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**Introduction to GPU programming**

*Practice 4. Numerical integration of heat equation*

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# Objectives

The objective of this practice is to develop a simple implementation of numerical integration of 2D heat equation using explicit scheme.

# Abstract

This practice is devoted to development a simple implementation of numerical integration of 2D heat equation using explicit scheme. We illustrate using 2D indexes in kernels and device functions.

# BRIEF OVERVIEW

We consider a Dirichlet problem for 2D heat equation:

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Simulation area is covered by uniform grid with steps ∆x, ∆y, and ∆t. Denote – grid value in point at -th time moment. We will use explicit scheme for numerical integration:

Computation is performed iteration by iteration in time, computations within one time iteration are independent. For simplicity we assume boundary conditions are stationary (do not depend on time), in this case on each time step we can only update internal values.

Implementation on CPU is as follows:

float f(float x, float y) {

return 0.05f \* x \* x + 0.001f \* y \* y \* y;

}

void step(int nx, int ny, float dx, float dy, float dt, float a, float c,

const float \* v, float \* newV) {

for (int i = 1; i < nx - 1; ++i)

for (int j = 1; j < ny - 1; ++j) {

float x = a + i \* dx;

float y = c + j \* dy;

newV[i \* ny + j] = v[i \* ny + j] + dt \* (f(x, y) +

(v[(i + 1) \* ny + j] - 2.0f \* v[i \* ny + j] + v[(i - 1) \* ny + j]) / (dx \* dx) +

(v[i \* ny + (j + 1)] - 2.0f \* v[i \* ny + j] + v[i \* ny + (j - 1)]) / (dy \* dy));

}

}

Two loops with independent iterations are easily converted to a kernel with 2D indexes. Note that we can not call right-hand side function f from a kernel, as it is not declared with \_\_device\_\_ qualifier. A solution is to either declare it with \_\_host\_\_ and \_\_device\_\_ qualifiers, or to create a copy with \_\_device\_\_ qualifier. We choose the latter:

\_\_device\_\_ float f\_gpu(float x, float y) {

return 0.05f \* x \* x + 0.001f \* y \* y \* y;

}

\_\_global\_\_ void kernel(int nx, int ny, float dx, float dy, float dt, float a,

float c, const float \* v, float \* newV) {

int i = 1 + blockIdx.x \* blockDim.x + threadIdx.x;

int j = 1 + blockIdx.y \* blockDim.y + threadIdx.y;

if ((i < nx - 1) && (j < ny - 1)) {

float x = a + i \* dx;

float y = c + j \* dy;

newV[i \* ny + j] = v[i \* ny + j] + dt \* (f\_gpu(x, y) +

(v[(i + 1) \* ny + j] - 2.0f \* v[i \* ny + j] + v[(i - 1) \* ny + j]) / (dx \* dx) + (v[i \* ny + (j + 1)] - 2.0f \* v[i \* ny + j] + v[i \* ny + (j - 1)]) / (dy \* dy));

}

}

Note that .x index is used for rows and .y is used for columns. Conditions in the kernel correspond to conditions in the loop in the implementation for CPUs. A function to invoke the kernel is as follows:

void step\_gpu(int nx, int ny, float dx, float dy, float dt, float a, float c, const float \* v\_gpu, float \* newV\_gpu) {

const int threadsX = 16;

const int threadsY = 16;

dim3 blocks(threadsX, threadsY);

dim3 grid(((nx - 2) + (threadsX - 1)) / threadsX,

((ny - 2) + (threadsY - 1)) / threadsY);

kernel<<<grid, blocks>>>(nx, ny, dx, dy, dt, a, c, v\_gpu, newV\_gpu);

}

We fix block size for both dimensions and compute number of blocks for each dimension using the same scheme as for vector addition.

# FOR STUDENTS

CUDA C language is described in [1, 2].

# References

1. Sanders J., Kandrot E. CUDA by Example: An Introduction to General-Purpose GPU Programming. – Addison-Wesley Professional, 2010. – 312 p.
2. NVIDIA CUDA C Programming Guide. [http://docs.nvidia.com/cuda/cuda-c-programming-guide/].

# Individual work

1. Create implementation of the kernel using 1D indexes.
2. Implement a version with non-stationary boundary condition.