# Decoupled Access/Execute Metaprogramming for GPU-Accelerated Systems

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#### I. INTRODUCTION

We describe the evaluation of several implementations of a simple image processing filter on an NVIDIA GTX 280 card. Our experimental results show that performance depends significantly on low-level details such as data layout and iteration space mapping which complicate code development and maintenance.

We propose extending a CUDA<sup>1</sup> or OpenCL<sup>2</sup> like model with decoupled Access/Execute ("Æcute" [1]) metadata, describing its iteration space ordering and partitioning (*execute* metadata) and its memory access patterns (*access* metadata).

We believe that using Æcute metadata will make software engineering for accelerated systems more disciplined and productive, by separating algorithm representation from lowlevel mapping and tuning.

#### II. MOTIVATING EXAMPLE

We consider a vertical mean image filter, for which the output pixel at position (x, y) is given by the formula

$$\mathbf{O}_{x,y} = \frac{1}{D} \sum_{k=0}^{D-1} \mathbf{I}_{x,y+k}, \text{ where}$$
(1)

- I is a  $W \times H$  grey-scale input image;
- **O** is a  $W \times (H D)$  grey-scale output image;
- *D* is the *diameter* of the filter, *i.e.* the number of input pixels over which the mean is computed (typically, *D* ≪ *H*);
- $0 \le x < W, 0 \le y < H D.$

Let N be the number of output pixels:  $N = W \times (H - D)$ . A naïve parallel algorithm can run N threads, each producing a single output pixel, which requires  $\Theta(ND)$  reads and arithmetic operations. A good parallel algorithm, however, must be efficient and scalable [2].

#### A. Scalable algorithm

The algorithm in Listing 1 *strips* the computation in the vertical dimension, where up to T outputs in the same strip are computed serially in two *phases*. The first phase in lines 6–10 computes  $\mathbf{O}_{x,y0}$  according to (1). The second phase in lines 12–19 computes  $\mathbf{O}_{x,y}$  for  $y \ge y0 + 1$  as  $\mathbf{O}_{x,y-1} + (\mathbf{I}_{x,y+D-1} - \mathbf{I}_{x,y-1})/D$ .

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// for each column	1
for (int $x = 0$ ; $x < W$ ; $++x$ )	2
{ // for each strip of rows	3
for (int y0 = 0; y0 < H-D; y0 += T)	4
{	5
<pre>// first phase: convolution</pre>	6
<pre>float sum = 0.0f;</pre>	7
<b>for</b> ( <b>int</b> $k = 0$ ; $k < D$ ; ++k)	8
sum += I[(y0+k)*W + x];	9
O[y0*W + x] = sum / (float)D;	10
-	11
<pre>// second phase: rolling sum</pre>	12
<pre>for(int dy = 1; dy &lt; min(T,H-D-y0); ++dy)</pre>	13
{	14
int $y = y0 + dy;$	15
sum -= I[(y-1)*W + x];	16
sum += I[(y-1+D) *W + x];	17
$O[y \star W + x] = sum / (float)D;$	18
}	19
}	20
}	21

Listing 1: Vertical mean image filter algorithm in C.

This algorithm performs  $\Theta(N + ND/T)$  reads and arithmetic operations, significantly reducing memory bandwidth and compute requirements for  $T \gg D$ . Since the x and y0 loops carry no dependences, up to  $\lceil N/T \rceil$  threads can run in parallel.

Note that since the order of arithmetic operations is undefined in (1), both the naïve and scalable algorithms are functionally, if not arithmetically, equivalent.

Clearly, the optimal value of T depends on problem parameters (W, H and D), and device parameters (*e.g.* the number of cores and memory partitions). Thus, in §II-C we use the iterative compilation approach to find the optimum.

#### B. Efficient implementation

Implementing the vertical mean filter efficiently on a GPU requires mapping the iteration space onto threads, which are grouped into blocks located in a grid.

The most natural iteration space mapping is into thread blocks on a 2D grid, with each block producing a rectangular WPBX  $\times$  WPBY section of the output image. However, if the image width is not a multiple of WPBX, significant portions of thread blocks covering the right edge of the image may be unused, as illustrated by Figure 1a.

This issue can be alleviated by mapping into thread blocks on a 1D grid that covers the image by wrapping around the right edge, as illustrated by Figure 1b. As we show in §II-C, a mapping that maximises thread utilisation suffers from misalignment, if the image width is not a multiple of the size

<sup>&</sup>lt;sup>1</sup>http://www.nvidia.com/cuda

<sup>&</sup>lt;sup>2</sup>http://www.khronos.org/opencl



(a) A 2D grid mapping loses efficiency from unused threads off the right image edge.



(b) A 1D grid mapping uses its threads more efficiently by wrapping around the right image edge. For efficiency, it must take into account alignment, which complicates both memory access and iteration.

Fig. 1: Different mapping strategies result in different utilisation of threads. Light and dark regions of blocks denote used and unused threads, respectively.

of the SIMD unit (*warp* in NVIDIA's terminology); a better mapping takes alignment into account by wasting a small number of threads on the right of the image, thus ensuring that the first pixel of each row is handled by the first thread in a SIMD unit.

#### C. Experimental results

Figure 2 presents experimental results obtained on a dualcore 3GHz Intel Core 2 Duo E8400 system with 2GiB RAM, equipped with an NVIDIA GTX 280 card, running 64-bit Linux Ubuntu 8.04. Code is compiled using CUDA SDK 2.2 and GCC 4.2.4 with the "-O3" optimisation settings. We measure the kernel execution time only and get the best throughput out of 50 runs.

In all the experiments, we fix the number of threads per block at 128 (128×1), as we nearly achieve the peak memory efficiency with this setting:  $\approx 10$  Mpixel/s × 4 bytes/pixel × (2 reads + 1 write) = 120 GB/s (close to the bandwidth of aligned copy on this card). Thus, WPBX = 128 and WPBY = T.

Figure 2a shows that the 1D and 2D grid versions are similar in throughput when applied to a  $5120 \times 3200$  image, where 5120 is a multiple of 128 pixels. The throughput is below 1000 Mpixel/s when each thread produces a single pixel, climbs fast with increasing serial effiency, achieving (by the 1D grid version) the peak throughput of 9884 Mpixel/s when T = 355, and then declines with decreasing parallelism.

When applied to a  $5121 \times 3200$  image, however, the 2D grid version only achieves 7017 Mpixel/s, as shown by the bottom line in Figure 2b. Whilst we allocate memory using the cudaMallocPitch function, which pads the image to a multiple of 16 pixels to enable global memory access coalescing (5136 pixels in this case), such allocation leads to DRAM partition conflicts. We remedy the conflicts by manually padding the image to a multiple of 32, 64 and 128.

As the results of padding to a multiple of 64 and 128 are barely distinguishable, we fix the image padding at a multiple of 64 (5184 pixels) for all subsequent experiments.

Figure 2c shows that the 1D grid mapping that maximises thread utilisation by wrapping on 5121 pixels only achieves 5998 Mpixel/s, whilst wrapping on the image padding of 5184 pixels performs worse than wrapping on the warp size multiple of 5152 pixels.

Figure 2d summarises the throughput for the misaligned image padded to 5184 pixels: the 1D grid version wrapped on 5152 pixels achieves 9575 Mpixel/s at T = 396, whilst the 2D grid version achieves only 9056 Mpixel/s at T = 409; thus, the 1D grid version performs 6% better than the 2D grid one.

### III. TOWARDS METAPROGRAMMING

To ease the programmer's burden of mapping and tuning computation kernels to GPU architectures, we propose extending a kernel's description with decoupled Access/Execute metadata. Execute metadata for a kernel describes its iteration space ordering and partitioning. Access metadata for a kernel describes memory locations the kernel may access on each iteration.

// Array descriptors (C array wrappers)	1
Array2D< <b>float</b> > arrayI(&I[0][0], W, H);	2
Array2D< <b>float</b> > arrayO(&O[0][0], W, H-D);	3
	4
// Execute metadata: parallel iteration space	5
<pre>IterationSpace1D x(0,W);</pre>	6
<pre>IterationSpace1D y(0,H-D);</pre>	7
<pre>IterationSpace2D iterXY(x,y);</pre>	8
	9
// Access metadata: iteration space -> memory	10
<pre>VerticalStrip2D_R accessI(iterXY, arrayI, D);</pre>	11
<pre>Point2D_W accessO(iterXY, arrayO);</pre>	12

Listing 2: Æcute metadata for the vertical mean image filter.

We give an example of  $\mathcal{E}$ cute metadata for the vertical mean image kernel in Listing 2. In lines 1–3 we wrap accesses to plain C arrays I[W][H] and O[W][H–D] into  $\mathcal{E}$ cute array descriptors arrayI and arrayO to cleanse the kernel of uncontrolled side-effects. In lines 5–8 we construct a 2D iteration space descriptor iterXY from 1D descriptors x and y, having the same bounds as the output image dimensions. By default, an iteration space is parallel in every dimension. Finally, in lines 10–12 we specify that on each iteration of the 2D iteration space the kernel reads a vertical strip of D pixels from arrayI and writes a single pixel to arrayO.

Similar to Stanford's Sequoia language [3], we target systems with software-managed memory hierarchies and seek to separate a high-level algorithm representation from a systemspecific mapping. Unlike Sequoia, we base our mapping on partitioning (manually or automatically) an iteration space into disjoint subspaces and infer memory access of subspaces from Æcute metadata.

For example, for GPU-accelerated systems, a hierarchy of iteration space partitions can specify subspaces to be executed:

- at the lowest level, by individual threads:
  - //  $1 \mathrm{xT}$  outputs per thread
  - iterXY.partitionThreads(1,T);









(c)  $5121 \times 3200$  image. Data padded 5184 (a multiple of 64) pixels.(d)  $5121 \times 3200$  image. Data padded to 5184 (a multiple of 64) pixels.1D grid wrapped on the image width and multiples of 16, 32 and 64 pixels.2D grid; 1D grid wrapped on 5152 (a multiple of 32) pixels.

Fig. 2: Comparison of different mappings with various image sizes, data padding and thread wrapping alignment.

• at the middle level, by blocks of possibly cooperating threads:

```
// 128xT outputs per block
iterXY.partitionBlocks(128,T);
```

• at the highest level, by possibly cooperating compute devices:

```
// (W/2)x(H-D) outputs per device
iterXY.partitionDevices(W/2,H-D)
```

## IV. WORK IN PROGRESS

We are working on a tool that will take a high-level algorithm representation and generate efficient device-specific OpenCL code. The representation will be kept similar to C++, *e.g.* as in Listing 1 with accesses to C arrays replaced with accesses to Æcute array descriptors as in Listing 2. Code generation will be particularly oriented towards effectively orchestrating data movement in software-managed memory hierarchies, including automatically handling such low-level details as data alignment and padding.

#### REFERENCES

- L. W. Howes, A. Lokhmotov, A. F. Donaldson, and P. H. Kelly, "Deriving efficient data movement from decoupled Access/Execute specifications," in *Proceedings of the 4th International conference on High-Performance Embedded Architectures and Compilers (HiPEAC)*, ser. LNCS, vol. 5409. Springer, 2009, pp. 168–182.
- [2] C. Lin and L. Snyder, *Principles of Parallel Programming*, 1st ed. Boston, MA, USA: Addison-Wesley, 2008.
- [3] K. Fatahalian, D. R. Horn, T. J. Knight, L. Leem, M. Houston, J. Y. Park, M. Erez, M. Ren, A. Aiken, W. J. Dally, and P. Hanrahan, "Sequoia: programming the memory hierarchy," in *Proceedings of the ACM/IEEE conference on Supercomputing*, 2006, p. 83.