

Applying GPUs for massive calculations Determination of suitable applications

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Personal introduction

Experience within Philips (focused on technical part)

- Professional systems (mostly PC based)
- Real-time embedded consumer systems
- IC-design SW, NP-complete algorithms

Teaching:

- Execution Architecture course
- SW reliability
- · 3TU / OOTI embedded computing platform

Leading the Apptech Performance team

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Introduction

• Main GPU experience: nVidia, G80, Tesla, CUDA.

→ Advanced GPU optimization course for performance team

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- · This talk is 25 minutes: select and focus
 - GPU, GPGPU: General Purpose is an overstatement, so
 - Determination of suitable software \rightarrow suitable algorithm
 - A suitable algorithm is fast and energy efficient.
- · Skip a lot of GPU introduction, given previous speakers presentations.

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Suitable software



- A GPU is a parallel co-processor for a CPU, doing one task at a time.
- A GPU consists of ~32 multi-processors (MP)
 →Single Tasking Multiprocessing
- A multi-processor is a SIMD processor
 - 8x in HW, 32x fault in SW (4-stage pipeline).
- 1 algorithm applied in parallel to multi-data sets
 - Multi-threading: all threads execute the same code.
 - Vector processor is therefore a better description.
- Data-dependent if-then-else : MP serializes then + else (unless)
 Sorting / filtering based on ITE less suitable (alike control SW) →Number crunching

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Contents

- 1. Introduction
- 2. Suitable software
- 3. Suitable parallelism
- 4. Global modeling of algorithms
 - Intro & example
 - Steps, targets, how-to
 - Practice
- 5. Optimization, recognizing a bottleneck
- 6. Conclusion

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Introduction - overview



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Suitable parallelism

Estimation of degree of suitable number of data sets

- HW-Theoretical: 32 x 32 ≈ 1000
- · Overhead sources are relatively large (but undocumented), e.g.,
 - GPU data loading
 - GPU code loading (kernel) and kernel invocations
 - PC driver cost (~ 10 30 us = 30 MFLOP on Tesla)
- Practical required degree of parallelism: >> 1.000, e.g. 30.000+
 Depends on typical gpu-job time

→Massive parallel greedy number crunching jobs

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Global Modeling of Algorithms – simple example

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• Is a GPU suitable to calculate matrix C = A+B ?



• Exploiting 10⁻⁶ of the FLOP potential only.

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Global Modeling of Algorithms: intro and example

Convention: device = graphics card, host = PC cpu.

GPU (Tesla C1060) hardware characteristics

- 1 TFLOPS max (~87 GFLOPS for double precision)
 20 DVDs per
 second
- Data rate ~100 GB/s max (device transfers)
- Data rate ~3 GB/s max (device-host transfers)

Simple example: C = A + B (large n x n matrices, massive parallel)3 device-host word transfers, 3 word device transfers, 1 FLOP per word.Device – host transfer utilization: 1, i.e., 0.75 G C_{ij} per second.Device transfer utilization: 0.033FLOPS utilization: 0.75 x10⁻⁶

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Global Modeling - single precision

GPU characteristic for single precision:

- ~1 TFLOPS max
- Data rate ~25 G words/s max (device transfers)
- Data rate ~0.75 G words/s max (device-host transfers)

Key is the number of times a single data element is used in calculations

· When on device: ~40 calculations per data element required,

• When on host: ~1300 calculations per data element required .. before the calculation capacity becomes the bottleneck.

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Global Modeling: boundaries



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Global modeling – steps and target

Target
>> 1000
~TFLOPS
~40 dev, ~1300 host

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Global modeling - steps - how to

- Derive essential parallelism
 Outer loops may be parallelized
- Computations (limited focus in this phase)
- Data reference patterns
 optimize dataref reuse

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13

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Global modeling – data reference optimization

- Data reference pattern optimization:
- Reuse coefficients
- Reuse power calculations
- Reuse cross-terms
- This is the reason why per processor thousands registers are available, next to the 'shared memory'.

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Data referencing - examples

Algorithm:	host-dev	device
Matrix addition, subtraction, transposition:	no advantage	
 Matrix multiplication O(n³) → n² per element 	n > ~ 20	n > ~ 5
 Solve linear equations O(n³) → n² per element 	n > ~ 20	n > ~ 5
 2D FFT O(n²log n) → 5 n log n per element 	n > ~ 32	n > ~ 4

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The overhead per invocation is so large, that the first part of the range only applies when multiple matrices are handled per invocation.

Complex number calculations are more suitable: mult/div uses data twice.

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Summary of candidate SW

- Massive Parallel (>> 1000, e.g. 30.000)
- Computations (~ Teraflops)
- · Data transfer is typically the major bottleneck, suitable algorithms have
- many calculations per data element (~1000 per i/o, ~30 per device elt)
- GPU Jobsize should be substantial w.r.t. invocation overhead.
- · Set of algorithms of which data resides on the device increases suitability

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Practice: Sequence of algorithms

- Ineffective GPU alg's can be useful to prevent dev-host communication.
- Despite multiple kernel invocations
- Real iterating algorithms enhance greediness and limit dev-host communication.

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Next phase: optimization



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17

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14

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Optimization

- This part requires explaining the 10 bottlenecks of a GPU.
- · How to bypass them.
- A bit too much for now. It is in the optimization course.
- How to recognize a bottleneck is relatively simple, after the suitability of the algorithm has been investigated.

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Optimizations: how to recognize a bottleneck



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Optimization - some remarks

- Direct C to Cuda translation is easy, but is non-optimized.
 - Marketing of how easy and generic CUDA is, shoots in the foot.
 - Available numerical libraries quality is work-in-progress.
 - Positive impression of the support by nVidia.
- Potentially suitable algorithm typically required a complete reordering / rewriting, resulting in a speedup of factors ~100 w.r.t. first version, both for single- and double precision.

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21

Conclusions

- GPU is suitable for running extremely greedy calculations that run massively parallel, with limited cpu to gpu memory transfers, and lots of calculations per data element.
- Global modeling of an algorithm gives potential suitability, without knowledge of GPU internals.
- Though CUDA hardly contains hardware specifics, real performance requires in-depth understanding how CUDA runs in hardware, what are the bottlenecks.

22

• It's easy to notice if one suffers from a bottleneck.

