

# OpenMP for GPU

Kenjiro Taura

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- recent OpenMP supports offloading to GPU (`target` directive)
- official home page: <http://openmp.org/>
- specification:  
<http://openmp.org/wp/openmp-specifications/>
- latest version is 5.0  
(<https://www.openmp.org/spec-html/5.0/openmp.html>)
- section numbers below refer to those in OpenMP spec 5.0

# Compiling OpenMP programs for GPUs

- LLVM (clang/clang++) : compile with `-fopenmp -fopenmp-targets=nvptx64`

```
1 $ clang -Wall -fopenmp -fopenmp-targets=nvptx64 program.c
```

- you get a warning: “CUDA version is newer than the latest supported version 11.5” and `-Wunknown-cuda-version` suppresses it

- NVIDIA HPC SDK (nvc/nvc++) : compile with `-mp -target-gpu`

```
1 $ nvc -Wall -mp -target=gpu program.c
```

# Directives overview

- 1 move control
  - `target` : moves the execution to GPU
- 2 parallelize
  - `teams` and `distribute`
    - `teams` : creates a number of teams executing the same statement ( $\approx$  `parallel` pragma)
    - `distributed` : distribute iterations of a for loop among teams ( $\approx$  `for` pragma)
  - `parallel` and `for`
    - `parallel` : creates a number of threads executing the same statement in a team
    - `for` : distribute iterations of a for loop among threads in a team
  - think of `teams` + `distributed` another layer outside `parallel` + `for`
- 3 move (or sync) data
  - `target data` : move/sync data between CPU and GPU

# Implementation note

- while not specified anywhere in the spec (and there are cases they behave differently to below), you can think of
  - a team  $\sim$  a thread block
  - a thread  $\sim$  a CUDA thread
- it at least helps you understand why things look so redundant ...

# Frequently-used combined idioms

- all combined

```
1 #pragma omp target teams distribute parallel for
2 for (int i = start; i < end; i += incr) {
3     S
4 }
```

- teams + distributed to outer loop and parallel + for to outer loop

```
1 #pragma omp target teams distribute
2 for (int i = start; i < end; i += incr) {
3     #pragma omp parallel for
4     for (int j = start'; j < end'; j += incr') {
5         S
6     }
7 }
```

- similar to launching a kernel doing  $S$ , but
  - you don't have to adjust thread block size
  - the program is orthogonal to thread count

# Data mapping

a major headache when programming in CUDA is data management

- the only “transparent” data transfer is argument passing

```
1 f<<<nb,bs>>>(a, b, c, ...);
```

- getting the result back from GPU is already painful

```
1 cudaMalloc(&r_dev, ...);  
2 f<<<nb,bs>>>(a, b, c, ..., r);  
3 cudaMemcpy(r, r_dev, ...);
```

- for persistent data,
  - maintain two pointers to logically same data (CPU version and GPU version)
  - get them synched when necessary (before and after a kernel launch)

“data mapping” of OpenMP alleviates the pain

# Data mapping example

```
1 #pragma target data map(to: a[b:c]) map(from: x)
2   S
```

- send the array range  $a[b:c]$  ( $a[b]$ ,  $a[b+1]$ , ...,  $a[c-1]$ ) to GPU before  $S$
- send  $x$  from GPU after  $S$
- you can combine **to:** and **from:** into **tofrom:**
- somewhat “declarative” way of understanding this
  - expressions  $a[i]$  ( $b \leq i < c$ ) become valid (“mapped”) on GPU during  $S$
  - expressions  $x$  become valid on CPU after  $S$
- note: you can specify **map** clauses as part of **target** (not **target data**) directive, too
- learn details with tht notebook