OpenMP for GPU

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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`

```
$ 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`

```
$ 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, ...);
```

- gettint 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 launche)

"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