Performance and profiling in Julia

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Outline

- 1. Write code with performance in mind (think like a compiler)
- 2. Profile your code
- 3. Optimize your code

Step one, put your code in functions

A global variable might have its value and its type change at any point. This makes it difficult for the compiler to optimize code using global variables.



Any code that is performance critical or being benchmarked should be inside a function.

```
function foo()
                                             a = 1
a = 1
                                             for i = 1:100 000
\texttt{@btime for i} = 1:100 000
                                                  a += 1
    qlobal a
                                             end
    a += 1
                                             а
end
                                         end
    1.543 ms
                                         @btime foo()
    (100000 allocations: 1.53 MiB)
                                             1.270 ns
                                              (0 allocations: 0 bytes)
```

Avoid global variables (unless declared const or type annotated)

```
PRINT = false
                                    const PRINT = false
function foo()
                                    function foo()
  for i = 1:1 000 000 000
                                      for i = 1:1 000 000 000
    if PRINT
                                        if PRINT
      print(i)
                                          print(i)
    end
                                        end
  end
                                      end
end
                                    end
@time foo()
                                   @time foo()
  0.795711 seconds
                                      0.000002 seconds
```

PRINT is false in both cases, but the compiler can rely on it in the second case

Type declarations, type stability

Useful as assertion for debugging, but does not make the code faster (unless mitigating type instability).

Exception: Declare specific types for fields of composite types so that the compiler knows the memory layout

```
struct Foo struct Foo field::Type end end
```

It is in general bad for performance when the type of a variable can be changed at runtime, type annotation will prevent this.

Type stability

runtime.

An example of type instability

```
stable(i) = rand() > .5 ? 1 : -1
                                     unstable(i) = rand() > .5 ? 1. : -1
function foo()
                                     function bar()
    a = 1
                                         a = 1
    for i = 1:100 000
                                         for i = 1:100 000
        a += stable(i)
                                             a += unstable(i)
    end
                                         end
    а
                                         а
end
                                     end
@btime foo()
                                     @btime bar()
  162.940 us
                                       684.704 us
In bar the compiler does not know ahead of time which +
method to call (Int or Float64) and dispatch happens at
```

Type stability

Now the compiler knows that the type is Int after the type assertion, and dispatch can be determined at compile time, even though the return type of quasistable is set valued.

Julia uses column major convention

```
function foo()
                                          function bar()
    x = Matrix{Float64}(undef, 4096, 4096)
                                              x = Matrix{Float64}(undef, 4096, 4096)
    for i = 1:size(x.1)
                                              for j = 1:size(x,2)
        for j = 1:size(x,2)
                                                   for i = 1:size(x,1)
            x[i.i] = i*i
                                                      x[i.i] = i*i
        end
                                                   end
    end
                                              end
end
                                          end
@btime foo()
                                          @btime bar()
102.165 ms (2 allocations: 128.00 MiB) 19.745 ms (2 allocations: 128.00 MiB)
```

Think about this when you are choosing how to store your data!

Avoid unnecessary memory allocation

Julia passes arrays as references.¹ Use this to re-use already allocated memory.

```
function food()
    A = Matrix(Int64) {undef,100,100}
    for i = eachindex(A)
        A[i] = i
    end
    return A
end

function eat()
    for i = 1:10_000
        chicken = food()
        sum(chicken)
    end
end

@btime eat()
144.466 ms (20000 allocations: 763.70 MiB)
```

New plate every time, lots of time to clean! (garbage collect)

```
function beer!(A)
    for i = eachindex(A)
        A[i] = i
    end
end

function drink()
    glass = Matrix{Int64}{undef,100,100}
    for i = 1:10_000
        beer!(glass)
        sum(glass)
    end
end
end
end
ebtime drink()
54.119 ms (2 allocations: 78.20 KiB)
```

Use the same glass every time, drink beer faster!



¹Technically, by sharing

Try to keep structs immutable

- ▶ Mutable structs (typically) end up on the heap.
- isbits structs (typically) end up on the stack.
- ▶ Mutability is bug prone.

```
julia> struct Foo x::Float64 end

julia> isbitstype(Foo)
true

julia> mutable struct Bar x::Float64 end
julia> isbitstype(Bar)
false
```

Profile your code

Profiling

Profiling

Your goto-tool is always @btime from BenchmarkTools.jl, watch memory allocation and GC-time

- ► Type instability
- ► Allocations
- ▶ Do not benchmark in global scope
- ▶ Do not benchmark compilation time (unless it gets insanely long)
- ▶ Interpolate global variables @btime testfun(\$a)
- **©btime** runs the expression several times and reports the minimum

Profiling

```
Julia has built in profiling capabilities
julia> @profile foo()
julia> Profile.print()
     23 client.jl; start; line: 373
       23 client.jl; run_repl; line: 166
          23 client.jl; eval_user_input; line: 91
             23 profile.jl; anonymous; line: 14
                8 none; myfunc; line: 2
                 8 dSFMT.jl; dsfmt gv fill array close open!; line: 128
                15 none; myfunc; line: 3
                 2 reduce.jl; max; line: 35
                 2 reduce.jl; max; line: 36
                 11 reduce.jl; max; line: 37
```

Vscode profile view

The profile viewer in Vscode is nicer

@profview bode(P)

```
we were a significant content of the fragment regime of system 'sys' at the fragment of system 'sys' at the size (linguistic), series (
```

See also ProfileView.jl and StatProfilerHTML.jl

Profiling tools

@profview_allocs

An allocation profiler, similar to the time profiler.

@code_warntype

Ask the compiler what it thinks about your types

```
julia> naive_relu(x) = sum(x < 0 ? 0 : x)
```

```
iulia> @code_warntype naive_relu(1.0)
MethodInstance for naive_relu(::Float64)
  from naive_relu(x) in Main at REPL[39]:1
Arguments
  #self#::Core.Const(naive_relu)
  x::Float64
  @_3::Union{Float64, Int64}
Body::Union{Float64, Int64}
         goto #3 if not %1
         goto #4
   %6 = (0_3::Union{Float64, Int64}
    %7 = Main.sum(%6)::Union{Float64, Int64}
         return %7
```

Benchmarking

- ▶ Put your code in functions
- Let the function compile before timing (or use @btime)
- ▶ Watch out for unexpected memory allocation
- ▶ Read the performance tips!

Optimize your code

Optimization

Optimize your code

- ▶ Write a test before you start optimizing to make sure you are still calculating the same thing
- ▶ Use the result of @profview, @profview_allocs, @btime
- ▶ If your code spends 50% doing garbage collection, you can sometimes reduce your running time with more than 50% by better memory management.

Optimize your code

```
function slowfunc(x)
    a = foo(x)
    for i = 1:1000
        b = bar(a, randn(2))
        for j = 1:1000
            c = randn(5)
            b += baz(c)
        end
        a += b
    end
    a
end
```

Where do you start looking?

SIMD

Single Instruction Multiple Data

```
a64 = randn(1000000)
a32 = randn(Float32, 1000000)
function regular sum(x)
                                  function simd sum(x)
   s = zero(eltype(x))
                                       s = zero(eltype(x))
   for i = eachindex(x)
                                       @inbounds @simd for i = eachindex(x)
       s += x[i]
                                          s += x[i]
   end
                                       end
   S
                                       S
end
                                   end
@btime regular sum($a64)
@btime simd sum($a64)
646.580 μs (0 allocations: 0 bytes) 62.899 μs (0 allocations: 0 bytes)
@btime regular sum($a32)
@btime simd sum($a32)
610.543 µs (0 allocations: 0 bytes) 31.910 µs (0 allocations: 0 bytes)
```

SIMD

${\bf Loop Vectorization. jl}$

```
using LoopVectorization
function turbo_sum(x)
    s = zero(eltype(x))
    @tturbo for i = eachindex(x)
        s += x[i]
    end
    S
end
@btime turbo_sum($(a64))
12.624 µs (0 allocations: 0 bytes)
@btime turbo_sum($(a32))
11.332 µs (0 allocations: 0 bytes)
```

StaticArrays

One of the biggest speed-ups (after choosing the right algorithm) if often to use StaticArrays where available

- ▶ Size known at compile time
- Optimized operations
- ► Stack allocated (as opposed to heap allocated)

Benchmarks for 3×3 Float64 matrices

```
Matrix multiplication -> 8.2x speedup
Matrix multiplication (mutating) -> 3.1x speedup
Matrix addition -> 45x speedup
Matrix addition (mutating) -> 5.1x speedup
Matrix determinant -> 170x speedup
Matrix inverse -> 125x speedup
Matrix symmetric eigendecomposition -> 82x speedup
Matrix Cholesky decomposition -> 23.6x speedup
```

StaticArrays

Details

- ➤ Size and type hard coded, known at compile time
- ▶ VectorSVector has same memory layout as Matrix

Misc.

end

```
FillArrays.il Represent special arrays efficiently
repmat, repeat If you use these to force your problem into a vectorized
              form, you need to de-matlabify yourself
collect(1:10) You often do not need to collect.
Avoid allocating slices
              A[:,i] allocates and copies data, @view(A[:,i]) doesn't.
              (A[i,:] might however be worth it.)
     Parallel Threading and distributed computing
   dot-fusion R = sin.(exp.(A.^2)) compiles into a single loop
                No temporary arrys
                ► Single pass over data
    R = similar(a)
    for i in eachindex(a)
        R[i] = sin(exp(a[i]^2))
```

Other resources

▶ I would first and foremost recommend the performance tips section in the manual, it's quite comprehensive and readable: https:

//docs.julialang.org/en/v1/manual/performance-tips/index.html

Chris Rackaukas has some tutorials on solving ODEs and PDEs in Julia. He highlights a lot of neat Julia functionality and goes through a lot of performance optimizations that extend also outside the realm of ODEs and PDEs https://youtu.be/KPEqYtEd-zY Watch around minute 49 for performance optimization https://youtu.be/okGybBmih0E

An introduction to high performance custom arrays | Matt Bauman https://www.youtube.com/watch?v=jS9eouMJf_Y&t= 1831s&list=PLP8iPy9hna6Qsq5_-zrg0NTwqDSDYtfQB&index=82

Homework

Monte-Carlo simulation of a bootstrap particle filter

- ▶ I provide the baseline code
- ▶ My code provides a decent particle filter implementation
- ▶ The code is bad from a julia-performance point of view
- ▶ Your job is to optimize it
- Optimized code has to be equivalent (do not implement different algorithm)

$$\begin{split} x^+ &= 0.5x + \frac{25x}{1+x^2} + 8\cos(1.2(t-1)) + w \\ y &= 0.05x^2 + v \\ w, v &\sim \mathcal{N}(0, \sigma_w), \ \mathcal{N}(0, \sigma_v) \quad E(wv^\top) = 0 \end{split}$$

The particle filter

```
for t = 2:T # Main loop
    # Resample
    j = resample(w[t-1,:])
    # Time update
    xp[t,:] = f(xpT,t-1) + ow*randn(1,N)
    # Measurement update
    w[t,:] = wT + g(y[t]-0.05xp[t,:].^2)
    # Normalize weights
    w[t,:] -= log(sum(exp(w[t,:])))
end
```

The Monte-Carlo simulation

```
particle count = [5 10 20 50 100 200 500 1000 10 000]
time steps = [20, 200, 2000]
for (Ti,T) in enumerate(time steps)
  for (Ni, N) in enumerate(particle count)
    # Calculate how many Monte-Carlo runs to perform for the current
    # T,N configuration
    montecarlo runs =
        maximum(particle_count)*maximum(time_steps) / T / N
    for mc iter = 1:montecarlo runs
      for t = 1:T-1 # Simulate one realization of the model
        x[t+1] = f(x[t],t) + \sigma w*randn()
        y[t+1] = 0.05x[t+1]^2 + \sigma v*randn()
      end # t
      xh = pf(y, N, q, f, \sigma w0) # Run the particle filter
      RMS += rms(x-xh) # Store the error
    end # MC
```