井 HPC.NRW

SEVERAL WAYS TO SAXPY

Julia + CUDA.jl

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THE COMPETENCE NETWORK FOR HIGH-PERFORMANCE COMPUTING IN NRW.

AGENDA



- The SAXPY problem
- CUDA.jl
 - CPU and GPU memory allocations
 - memory transfer
- SAXPY with CUDA.jl
 - GPU broadcasting (higher-order abstraction)
 - GPU kernel (the manual way)
 - CUBLAS (the library way)
- Summary



THE SAXPY PROBLEM

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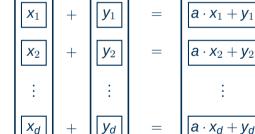
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- linear combination of two float arrays
- results written to third array



A typical Julia code:

<pre># define constants const dim = 100_000_000 const a = 3.1415</pre>
<pre># allocate vectors x = ones(Float32, dim) y = ones(Float32, dim) z = zeros(Float32, dim)</pre>
<pre># perform SAXPY z .= a .* x .+ y</pre>



CUDA.JL https://github.com/JuliaGPU/CUDA.jl

Programming interface for working with NVIDIA CUDA GPUs in Julia.

- high-level array abstractions
- tools for writing CUDA kernels
- wrappers for various CUDA libraries (e.g. cuFFT)

julia>] add CUDA

using CUDA
CUDA.versioninfo()





JULIA + CUDA.JL memory allocation and transfer

CPU memory (Array)

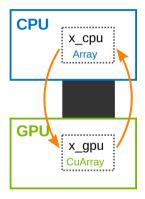
x_cpu = ones(Float32, dim)

- GPU memory (CuArray)

x_gpu = CUDA.ones(Float32, dim)

- type conversion triggers memory transfer

x_gpu = CuArray(x_cpu) # CPU -> GPU x_cpu = Array(x_gpu) # GPU -> CPU





SAXPY WITH CUDA.JL

GPU broadcasting



using CUDA

```
# define constants
const dim = 100_000_000
const a = 3.1415
```

```
# allocate vectors on the GPU
x = CUDA.ones(Float32, dim)
y = CUDA.ones(Float32, dim)
z = CUDA.zeros(Float32, dim)
```

```
# perform SAXPY
CUDA.@sync z .= a .* x .+ y
```

- almost the same code as for the CPU
- vectors x, y, z are initialized as CuArrays directly on the GPU
- CUDA.@sync make CPU wait until the GPU finishes SAXPY (blocking)

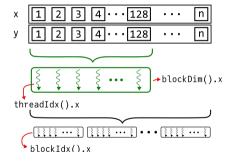


SAXPY WITH CUDA.JL

GPU kernel

```
# define GPU kernel
function saxpy_gpu_kernel!(z,a,x,y)
    i = (blockIdx().x - 1) * blockDim().x +
    threadIdx().x
    if i <= length(z)
       @inbounds z[i] = a * x[i] + y[i]
    end
    return nothing
end</pre>
```

- z[i] = a * x[i] + y[i] as if surrounded by implicit for-loop
- threadIdx().x, blockDim().x, blockIdx().x built-in variables that identify the thread global index

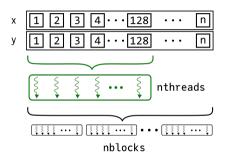




SAXPY WITH CUDA.JL

launching the kernel

```
using CUDA
# [define GPU kernel, constants, and
# allocate vectors on the GPU ...]
# define GPU execution parameters
nthreads = CUDA.attribute(
    device(),
    CUDA. DEVICE ATTRIBUTE MAX THREADS PER BLOCK
nblocks = cld(dim, nthreads)
CUDA.@sync @cuda(
  threads=nthreads,
  blocks=nblocks,
  saxpy_gpu_kernel!(z,a,x,y)
```



 @cuda macro launches kernel on the GPU with the given launch configuration.



SAXPY WITH CUDA.JL CUBLAS



using CUDA using CUDA.CUBLAS

```
# define constants
const dim = 100_000_000
const a = 3.1415
```

```
# allocate vectors on the GPU
x = CUDA.ones(Float32, dim)
y = CUDA.ones(Float32, dim)
```

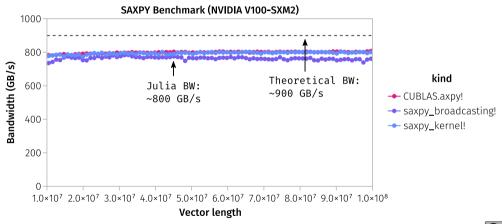
perform SAXPY
(y is overwritten with the result)
CUDA.@sync CUBLAS.axpy!(dim, a, x, y)

- CUDA.jl provides low-level wrappers of the vendor library CUBLAS.
- CUBLAS.axpy! for Float32 input directly calls the cuBLAS function cublasSaxpy_v2.
- CUDA.@sync make CPU wait until the GPU finishes SAXPY (blocking)



SAXPY PERFORMANCE







INNOVATION THROUGH COOPERATION.





- Julia + CUDA.jl
 - NVIDIA CUDA interface: higher-order abstractions, library wrappers
- SAXPY with CUDA.jl
 - GPU broadcasting: cuArray moves computation to the GPU
 - zeros(Float32, N) VS. CUDA.zeros(Float32, N)
 - SAXPY: essentially the same code as for the CPU
 - GPU kernel
 - definition: blockIdx(), blockDim(), threadIdx()
 - launching: CUDA.@sync @cuda threads=T blocks=B kernel()
 - CUBLAS: CUDA.@sync CUBLAS.axpy!()
 - transfer GPU result to CPU: result_cpu = Array(result_gpu)
 - memory release: result_gpu = nothing; GC.gc(true)
- SAXPY performance on NVIDIA V100: 800 GB/s

