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## Lecture 12

Reduction

## A few more CUDA issues

## Sorting on GPU

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## Last time

- Coalescing
- Constant memory
- Texture memory

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## The world's simplest demo on texture memory?

texobjdemo.cu

Simple texture memory example.

Array of numbers, accessed at non-integer coordinates.

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## Upcoming and ongoing labs

This week: Lab 4: Intro to CUDA, Mandelbrot then

Lab 5: Image filtering.
Shared memory in focus!
Lab 6: Reduction and sorting with OpenCL.

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## Lecture questions

1) How can you efficiently compute the average of a dataset with CUDA?
2) In what way does bitonic sort fit the GPU better than many other sorting algorithms?
3) What is the reason to use pinned memory?
4) What problem does atomics solve?

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## Reduction

Parallelizing problems of limited parallel nature Problem seen in Kessler 1.3.1.4 and 1.5.2-1.5.4 Global sum.

## Examples of reduction problems

Extracting small data from larger

- Finding max or min
- Calculating median or average
- Histograms

Common problems!

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## Sequentially trivial

Loop through data
Add/min/max, accumulate results
Fits badly in massive parallelism!

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## Tree-based approach



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## In 2D, typically 4-to-1 per level

## Pyramid hierarchy



## Tree-based approach

# Each level parallel! Can be split onto large numbers of threads 

## but

the parallelism is reduced for each level, and the results need to be reorganized to a smaller number of threads!

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etc
16
8

4

2
1


# Multiple kernel runs for varying size! 

## For $\mathbf{n}=\mathbf{k}$ downto 0 do Launch $2^{\text {n }}$ kernels

Multiple levels can be merged into one - but not all of them!

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# Again: You can not synchronize between blocks! 

## Not all blocks are simultaneously active

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## Multiple levels per kernel run for avoiding overhead



## Doubly interesting due to study with many optimizations:

## Many possibilities:

- Avoid "if" statements, divergent branches - Avoid bank conflicts in shared memory
- Loop unrolling to avoid loop overhead (classic old-style optimization!)


# Huge speed difference reported by Harris 

|  | Time ( $\mathbf{2}^{22}$ ints) | Bandwidth | Step Speedup | Cumulative Speedup |
| :---: | :---: | :---: | :---: | :---: |
| Kernel 1: <br> interleaved addressing with divergent branching | 8.054 ms | 2.083 GB/s |  |  |
| Kernel 2: <br> interleaved addressing with bank conflicts | 3.456 ms | 4.854 GB/s | 2.33x | 2.33x |
| Kernel 3: sequential addressing | 1.722 ms | 9.741 GB/s | 2.01x | 4.68x |
| Kernel 4: <br> first add during global load | 0.965 ms | 17.377 GB/s | 1.78x | 8.34x |
| Kernel 5: <br> unroll last warp | 0.536 ms | 31.289 GB/s | 1.8x | 15.01x |
| Kernel 6: <br> completely unrolled | 0.381 ms | 43.996 GB/s | 1.41x | 21.16x |
| Kernel 7: <br> multiple elements per thread | 0.268 ms | 62.671 GB/s | 1.42x | 30.04x |

However, some of these optimizations are no longer valid.

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## Alternative: Reduction in many levels, but making sure idle threads are dense!

With every other thread idle/finished half the performance.

With every other warp idle finished good performance!

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## Skip every other thread over and over in same kernel - waste!



## Keep active threads together - better!



Threads and memory both behave like this for coalescing.

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## Divergent branching =

 "if" statements:
## Branches can be bad in GPU code!

Why?

## Divergent branching in SIMD:

All branches execute all code! Data masked with result of "if".

Warp-level problem!
Can not be avoided within warps if a single thread gets a different result from others. Can be avoided if all threads in warp take same branch

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## Divergent warp

if $X$ then 10010110
I
I and with 10010110
else
I
I and with 01101001
I
endif

Non-divergent warp
if $X$ then 11111111
I
|
else
I
I
endif


## Conclusions:

- Multiple kernel runs for varying problem size
- Multiple kernel runs for synchronizing blocks
- Optimizing matters! Not only shared memory and coalescing!

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# More memory <br> Managed memory 

Atomics
Pinned memory

## Managed memory

## Makes read/write memory as easy as constant!

New, simpler Hello World!

```
#include <stdio.h>
const int N = 16;
const int blocksize = 16;
__global_
void hello(char *a, int *b)
{
a[threadldx.x] += b[threadIdx.x];
}
```

```
int main()
```

int main()
{
{
printf("%s", a);
printf("%s", a);
dim3 dimBlock( blocksize, 1 );
dim3 dimBlock( blocksize, 1 );
dim3 dimGrid( 1, 1 );
dim3 dimGrid( 1, 1 );
hello<<<dimGrid, dimBlock>>>(a, b);
hello<<<dimGrid, dimBlock>>>(a, b);
cudaDeviceSynchronize(); // Synchronize
cudaDeviceSynchronize(); // Synchronize
printf("%s\n", a);
printf("%s\n", a);
return EXIT_SUCCESS;
return EXIT_SUCCESS;
}
}
__managed
char a[N] = "Hello \0\0\0\0\0\0";
__managed__ int b[N] = {15,10,6,0,-11, 1, 0, 0, 0,0,0,0,0,0,0,0};

```

\section*{Managed memory}

Managed memory must be declared __managed
Memory accessible both from CPU and GPU. Risk for racing!

Do not expect performance penalty (but always be ready for surprises).

Not supported everywhere.

\section*{Atomic operations}

A special memory access method, for avoiding conflicts and race conditions.

Available in CUDA from Compute model 1.1.
To use it, specify model with
-arch compute_11
(or higher)

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\section*{Example: Histogram}

Simple method for gathering statistics about a set of data. Much data in, little out.

Common in image processing.

for all elements in a[]
\(h[a[i]]+=1\)


\section*{Histogram memory conflicts}

If you try to parallelize these operations, multiple threads will write simultaneously at the same item

Non-atomic operations will read \(\mathrm{h}[\mathrm{a}[\mathrm{i}]\) ], add 1, and write back.
\begin{tabular}{|l|l|l|}
\hline Read & -10 & \(\rightarrow\) Read \\
\hline Add 1 & & Add 1 \\
\hline Write back & \(\rightarrow 11\) & \(\rightarrow\) Write back \\
\hline
\end{tabular}

Unknown write order
\begin{tabular}{|c|c|c|}
\hline Read & 10 & \\
\hline Add 1 & 10 & Read \\
\hline Write back & 11 & Add 1 \\
\hline & 11 & - Write back \\
\hline
\end{tabular}

Write unsynchronized values in sequence

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\section*{Solution: Atomics}

Read - modify - write in one operation
Guaranteed not to be subject to racing
atomicAdd, atomicSub, atomicExch, atomicMin, atomicMax, atomicInc, atomicDec, atomicCAS, atomicAND, atomicOR, atomicXor

More types in Fermi and up

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\section*{But it comes for a cost!}

\section*{Slower than other operations}

Global memory only as of Compute Capability 1.1
Shared memory atomics in modern GPUs.
Simpler but slower than reduction solutions!

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\section*{Example: Find maximum}
for all elements \(i\) in \(a[]\) \(\operatorname{maxValue}=\max (\operatorname{maxValue}, \mathrm{a}[\mathrm{i}])\)

\section*{Easy? Yes! Parallel? No!}

All threads will write to the same memory element!

Use atomics? Very slow! All write at the same time, must wait -> sequential performance!

Solution: Use reduction instead!

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\section*{Atomic conclusions}

\section*{Simplifies some operations}

Serializes conflicting operations
Can hurt performance! Don't overuse!

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\section*{More exotic optimizations and tools}

Pinned memory
Multiple streams
Not where you start but let's not ignore the options.

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\section*{Pinned memory}

Can boost performance for memory transfer
Page-locked memory
So far: malloc() and cudaMalloc()
New call: cudaHostAlloc()
Allocated page-locked memory! Fixed physical location!

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\section*{Pinned memory}

Page-locked memory is a limited resource!
For non-pinned memory, CUDA copies it internally to pagelocked memory, then DMA to GPU. Transfer time goes up!

Picture based on an NVidia article


Normal, pageable data transfer


Pinned data transfer

\section*{Pinned memory, streams, overlapping computation}

> Pinned memory is part of an optimization approach with overlapping computations

No longer just a slight speedup of data transfer!
cudaMemCpyAsynch() can copy locked memory asynchronously!

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\section*{Multiple streams}

CUDA commands are placed in a queue, a stream!
These are the same queues as you can post CUDA events to.

We usually only use the default CUDA stream.
Multiple CUDA streams can be used to overlap work especially computing and data transfers!

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\section*{Single stream computation}

The kernel can not run until the data is transferred.

For this example, 2/3 data transfer, 1/3 computation
\begin{tabular}{|l|}
\hline Copy data to GPU \\
\hline Run kernel \\
\hline Copy result to CPU \\
\hline Copy data to GPU \\
\hline Run kernel \\
\hline Copy result to CPU \\
\hline
\end{tabular}

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\section*{Dual stream computation}

While one stream runs a kernel, the other stream performs data copying,

More time for computing, in this example kernels are running \(1 / 2\) of the time instead of \(1 / 3\).
\begin{tabular}{|l|l|}
\hline \multicolumn{2}{|c|}{ Copy data to GPU } \\
\multicolumn{2}{|c}{} \\
\hline Run kernel & Copy data to GPU \\
\hline Copy result to CPU & Run kernel \\
\hline Copy data to GPU & - \\
\hline Run kernel & Copy result to CPU \\
\hline- & Copy data to GPU \\
\hline Copy result to CPU & Run kernel \\
\hline & - \\
\cline { 2 - 2 } & Copy result to CPU \\
\cline { 2 - 3 } &
\end{tabular}

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\section*{Not all devices...}

Asynchronous data copying as well as concurrent execution is not guaranteed...
so make a device query!
CU_DEVICE_ATTRIBUTE_ASYNCH_ENGINE_COUNT: Can we copy memory asynch?

CU_DEVICE_ATTRIBUTE_CONCURRENT_KERNELS: Can we run multiple kernels?

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\section*{Debugging CUDA}

Let's get a bit more efficient when your code doesn't work
- Catch error codes
- printf() from kernels
- cudagdb

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\section*{Catch those error codes}
```

// Check for errors everywhere
err = cudaMalloc( (void**)\&ad, csize );
// If the GPU won't even take our data we are toasted
if (err) printf("cudaMalloc %d %s\n", err, cudaGetErrorString(err));
dim3 dimBlock( blocksize, 1 );
dim3 dimGrid( 1, 1 );
hello<<<dimGrid, dimBlock>>>(ad, bd);
// Most important thing to check? Did the kernel run at all?
err = cudaPeekAtLastError();
if (err) printf("cudaPeekAtLastError %d %s\n", err, cudaGetErrorString(err));

```
and pass them to cudaGetErrorString() for an explanation

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\section*{printf() from kernels}

Yes - printf() if legal in a kernel since Compute Capability 2.0

But don't try to print 100000 messages per second...

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\section*{More advanced debugger tools}

\section*{There are more tools to help you out there!}

\author{
cudagdb
}

Variant of the GDB debugger
Allows breakpoints and single-stepping
CUDA kernels!

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\section*{Sorting on GPUs}

Revisiting some algorithms from lecture 6:
Some not-so-good sorting approaches
Bitonic sort
QuickSort
Concurrent kernels and recursion

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\section*{Adapt to parallel algorithms}

Many sorting algorithms are highly sequential
Suitable for parallel implementation?
- Data driven execution
- Data independent execution

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\section*{Data driven execution}

Computing pattern depends on data
Usually harder to parallellize!
Example: QuickSort.

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\section*{Data independent execution}

Known computing pattern
Easier to parallellize - always the same plan
Example: Bitonic sort

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\section*{Bubble sort}

Loop through data, compare neighbors

\section*{Extremely sequential}

Inefficient
Parallel version: Bubble sort with odd-even transposition method
Compare all items pairwise
Two phases, "odd phase" and "even phase" (shifted one step)

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\section*{Bubble sort, parallel version}

Bubble sort with odd-even transposition method
Compare all items pairwise
Two phases, "odd phase" and "even phase" (shifted one step)
Fully sorted after n phases


Even phase
Odd phase
\(\mathrm{O}\left(\mathrm{n}^{2}\right)\)

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\section*{Suitable for GPU?}

Not as bad as it seems at first look:
- Data independent
- Excellent locality
- Appears to have possibilities to use shared memory but with some costly transfers at edges between blocks.
- But certainly not optimal at very large sizes
"Better" algorithms don't necessary beat this all that easily!

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\section*{Rank sort}

Count number of items that are smaller

\section*{Easy to parallelize:}
- One thread per item
- Loop through entire data
- Store in index decided from count of number of smaller items.

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\section*{Suitable for GPU?}

Again, not as bad as it seems at first look:
- Data independent
- Excellent locality - especially good for broadcasting (e.g. constant memory). Also suitable for shared memory.
- Again, \(O\left(n^{2}\right)\) : Will grow at very large sizes

Two bad ones that are not quite as bad as they seem.
N parallel iterations may beat NlogN sequential ones!

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\section*{Bitonic sort}
(As described in Kessler 2.3)
Bitonic set: Two monotonic parts in different direction.


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\section*{Bitonic sort}
(According to Batcher:) Let a be a bitonic set with a maximum at k , consisting of two monotonic parts, one increasing, \(\mathrm{a}^{-}\)(from item 1 to \(k\) ) and one decreasing, \(a^{+}(k+1\) to \(n\) )

Then two new sets can be constructed as
\[
\begin{aligned}
a^{\prime} & =\min \left(a_{1}, a_{k+1}\right), \min \left(a_{2}, a_{k+2}\right) \ldots \\
a^{\prime \prime} & =\max \left(a_{1}, a_{k+1}\right), \max \left(a_{2}, a_{k+2}\right) \ldots
\end{aligned}
\]

These two sets are also bitonic and \(\max \left(a^{\prime}\right) \leq \min \left(\mathrm{a}^{\prime \prime}\right)\) !


\section*{Bitonic sort by divide-andconquer}

Bitonic sort works on a bitonic sequence: partially sorted

The parts must be sorted. Sort them by bitonic sort!

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\section*{Bitonic sort example}
\begin{tabular}{|c|c|c|c|c|c|c|c|c|c|c|c|c|c|c|c|}
\hline 7 & & & & & & & & & & & & & & & \\
\hline & & 7 & & & \(\nabla\) & 3 & & & & 3 & & & & & 2 \\
\hline 8 & & 8 & & 8 & & 7 & & & & 4 & & & & & 3 \\
\hline 3 & & & & 7 & \(\downarrow\) & 8 & & & & 2 & & & & & 4 \\
\hline 5 & & & & \[
5
\] & \[
\Delta
\] & 6 & V & & & 6 & & & & & 5 \\
\hline 6 & \[
\dagger
\] & & & \[
6
\] & & \[
5
\] & & & & \[
5
\] & & & & & \\
\hline 2 & 4 & & & & & & & 1 & - & 7 & & \(V\) & 7 & & 7 \\
\hline 4 & & & & 2 & & 2 & & & & 8 & \(\downarrow\) & & 8 & \(\nabla\) & \\
\hline
\end{tabular}

Bitonic sort of smaller parts

Bitonic sort of main part
Reverse parts
(bitonic merge)
Reverse parts
(bitonic merge)

\section*{Bigger example}

\section*{The problem scales nicely, uniformly}


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\title{
Get those steps right
}

> Step length

\section*{Step direction}

Comparison direction
Calculated from stage number and stage length

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\section*{Code examples}

\section*{Sequential:}

Recursive example
Iterative example

\author{
Parallel:
}

CUDA example (not optimized)

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\section*{Bitonic sort}
- Data independent, no worst case
- Fast: O(n•昚2n) (Why?)
- Good locality in some parts
but
- Big leaps in addressing for some parts

\section*{What about those big leaps?}

Small leaps: Can be computed within one block. Shared memory friendly.

Big leaps (>number of threads/block): No synchronization possible between blocks!

But we must synchronize!
-> multiple kernel runs!

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\section*{QuickSort}

\section*{Very popular algorithm for sequential implementation}


Data driven, data dependent reorganization, non-uniform
Fancy name - nobody expect QuickSort to be nothing but optimal

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\section*{QuickSort is}

Fast: \(O(n \cdot \log n)\) in typical cases
\(O\left(n^{2}\right)\) in the worst case
Data driven, data dependent reorganization, non-uniform

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\section*{QuickSort on GPU}

Initially ignored as impractical
CUDA implementations exist
Data driven approaches increasingly suitable as GPUs become more flexible

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\section*{Parallel QuickSort}

Several stages to consider:
- Pivot selection. Usually just grab one.
- Comparisons
- Partitioning
- Concatenate result

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\section*{Pivot selection}

If we could always pick a pivot that splits the data in half...


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\section*{but you can't do that without sorting! (Or a histogram.) But how about a random one?}


There is a worst case caused by bad pivots. Live with it!

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\section*{Comparisons}

\section*{Easy to parallelize}

One thread per comparison not unreasonable! (GPUs don't have a problem with many threads!)

No problem!

\title{
Partitioning
}

The big problem!

\section*{Sequential partitioning: Bad!}

Parallel partitioning 1: Atomic fetch \& increment. (GPUs have atomics!)

Parallel partitioning 2: Divide and conquer

\section*{In-place sorting not feasible}

Split to two list of same size as original. Massive number of threads!

Then we must pack to smaller size.


\title{
Packing to smaller size not trivial
}

\section*{Data dependent}

Use parallel prefix sum to create a look-up table for addressing. (Kessler 1.6.3)

Computes sum of all previous items.

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\section*{Parallel prefix sum}

\section*{Similar to reduction but full output.}

\begin{tabular}{|l|l|l|l|l|l|l|l|}
\hline\(\# 1\) & \(\# 1+2\) & \(\# 3\) & \(\# 1 . .4\) & \(\# 5\) & \(\# 5+6\) & \(\# 7\) & \(\# 1 . .8\) \\
\hline
\end{tabular}


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\section*{Parallel prefix sum}

\section*{Example}


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\section*{For sorting: Binary parallel prefix sum}


\section*{Parallel prefix sum on GPU}
- No reason to use few threads. Use as many as you have output items.
- Multiple kernel runs to adapt to problem size variation.
- As described above, non-coalesced. Pack intermediate values for coalescing. If using shared memory, risk of bank conflicts. [Capannini]

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Thus, QuickSort is not impossible, but more complex than before.

Note:
GPUs have Compare-And-Swap atomics!
GPUs favor massive numbers of threads. One thread per comparison is more than OK!

Implementations available. Example:
https://sourceforge.net/projects/cuda-quicksort/

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\section*{Recursion}

GPUs can't do recursion efficiently... or can they?
Since Kepler we have concurrent kernels
Not only a matter of launching kernels from CPU!

\section*{A kernel can spawn new kernels!}

Do recursion by spawning new kernels!

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\section*{Concurrent kernels, Dynamic Parallelism}

Less work for the CPU to manage the computation.

CPU



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\section*{Recursion can look like this:}
```

_global__ void quicksort(int *data, int left, int right)
{
int nleft, nright;
cudaStream_t s1, s2;
// Partitions data based on pivot of first element.
// Returns counts in nleft \& nright
partition(data+left, data+right, data[left], nleft, nright);
// If a sub-array needs sorting, launch a new grid for it.
// Note use of streams to get concurrency between sub-sorts
if(left < nright) {
cudaStreamCreateWithFlags(\&s1, cudaStreamNonBlocking);
quicksort<<<<..., sl >>>(data, left, nright);
}
if(nleft < right) {
cudaStreamCreateWithFlags(\&s2, cudaStreamNonBlocking);
quicksort<<< ..., s2 >>>(data, nleft, right);
}
}
_host__ void launch_quicksort(int *data, int count)
quicksort<<<< ... >>>(data, 0, count-1);
}

```

Source: http://blogs.nvidia.com/blog/2012/09/12/how-tesla-k20-

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\section*{Advantages}
- Less work for CPU
- Less synchronizing (from CPU side)
- Easier programming!

They claim it matters this much (but your milage will vary)


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\section*{Recursive CUDA kernels, a significant improvement}

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\title{
Other non-trivial algorithms
}

\author{
FFT, Fast Fourier Transform
}

Distance transform
Fractal Brownian Motion

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\section*{Fast Fourier Transform}

Based on a sequence of "butterflies"
Similarily to Bitonic sort, can be computed several stage in one run for the "smaller" stages


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\section*{Distance transform}

Fast and simple version by Danielsson 1980: "Jump flooding"

Makes "jumps" of various length


Every "jump" need to be one kernel run!

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\section*{Fractal Brownian Motion}

Used for e.g. realistic looking procedural terrains
Among other methods:
- Diamond-square
- Multi-pass Perlin noise

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\section*{Diamond-square algorithm}
1) Midpoint from corners
2) Edge from corners and midpoints


Repeat to desired resolution

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\section*{Multi-pass Perlin noise}

\section*{Theoretically slower than Diamond-square}

\section*{BUT}
can be computed by independent threads! One kernel run!


Single octave
Needs log \(\mathbf{N}\) passes of different frequency

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\section*{Conclusion}

Algorithms with dependency in computed data often need multiple kernel runs.

This is an extra cost!
Does it pay when the computational complexity is lower?

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\section*{That's all folks!}```

