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Mixing PLASMA and MAGMA with StarPU

- « Super-Linear » efficiency in QR?
 - Kernel efficiency
 - sgeqrt
 - CPU: 9 Gflops GPU: 30 Gflops (Speedup : ~3)
 - stsqrt
 - CPU: 12Gflops GPU: 37 Gflops (Speedup: ~3)
 - somqr
 - CPU: 8.5 Gflops GPU: 227 Gflops (Speedup: ~27)
 - Sssmqr
 - CPU: 10Gflops GPU: 285Gflops (Speedup: ~28)
 - Task distribution observed on StarPU
 - sgeqrt: 20% of tasks on GPUs
 - Sssmqr: 92.5% of tasks on GPUs
 - Taking advantage of heterogeneity !
 - Only do what you are good for
 - Don't do what you are not good for



Scheduling parallel tasks



Parallel tasks

Why do we need parallel tasks ?

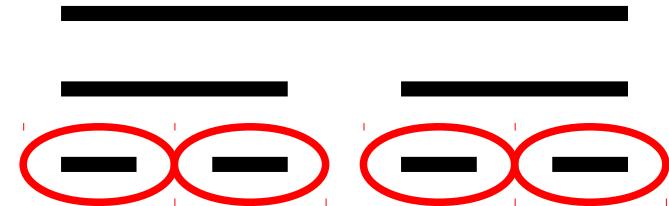
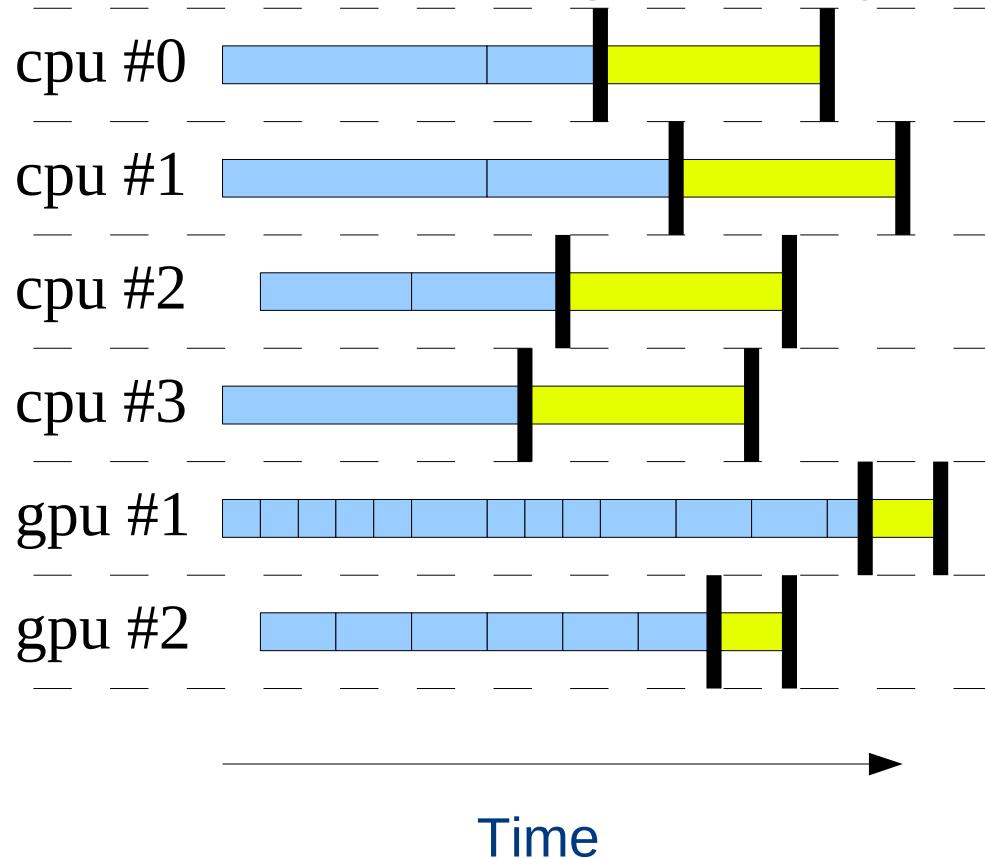
- Take advantage of multicore architectures
- Task parallelism may not be suited at a fine grain
 - Use other parallel paradigms
 - eg. OpenMP
 - Use existing parallel libraries
 - eg. do not reimplement parallel BLAS ...
- Alleviate granularity concerns
 - Less tasks
 - Large enough tasks (suited for the GPU)



Parallel tasks

Scheduling parallel tasks

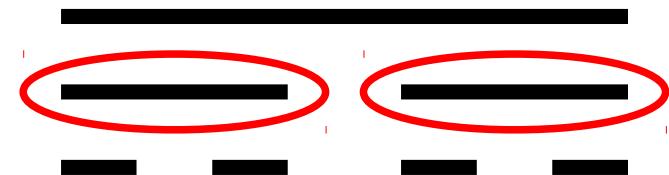
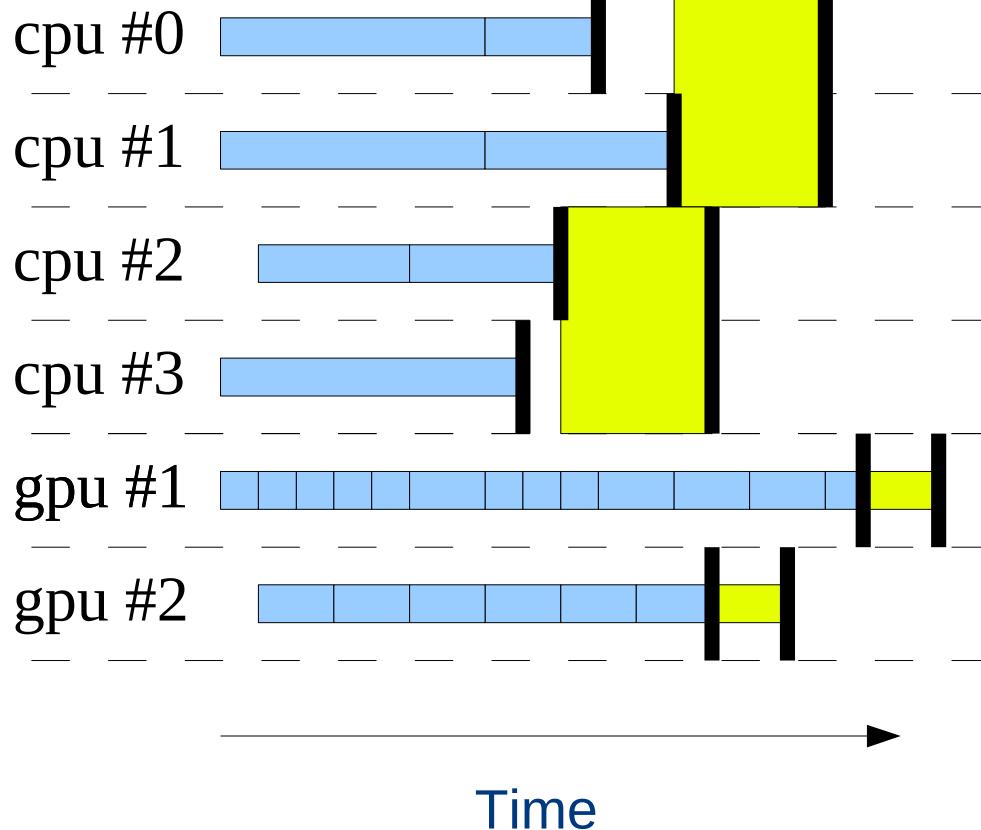
- StarPU allocates processing units



Parallel tasks

Scheduling parallel tasks

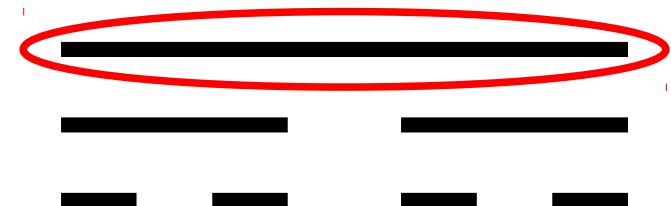
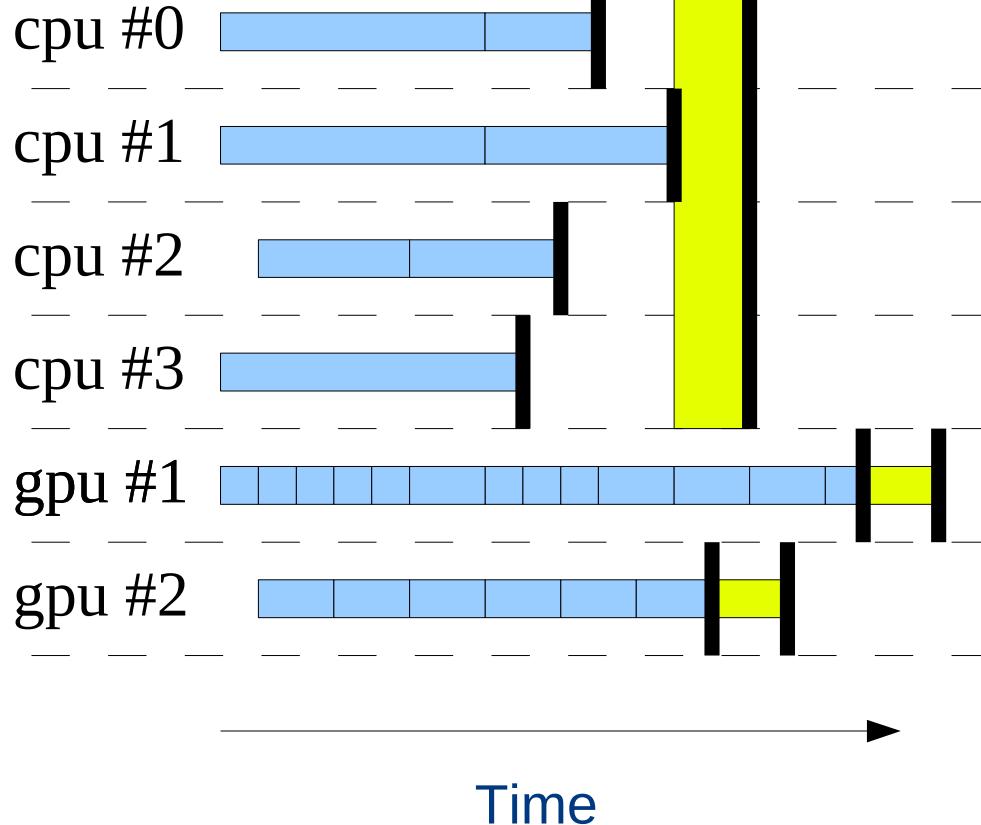
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Parallel tasks

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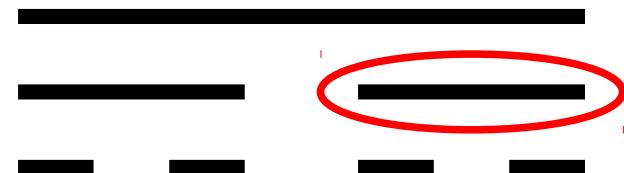
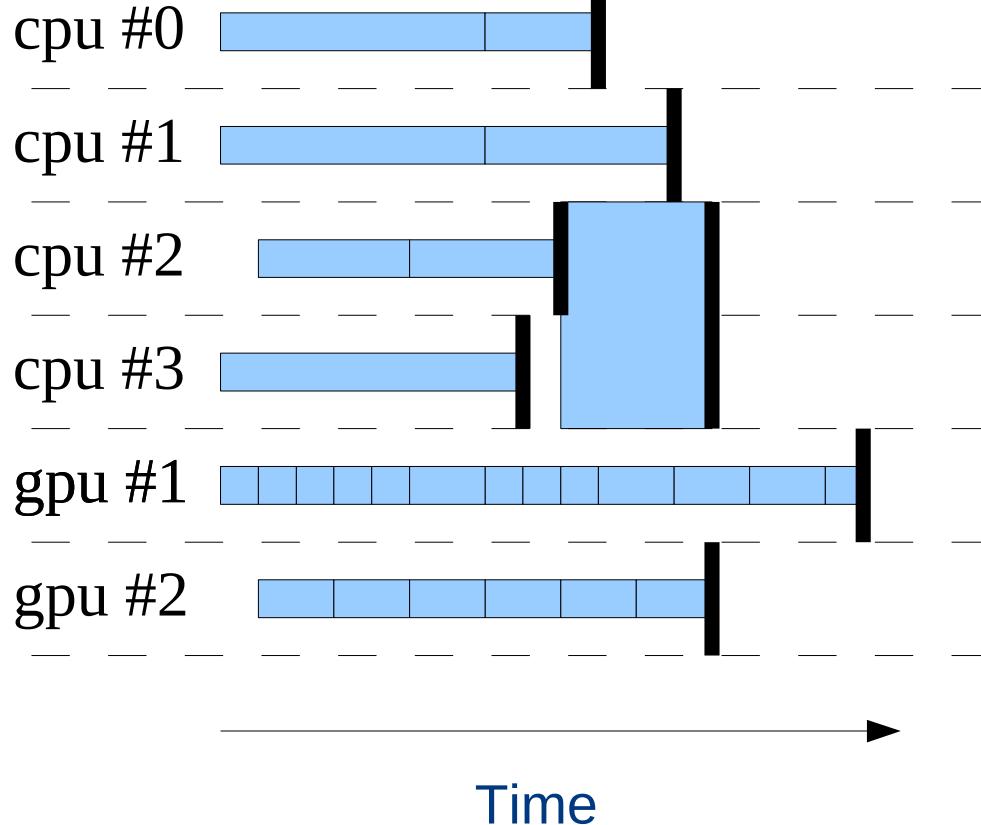
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Parallel tasks

Scheduling parallel tasks

- StarPU allocates processing units



Adding support for MPI in StarPU



Accelerating MPI applications with StarPU

- Keep MPI SPMD style
 - Static distribution of data
 - Scheduling within the node only
 - No load balancing between MPI processes
- Inter-process data dependencies
 - MPI communications triggered by StarPU data availability
 - Support from StarPU's memory management
 - Automatically construct MPI datatype



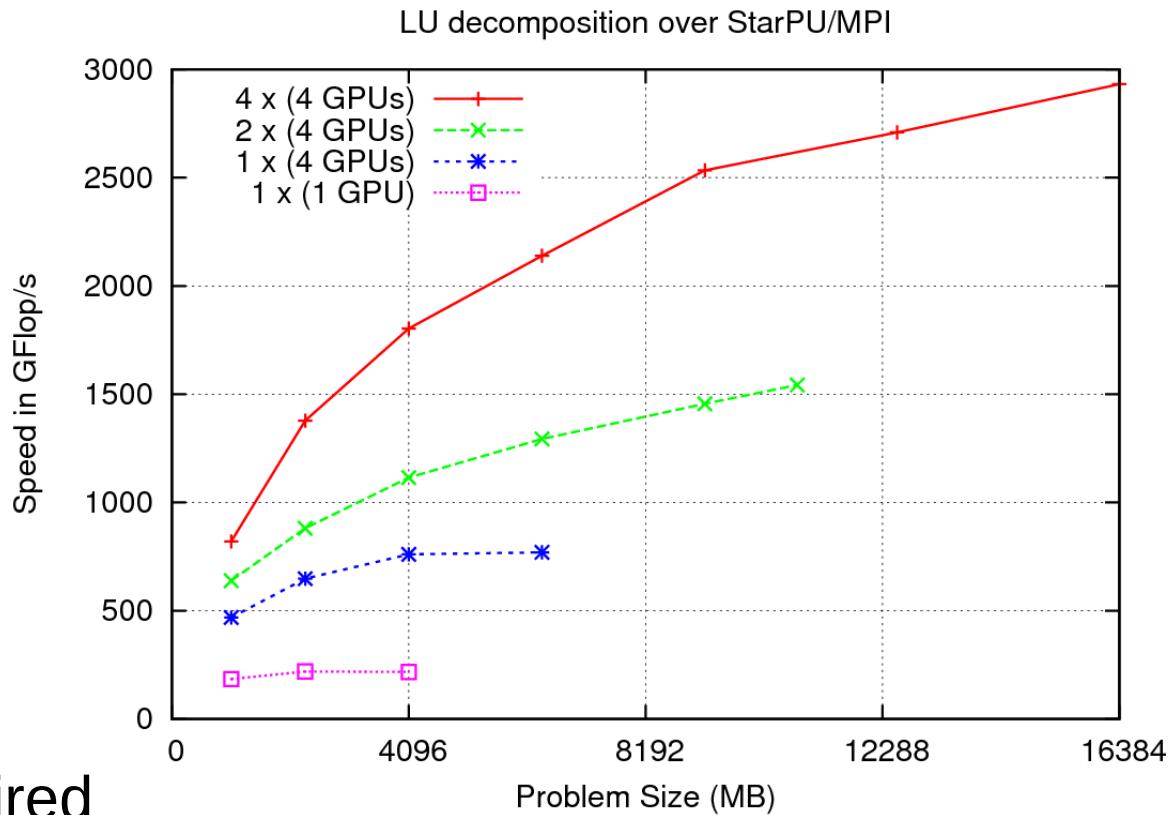
Accelerating MPI applications with StarPU

- Provided API
 - `starpu_mpi_{send,recv}`
 - `starpu_mpi_{isend,irecv}`
 - `starpu_mpi_{test,wait}`
 - `starpu_mpi_{send,recv}_detached`
 - `starpu_mpi_*_array`
- Detached calls
 - No need to explicitly test/wait for the request
 - Automatic progression
- Automatic data dependencies
 - MPI transfers ~ StarPU tasks
 - Accelerating legacy codes



Accelerating LU/MPI with StarPU

- LU decomposition
 - MPI+multiGPU
- Static MPI distribution
 - 2D block cyclic
 - ~SCALAPACK
 - No pivoting !
- Algorithmic work required
 - Collaboration with UTK



Conclusion



Conclusion

Summary

- StarPU

- Freely available under LGPL
- Available on Linux, OS/X, Windows
- Open to external contributors!

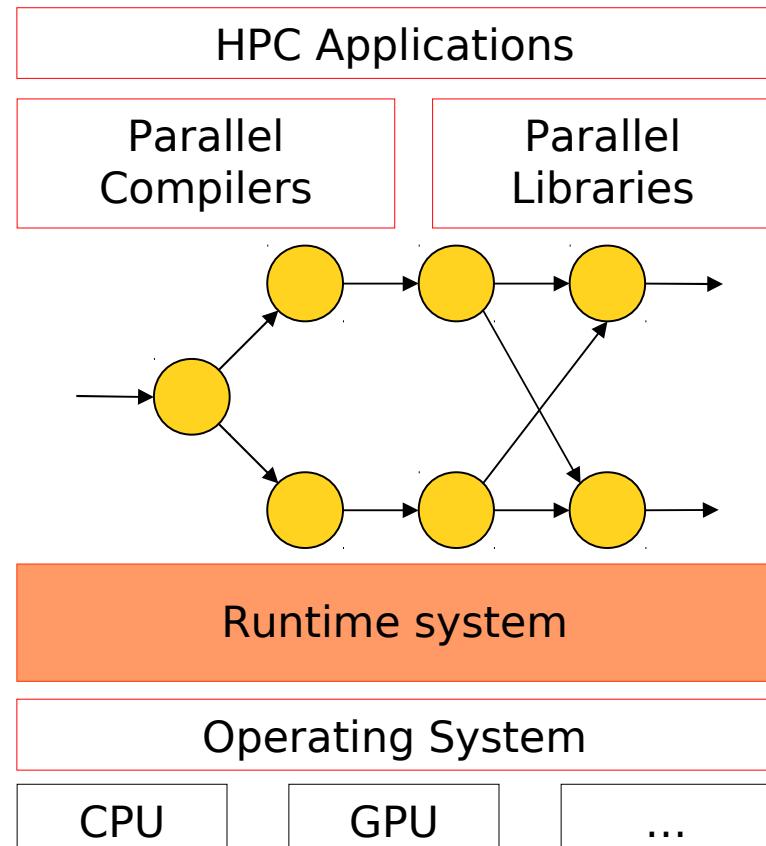
- Task Scheduling

- Required on hybrid platforms
- Auto-tuned performance models

- Combined PLASMA and MAGMA

- Parallel tasks

- MPI extensions



Conclusion

Future work

- Implement more algorithms

- LU, Hessenberg
- Communication Avoiding algorithms
- Hybrid Scalapack
- Provide higher level constructs (eg. reductions)

HPC Applications

Compiling environment

Specific libraries

- Provide a back-end for compilers
- StarSS, XscalableMP, HMPP

Runtime system

- Support new architectures
- Intel SCC, Fermi cards, ...

Operating System

- Dynamically adapt granularity
- Divisible tasks

Hardware



Conclusion

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Thanks for your attention !
Any question ?





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Performance Models

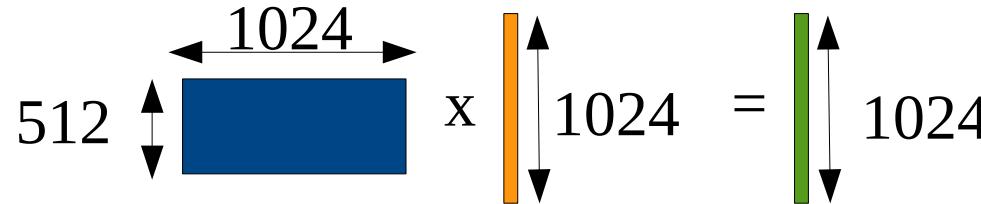
Our History-based proposition

- Hypothesis

- Regular applications
- Execution time independent from data content
 - Static Flow Control

- Consequence

- Data description fully characterizes tasks
- Example: matrix-vector product



- Unique Signature : $((1024, 512), \text{CRC}(1024), \text{CRC}(1024))$
- Per-data signature
 - $\text{CRC}(1024, 512) = 0x951ef83b$
- Task signature
 - $\text{CRC}(\text{CRC}(1024, 512), \text{CRC}(1024), \text{CRC}(1024)) = 0x79df36e2$

Performance Models⁶²

Our History-based proposition

- Generalization is easy

- Task $f(D_1, \dots, D_n)$
- Data
 - $\text{Signature}(D_i) = \text{CRC}(p_1, p_2, \dots, p_k)$
- Task \sim Series of data
 - $\text{Signature}(D_1, \dots, D_n) = \text{CRC}(\text{sign}(D_1), \dots, \text{sign}(D_n))$

- Systematic method

- Problem independent
- Transparent for the programmer
- Efficient



Evaluation⁶³

Example: LU decomposition

Speed (GFlop/s)		
	(16k x 16k)	(30k x 30k)
ref.	89.98 \pm 2.97	130.64 \pm 1.66
1 st iter	48.31	96.63
2 nd iter	103.62	130.23
3 rd iter	103.11	133.50
\geq 4 iter	103.92 \pm 0.46	135.90 \pm 0.00

- Faster
- No code change !
- More stable
- Dynamic calibration
- Simple, but accurate

