Julia: From multicore to GPU parallelism

presented by Steffen Haug

WHAT IS JULIA?

(Briefly)

1. High level programming language



- 1. High level programming language
- 2. Dynamically typed



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- 3. For HPC



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- 2. Dynamically typed
- 3. For HPC
- 4. Novel compilation infrastructure (Central on CPU and GPU)

Untyped AST specialized at runtime



- 1. High level programming language
- 2. Dynamically typed
- 3. For HPC
- 4. Novel compilation infrastructure (Central on CPU and GPU)

Untyped AST specialized at runtime

"Looks like Python - runs like Fortran"



```
function square!(xs)
  for i in eachindex(xs)
    xs[i] ^= 2
  end
end

V = rand(10000);
@code_native syntax=:intel square(V)
```

Function that squares (mutates) a vector.

```
function square!(xs)
  for i in eachindex(xs)
    xs[i] ^= 2
  end
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```
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; [LILL
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veries year, year, year
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        .nie julia_nquere_bi, .idum_emilijulia_nquere_bi
.nii mierus
```

.amtion ".mate.DE-stant", " , (projetts



```
square:
 mov rax, gword ptr [rdi + 8]
 test pax, pax
 ie IRRR 7
 mov rex. award ptr [rdi]
 nov edx, 1
 стр гах. 16
 ib .LBB0 S
 mov rsi, rax
 and rsi. -16
 lea rdx, [rsi + 1]
 xor edi. edi
LBB0 3:
 vmovups xmm8, xmmword ptr [rcx + 8*rdi]
 vmovups xmm1, xmmword ptr [rcx + 8*rdi + 32]
 vmovups xmm2, xmmword ptr [rcx + 8*rdi + 64]
 vmovups xmm3, xmmword ptr [rcx + 8*rdi + 96]
 vinsertf128 vmm8, vmm8, xmmword ptr [rcx + 8*rdi + 16], 1
 vinsertf128 ymm1, ymm1, xmmword ptr [rcx + 8*rdi + 48], 1
 vinsertf128 vmm2, vmm2, xmmword ptr [rcx + 8*rdi + 88], 1
 vinsertf128 vmm3, vmm3, xmmword ptr [rcx + 8*rdi + 112], 1
 vmulpd ymm0, ymm0, ymm0
 vmulpd vmm1, vmm1, vmm1
 vmulpd ymm2, ymm2, ymm2
 vmulpd ymm3, ymm3, ymm3
 vextractf128 xmmword ptr [rcx + 8*rdi + 16], vmm0, 1
 vmovupd xmmword ptr [rcx + 8*rdi], xmm8
 vextractf128 xmmword ptr [rcx + 8*rdi + 48], vmm1, 1
 vmovupd xmmword ptr [rcx + 8*rdi + 32], xmm1
 vextractf128 xmmword ptr [rcx + 8*rdi + 80], ymm2, 1
 vmovupd xmmword ptr [rcx + 8*rdi + 64], xmm2
 vextractf128 xmmword ptr [rcx + 8*rdi + 112], ymm3, 1
 vmovupd xmmword ptr [rcx + 8*rdi + 96], xmm3
 add rdi, 16
 cmp rsi, rdi
 ine .LBB0 3
 cmp rax. rsi
 je .LBB0_7
LBB0 5:
 dec rdx
LBB0 6:
 vmovsd xmm8, gword ptr [rex + 8*rdx]
 vmulsd xmm0, xmm0, xmm0
 vmovsd gword ptr [rex + 8*rdx], xmm0
 inc rdx
 cmp rdx, rax
 ib .LBB0 6
.LBB0_7:
 vzeroupper
 ret
```



Remove some

```
.LBB0 3:
 vmovuns xmm0, xmmword ntr [rcx + 8*rdi]
 vmovups xmm1, xmmword ptr [rcx + 8*rdi + 32]
 vmovups xmm2, xmmword ptr [rcx + 8*rdi + 64]
 vmovuns xmm3, xmmword ntr [rcx + 8*rdi + 96]
 vinsertf128 vmm0, vmm0, xmmword ptr [rcx + 8*rdi + 16], 1
 vinsertf128 ymm1, ymm1, xmmword ptr [rcx + 8*rdi + 48], 1
 vinsertf128 vmm2, vmm2, xmmword ptr [rcx + 8*rdi + 80], 1
 vinsertf128 ymm3, ymm3, xmmword ptr [rcx + 8*rdi + 112], 1
 vmulpd ymm0, ymm0, ymm0
 vmulpd vmm1, vmm1, vmm1
 vmulpd ymm2, ymm2, ymm2
 vmulpd vmm3, vmm3, vmm3
 vextractf128 xmmword ptr [rcx + 8*rdi + 16], ymm0, 1
 vmovupd xmmword ptr [rcx + 8*rdi], xmm0
 vextractf128 xmmword ptr [rcx + 8*rdi + 48], ymm1, 1
 vmovupd xmmword ptr [rcx + 8*rdi + 32], xmm1
 vextractf128 xmmword ptr [rcx + 8*rdi + 80], ymm2, 1
 vmovund xmmword ntr [rcx + 8*rdi + 64], xmm2
 vextractf128 xmmword ptr [rcx + 8*rdi + 112], vmm3, 1
 vmovupd xmmword ptr [rcx + 8*rdi + 96], xmm3
 add rdi. 16
 cmp rsi, rdi
 ine .LBB0 3
```

Julia has unrolled the loop and used SIMD-instructions!

```
void square(double *xs, int N) {
 for (int i = 0; i < N; i++) {
   xs[i] *= xs[i];
.L1:
  vmovupd vmm0.YMMWORD PTR [rdi+rsi*1]
  vmovupd ymm1,YMMWORD PTR [rdi+rsi*1+0x20]
  vmovupd vmm2.YMMWORD PTR [rdi+rsi*1+0x40]
  vmovupd ymm3,YMMWORD PTR [rdi+rsi*1+0x60]
  vmulpd ymm0,ymm0,ymm0
  vmulpd vmm1.vmm1.vmm1
  vmulpd ymm2,ymm2,ymm2
  vmulpd vmm3, vmm3, vmm3
  vmovupd YMMWORD PTR [rdi+rsi*1],ymm0
  vmovupd YMMWORD PTR [rdi+rsi*1+0x20].vmm1
  vmovupd YMMWORD PTR [rdi+rsi*1+0x40],ymm2
  vmovupd YMMWORD PTR [rdi+rsi*1+0x60],ymm3
        rsi.0xffffffffffff80
  sub
        rdx,rsi
  cmp
  ine
        .L1
```

This is what Clang does.

(Slightly better due to Julia-arrays being strided)

Julia uses LLVM



Julia can do the same optimizations we expect in C/C++

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Julia can do the same optimizations we expect in C/C++

 $... because \ it \ is \ literally \ the \ same \ optimizer!$

But there are some <u>huge</u> caveats

1. There is some *slight* compilation overhead at startup

However: As problem size grows, relative compilation time decreases.

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BUT THERE ARE SOME HUGE CAVEATS

- 1. There is some slight compilation overhead at startup
- 2. Julia is garbage-collected
- 3. We have no explicit control over memory (Allocation, copying, ...)

Accidental heap operations can really hurt!

But there are some <u>huge</u> caveats

- 1. There is some *slight* compilation overhead at startup
- 2. Julia is garbage-collected
- 3. We have no explicit control over memory (Allocation, copying, ...)
- 4. The whole thing breaks if Julia can't figure out the types

There are many ways to shoot yourself in the foot! \cdots

```
Δx = 0.1

function riemann!(X, x)

N = Length(x)

X[0] = 0.0

for i in 2:N

X[i] = X[i - 1] + Δx * X[i]

end

v = rand(100000)

V = zeros(100000)

Qtime riemann!(V, v)
```

```
const \( \Delta \times = 0.1 \)

function riemann!(X, x)

N = length(x)

X[0] = 0.0

for i in 2:N

X[i] = X[i - 1] + \( \Delta \times X[i] \)

end

end

v = rand(100000)

V = zeros(100000)

@time riemann!(V, v)
```

Two seemingly quite similar programs... $\,$

```
Δx = 0.1

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    end
end

v = rand(100000)
v = zeros(100000)
0time riemann!(V, v)
```

0.000175 seconds

One is $\sim 60 \times$ faster than the other!

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        X[i] = X[i - 1] + \Delta x * X[i]
    end
end

v = rand(100000)
v = zeros(100000)
@time riemann!(V, V)
```

0.000175 seconds

0.010550 seconds (499.49 k allocations)

And 500K allocations? 🤯

```
Δx = 0.1

function riemann!(X, x)

N = length(x)

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for i in 2:N

X[i] = X[i - 1] + Δx * X[i]

end

end

v = rand(100000)

V = zeros(100000)

Qtime riemann!(V, v)
```

```
const \( \Delta x = 0.1 \)
function riemann!(X, \( x \))
    N = length(x)
    X[0] = 0.0
    for i in 2:N
        X[i] = X[i - 1] + \( \Delta x \) X[i]
    end
end

v = rand(100000)
v = zeros(1000000)
@time riemann!(V, \( v \))
```

0.000175 seconds

And 500K allocations? 🔯

 Δx could be reassigned \implies Julia can't assume its type!

```
Δx = 0.1

function riemann!(X, x)

N = length(x)

X[0] = 0.0

for i in 2:N

X[i] = X[i - 1] + Δx * X[i]

end

end

v = rand(100000)

V = zeros(100000)

@time riemann!(V, v)
```

```
const \( \Delta x = 0.1 \)
function riemann!(X, X)
    N = length(X)
    X[e] = 0.0
    for in 2:N
        X[i] = X[i - 1] + \( \Delta x * X[i] \)
    end
end

v = rand(100000)
v = zeros(1000000)
@time riemann!(V, v)
```

0.000175 seconds

And 500K allocations? 🔯

Leads to runtime type checks and dynamic dispatching...

```
Δx = 0.1

function riemann!(X, x)

N = length(x)

X[0] = 0.0

for i in 2:N

X[i] = X[i - 1] + Δx * X[i]

end

end

v = rand(100000)

V = zeros(100000)

@time riemann!(V, v)
```

0.000175 seconds

And 500K allocations? 🔯

 \ldots and forces intermediate calculations onto the heap!

```
\Delta x = 0.1
function riemann!(X, x)
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v = rand(100000)
V = zeros(100000)
@time riemann!(V, v)
```

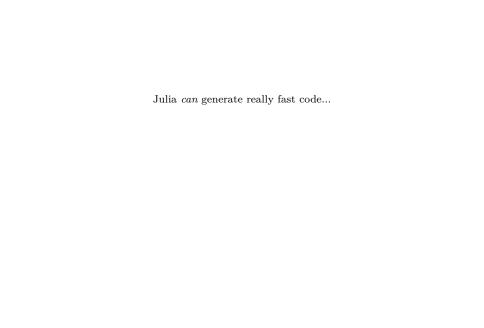
0.000175 seconds

And 500K allocations?



...and forces intermediate calculations onto the heap!







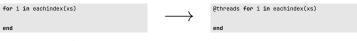
Julia can generate really fast code...

But the *lack of explicit control* requires a lot of attention.

Even more so when we parallelize our programs.

Multicore

using .Threads: @threads



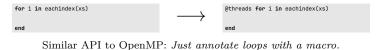
Similar API to OpenMP: Just annotate loops with a macro.



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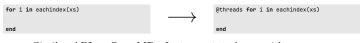
The normal caveats apply

1. Spawning threads involve some overhead



THE NORMAL CAVEATS APPLY

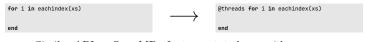
- 1. Spawning threads involve some overhead
- 2. Memory is shared (Race conditions, deadlocks)



Similar API to OpenMP: Just annotate loops with a macro.

THE NORMAL CAVEATS APPLY

- 1. Spawning threads involve some overhead
- 2. Memory is shared (Race conditions, deadlocks)
- 3. Possible speedup severely limited by core count



Similar API to OpenMP: $Just\ annotate\ loops\ with\ a\ macro.$

The normal caveats apply

- 1. Spawning threads involve some overhead
- 2. Memory is shared (Race conditions, deadlocks)
- 3. Possible speedup severely limited by core count
- 4. Domain decomposition sometimes necessary to get speedup on simple problems

GPU PARALLELISM



LEVEL 1: "VECTORIZED" ARRAY API

(Vectorized in the sense that you only operate on arrays, Numpy style)

```
# In-place operations.
v = CUDA.rand(10000000)
v *= 10.0
v ^= 2
v •= sqrt.(v)

# Parallel scan.
w = CUDA.zeros(10000000)
accumulate!(max, w, v)
```

v, w of type CuArray that manages a device-side buffer.

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v *= 10.0
v ^= 2
v •= sqrt.(v)

# Parallel scan.
w = CUDA.zeros(10000000)
accumulate!(max, w, v)
```

Block- and gridsizes automatically determined.

```
# FFT (10M-point FFT takes 200ms on 980)
using CUDA.CUFFT
\( \omega = \text{fft}(v) \)

# Kernels get fused
\( \omega = \omega \cdot \cdot
```

Bindings to CUDA libraries like cuFFT and cuBLAS (Also cuSPARSE, cuSOLVER, ...)

```
# FFT (10M-point FFT takes 200ms on 980)
using CUDA.CUFFT
\omega = fft(v)

# Kernels get fused
w •= w.^2 .+ v

# Check out what happens on the GPU
@device_code_sass w •= w.^2 .+ v
```

GPU code generated on the fly using LLVMs PTX/SASS backend... (Same sort of infrastructure as specializing Julia functions)

```
# FFT (10M-point FFT takes 200ms on 980)
using CUDA.CUFFT
ω = fft(v)

# Kernels get fused
w •= w.^2 .+ v

# Check out what happens on the GPU
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```

Enables kernel fusion and global optimization (!)

```
# FFT (10M-point FFT takes 200ms on 980)
using CUDA.CUFFT

\omega = fft(v)

# Kernels get fused
w \cdot = w.^2 .+ v

# Check out what happens on the GPU
@device_code_sass w \cdot = w.^2 .+ v
```

Like with CPU code, you can inspect the generated GPU code. (SASS/PTX code is too long/ugly to put on a slide)

LEVEL 2: JULIA ON THE GPU

```
function cusquare!(xs)
    i = (blockIdx().x - 1) * blockDim().x + threadIdx().x
    if i < length(xs)
        xs[i] *= xs[i]
    end
    return nothing
end

Nt = 1024
Nb = ceil(Int, length(v) / Nt)
@cuda threads=Nt blocks=Nb cusquare!(v)</pre>
```

cusquare! is defined as a regular Julia function!

(It won't run if you call it; block- and thread-ID is not defined on the CPU)

```
function cusquare!(xs)
    i = (blockIdx().x - 1) * blockDim().x + threadIdx().x
    if i < length(xs)
        xs[i] *= xs[i]
    end
    return nothing
end

Nt = 1024
Nb = ceil(Int, length(v) / Nt)
@cuda threads=Nt blocks=Nb cusquare!(v)</pre>
```

The Qcuda macro transforms typed Julia IR into GPU code!



```
function cusquare!(xs)
    i = (blockIdx().x - 1) * blockDim().x + threadIdx().x
    if i < length(xs)
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    end
    return nothing
end

Nt = 1024
Nb = ceil(Int, length(v) / Nt)
@cuda threads=Nt blocks=Nb cusquare!(v)</pre>
```

Basically, almost any sensible Julia can just work on the GPU

```
function cusquare!(xs)
    i = (blockIdx().x - 1) * blockDim().x + threadIdx().x
    if i < length(xs)
        xs[i] *= xs[i]
    end
    return nothing
end

Nt = 1024
Nb = ceil(Int, length(v) / Nt)
@cuda threads=Nt blocks=Nb cusquare!(v)</pre>
```

However, there is no escape hatch: The code must type check, and can't allocate

PITFALLS

PITFALLS

CuArray controls all memory transfer

PITFALLS

Accidentally reading back to the CPU is really painful!

Future Outlook

LEVEL 3: MULTI-GPU / MULTI-NODE

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CUDA.jl composes with DistributedArrays.jl

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DistributedArrays.jl: Domain decomposition and distribution algorithms

CUDA.jl: Manages device-side buffers and host-device communication

Both understands Julias abstract concept of an "Array", so they work together.

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CUDA.jl composes with DistributedArrays.jl

DistributedArrays.jl: Domain decomposition and distribution algorithms

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Both understands Julias abstract concept of an "Array", so they work together.

Rapid Software Prototyping for Heterogeneous and Distributed Platforms
Besard et. al.

Julia makes bold claims

Julia makes bold claims: "Looks like Python – runs like C/C++/Fortran"

Julia makes bold claims: "Only dynamic language to achieve petascale"

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It is a truth with a big asterisk.

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If there is one thing that is true, it is that you can't escape

LOCATION, LOCATION

Julia makes bold claims: "Only dynamic language to achieve petascale"

If there is one thing that is true, it is that you can't escape

LOCATION, LOCATION

Julia has real potential, but it is not a free lunch! (Especially on the GPU)

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If there is one thing that is true, it is that you can't escape

LOCATION, LOCATION

I spend a lot of time "fighting the abstraction"

Julia makes bold claims: "Only dynamic language to achieve petascale"

If there is one thing that is true, it is that you can't escape

LOCATION, LOCATION

Allocation, garbage-collection, copies, CUDA device $\leftrightarrow \! host$ transfers, etc. are all implicit!

Julia makes bold claims: "Only dynamic language to achieve petascale"

If there is one thing that is true, it is that you can't escape

LOCATION, LOCATION

Allocation, garbage-collection, copies, CUDA device $\leftrightarrow \! host$ transfers, etc. are all implicit!

As Julia takes away your control, you have to be careful.

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LOCATION, LOCATION

Allocation, garbage-collection, copies, CUDA device $\leftrightarrow \! host$ transfers, etc. are all implicit!

The abstraction leaks!

What niche does Julia even fill?

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If you don't understand the compilation process, Julia is basically unusable.

What niche does Julia even fill?

 ${\it Julia\ wont't\ replace\ lower-level\ languages}.$

That being said...

I quite like Julia

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Julia makes HPC a first-class citizen in the "Matlab workflow"

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Julia makes HPC a first-class citizen in the "Matlab workflow"

- Interactive REPL/Notebook-based workflow
- 2. Good for fucking around with data
- 3. Sensible package manager
- 4. "Math stuff" is easily accessible
- 5. Visualization is made easy

