Thread Layouts for 2D Computational Problems



- Many computational problems have a 2D domain (e.g., CV)
 - Many others have a 3D domain (e.g., fluids simulation)
- Solution: layout threads in 2D

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Simplifies index calculations a lot





- Definition:
 - For each $c \in \mathbb{C}$ consider the (infinity) sequence

$$z_{i+1} = z_i^2 + c$$
 , $z_0 = 0$

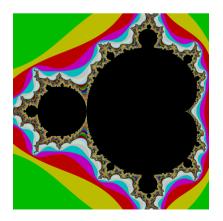
Define the Mandelbrot set

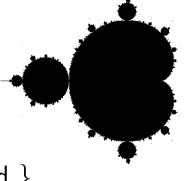
 $\mathbb{M} = \{ c \in \mathbb{C} \mid \text{sequence } (z_i) \text{ remains bounded } \}$

Theorem (w/o proof):

$$\exists t: |z_t| > 2 \implies c \notin \mathbb{M}$$

- Visualizing M nicely:
 - Color pixel c = (x,y) black, if |z| remains <2 after "many" iterations
 - Color c depending on the number of iterations necessary to make |z_t| > 2









- A few interesting facts about M
 (with which you can entertain people at a party ☺):
 - The length of the border of \mathbb{M} is infinite
 - ∭ is connected
 - (i.e., all "black" regions are connected with each other)
 - Mandelbrot himself believed $\ensuremath{\mathbb{M}}$ was disconnected
 - For each color, there is exactly one "ribbon" around M, i.e., there is exactly one ribbon of values *c*, such that |*z*₁| > 2, there is exactly one ribbon of values *c*, such that |*z*₂| > 2, etc. ...
 - Each such "iteration ribbon" reaches goes completely around M and it is connected (i.e., there are no "self intersections")
 - There is an infinite number of "mini Mandelbrot sets", i.e., smaller copies of M(self similarity)

Computing the Mandelbrot Set on the GPU

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- Embarrassingly parallel: each pixel computes their own zsequence, then sets the color
- Usual code for allocating memory, here a bitmap:

```
const unsigned int bitmap_size = img_size * img_size * 4;
h_bitmap = new unsigned char[bitmap_size];
cudaMalloc( (void**) &d_bitmap, bitmap_size );
```

Setup threads layout, here a 2D arrangement of blocks

```
dim3 threads( 16, 16 );
dim3 blocks( img_size/threads.x, img_size/threads.y );
```

- Here, we assume image size = multiple of 32
 - Simplifies calculation of number of blocks
 - Also simplifies kernel: we don't need to check whether thread out of range
 - See example code on web page how to ensure that





• Launch kernel:

```
mandelImage<<< blocks,threads >>>( d_bitmap, img_size );
```

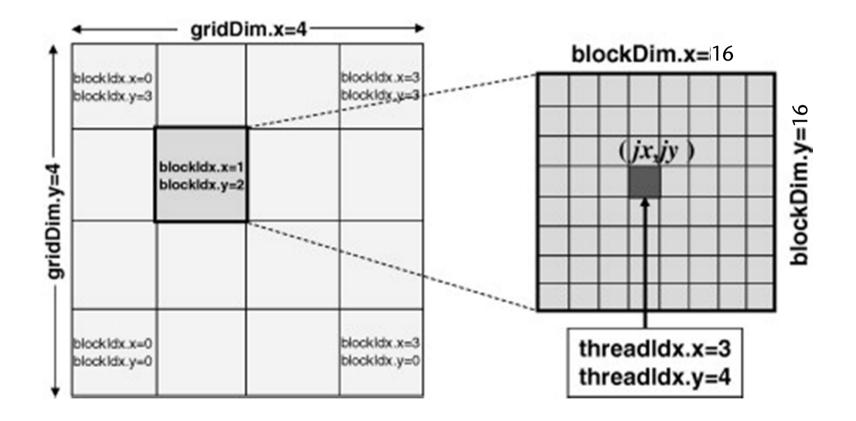
Implementation of the kernel (simplified):

```
__global__
void mandelImage( char4 * bitmap, const int img_size )
{
    int x = blockIdx.x * blockDim.x + threadIdx.x;
    int y = blockIdx.y * blockDim.y + threadIdx.y;
    int offset = x + y * (gridDim.x * blockDim.x); // x + y * width
    int isOutsideM = isPointInMandelbrot( x, y, img_size );
    bitmap[offset].x = 255 * isOutsideM; // red = outside
    bitmap[offset].y = bitmap[offset].z = 0;
    bitmap[offset].w = 255;
}
```



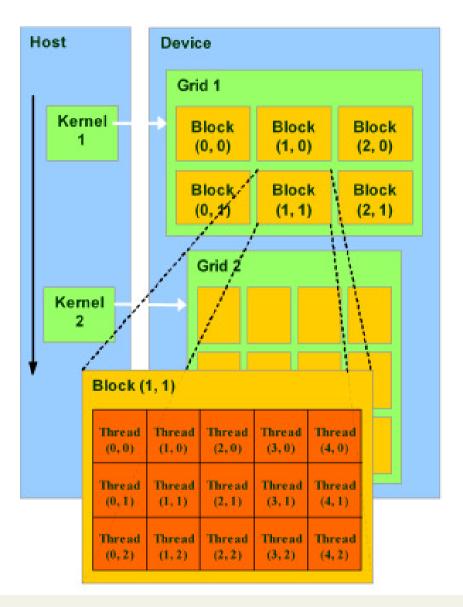


Visualization of our layout:





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- In general, the layout of threads can change from kernel to kernel:





A Word About dim3



Definition (done by CUDA):

```
struct dim3 // is actually a C++ class
{
    unsigned int x, y, z;
};
```

Usage:

dim3 layout (nx); = dim3 layout (nx, 1); = dim3 layout (nx, 1, 1);

dim3 layout (nx, ny); = dim3 layout (nx, ny, 1);

Launching a kernel like this: kernel<<<N, M>>>(...);

```
is equivalent: dim3 threads(M,1);
    dim3 blocks(N,1);
    kernel<<<blocks,threads>>>(...);
```



Implementation of the Kernel



```
device
int isPointInMandelbrot( int x, int y,
                         const int img_size, float scale )
 cuComplex c( (float) (x - img size/2) / (img size/2),
               (float) (y - img_size/2) / (img_size/2) );
 c *= scale;
 cuComplex z(0.0, 0.0); // z i of the sequence
 for ( int i = 0; i < 200; i ++ )</pre>
  {
   z = z \star z + c;
   if ( z.magnitude2() > 4.0f ) // |z|^2 > 2^2 -> outside
     return i;
  }
 return 0;
```





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```
struct cuComplex // define a class for complex numbers
{
  float r, i; // real and imaginary part
   device // constructor
  cuComplex(float a, float b) : r(a), i(b) {}
  device // |z|^2
  float magnitude2( void )
  ł
     return r * r + i * i;
  }
  device // z1 * z2
  cuComplex operator * (const cuComplex & a)
  {
     return cuComplex(r*a.r - i*a.i, i*a.r + r*a.i);
  }
  // for more: see example code on web page
};
```





		Executed on:	Only callable from:
<u>device</u> float De	<pre>viceFunc();</pre>	device	device
<u>global</u> void Ke	<pre>rnelFunc();</pre>	device	host
<u>host</u> float Ho	stFunc();	host	host

• Remarks:

- global defines a kernel function
- Each '___' consists of two underscore characters
- A kernel function must return void
- device and __host__ can be used together

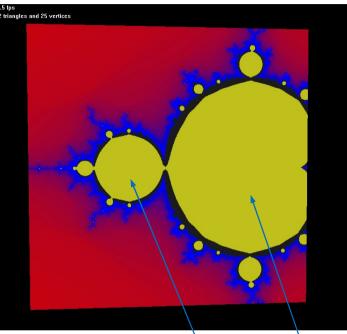


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- Example for the latter: make cuComplex usable on both device and host

```
struct cuComplex // define a class for complex numbers
{
   float r, i; // real, imaginary part
   __device_ __host__
   cuComplex( float a, float b ) : r(a), i(b) {}
   __device_ __host__
   float magnitude2( void )
   {
     return r * r + i * i;
   }
   // etc. ...
};
```



- An "Optimization":
 - The sequence of z_i can either converge towards single (complex) value,
 - or it can end up in a cycle of values,
 - or it can be chaotic.
- Idea:
 - Try to recognize such cycles;
 if you realize that a thread's is
 caught in a cycle, exit immediately
 (should happen much earlier in most cases)
 - Maintain an array of the k most recent elements of the sequence
- Last time I checked: 4x slower than the brute-force version!



All points here
convergeAll points here
convergetowards cycle
of length 2towards
fixed point

Querying the Device for its Capabilities



- How do you know how many threads can be in a block, etc.?
- Query your GPU, like so:

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For Your Reference: the Complete Table of the cudaDeviceProp



DEVICE PROPERTY	DESCRIPTION
<pre>char name[256];</pre>	An ASCII string identifying the device (e.g., "GeForce GTX 280")
<mark>size_t</mark> totalGlobalMem	The amount of global memory on the device in bytes
<pre>size_t sharedMemPerBlock</pre>	The maximum amount of shared memory a single block may use in bytes
int regsPerBlock	The number of 32-bit registers available per block
<pre>int warpSize</pre>	The number of threads in a warp
<pre>size_t memPitch</pre>	The maximum pitch allowed for memory copies in bytes
int maxThreadsPerBlock	The maximum number of threads that a block may contain
<pre>int maxThreadsDim[3]</pre>	The maximum number of threads allowed along each dimension of a block
<pre>int maxGridSize[3]</pre>	The number of blocks allowed along each dimension of a grid
<pre>size_t totalConstMem</pre>	The amount of available constant memory



DEVICE PROPERTY	DESCRIPTION
int major	The major revision of the device's compute capability
int minor	The minor revision of the device's compute capability
<pre>size_t textureAlignment</pre>	The device's requirement for texture alignment
int deviceOverlap	A boolean value representing whether the device can simultaneously perform a cudaMemcpy() and kernel execution
int multiProcessorCount	The number of multiprocessors on the device
<pre>int kernelExecTimeoutEnabled</pre>	A boolean value representing whether there is a runtime limit for kernels executed on this device
int integrated	A boolean value representing whether the device is an integrated GPU (i.e., part of the chipset and not a discrete GPU)
int canMapHostMemory	A boolean value representing whether the device can map host memory into the CUDA device address space
int computeMode	A value representing the device's computing mode default, exclusive, or prohibited
int maxTexture1D	The maximum size supported for 1D textures

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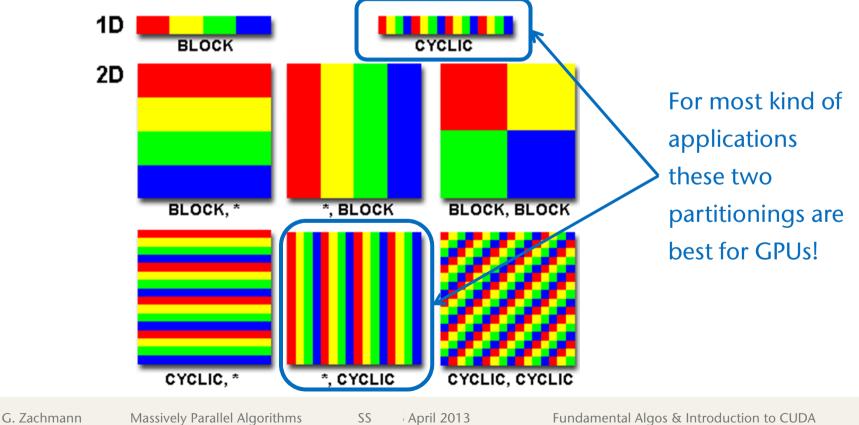
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DEVICE PROPERTY	DESCRIPTION
<pre>int maxTexture2D[2]</pre>	The maximum dimensions supported for 2D textures
<pre>int maxTexture3D[3]</pre>	The maximum dimensions supported for 3D textures
<pre>int maxTexture2DArray[3]</pre>	The maximum dimensions supported for 2D texture arrays
int concurrentKernels	A boolean value representing whether the device supports executing multiple kernels within the same context simultaneously





- Problem: your input, e.g. the vectors, is larger than the maximally allowed size along one dimension?
 - I.e., what if vec_len > maxThreadsDim[0] * maxGridSize[0]?
- Solution: partition the problem (color = thread ID)



Example: Adding Huge Vectors



- Vectors of size 100,000,000 are not uncommon in highperformance computing (HPC) ...
- The thread layout:

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```
dim3 threads(16,16); // = 256 threads per block
int n_threads_pb = threads.x * threads.y;
int n_blocks = (vec_len + n_threads_pb - 1) / n_threads_pb;
int nb_sqrt = (int)( ceilf( sqrtf( n_blocks ) ) );
dim3 blocks( nb_sqrt, nb_sqrt );
```

Kernel launch:

```
addVectors<<< threads, blocks >>>( d_a, d_b, d_c, n );
```

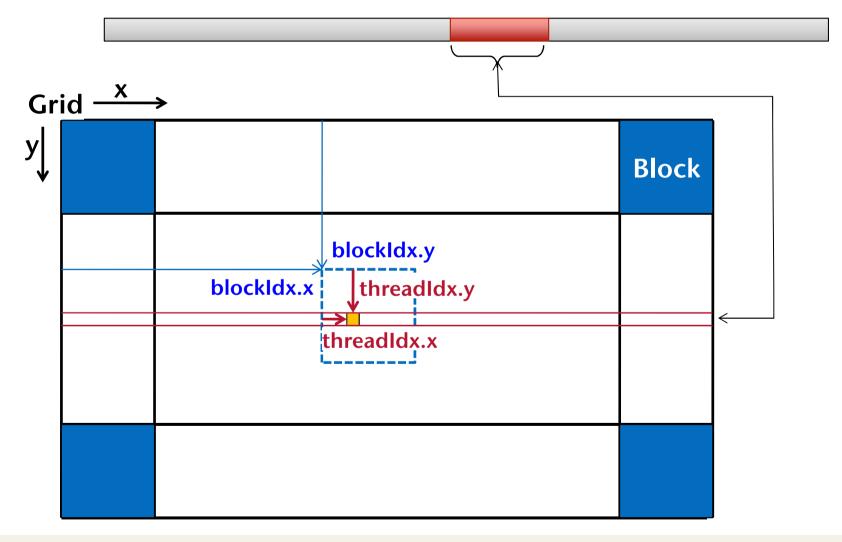
Index computation in the kernel:

```
unsigned int tid_x = blockDim.x * blockIdx.x + threadIdx.x;
unsigned int tid_y = blockDim.y * blockIdx.y + threadIdx.y;
unsigned int i = tid_y * (blockDim.x * gridDim.x) + tid_x;
```





Visualization of this index computation:





Constant Memory



- Why is it so important to declare constant variables/instances in C/C++ as const ?
- It allows the compiler to ...
 - optimize your program a lot
 - do more type-checking
- Something similar exists in CUDA → constant memory

Example: a Simple Raytracer



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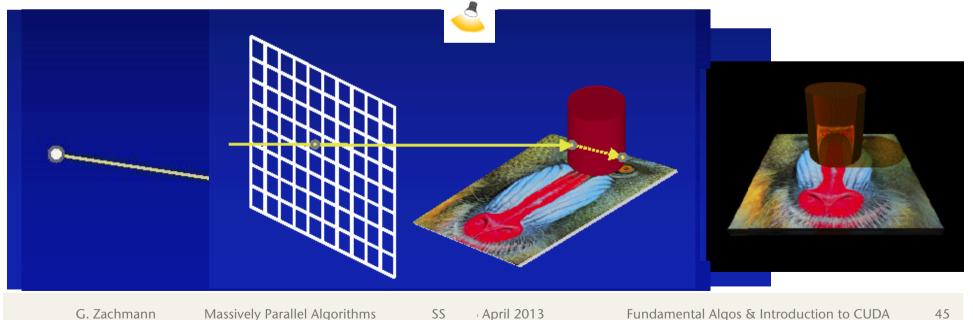
The ray-tracing principle:

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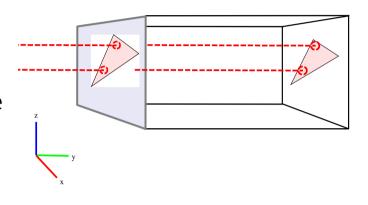
- 1. Shoot rays from camera through every pixel into scene (primary rays)
- 2. If the rays hits more than one object, then consider only the first hit
- 3. From there, shoot rays to all light sources (shadow feelers)
- 4. If a shadow feeler hits another obj \rightarrow point is in shadow w.r.t. that light source. Otherwise, evaluate a lighting model (e.g., Phong [see "Computer graphics"])
- 5. If the hit object is glossy, then shoot reflected rays into scene (secondary rays) \rightarrow recursion
- 6. If the hit object is transparent, then shoot refracted ray \rightarrow more recursion

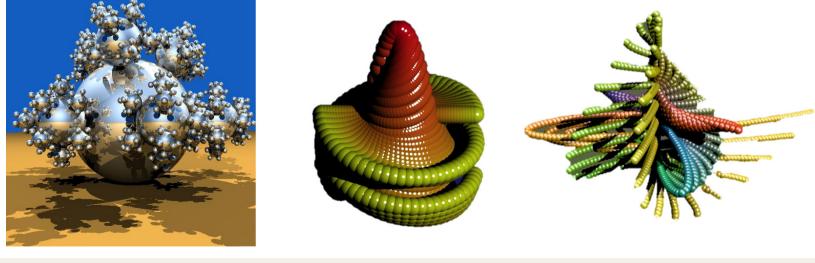


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- Simplifications (for now):
 - Only primary rays
 - Camera at infinity → primary rays are orthogonal to image plane
 - Only spheres
 - They are so easy, every ray tracer has them $\ensuremath{\textcircled{\sc o}}$







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• The data structures:

```
struct Sphere
{
    Vec3 center; // center of sphere
    float radius;
    Color r, g, b; // color of sphere
    __device__
    bool intersect( const Ray & ray, Hit * hit )
    {
        ...
    }
};
```





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• The mechanics on the host side:

```
int main( void )
{
   // create host/device bitmaps (see Mandelbrot ex.)
   . . .
   Sphere * h_spheres = new Sphere[n_spheres];
   // generate spheres, or read from file
   // transfer spheres to device (later)
   // generate image by launching kernel
   // assumption: img size = multiple of 16!
   dim3 threads (16, 16);
   dim3 blocks( img size/16, img size/16 );
   raytrace<<<blocks,threads>>>( d_bitmap );
   // display, clean up, and exit
};
```

The mechanics on the device side

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```
qlobal
void raytrace( unsigned char * bitmap ) {
   // map from thread id to pixel pos
   int x = blockIdx.x * blockDim.x + threadIdx.x;
   int y = blockIdx.y * blockDim.y + threadIdx.y;
   int offset = x + y * (gridDim.x * blockDim.x);
  Ray primary(x, y, camera); // generate primary ray
   // check intersection with scene, take closest one
  min dist = INF;
   int hit sphere = MAX INT;
  Hit hit;
   for ( int i = 0; i < n_spheres; i ++ ) {</pre>
      if ( intersect(ray, i, & hit) ) {
         if ( hit.dist < min dist ) {</pre>
           min dist = hit.dist; // found a closer hit
           hit sphere = i;
                                     // remember sphere; hit info
                                     // is already filled
         }
   // compute color at hit point (if any) and set in bitmap[offset]
```



Declaration & transfer



• Since it's constant memory, we declare it as such:

```
const int MAX_NUM_SPHERES 1000;
____constant___ Sphere c_spheres[MAX_NUM_SPHERES];
```

Transfer now works by a different kind of Memcpy:



- Access of constant memory on the device (i.e., from a kernel) works just like with any globally declared variable
- Example:

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```
constant Sphere c spheres[MAX NUM SPHERES];
 device
 bool intersect( const Ray & ray, int s, Hit * hit )
 {
    Vec3 m( c_spheres[s].center - ray.orig );
    float q = m*m - c_spheres[s].radius*c_spheres[s].radius;
    float p = \ldots
     solve pq( p, q, *t1, *t2 );
     . . .
                                                                m
(t \cdot \mathbf{d} - \mathbf{m})^2 = r^2 \Rightarrow t^2 - 2t \cdot \mathbf{md} + \mathbf{m}^2 - r^2 = 0
```