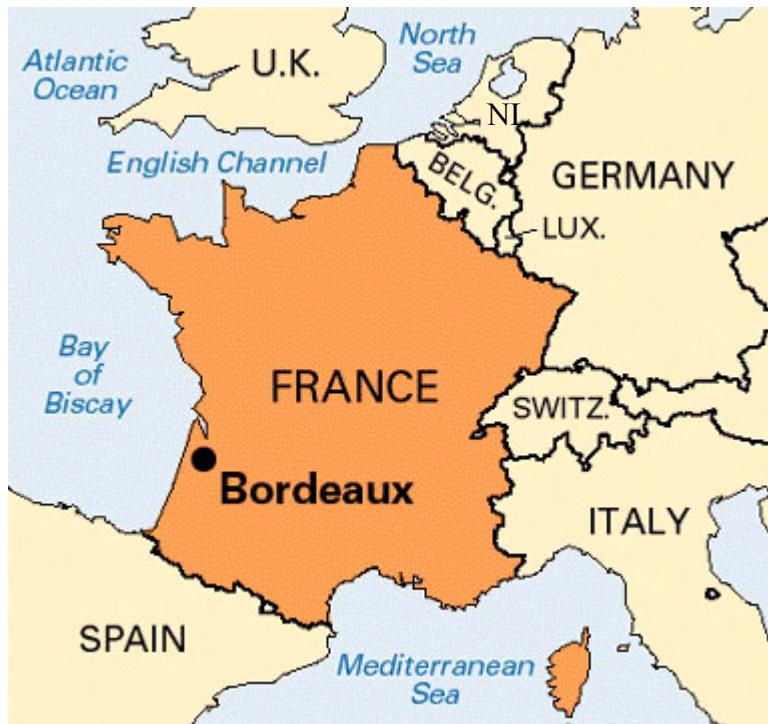




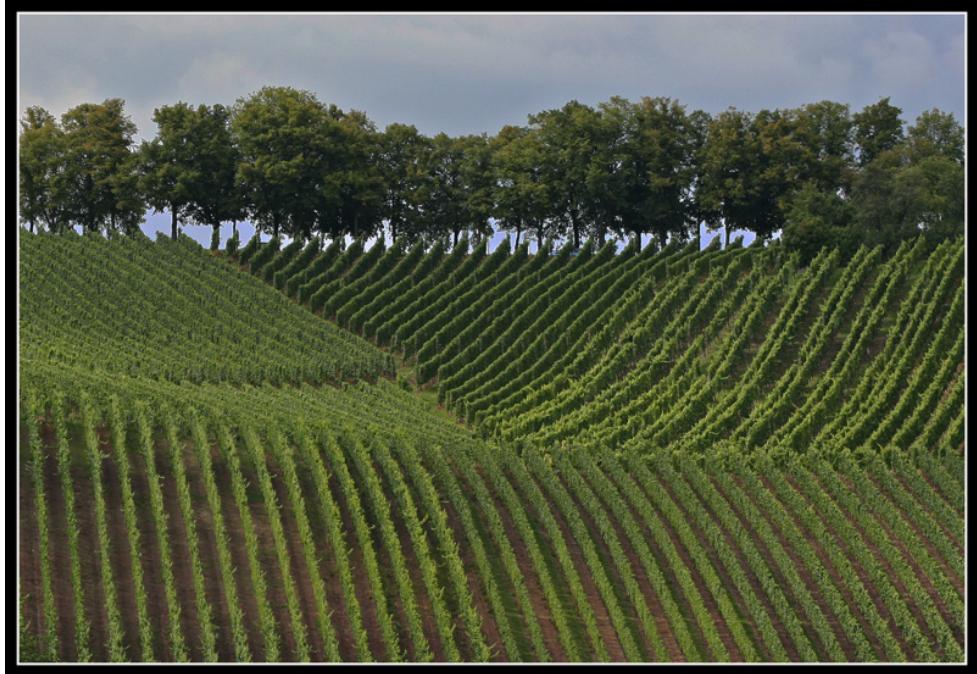
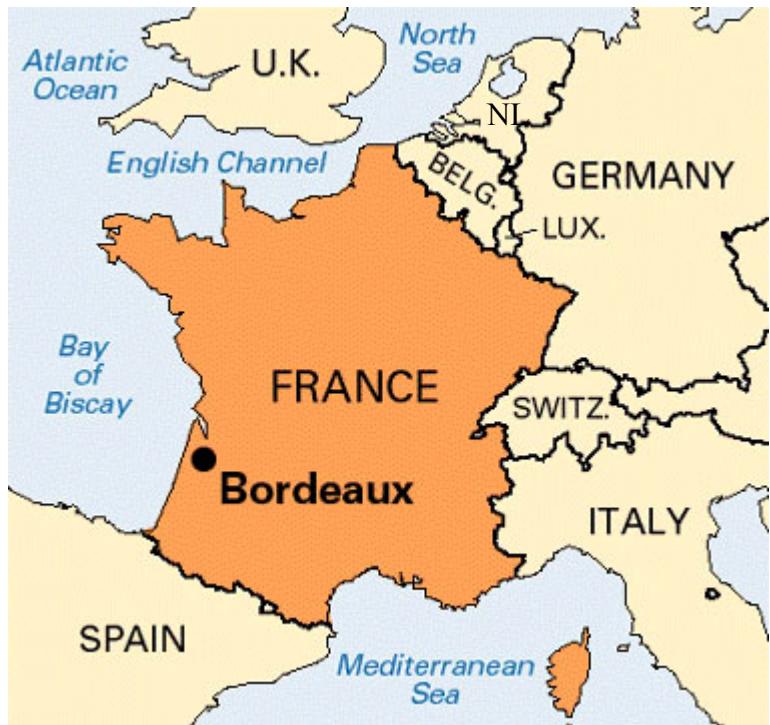
StarPU: seamless computations among CPUs and GPUs

Cédric Augonnet, Samuel Thibault,
Olivier Aumage, Nathalie Furmento
INRIA Runtime Team-Project

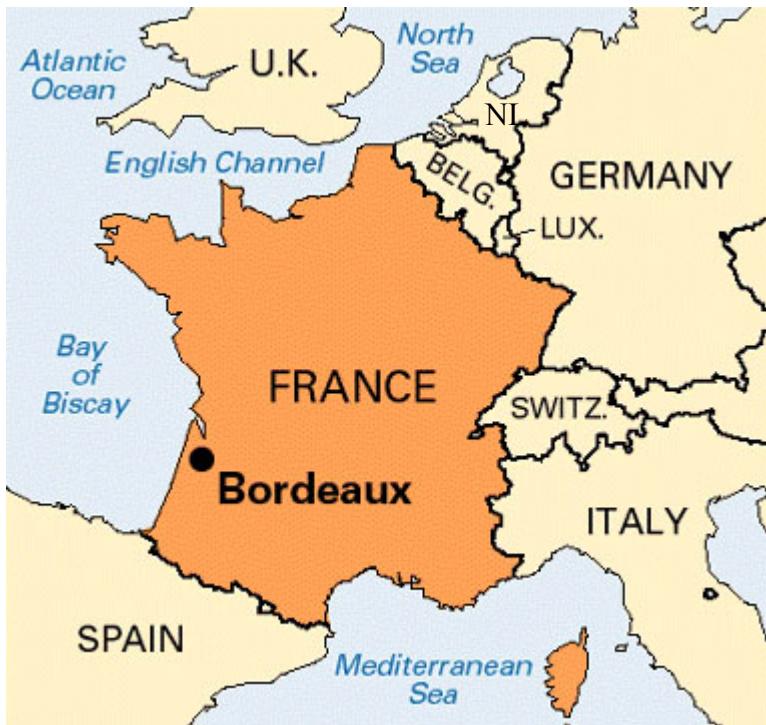
The RUNTIME Team



The RUNTIME Team



The RUNTIME Team

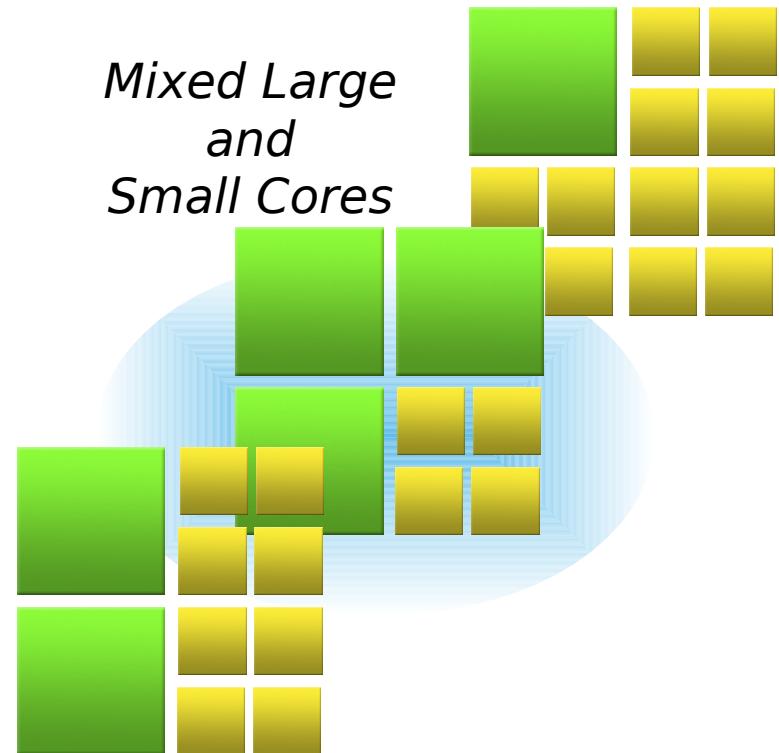


Doing Parallelism for centuries !

Introduction

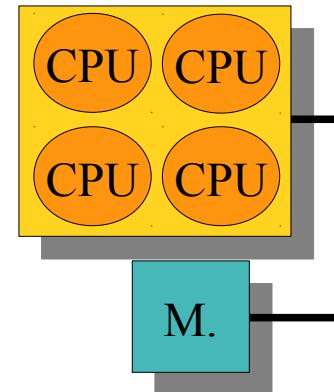
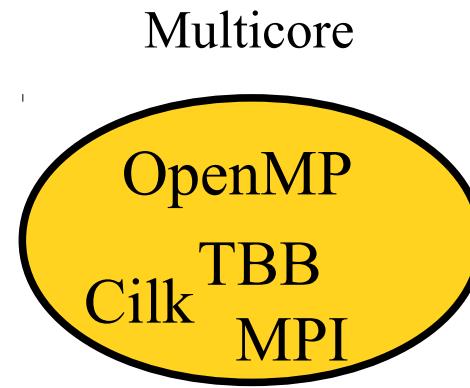
Toward heterogeneous multi-core architectures

- Multicore is here
 - Hierarchical architectures
 - Manycore
- Architecture specialization
 - Now
 - Accelerators (GPGPUs, FPGAs)
 - Coprocessors (Xeon Phi)
 - Fusion
 - DSPs
 - All of the above
 - In the near Future
 - Many simple cores
 - A few full-featured cores



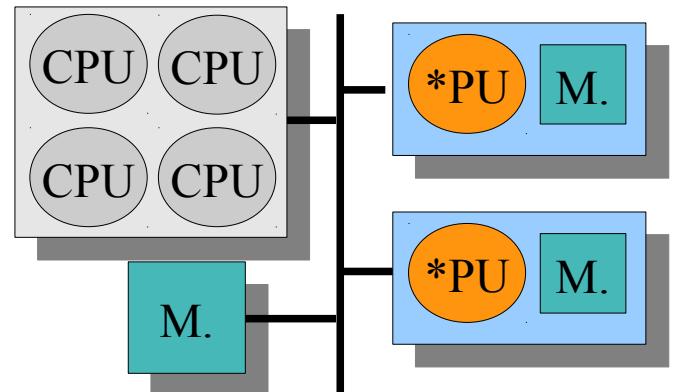
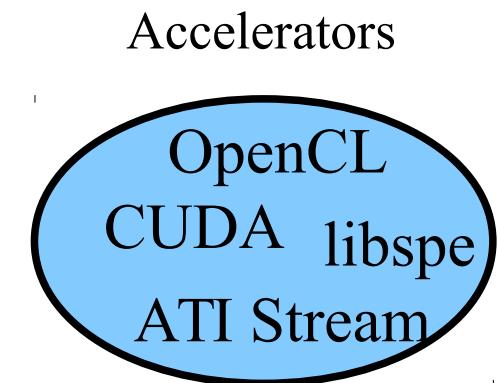
How to program these architectures?

- Multicore programming
 - pthreads, OpenMP, TBB, ...



How to program these architectures?

- Multicore programming
 - pthreads, OpenMP, TBB, ...
- Accelerator programming
 - Consensus on OpenCL/OpenACC?
 - (Often) Pure offloading model

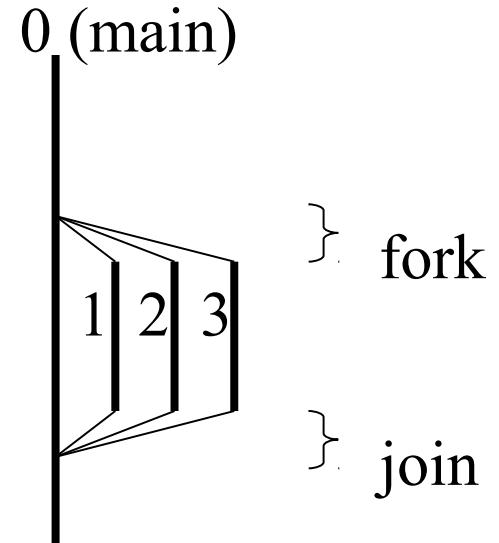


OpenMP

A portable approach to shared-memory programming

- Extension to existing languages
 - C, C++, Fortran
 - Set of programming directives
- Fork/join approach
 - Parallel sections
- Well suited to data-parallel programs
 - Parallel loops
- OpenMP 3.0 introduced *tasks*
 - Support for irregular parallelism

```
int matrix[MAX][MAX];  
...  
#pragma omp parallel for  
for (int i; i < 400; i++)  
{  
    matrix[i][0] += ...  
}
```



How to program these architectures?

Accelerator programming

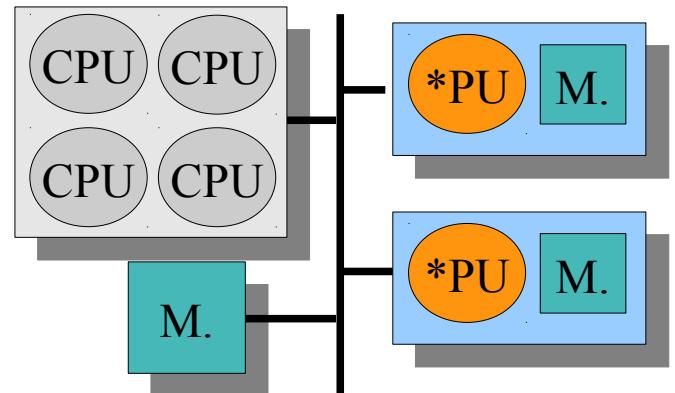
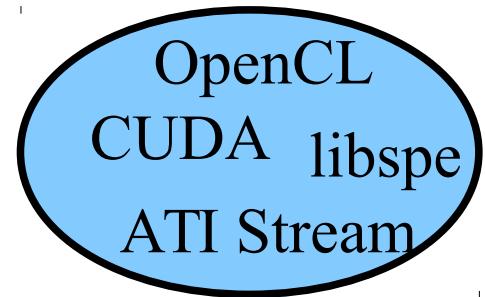
- OpenMP extension

```
int matrix[MAX][MAX];
...
#pragma omp target device(acc0)
map(matrix)

#pragma omp parallel for
for (int i; i < 400; i++)
{
    matrix[i][0] += ...
}
```

- Still quite hand-tuned

Accelerators



How to program these architectures?

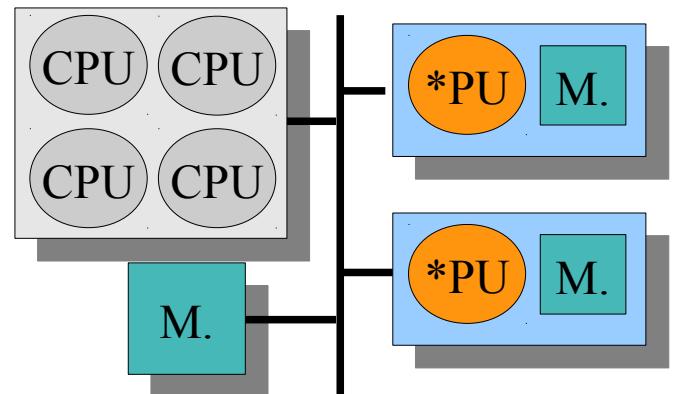
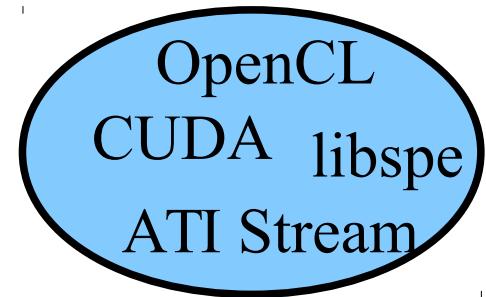
Accelerator programming

- OpenACC

```
int matrix[MAX] [MAX] ;
...
#pragma acc kernels copy(matrix)
for (int i; i < 400; i++)
{
    matrix[i][0] += ...
}
```

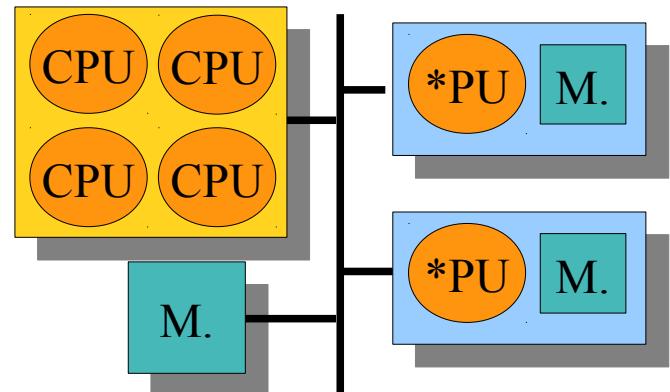
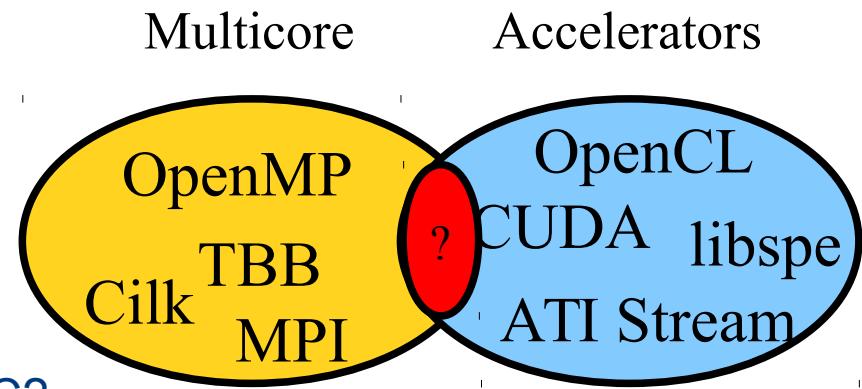
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Accelerators



How to program these architectures?

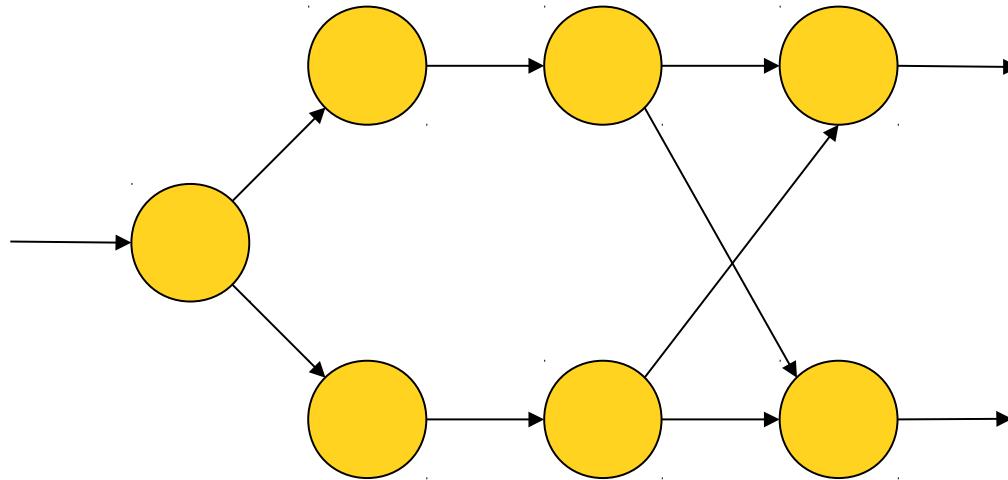
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 - pthreads, OpenMP, TBB, ...
- Accelerator programming
 - Consensus on OpenCL/OpenACC?
 - (Often) Pure offloading model
- Hybrid models?
 - Take advantage of all resources 😊
 - Complex interactions and distribution ☹



Task graphs

- Well-studied expression of parallelism
- Departs from usual sequential programming

Really ?



Task management

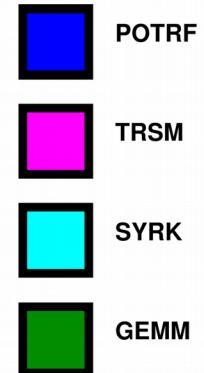
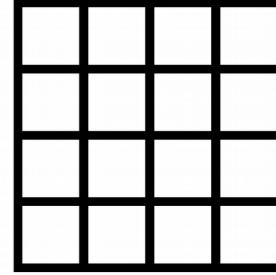
Implicit task dependencies

- Right-Looking Cholesky decomposition (from PLASMA)

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task_wait_for_all();

```



Task management

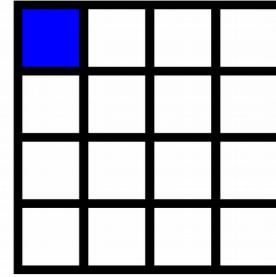
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	POTRF
	TRSM
	SYRK
	GEMM

Task management

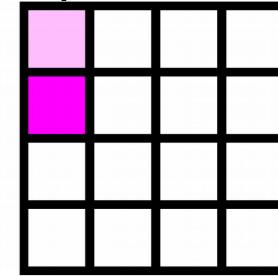
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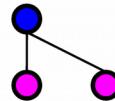
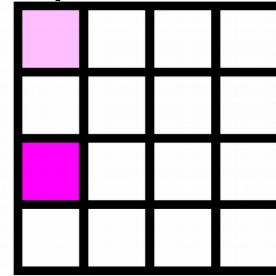
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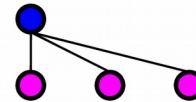
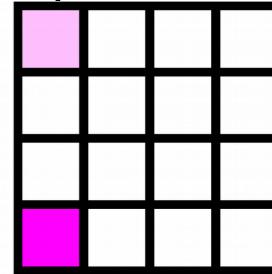
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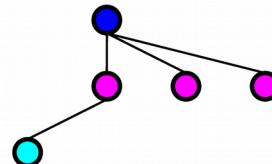
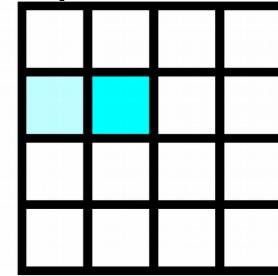
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Task management

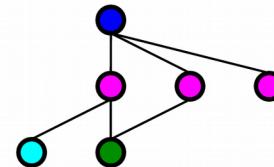
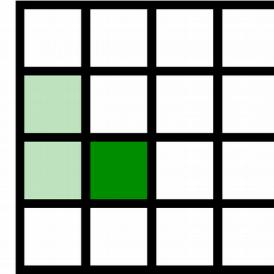
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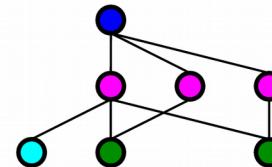
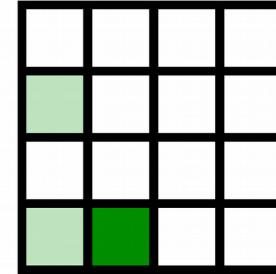
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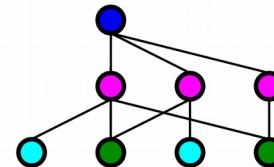
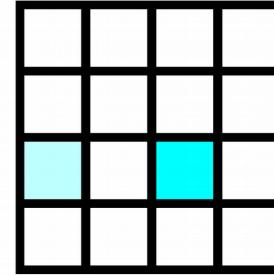
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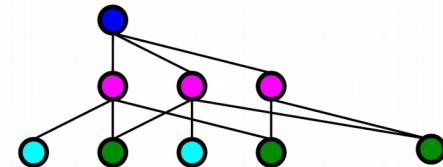
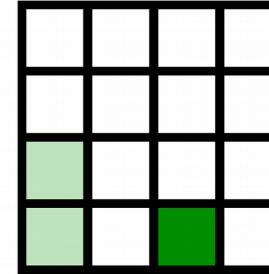
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Task management

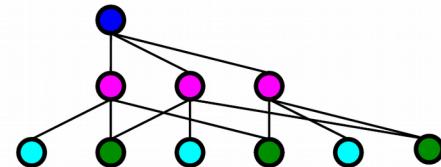
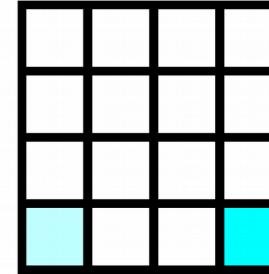
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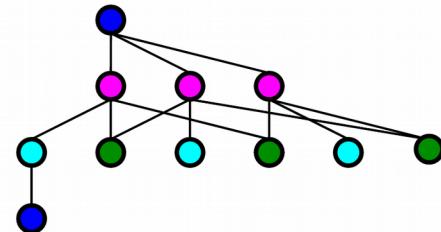
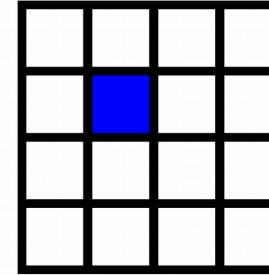
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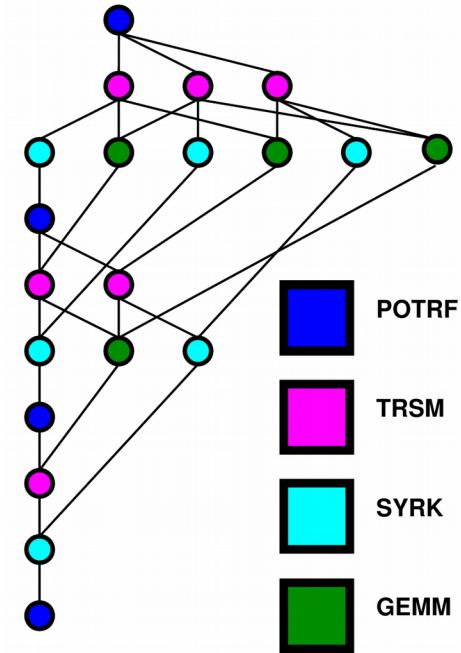
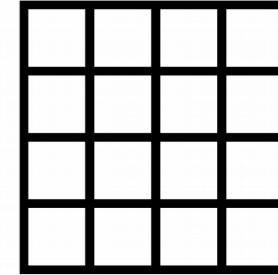
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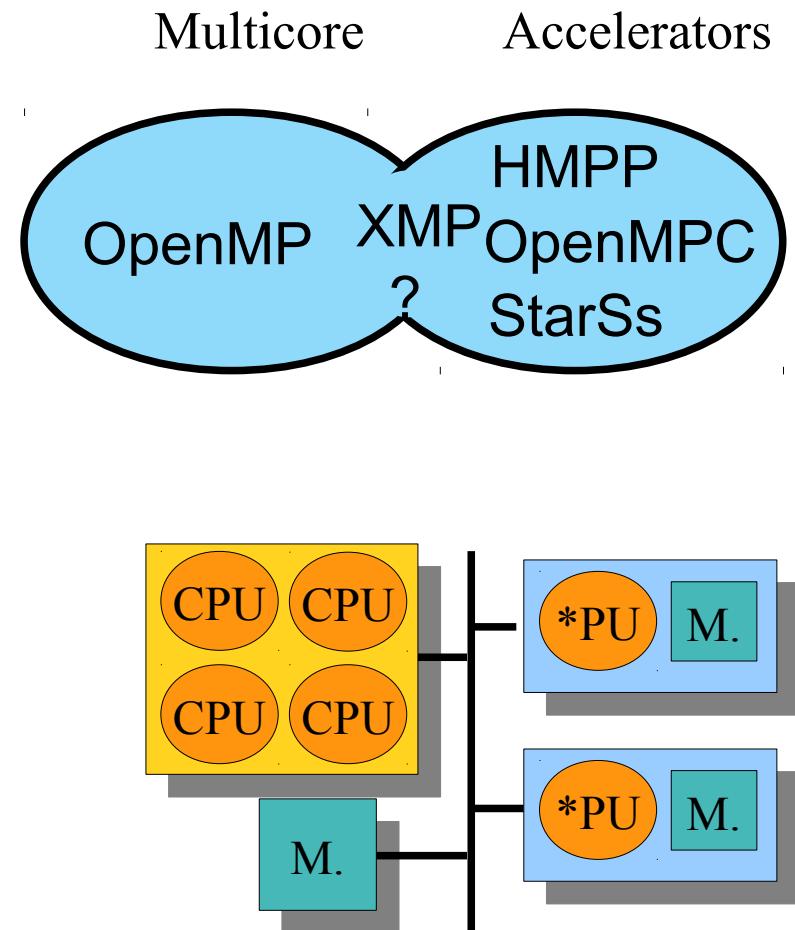
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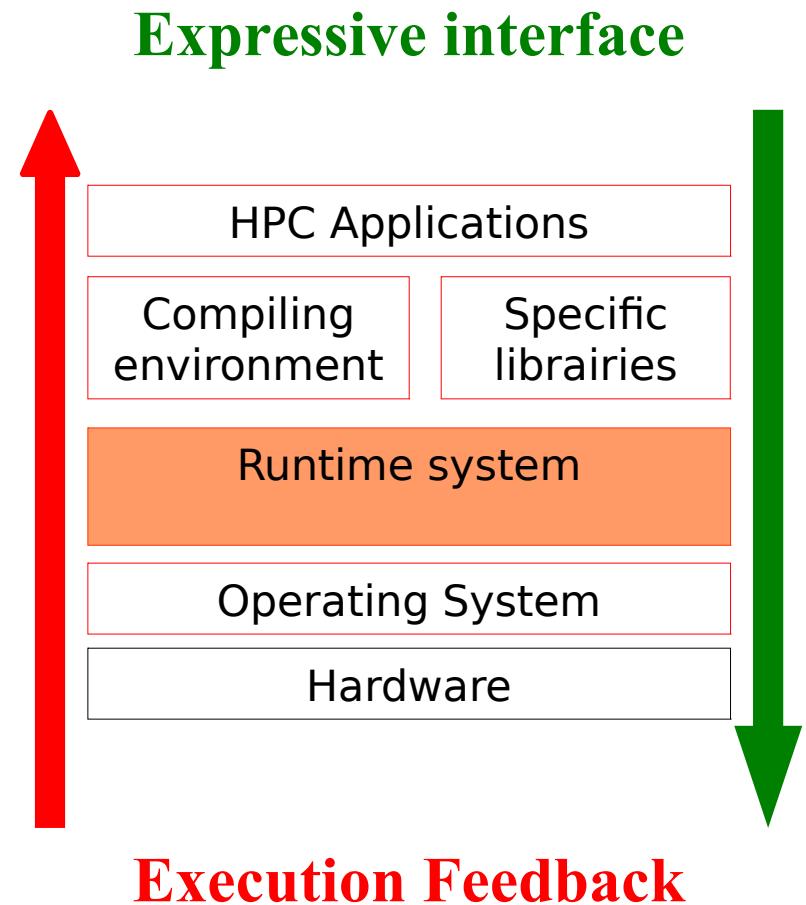
How to program these architectures?

- A uniform way
 - Use a single (or a combination of) high-level programming language to deal with network + multicore + accelerators
 - Increasing number of directive-based languages
 - Use simple directives... and good compilers!
 - XcalableMP
 - HMPP
 - StarSs
 - Much better potential for *composability*
 - If compiler is clever!



Challenging issues at all stages

- Applications
 - Programming paradigm
 - BLAS kernels, FFT, ...
- Compilers
 - Languages
 - Code generation/optimization
- Runtime systems
 - Resources management
 - Task scheduling
- Architecture
 - Memory interconnect

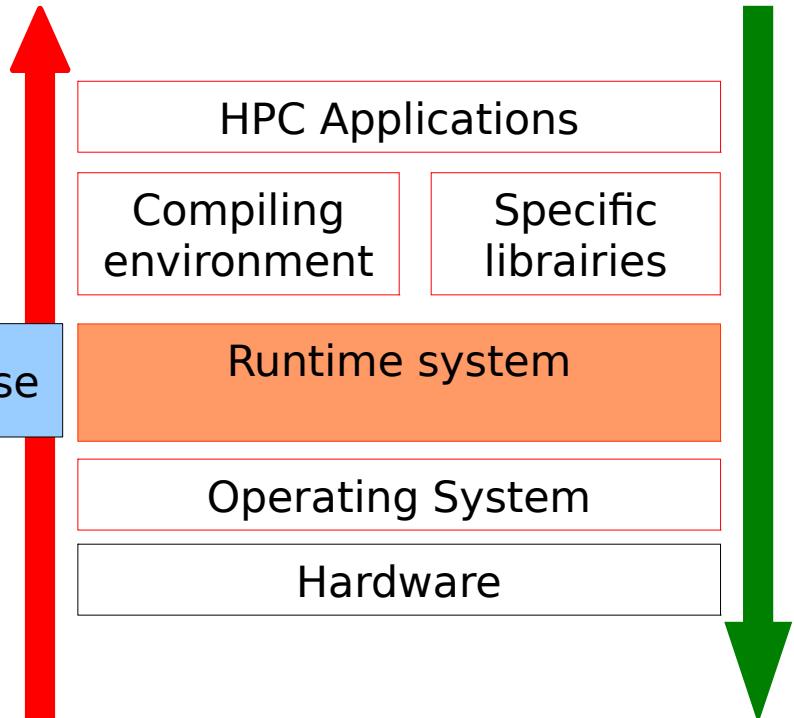


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Scheduling expertise

Expressive interface



Execution Feedback

Overview of StarPU

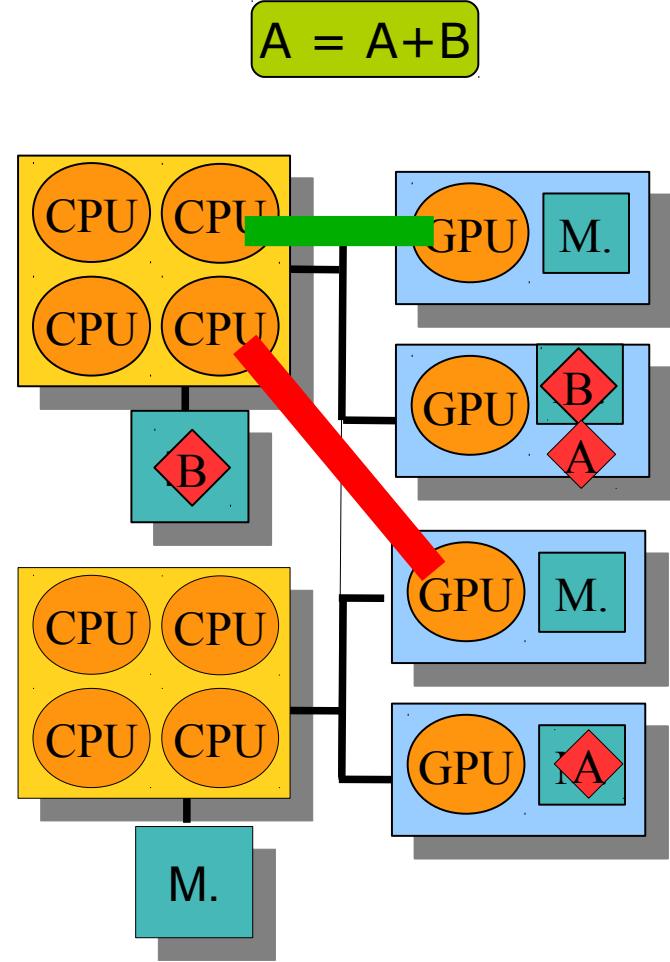
Rationale

Task scheduling

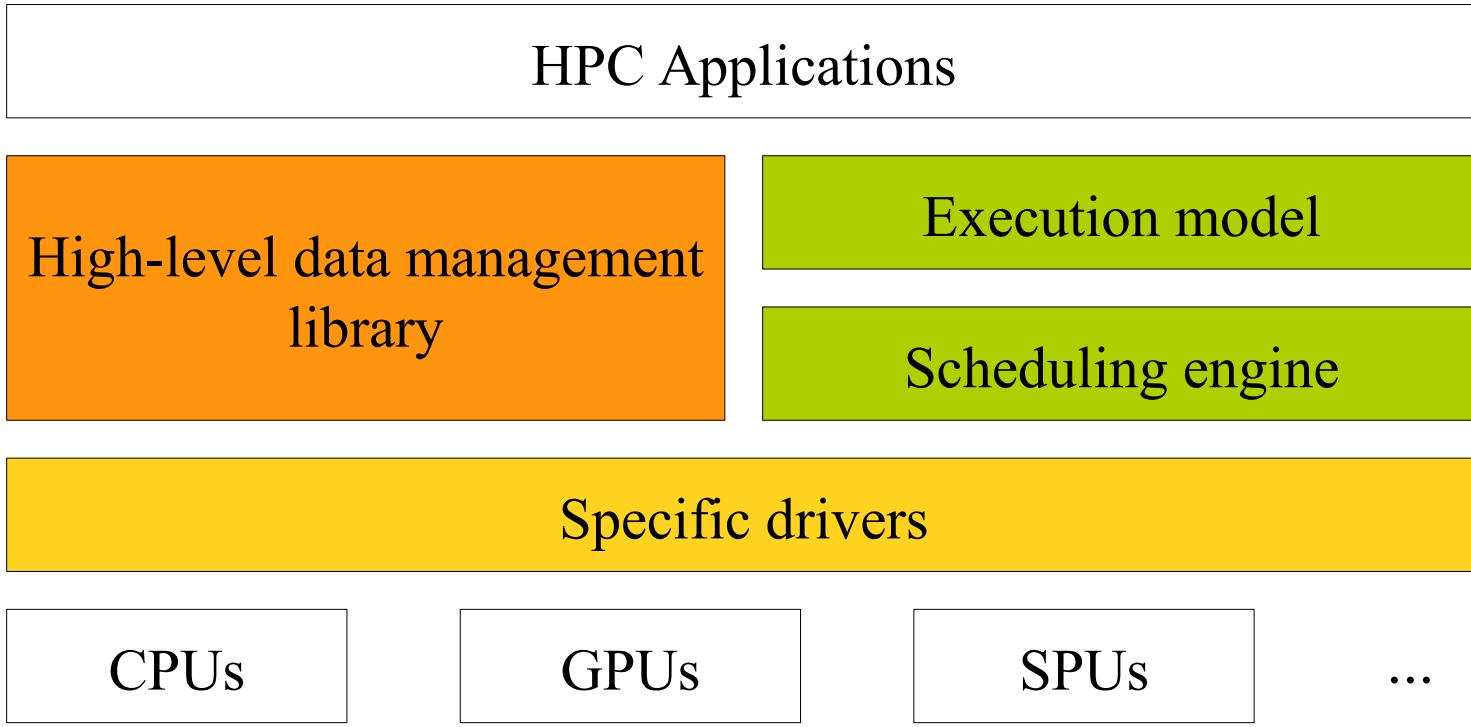
- Dynamic
- On all kinds of PU
 - General purpose
 - Accelerators/specialized

Memory transfer

- Eliminate redundant transfers
- Software VSM (Virtual Shared Memory)



The StarPU runtime system

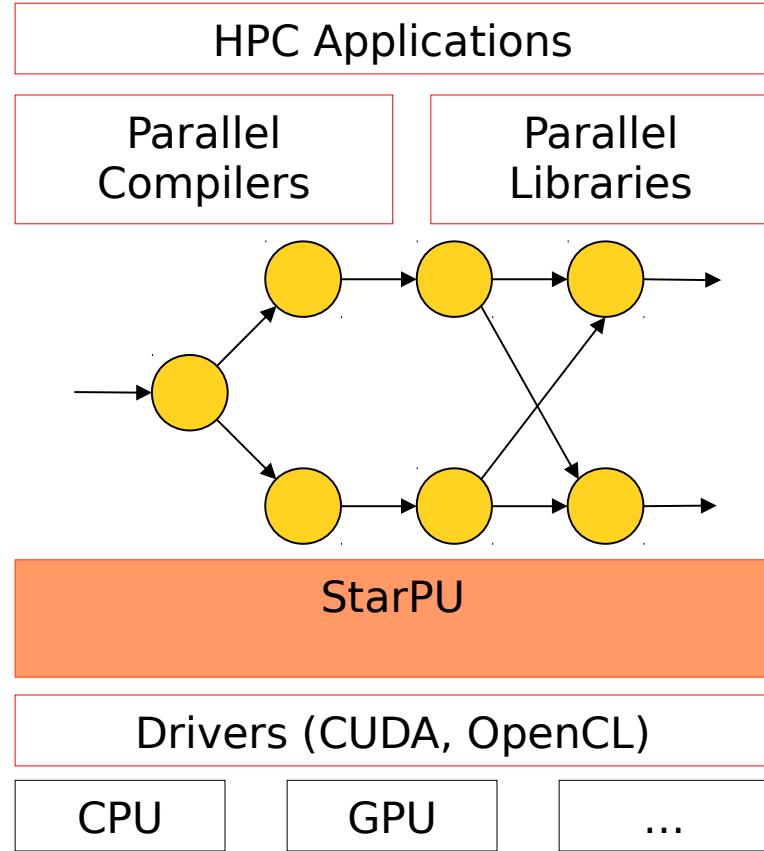


Mastering CPUs, GPUs, SPUs ... *PUs → **StarPU**

The StarPU runtime system

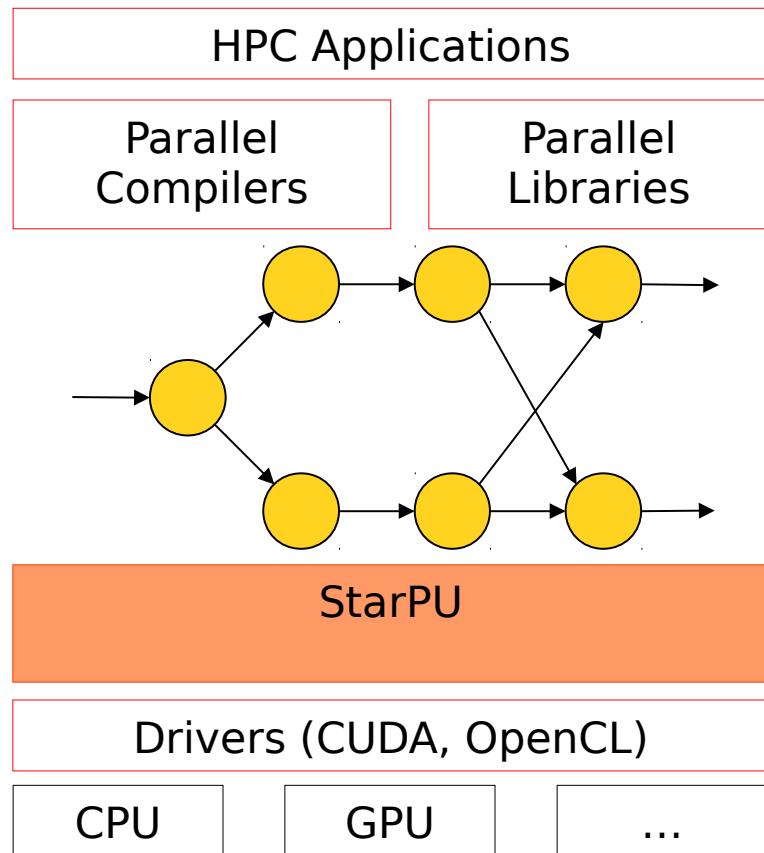
The need for runtime systems

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



Data management

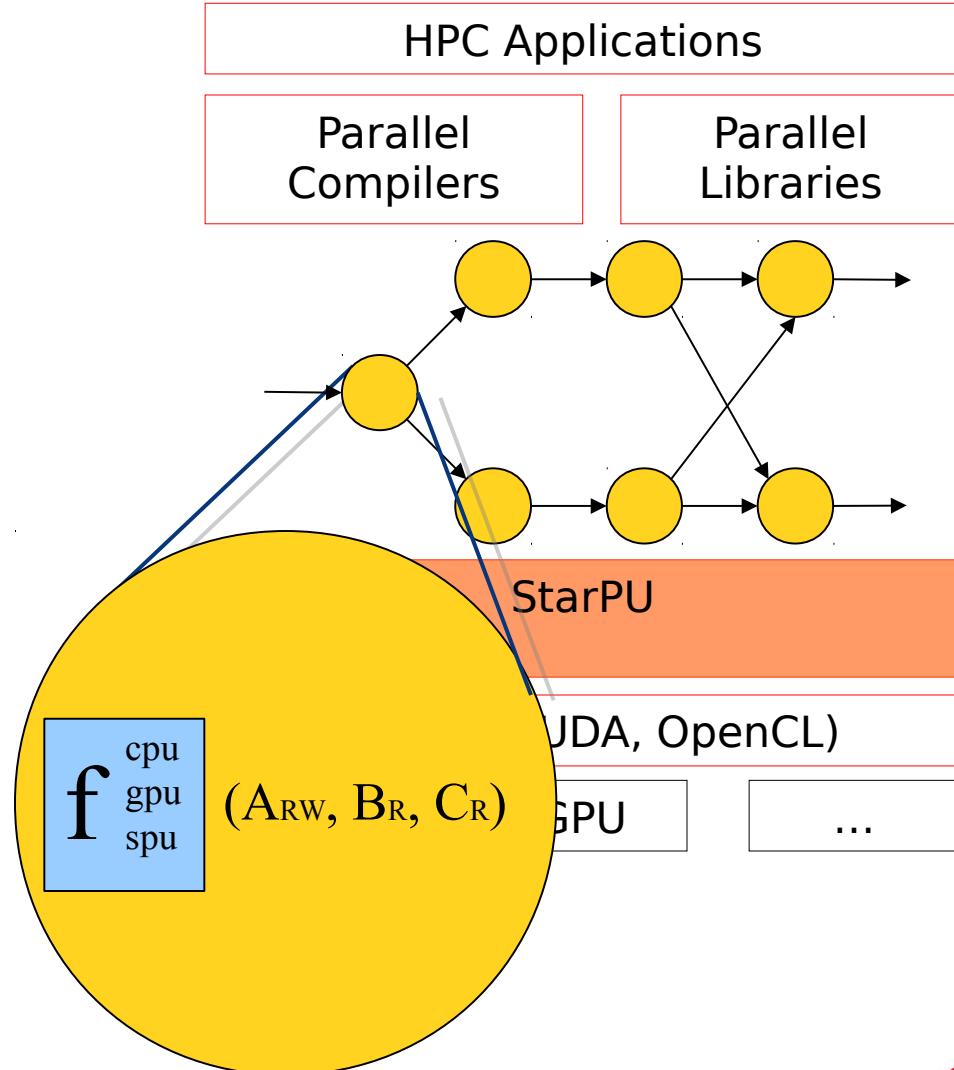
- StarPU provides a **Virtual Shared Memory (VSM)** subsystem
 - Replication
 - Weak consistency
 - Single writer
 - High level API
 - Partitioning filters
- Input & output of tasks = reference to VSM data



The StarPU runtime system

Task scheduling

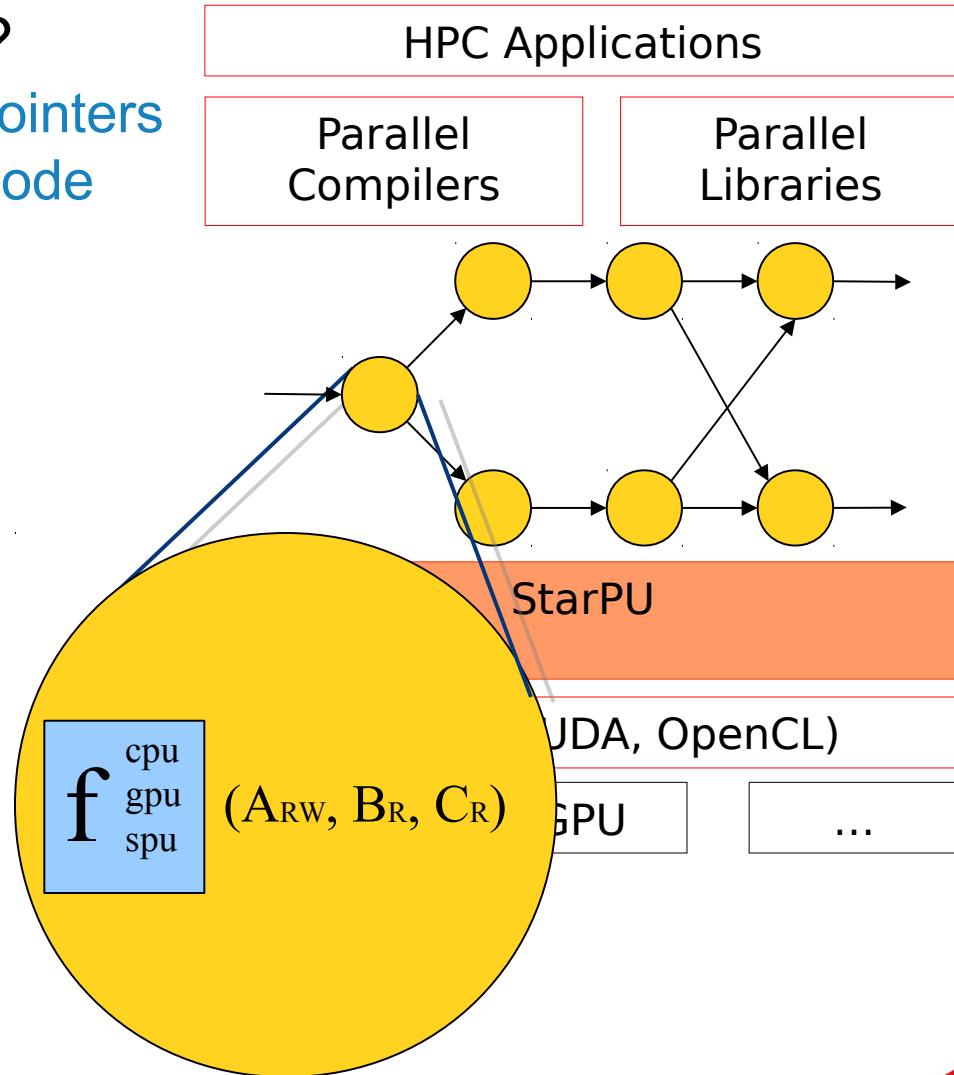
- Tasks =
 - Data input & output
 - Reference to VSM data
 - Multiple implementations
 - E.g. CUDA + CPU implementation
 - Non-preemptible
 - Dependencies with other tasks
 - Scheduling hints
- StarPU provides an **Open Scheduling platform**
 - Scheduling algorithm = plug-ins



The StarPU runtime system

Task scheduling

- Who generates the code ?
 - StarPU Task \approx function pointers
 - StarPU doesn't generate code
- Libraries era
 - PLASMA + MAGMA
 - FFTW + CUFFT...
- Rely on compilers
 - PGI accelerators
 - CAPS HMPP...



Task management

Implicit task dependencies

- Right-Looking Cholesky decomposition (from PLASMA)

For ($k = 0 \dots \text{tiles} - 1$)

{

POTRF($A[k,k]$)

for ($m = k+1 \dots \text{tiles} - 1$)

TRSM($A[k,k], A[m,k]$)

for ($m = k+1 \dots \text{tiles} - 1$)

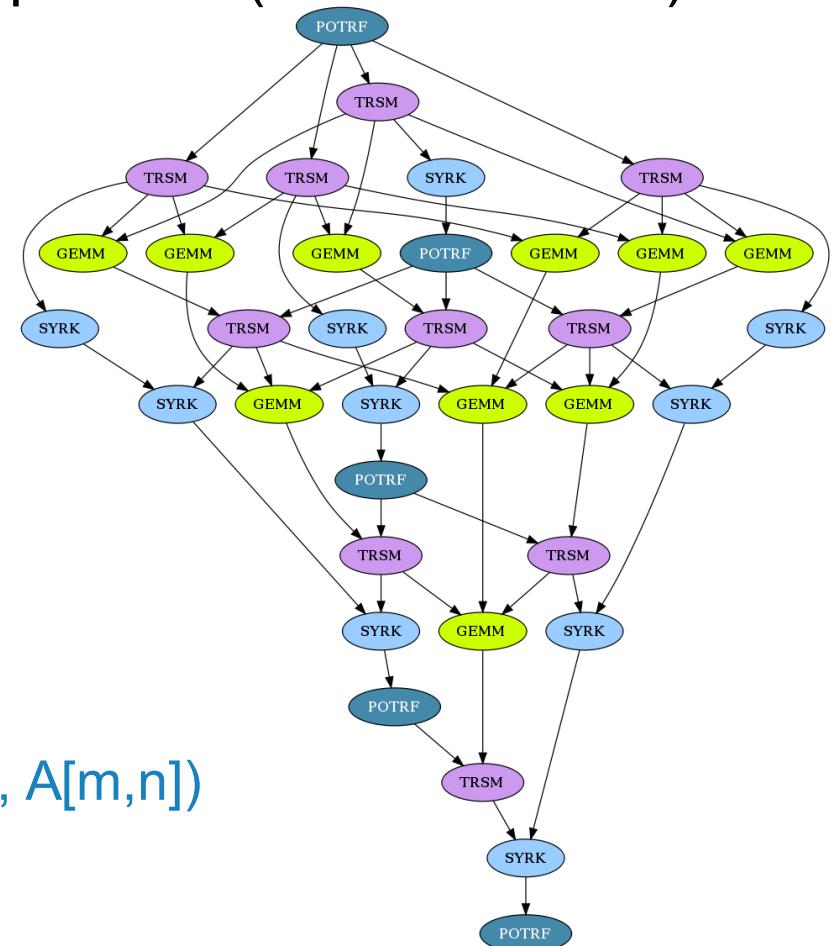
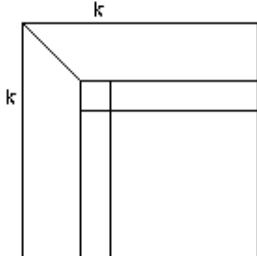
SYRK($A[m,k], A[m,m]$)

for ($m = k+1 \dots \text{tiles} - 1$)

for ($n = k+1 \dots m - 1$)

GEMM($A[m,k], A[n,k], A[m,n]$)

}



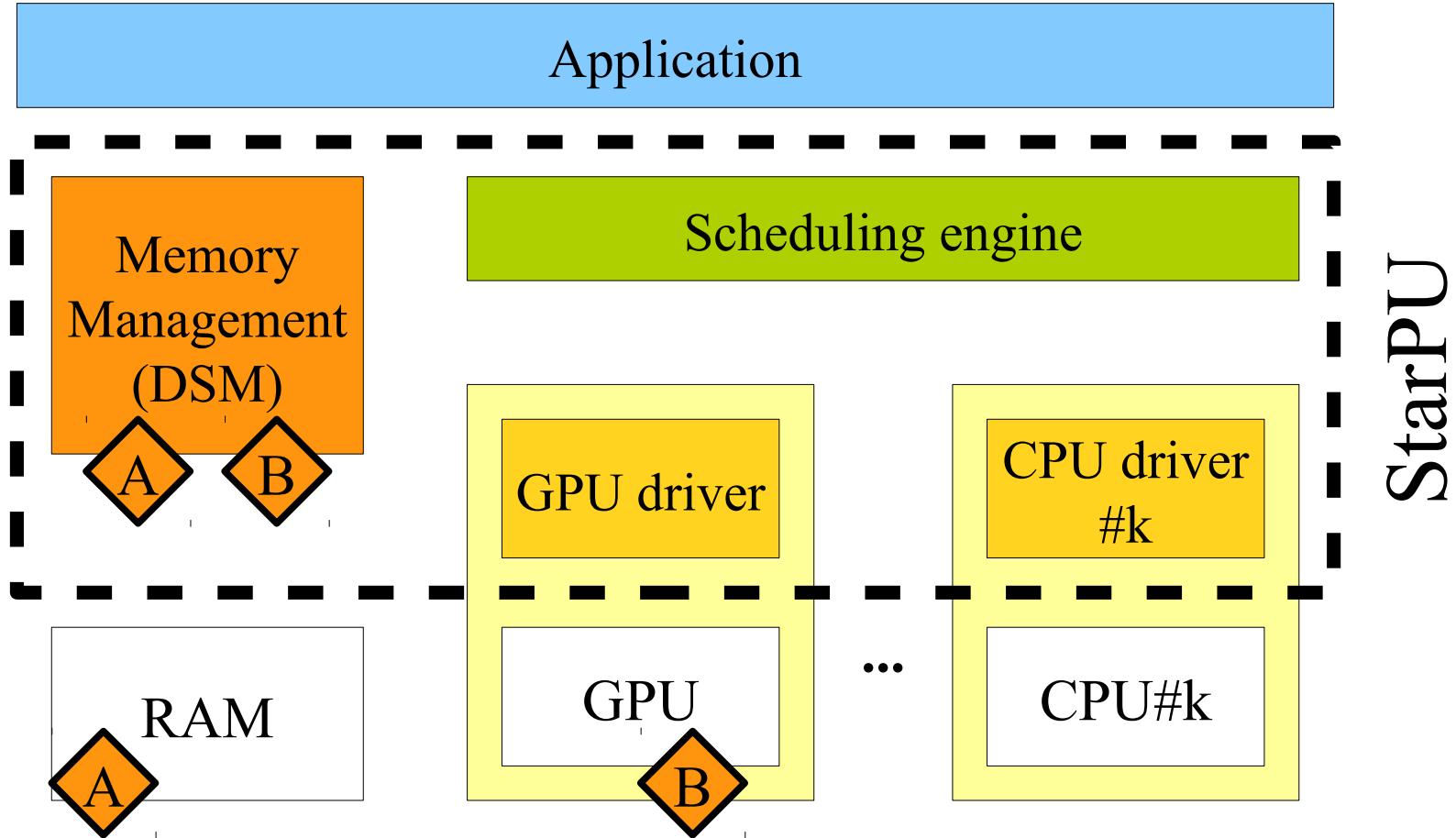
The StarPU runtime system

Development context

- History
 - Started about 6 years ago
 - PhD Thesis of Cédric Augonnet
 - StarPU main core ~ 40k lines of code
 - Written in C
- Open Source
 - Released under LGPL
 - Sources freely available
 - svn repository and nightly tarballs
 - See <http://runtime.bordeaux.inria.fr/StarPU/>
 - Open to external contributors
- [HPPC'08]
- [Europar'09] – [CCPE'11],... >400 citations

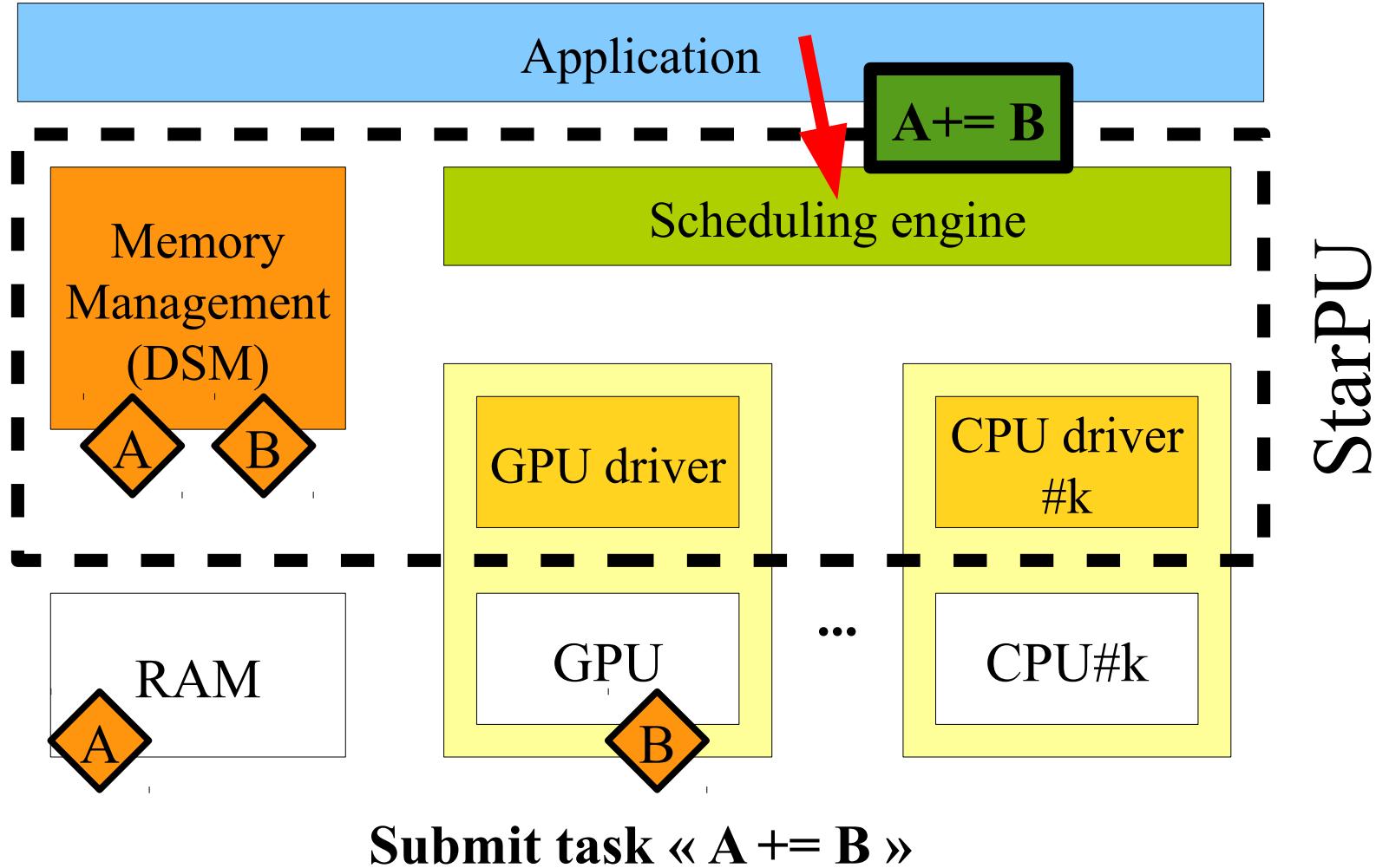
The StarPU runtime system

Execution model



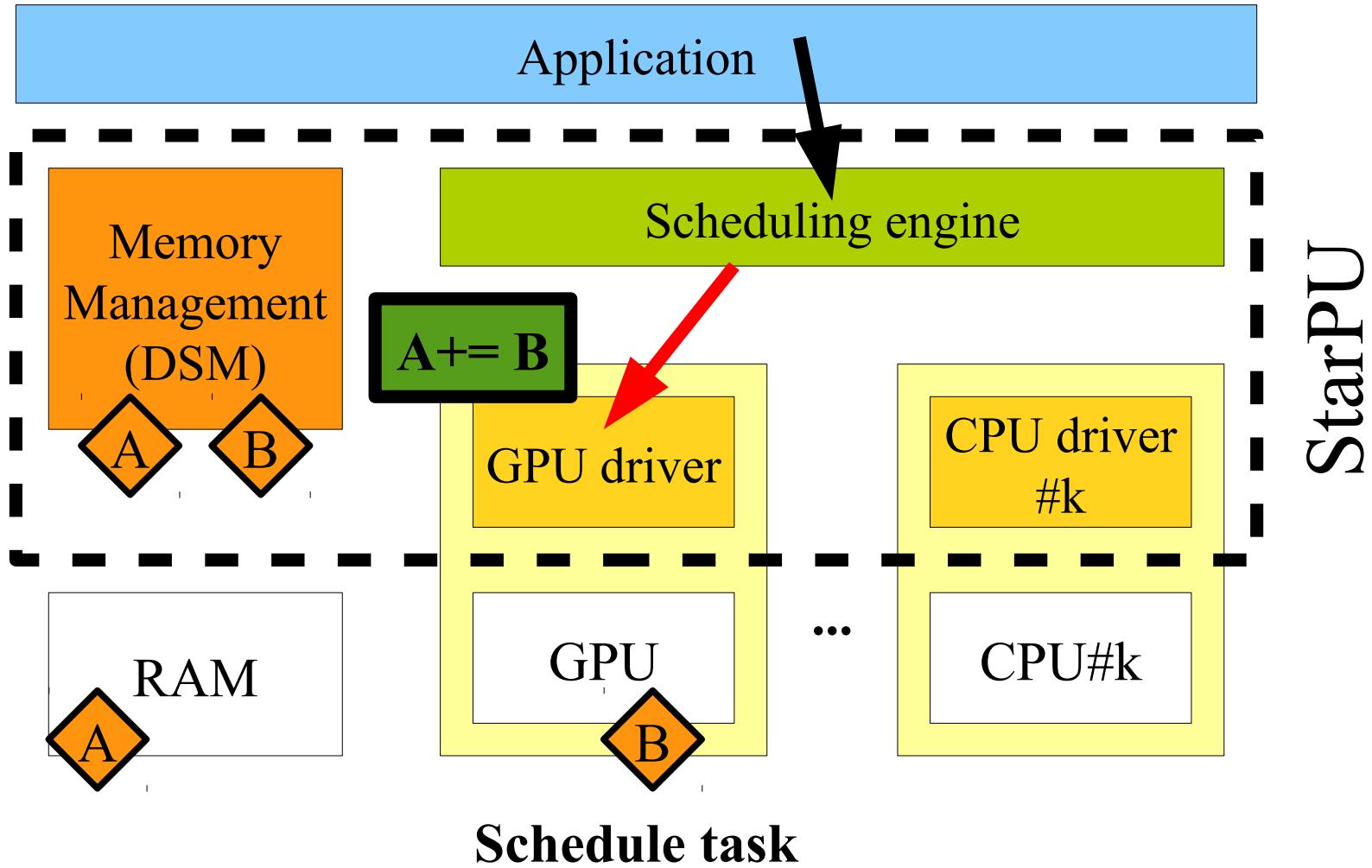
The StarPU runtime system

Execution model



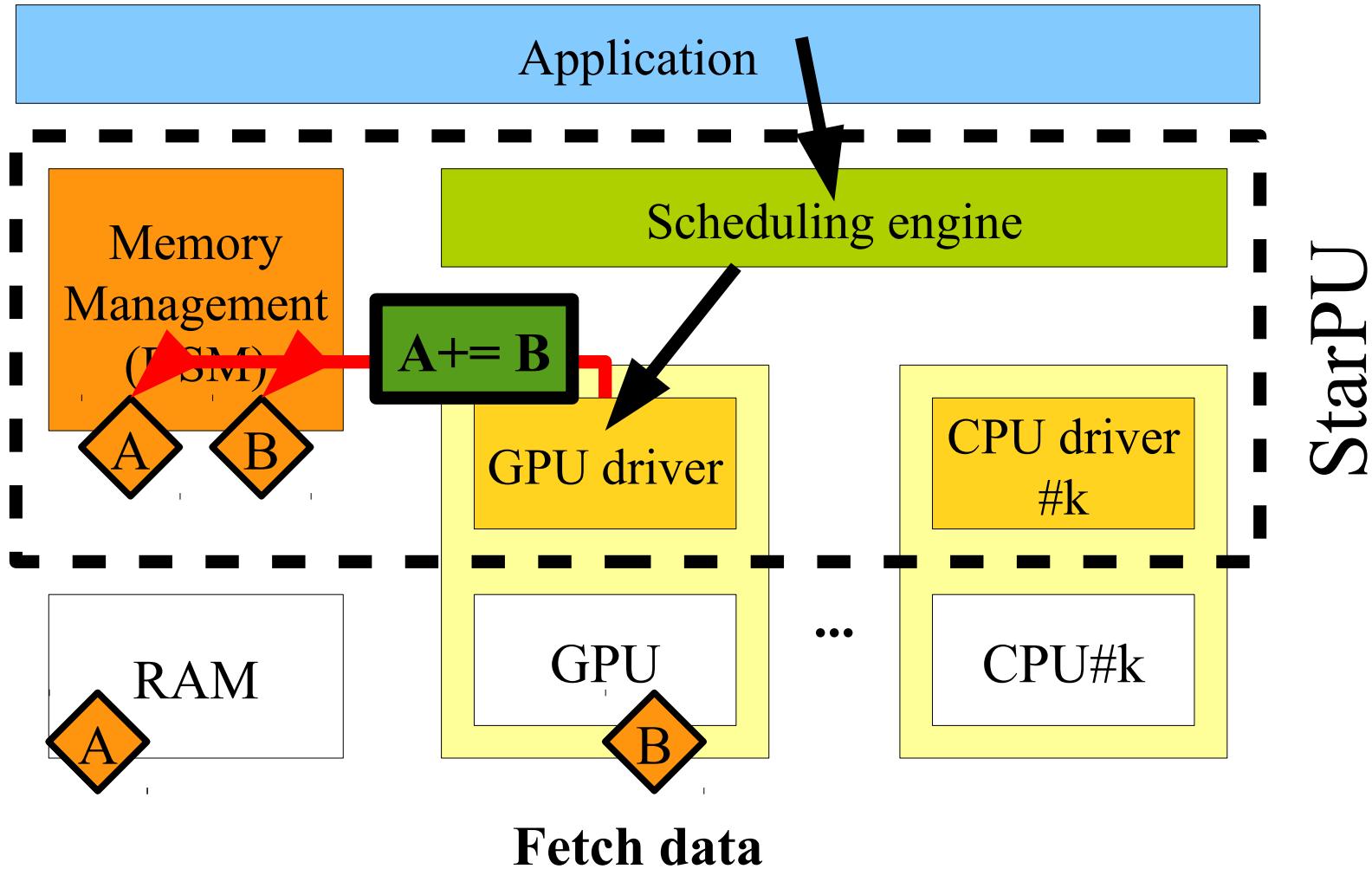
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Execution model



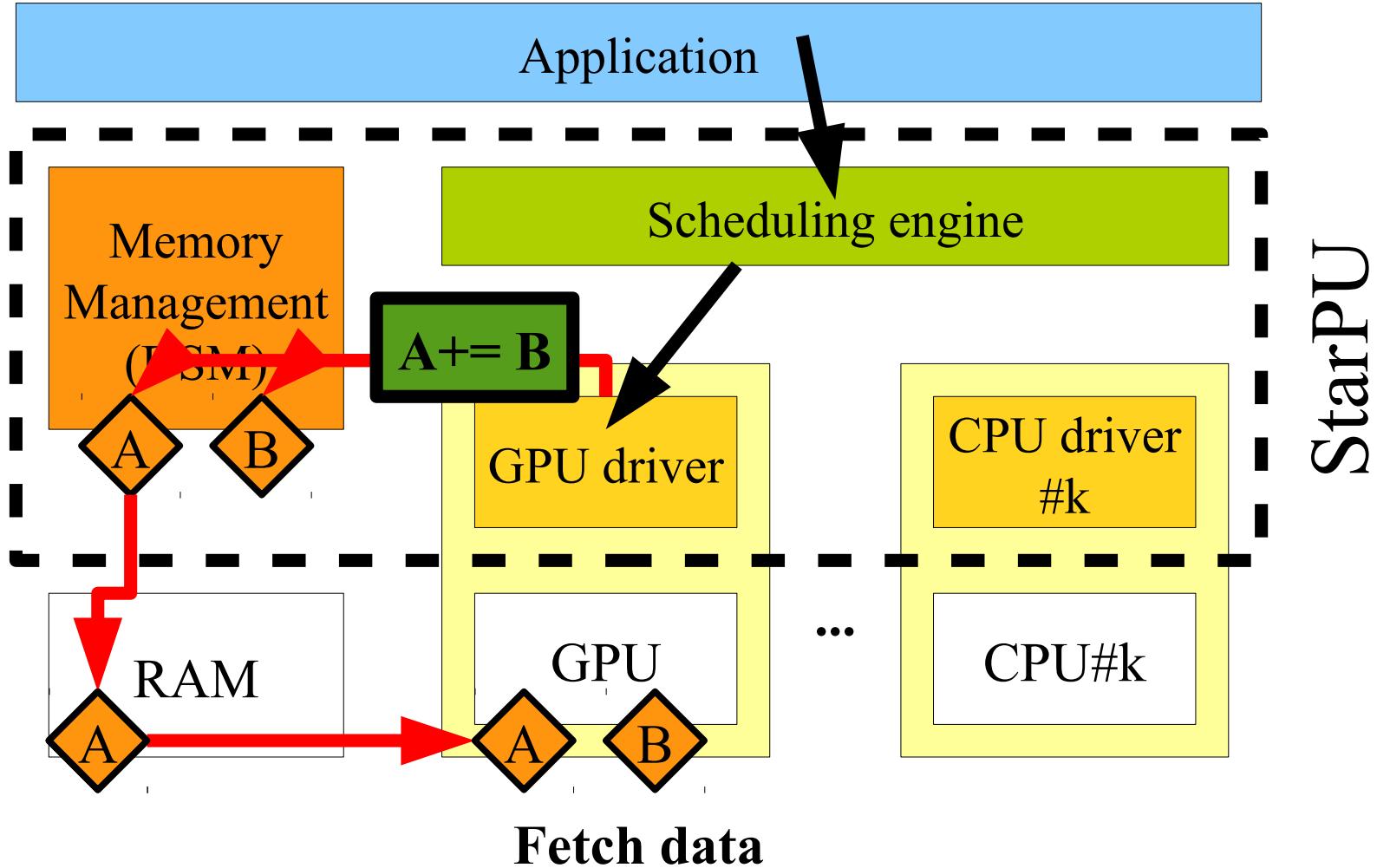
The StarPU runtime system

Execution model



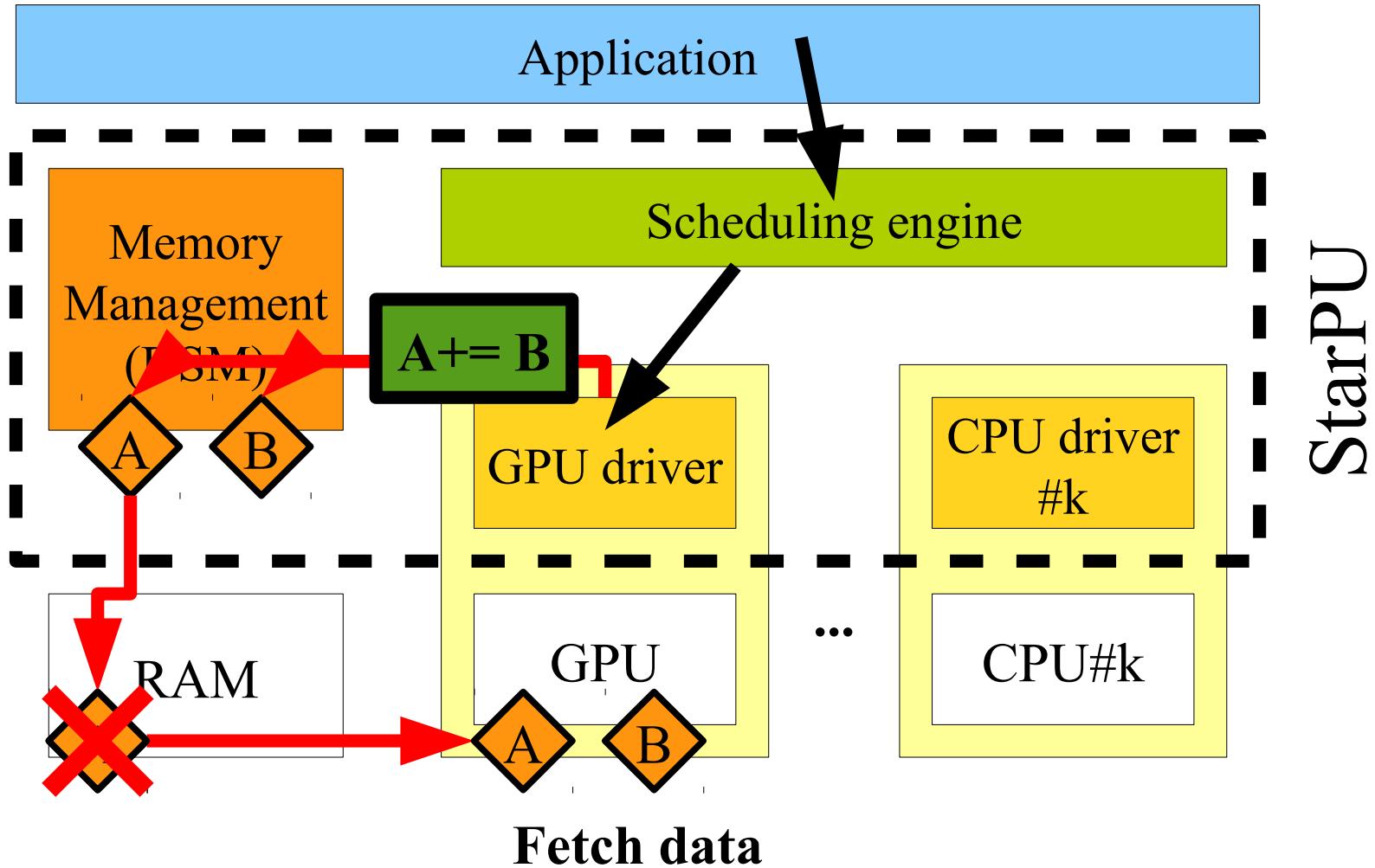
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Execution model



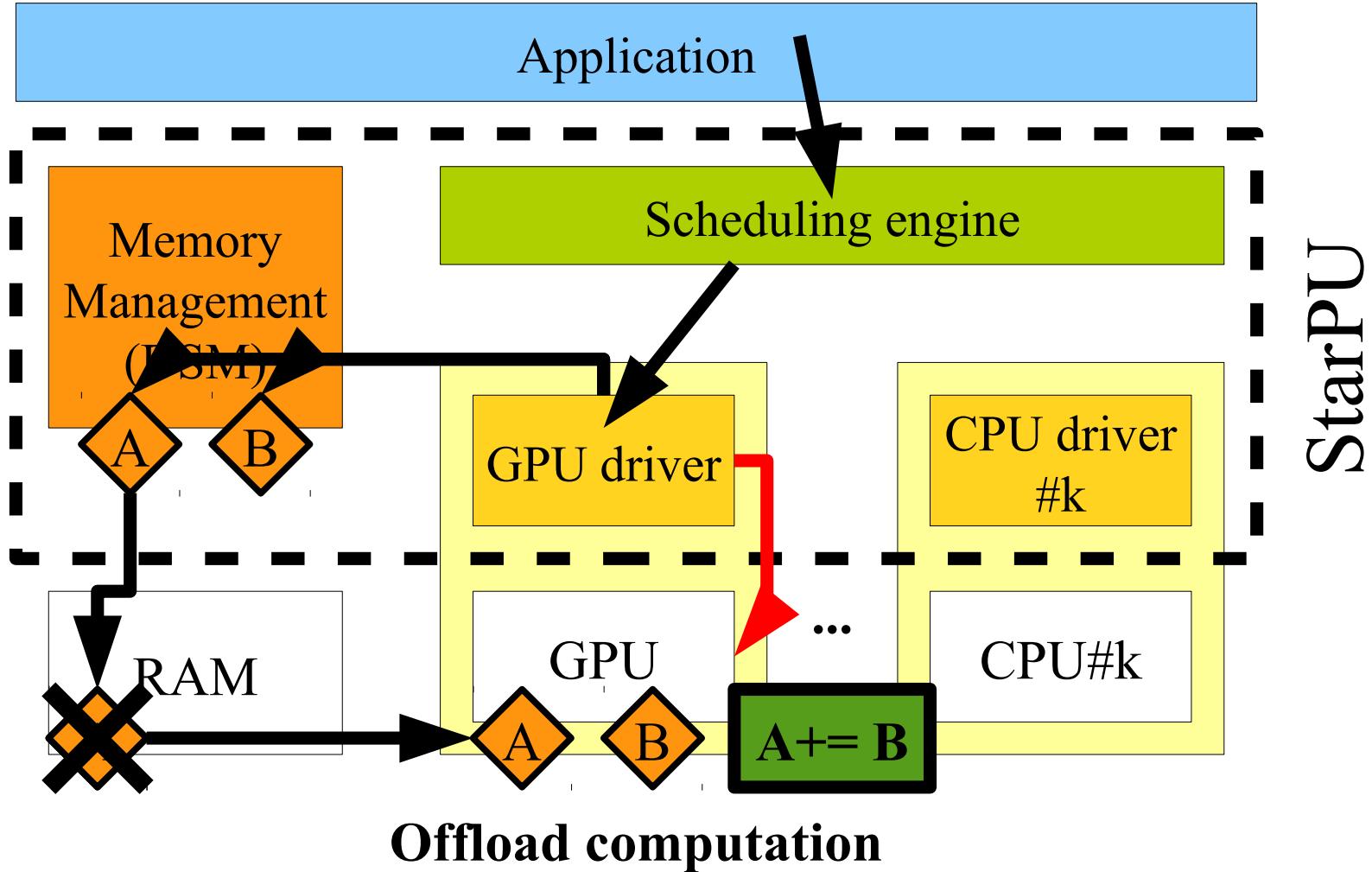
The StarPU runtime system

Execution model



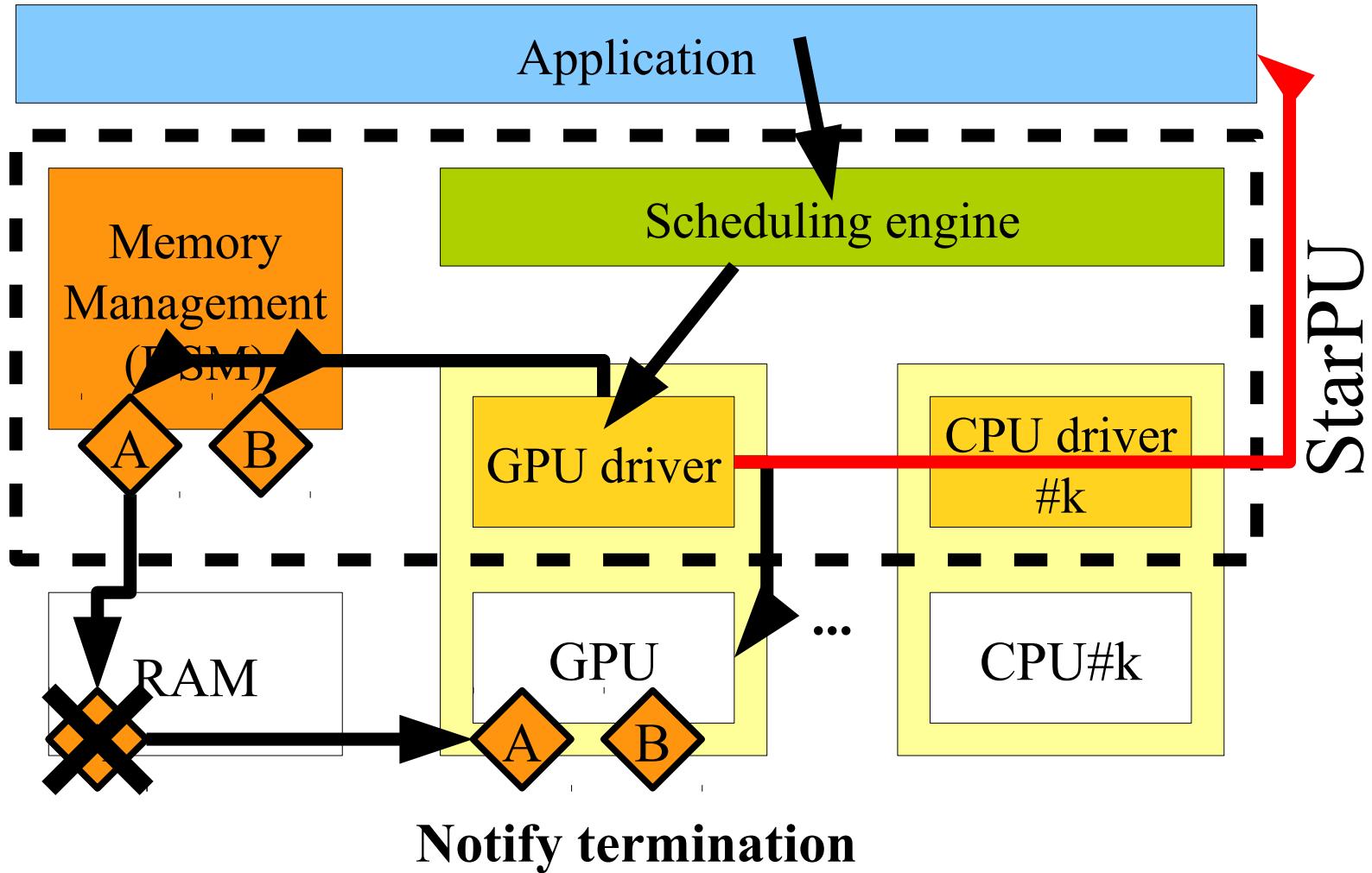
The StarPU runtime system

Execution model



The StarPU runtime system

Execution model



Optimizations

- Task pipelining
- Task execution / data transfer overlap
- GPU-GPU copies
- Data prefetch
- ...

Thus needs

- Asynchronous API with fine-grain synchronization
- Non-blocking API
- Pitched 2D copy & such
- Thread safety

Host memory mapped on GPU & vice-versa is useful, too

Programming interface

Scaling a vector

Data registration

- Register a piece of data to StarPU

- `float array[NX];`
`for (unsigned i = 0; i < NX; i++)`
`array[i] = 1.0f;`

```
starpu_data_handle vector_handle;  
starpu_vector_data_register(&vector_handle, 0,  
    array, NX, sizeof(vector[0]));
```

- Unregister data
 - `starpu_data_unregister(vector_handle);`

Scaling a vector

Defining a codelet (4)

- Codelet = multi-versionned kernel
 - Function pointers to the different kernels
 - Number of data parameters managed by StarPU

```
starpu_codelet scal_cl = {  
    .cpu_func = scal_cpu_func,  
    .cuda_func = scal_cuda_func,  
    .opencl_func = scal_opencl_func,  
    .nbuffers = 1,  
    .modes = STARPU_RW  
};
```

Scaling a vector

Defining a task

- Define a task that scales the vector by a constant

```
struct starpu_task *task = starpu_task_create();
task->cl = &scal_cl;
```

```
task->buffers[0].handle = vector_handle;
```

```
float factor = 3.14;
task->cl_arg = &factor;
task->cl_arg_size = sizeof(factor);
```

```
starpu_task_submit(task);
starpu_task_wait(task);
```

Scaling a vector

Defining a task, starpu_insert_task helper

- Define a task that scales the vector by a constant

```
float factor = 3.14;
```

```
starpu_insert_task(  
    &scal_cl,  
    STARPU_RW, vector_handle,  
    STARPU_VALUE,&factor,sizeof(factor),  
    0);
```

Scaling a vector

Defining a task, gcc plugin

```
void scale_vector(int size, float vector[size], float factor)
    __attribute__((task));
void scale_vector_cpu(int size, float vector[size], float factor)
    __attribute__((task_implementation("cpu", scale_vector))) ;
void scale_vector_cpu(int size, float vector[size], float factor)
{ ... }
int main(void) {
    static float input[NX];
#pragma starpu register input
    scale_vector(NX, input, 42);
#pragma starpu wait
#pragma starpu unregister input
}
```

Scaling a vector

Defining a task, gcc plugin

```
void scale_vector(int size, float vector[size], float factor)
    __attribute__((task));
void scale_vector_cpu(int size, float vector[size], float factor)
    __attribute__((task_implementation("cpu", scale_vector))) ;
void scale_vector_cpu(int size, float vector[size], float factor)
{ ... }
int main(void) {
    static float input[NX];
#pragma starpu register input
    scale_vector(NX, input, 42);
    frob_vector(NX, input, out1);
    shred_vector(NX, input, out2);
#pragma starpu wait
#pragma starpu unregister input
}
```

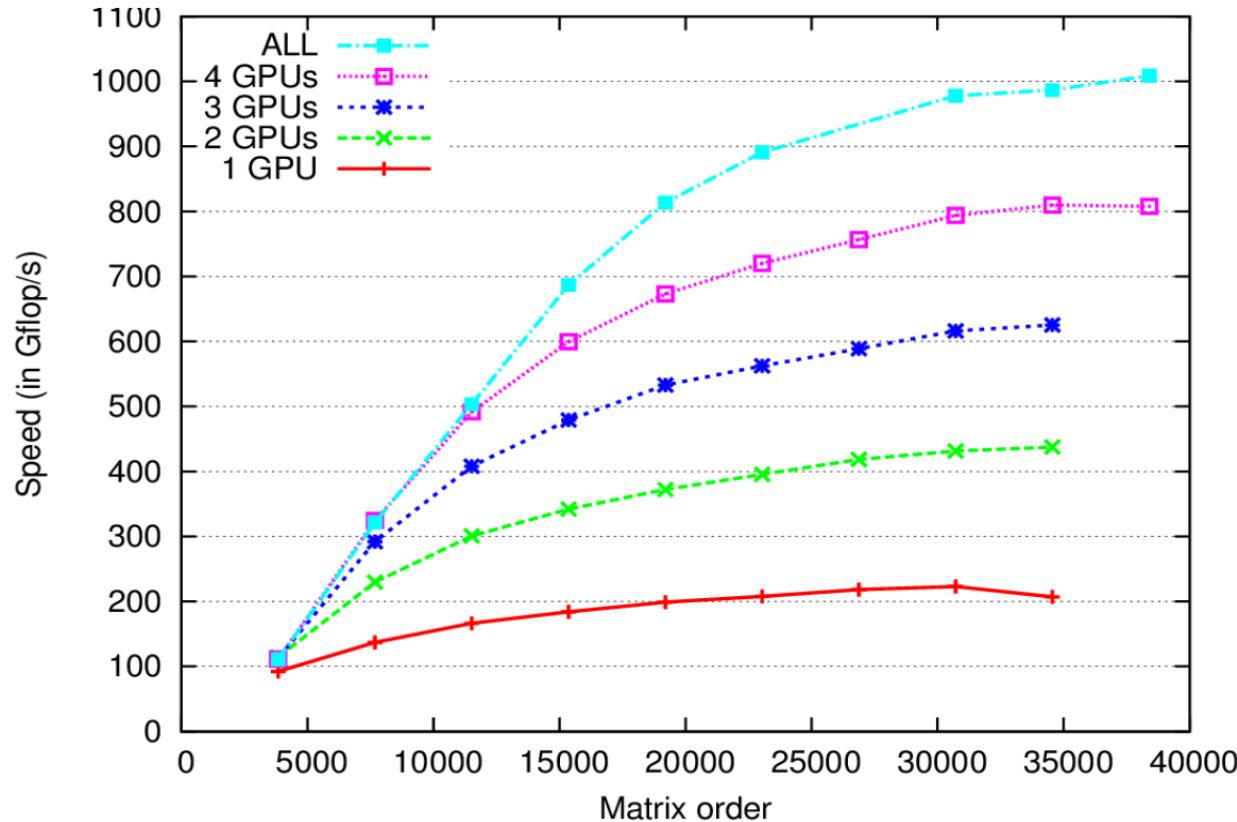
Scaling a vector

Defining a task, gcc plugin

```
void scale_vector(int size, float vector[size], float factor)
    __attribute__((task));
void scale_vector_opencl(int size, float vector[size], float factor)
    __attribute__((task_implementation("opencl", scale_vector))) ;
#pragma starpu opencl scale_vector_opencl \
    "vector-scale.cl" "vector_scal_kern" \
    group_size ngroups ;
```

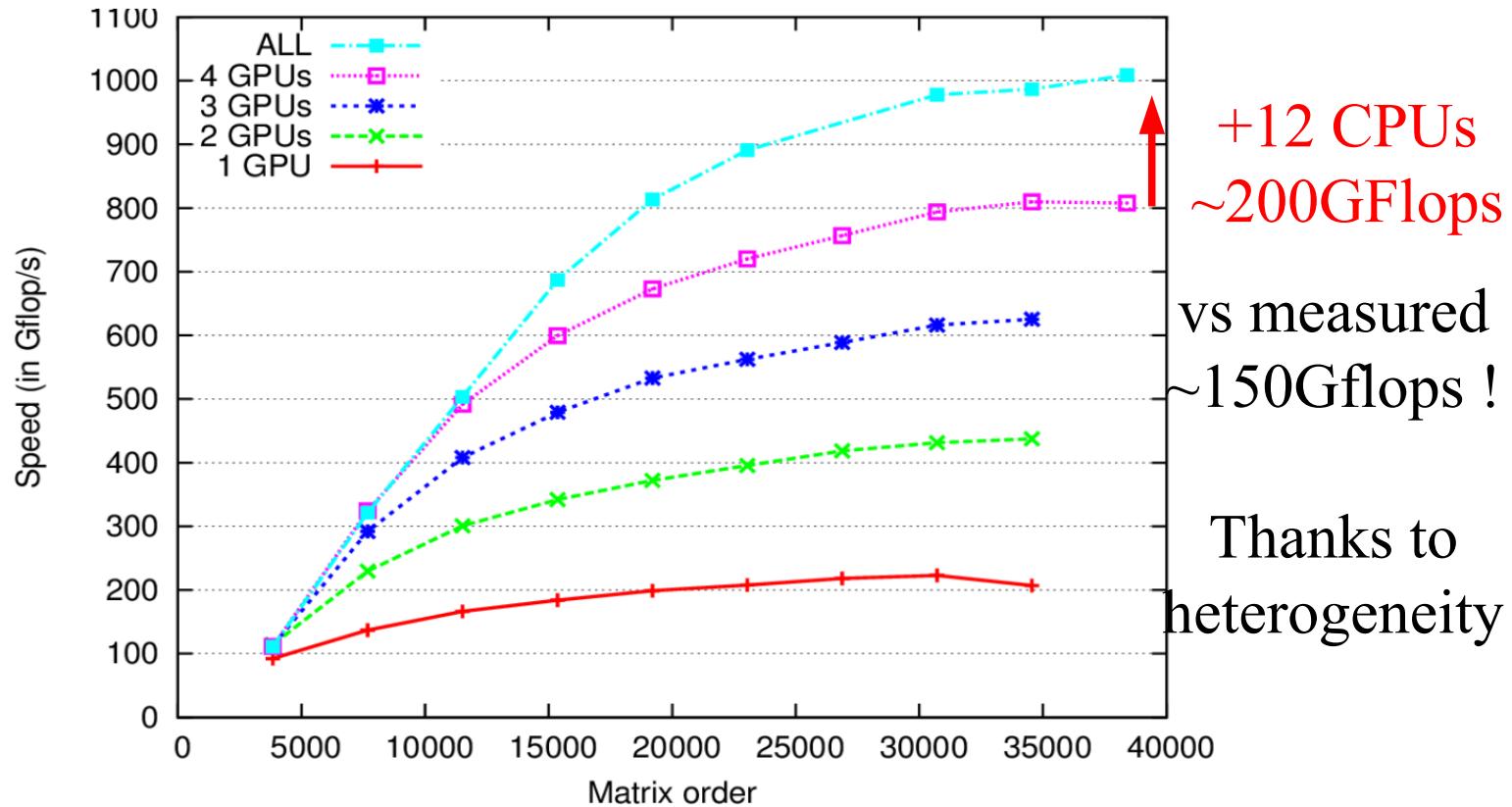
Mixing PLASMA and MAGMA with StarPU

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



Mixing PLASMA and MAGMA with StarPU

- QR decomposition
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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

Conclusion

Summary

Tasks

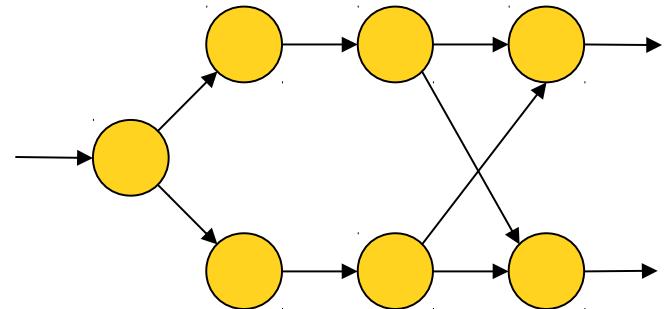
- Nice programming model
- Runtime playground
- Scheduling playground
- Algorithmic playground
- Used for various computations
 - Cholesky/QR/LU (dense/sparse), FFT, stencil, CG, FMM...
- <http://starpu.gforge.inria.fr>

Scheduling expertise

HPC Applications

Parallel
Compilers

Parallel
Libraries



Runtime system

Operating System

CPU

GPU

...

Conclusion

Summary

Scheduling researchers can experiment and tune various heuristics

- On actual applications
- Without even needing the hardware
 - And with fast experimentation time

Optimize

- Completion time
- Memory consumption
- Energy consumption
- ...
- <http://starpu.gforge.inria.fr>

Scheduling expertise

